

ARCADIS

Appendix A

Groundwater Sampling Logs

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Talleast	
WELL NO: DW-1	SAMPLE ID: DW-1	DATE: 02-01-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 82 feet to 92 feet	STATIC DEPTH TO WATER (feet): 13.65	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 0938	PURGING ENDED AT: 0954	TOTAL VOLUME PURGED (gallons): 1.6
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0948	1.0	1.0	0.1	20.21	7.95	28.04	123	0.51	97.1	clear	none
0951	0.3	1.3	0.1	20.51	8.01	28.03	127	0.47	97.1	clear	none
0954	0.3	1.6	0.1	20.75	8.02	27.97	135	0.49	95.1	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darin Johnson	SAMPLER(S) SIGNATURES: <i>Darin Johnson</i>	SAMPLING INITIATED AT: 0955	SAMPLING ENDED AT: 1000
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): < 100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: Y N	FIELD FILTERED: Y N	FILTER SIZE: _____ µm	DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
DW-1	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **transducers (2) removed @ 0935 / (2) reinserted @ 10:13**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: EW-UAFG-1	SAMPLE ID: EW-UAFG-1
DATE: 2-01-08	

PURGING DATA

WELL DIAMETER (inches): 6"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 98 feet to 108 feet	STATIC DEPTH TO WATER (feet): 15.61	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 108 feet) + 0.25 gallons = 0.53 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 103'	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 103'	PURGING INITIATED AT: 0930	PURGING ENDED AT: 0954	TOTAL VOLUME PURGED (gallons): 2.2

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0942	1.0	1.0	0.1	15.90	7.02	26.80	1.17	1.75	2.97	clear	none
0945	0.3	1.3	↓	15.91	7.02	26.80	1.17	1.58	1.39		
0948	↓	1.6	↓	15.92	7.02	26.86	1.17	1.37	1.27		
0951	↓	1.9	↓	15.92	7.02	26.79	1.17	1.35	1.19		
0954	↓	2.2	↓	15.92	7.02	26.80	1.17	1.34	1.43		
0957											

-241
-243
-245
-246
-246

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ann Coats / Arcadis	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 10:05	SAMPLING ENDED AT: 10:00
PUMP OR TUBING DEPTH IN WELL (feet): 103'	SAMPLE PUMP FLOW RATE (mL per minute): < 100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: (Y) N	FIELD-FILTERED: (Y) N	FILTER SIZE: _____ µm	DUPLICATE: Y (N)

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-UAFG-1	3	CG	40mL	HCl	—	—	8260	RFPP

MDS 2/11/08

REMARKS: **put tubing down well @ 9:03 - did not remove transducer**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Flow rate varied in beginning.
 1002 pump off - tubing back down well.
 total vol removed ~ 2.5 gallons

EW-101

DEP-SOP-001/01
Form FD 9000-24

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: EW-101	SAMPLE ID: EW-101
DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 6	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 30 feet	STATIC DEPTH TO WATER (feet): 3.93	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (.0026 gallons/foot X 35 feet) + .25 gallons = 0.34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): none	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 15	PURGING INITIATED AT: 12:30	PURGING ENDED AT: 12:45	TOTAL VOLUME PURGED (gallons): 1.5

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
12:36	.6	.6	0.1	4.20	5.32	29.59	1299	3.05	2.19	clear	none
12:39	.3	.9	0.1	4.23	5.31	29.67	1297	1.94	1.86	clear	none
12:42	.3	1.2	0.1	4.20	5.31	29.70	1299	1.90	1.90	clear	none
12:45	.3	1.5	0.1	4.24	5.31	29.72	1300	1.84	1.92	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darin Johnson	SAMPLER(S) SIGNATURES: Darin Johnson	SAMPLING INITIATED AT: 12:46	SAMPLING ENDED AT: 12:51
PUMP OR TUBING DEPTH IN WELL (feet): 15	SAMPLE PUMP FLOW RATE (mL per minute): 0.1	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N Filtration Equipment Type:	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-101	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

MDC
A/11/08

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Seasuta, FL
WELL NO: EW-102	SAMPLE ID: EW-102
	DATE: 1.31.08

PURGING DATA

WELL DIAMETER (inches): 6	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 31 feet to 36 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33.5				
FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33.5				
PURGING INITIATED AT: 1629		PURGING ENDED AT: 1647		TOTAL VOLUME PURGED (gallons): 1.44

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1638	0.12	0.72	0.08	8.42	6.90	25.05	1000	0.75	4.85	clear	NO
1641	0.24	0.96	0.08	8.42	6.91	25.02	1001	0.68	4.05	clear	NO
1644	0.24	1.20	0.08	8.42	6.91	24.99	1003	0.67	2.34	clear	NO
1647	0.24	1.44	0.08	8.42	6.96	24.99	1010	0.70	1.17	clear	NO

ORP
-188
-187
-186
-214

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0028; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/TH / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1648	SAMPLING ENDED AT: 1652
PUMP OR TUBING DEPTH IN WELL (feet): 33.5	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N	FILTRATION EQUIPMENT TYPE: <input checked="" type="radio"/> N	DUPLICATE: <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-102	3	CG	40mL	HCl	40mL	—	82608 82605im RFP	
FB 29	3	BCG	40mL	HCl	40mL	—	82605 82605im RFP	

1655

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC		SITE LOCATION: Tallevast	
WELL NO: EW-103	SAMPLE ID: EW-103	DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 6"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 5 feet to 20 feet	STATIC DEPTH TO WATER (feet): 3.14	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0020 gallons/foot X 10 feet) + 0.25 gallons = 0.28 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 5'	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 5'	PURGING INITIATED AT: 13:23	PURGING ENDED AT: 14:44	TOTAL VOLUME PURGED (gallons): 2.1 MDS 4/11/08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
14:32	0.1	0.1	0.1	3.92	6.03	23.7	423	0.80	37.3	clear	none
14:35	0.3	0.4	0.1	3.92	6.03	23.61	421	1.0	36.1	clear	none
14:38	0.3	0.7	0.1	3.93	6.04	23.59	420	1.89	31.1	clear	none
14:41	0.3	1.0	0.1	3.93	6.04	23.63	419	1.82	30.3	clear	none
14:44	0.3	1.3	0.1	3.93	6.04	23.58	420	1.78	31.0	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson		SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>		SAMPLING INITIATED AT: 14:45	SAMPLING ENDED AT: 14:50
PUMP OR TUBING DEPTH IN WELL (feet): 5		SAMPLE PUMP FLOW RATE (mL per minute): 500 MDS 4/11/08		TUBING MATERIAL CODE: pe	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> N		FILTER SIZE: _____ µm	
DUPLICATE: <input checked="" type="checkbox"/> N					

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-103	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABL Facility	SITE LOCATION: Tallevast
WELL NO: EW-104	SAMPLE ID: EW-104
DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 6	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 26 feet to 31 feet	STATIC DEPTH TO WATER (feet): 4.92	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= 0 gallons + (^{0.0210} **28** AC gallons/foot X ³³ **28** AC feet) + **0.25** gallons = **0.34** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 31	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 27	PURGING INITIATED AT: 14:20	PURGING ENDED AT: 14:41	TOTAL VOLUME PURGED (gallons): 2.1
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
14:29	0.1	0.1	0.1	5.48	5.53	24.91	565	2.80	44.5	red	none
14:32	0.3	1.2	0.1	5.54	5.54	24.9	566	2.13	42.4	reddish	none
14:35	0.3	1.5	0.1	5.58	5.56	24.87	567	1.88	43.2	red	none
14:38	0.3	1.8	0.1	5.59	5.59	24.84	568	1.79	40.3	red	none
14:41	0.3	2.1	0.1	5.59	5.6	24.84	569	1.79	38.5	red	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson / Arcadis	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 14:42	SAMPLING ENDED AT: 14:47
PUMP OR TUBING DEPTH IN WELL (feet): 28	SAMPLE PUMP FLOW RATE (mL per minute): <100 (MDS 4/11/08)	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	
Filtration Equipment Type: _____		DUPLICATE: Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-104	3	CG	40ml	HCl	—	—	8266	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

026
25

SITE NAME: Former ABC Facility SITE LOCATION: Tallevast, FL
WELL NO: EW-105 SAMPLE ID: EW-105 DATE: 1.30.08

PURGING DATA

WELL DIAMETER (inches): 6 TUBING DIAMETER (inches): 1/4 WELL SCREEN INTERVAL DEPTH: 5 feet to 25 feet STATIC DEPTH TO WATER (feet): 405 PURGE PUMP TYPE OR BAILER: PP

WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
= (5 feet - 405 feet) X 0.28 gallons/foot = 0.28 gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
= 0 gallons + (0.0526 gallons/foot X 10' feet) + 0.25 gallons = 0.28 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 5' FINAL PUMP OR TUBING DEPTH IN WELL (feet): 5' PURGING INITIATED AT: 11:40 PURGING ENDED AT: 12:07 TOTAL VOLUME PURGED (gallons): 2.3

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
11:55	1.1	1.1	0.1	4.12	6.42	28.73	1299	3.87	12.3	clear	none
11:58	0.3	1.4	0.1	4.14	6.42	28.71	1300	2.53	13.1	clear	none
12:01	0.3	1.7	0.1	4.14	6.41	28.66	1293	2.30	13.7	clear	none
12:04	0.3	2.0	0.1	4.16	6.39	28.74	1292	2.11	17.4	clear	none
12:07	0.3	2.3	0.1	4.16	6.39	28.71	1291	2.10	14.8	clear	none

116
109
108
105
103

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amey Coats / Arcadis SAMPLER(S) SIGNATURES: Amey Coats
PUMP OR TUBING DEPTH IN WELL (feet): 5 SAMPLE PUMP FLOW RATE (mL per minute): <100 TUBING MATERIAL CODE: teflon
FIELD DECONTAMINATION: Y N FIELD-FILTERED: Y N FILTER SIZE: 1µm DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-105	3	CG	10ml	HCl	←	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast
WELL NO: EW-106	SAMPLE ID: EW-106 DATE: 1-30-08

PURGING DATA

WELL DIAMETER (inches): 6"	TUBING DIAMETER (inches): 1 1/4"	WELL SCREEN INTERVAL DEPTH: 20 feet to 31 feet	STATIC DEPTH TO WATER (feet): 560	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 32 feet) + 0.25 gallons = 0.33 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 28	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 28	PURGING INITIATED AT: 11:38	PURGING ENDED AT: 12:02	TOTAL VOLUME PURGED (gallons): 2.4
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
11:50	1.2	1.2	0.1	6.53	5.57	28.42	463	.46	42.4	clear	none
11:53	.3	1.5	0.1	6.60	5.58	28.33	462	.45	20.6	clear	none
11:56	.3	1.8	0.1	6.62	5.57	28.42	461	.43	15.8	clear	none
11:59	.3	2.1	0.1	6.68	5.57	28.38	460	.41	12.8	clear	none
12:02	.3	2.4	0.1	6.68	5.56	28.43	460	.29	14.4	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amy Coats / Arcadis	SAMPLER(S) SIGNATURES: <i>Amy Coats</i>	SAMPLING INITIATED AT: 12:03	SAMPLING ENDED AT: 12:08
PUMP OR TUBING DEPTH IN WELL (feet): 28	SAMPLE PUMP FLOW RATE (mL per minute): < 100	TUBING MATERIAL CODE: Teflon	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-106	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Seasota, FL
WELL NO: EW-07	SAMPLE ID: EW-107
DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 6	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 7 feet to 27 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.026 gallons/foot X 30 feet) + 0.25 gallons 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 9.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 9.5	PURGING INITIATED AT: 1540	PURGING ENDED AT: 1555	TOTAL VOLUME PURGED (gallons): 1.65

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1546	0.66	0.66	0.11	3.41	6.62	28.39	285	1.19	2.71	clear	NO
1552	0.33	0.99	0.11	3.41	6.54	28.34	280	0.83	2.32	clear	NO
1552	0.33	1.32	0.11	3.41	6.47	28.32	279	0.68	2.37	clear	NO
1555	0.33	1.65	0.11	3.41	6.42	28.25	278	0.69	2.34	clear	NO

1549
1552
1555

OLP
-274
-273
-284
-285

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW, TH / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1556	SAMPLING ENDED AT: 1600
PUMP OR TUBING DEPTH IN WELL (feet): 9.5	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-107	3	CG	40ml	HCl	40ml	---	8260B, 8260SIM	RFPP
DUP 29	3	CG	40ml	HCl	40ml	---	8260B, 8260SIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Sarasota, FL
WELL NO: EW-108	SAMPLE ID: EW-108
DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 6	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 43.5 feet + 0.25 gallons = 0.36 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 1423	PURGING ENDED AT: 1443	TOTAL VOLUME PURGED (gallons): 1.40

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1434	0.77	0.77	0.07	8.51	7.95	29.59	970	0.65	4.07	clear	NO	ORP -362
1437	0.21	0.98	0.07	8.52	7.99	29.61	977	0.53	3.73	clear	NO	-370
1440	0.21	1.19	0.07	8.51	8.04	29.62	1002	0.52	3.90	clear	NO	-371
1443	0.21	1.40	0.07	8.53	8.05	29.63	1026	0.50	3.32	clear	NO	-382

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: TE/SW/ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1444	SAMPLING ENDED AT: 1447
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 2/1/08	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N Filtration Equipment Type: <input checked="" type="radio"/> <input type="radio"/>	FILTER SIZE: _____ µm	
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N MDS 2/1/08	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-108	3	CG	40ml	HCl	40ml	—	82608+82603im RFPF	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallevast, FL	
WELL NO: EW-109	SAMPLE ID: EW-109	DATE: 1.30.08	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: 5 feet to 20 feet	STATIC DEPTH TO WATER (feet): 3.27	PURGE PUMP TYPE OR BAILER: PD
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 5'	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 5'	PURGING INITIATED AT: 15:15	PURGING ENDED AT: 15:32	TOTAL VOLUME PURGED (gallons): 1.7
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
15:23	0.8	0.8	0.1	3.50	6.34	24.67	1.51	2.88	5.01	clear	none
15:26	0.3	1.1	0.1	3.53	6.29	24.58	1.51	2.07	4.06	clear	none
15:29	0.3	1.4	0.1	3.53	6.34	24.47	1.51	1.96	2.96	clear	none
15:32	0.3	1.7	0.1	3.53	6.35	24.49	1.51	1.90	3.10	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson/Aradis	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 15:33	SAMPLING ENDED AT: 15:38
PUMP OR TUBING DEPTH IN WELL (feet): 5'	SAMPLE PUMP FLOW RATE (mL per minute): <100 <i>MDS 4/11/08</i>	TUBING MATERIAL CODE: Teflon	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-109	3	CG	40ml	HCl	—	—	8260	RFPP <i>MDS 4/11/08</i>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: EW-110	SAMPLE ID: EW-110 DATE: 1.30.08

PURGING DATA

WELL DIAMETER (inches): 6"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 30 feet to 35 feet	STATIC DEPTH TO WATER (feet): 35.3	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable) = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = **0** gallons + (**0.0026** gallons/foot X **40** feet) + **0.25** gallons = **0.35** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 28'	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 28'	PURGING INITIATED AT: 15:25 15:25	PURGING ENDED AT: 15:30	TOTAL VOLUME PURGED (gallons): 1.5
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
15:24	0.1	0.1	0.1	4.16	5.01	24.77	1.35	1.51	6.40	clear	none
15:27	0.3	1.2	0.1	4.19	5.01	24.72	1.34	1.46	6.16	clear	none
15:30	0.3	1.5	0.1	4.22	5.0	24.74	1.32	1.32	5.94	clear	none
15:29											

15:24
15:27
15:30
~~15:29~~

67
65
66

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ami Gade / Arcadis	SAMPLER(S) SIGNATURES: <i>Ami Gade</i>	SAMPLING INITIATED AT: 15:31	SAMPLING ENDED AT: 15:36
PUMP OR TUBING DEPTH IN WELL (feet): 28' ^{MB 4/1/08}	SAMPLE PUMP FLOW RATE (mL per minute): <100 ^{MB 9/11/08}	TUBING MATERIAL CODE: Teflon	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
EW-110	3	CG	40mL	HCl	—	—	8260	RFPP

^{MB 9/11/08}

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: IWI-1	SAMPLE ID: IWI-1 DATE: 2-1-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 100 feet to 110 feet	STATIC DEPTH TO WATER (feet): 15.8	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 109.65 feet + 0.25 gallons = gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 104	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 104	PURGING INITIATED AT: 8:31	PURGING ENDED AT: 9:10	TOTAL VOLUME PURGED (gallons): 3.9

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
8:40	0.9	0.9	0.1	16.15	6.85	27.3	1.75	1.27	>999	gray	none	-156
8:45	0.5	1.4	0.1	16.15	6.90	27.3	1.70	1.01	300	"	"	-165
8:50	0.5	1.9	0.1	16.15	6.91	27.4	1.79	0.87	48	"	"	-175
8:55	0.5	2.4	0.1	16.20	6.90	27.5	1.83	0.76	46	"	"	-190
9:00	0.5	2.9	0.1	16.2	6.90	27.4	1.90	0.69	15	"	"	-206
9:05	0.5	3.4	0.1	16.2	6.90	27.4	2.02	0.65	16	"	"	-216
9:10	0.5	3.9	0.1	16.2	6.89	27.5	2.10	0.61	26	"	"	-220

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 9:11	SAMPLING ENDED AT: 9:15
PUMP OR TUBING DEPTH IN WELL (feet): 104 MDS 4/11/08	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/11/08	TUBING MATERIAL CODE: T MDS 4/11/08	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
IWI-1	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: IWI-2	SAMPLE ID: IWI-2 DATE: 2.1.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet): 19.1	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.026 gallons/foot X 169.12 feet) + 0.25 gallons = 0.69 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 165	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 165	PURGING INITIATED AT: 9:28	PURGING ENDED AT: 10:20	TOTAL VOLUME PURGED (gallons): 2.60

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
9:45	0.85	0.85	0.05	22.7	7.21	27.5	0.875	0.79	7999	gray	none	-243
9:50	0.25	1.10	0.05	23.8	7.21	27.5	0.940	0.72	190	"	"	-249
9:55	0.25	1.35	0.05	23.8	7.22	27.5	1.06	0.69	120	"	"	-250
10:00	0.25	1.60	0.05	23.8	7.21	27.4	1.06	0.64	754	"	"	-252
10:05	0.25	1.85	0.05	25.8	7.22	27.7	1.05	0.62	257	"	"	-252
10:10	0.25	2.10	0.05	25.8	7.21	27.61	1.04	0.59	36	"	"	-251
10:15	0.25	2.35	0.05	25.8	7.22	27.4	1.02	0.59	23	"	"	-251
10:20	0.25	2.60	0.05	25.8	7.21	27.31	1.02	0.57	21	"	"	-250

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:21	SAMPLING ENDED AT: 10:25
PUMP OR TUBING DEPTH IN WELL (feet): 165 ^{MDS 9/11/08}	SAMPLE PUMP FLOW RATE (mL per minute): <100 ^{MDS 9/11/08}	TUBING MATERIAL CODE: T	^{MDS 9/11/08}
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
IWI-2	3	CG	40ml	HCl	—	—	8260	RFPP ^{MDS 9/11/08}

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellovert</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>PZ-LSAS-1</u>	SAMPLE ID: <u>PZ-LSAS-1</u> DATE: <u>1/31/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>30.4</u> feet to <u>35.46</u> feet	STATIC DEPTH TO WATER (feet): <u>5.62</u>	PURGE PUMP TYPE OR BAILER: <u>(PP) peristaltic pump</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)

= gallons + (0.0026 gallons/foot X 39 feet) X 0.25 gallons = 0.35 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~32</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~32</u>	PURGING INITIATED AT: <u>1350</u>	PURGING ENDED AT: <u>1402</u>	TOTAL VOLUME PURGED (gallons): <u>1.80</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1353	0.45	0.45	0.15	6.65	6.22	29.79	1.71	0.74	14.6	clear	yes	-106
1356	0.45	0.90	0.15	6.68	6.24	29.74	1.74	0.63	6.41	clear	yes	-112
1359	0.45	1.35	0.15	6.68	6.25	29.75	1.74	0.62	3.9	clear	yes	-113
1402	0.45	1.80	0.15	6.70	6.27	29.71	1.75	0.61	2.31	clear	yes	-114
<u>Neil Smith 1/31/08</u>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith TARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil Smith</u>	SAMPLING INITIATED AT: <u>1405</u>	SAMPLING ENDED AT: <u>1410</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~32</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>~100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
PZ-LSAS-1	3	CG	40ml	HCl	1 -	~2	RFP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: P2-LSAS-2	SAMPLE ID: P2-LSAS-2 DATE: 1.31.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 30 feet to 35 feet	STATIC DEPTH TO WATER (feet): 5.7	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + 0.0026 gallons/foot X 35 feet + 0.25 gallons = 0.34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33	PURGING INITIATED AT: 11:37	PURGING ENDED AT: 12:10	TOTAL VOLUME PURGED (gallons): 1.65

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
11:45	0.4	0.4	0.05	6.9	6.64	29.56	1.20*	0.66	35	clear	none	-124
11:50	0.25	0.65	0.05	6.75	6.67	29.53	1.25*	0.47	14	"	"	-129
11:55	0.25	0.90	0.05	6.71	6.6	29.59	1.27*	0.43	11	"	"	-129
12:00	0.25	1.15	0.05	6.72	6.7	29.32	3.79*	0.41	1	"	"	-129
12:05	0.25	1.40	0.05	6.7	6.71	29.39	19.6*	0.39	2	"	"	-124
12:10	0.25	1.65	0.05	6.7	6.70	29.4	99.9	0.15	0	"	"	-122

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Nicah Forbes ARCADES		SAMPLER(S) SIGNATURES: <i>[Signature]</i>		SAMPLING INITIATED AT: 12:11	SAMPLING ENDED AT: 12:15
PUMP OR TUBING DEPTH IN WELL (feet): 33		SAMPLE PUMP FLOW RATE (mL per minute): <100		TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>		FIELD-FILTERED: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>		DUPLICATE: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
P2-LSAS-2	3	CG	40mL	HCl	—	—	8260	RFPP

REMARKS: ***Conductivity readings very strange - may be malfunctioning**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC Facility</u>		SITE LOCATION: <u>Tallevast, FL</u>	
WELL NO: <u>PZ-LSAS-3</u>	SAMPLE ID: <u>PZ-LSAS-3</u>	DATE: <u>1-31-08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>30</u> feet to <u>35</u> feet	STATIC DEPTH TO WATER (feet): <u>6.5</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>35</u> feet) + <u>0.25</u> gallons = <u>0.34</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>32.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>32.5</u>	PURGING INITIATED AT: <u>12:36</u>	PURGING ENDED AT: <u>13:15</u>	TOTAL VOLUME PURGED (gallons): <u>2.00</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	GRP
12:45	0.5	0.5	0.05	6.8	6.48	29.61	1.28	0.67	10	clear	none	-137
12:50	0.25	0.75	0.05	6.8	6.34	29.8	1.37	0.46	9	"	"	-131
12:55	0.25	1.00	0.05	6.8	6.28	29.4	1.45	0.42	1	"	"	-128
13:00	0.25	1.25	0.05	6.8	6.32	29.4	2.05	0.41	0	"	"	-128
13:05	0.25	1.50	0.05	6.8	6.36	29.5	1.35	0.39	0	"	"	-126
13:10	0.25	1.75	0.05	6.8	6.37	29.7	1.35	0.38	0	"	"	-127
13:15	0.25	2.00	0.05	6.8	6.37	29.7	1.35	0.38	0	"	"	-127

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nicah Forbes / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>13:16</u>	SAMPLING ENDED AT: <u>13:20</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>32.5</u> ^{MBS 4/11/08}	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u> ^{MBS 4/11/08}	TUBING MATERIAL CODE: <u>PE</u> ^{MBS 4/11/08}	
FIELD DECONTAMINATION: <u>Y</u> ⁽¹⁾ N	FIELD-FILTERED: <u>Y</u> N	FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
PZ-LSAS-3	3	CG	ADM	Hel	—	—	8260	RFPP

REMARKS: * conductivity meter may be malfunctioning

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Tellevis</i>		SITE LOCATION: <i>Sarasota, FL</i>	
WELL NO: <i>PZ-LSAS-4</i>	SAMPLE ID: <i>PZ-LSAS-4</i>	DATE: <i>1/21/08</i>	

PURGING DATA

WELL DIAMETER (inches): <i>2</i>	TUBING DIAMETER (inches): <i>1/4</i>	WELL SCREEN INTERVAL DEPTH: <i>30.5</i> feet to <i>35.5</i> feet	STATIC DEPTH TO WATER (feet): <i>16.19</i>	PURGE PUMP TYPE OR BAILER: <i>PP</i> <i>MDS 7/11/08</i>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<i>0.0026</i> gallons/foot X <i>38</i> feet) + <i>0.25</i> gallons = <i>0.25</i> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <i>1150</i>	PURGING ENDED AT: <i>1211</i>	TOTAL VOLUME PURGED (gallons): <i>2.52</i>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<i>3 1153</i>	<i>0.36</i>	<i>0.36</i>	<i>0.12</i>	<i>7.37</i>	<i>6.36</i>	<i>30.02</i>	<i>1.49</i>	<i>0.86</i>	<i>41.1</i>	<i>clear</i>	<i>yes</i>	<i>-126</i>
<i>1156</i>	<i>0.36</i>	<i>0.72</i>	<i>0.12</i>	<i>7.45</i>	<i>6.33</i>	<i>28.88</i>	<i>1.64</i>	<i>0.73</i>	<i>42.2</i>	<i>clear</i>	<i>yes</i>	<i>-134</i>
<i>1159</i>	<i>0.36</i>	<i>1.08</i>	<i>0.12</i>	<i>7.46</i>	<i>6.33</i>	<i>29.81</i>	<i>1.68</i>	<i>0.66</i>	<i>26.8</i>	<i>clear</i>	<i>yes</i>	<i>-41</i>
<i>1202</i>	<i>0.36</i>	<i>1.44</i>	<i>0.12</i>	<i>7.46</i>	<i>6.33</i>	<i>29.86</i>	<i>1.68</i>	<i>0.63</i>	<i>22.0</i>	<i>clear</i>	<i>yes</i>	<i>-130</i>
<i>1205</i>	<i>0.36</i>	<i>1.80</i>	<i>0.12</i>	<i>7.46</i>	<i>6.34</i>	<i>29.73</i>	<i>1.65</i>	<i>0.60</i>	<i>19.4</i>	<i>clear</i>	<i>yes</i>	<i>-142</i>
<i>1208</i>	<i>0.36</i>	<i>2.16</i>	<i>0.12</i>	<i>7.46</i>	<i>6.35</i>	<i>29.73</i>	<i>1.64</i>	<i>0.58</i>	<i>16.7</i>	<i>clear</i>	<i>yes</i>	<i>-143</i>
<i>1211</i>	<i>0.36</i>	<i>2.52</i>	<i>0.12</i>	<i>7.46</i>	<i>6.35</i>	<i>29.77</i>	<i>1.63</i>	<i>0.58</i>	<i>13.4</i>	<i>clear</i>	<i>yes</i>	<i>-144</i>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Ned Smith / ARCADIS</i>	SAMPLER(S) SIGNATURES: <i>Ned Smith</i>	SAMPLING INITIATED AT: <i>1215</i>	SAMPLING ENDED AT: <i>1220</i>
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <i>±100</i>	TUBING MATERIAL CODE: <i>PE</i>	
FIELD DECONTAMINATION: <i>Ⓞ</i>	FIELD-FILTERED: <i>Y</i> / <i>N</i> FILTER SIZE: _____ µm	DUPLICATE: <i>Y</i> / <i>N</i>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>PZ-LSAS-4</i>	<i>3</i>	<i>CG</i>	<i>40mL</i>	<i>HCL</i>	<i>-</i>	<i>6.2</i>	<i>826B + 160S in RFP</i>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallahassee</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>PZ-LSAS-5</u>	SAMPLE ID: <u>PZ-LSAS-5</u> DATE: <u>1/31/02</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>27.87</u> feet to <u>32.87</u> feet	STATIC DEPTH TO WATER (feet): <u>6.27</u>	PURGE PUMP TYPE OR BAILER: <u>PP peristaltic pump</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable
 = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = gallons + 0.0012 gallons/foot X 35 feet + 0.25 gallons = 0.34 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~30</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~30</u>	PURGING INITIATED AT: <u>1305</u>	PURGING ENDED AT: <u>1320</u>	TOTAL VOLUME PURGED (gallons): <u>1.80</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1208	0.36	0.36	0.12	7.26	6.51	30.02	1.42	1.21	16.7	clear	Yes	-116
1211	0.36	0.72	0.12	7.26	6.48	30.16	1.58	1.04	13.7	clear	Yes	-120
1314	0.36	1.08	0.12	7.26	6.28	30.14	1.77	1.00	11.6	clear	Yes	-113
1317	0.36	1.44	0.12	7.29	6.19	30.15	1.82	0.97	18.0	clear	Yes	-111
1320	0.36	1.80	0.12	7.29	6.19	30.19	1.80	0.91	16.4	clear	Yes	-111
<u>NO DATA 1/31/02</u>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nick Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Nick Smith</u>	SAMPLING INITIATED AT: <u>1321</u>	SAMPLING ENDED AT: <u>1325</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~30</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>1100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>X</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> <u> </u> µm	DUPLICATE: <u>Y</u> <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
PZ-LSAS-5	3	CG	40ml	HCL	-	6.2	8266666666	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: PZ-LSAS-6	SAMPLE ID: PZ-LSAS-6 DATE: 1.31.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 3 feet to 35 feet	STATIC DEPTH TO WATER (feet): 7.3	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 35 feet + 0.25 gallons = 0.34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33	PURGING INITIATED AT: 13:53	PURGING ENDED AT: 14:20	TOTAL VOLUME PURGED (gallons): 1.35

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
14:00	0.35	0.35	0.05	8.05	6.60	28.13	1.42	2.86	3	clear	MR
14:05	0.60	0.60	0.05	8.05	6.50	27.9	1.44	1.60	22	"	"
14:10	0.25	0.85	0.05	8.30	6.51	28.04	1.42	1.78	30	"	"
14:15	0.25	1.10	0.05	8.30	6.51	27.92	1.45	1.70	28	"	"
14:20	0.25	1.35	0.05	8.30	6.51	27.90	1.47	1.69	25	"	"

ORP
-92
-95
-95
-95
-95

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 14:21	SAMPLING ENDED AT: 14:25
PUMP OR TUBING DEPTH IN WELL (feet): 33	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y	FIELD-FILTERED: Y FILTER SIZE: _____ µm	DUPLICATE: Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
PZ-LSAS-6	3	CG	100 mL	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2-mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellico	SITE LOCATION: Shawnee, FL
WELL NO: P2-LSAS-7	DATE: 1/31/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 30.75 set to 35.75 set	STATIC DEPTH TO WATER (feet): 6.59	PURGE PUMP TYPE OR BAILER: (PP) geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0022 gallons/foot X 38 feet) + 0.05 gallons = 0.34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~32.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~32.5	PURGING INITIATED AT: 1230	PURGING ENDED AT: 1242	TOTAL VOLUME PURGED (gallons): 156

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1233	0.39	0.39	0.13	7.29	5.86	27.53	0.98	4.26	12.2	clear	yes
1236	0.37	0.76	0.13	7.34	5.88	27.54	0.99	4.13	5.96	clear	yes
1239	0.39	1.17	0.13	7.34	5.90	27.51	1.00	4.06	4.78	clear	yes
1242	0.39	1.56	0.13	7.34	5.92	27.57	1.02	3.99	5.82	clear	yes
<i>well shut</i>											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / DEP			SAMPLER(S) SIGNATURES: <i>Neil Smith</i>			SAMPLING INITIATED AT: 1245		SAMPLING ENDED AT: 1250	
PUMP OR TUBING DEPTH IN WELL (feet): ~32.5			SAMPLE PUMP FLOW RATE (mL per minute): ~100			TUBING MATERIAL CODE: PE			
FIELD DECONTAMINATION: N			FIELD-FILTERED: Y FILTER SIZE: _____ µm			DUPLICATE: Y			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
P2-LSAS-7	3	CG	40mL	HCL	—	~2	8200ft 8260sm	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellerant</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>RW-1</u>	SAMPLE ID: <u>RW-1</u> DATE: <u>4/2/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>6</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL (MPS) DEPTH: <u>15</u> feet to <u>19</u> feet	STATIC DEPTH TO WATER (feet): <u>5.49</u>	PURGE PUMP TYPE OR BAILER: <u>APP gasper</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<u>0.0026</u> gallons/foot X <u>24</u> feet) + <u>0.25</u> gallons = <u>0.30</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>14</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>14</u>	PURGING INITIATED AT: <u>0850</u>	PURGING ENDED AT: <u>0903</u>	TOTAL VOLUME PURGED (gallons): <u>1.12</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle % or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0854	0.28	0.32	0.08	5.51	6.10	27.85	0.482	1.59	13.2	clear	NO	ORP
0857	0.24	0.56	0.08	5.52	6.06	28.11	0.482	1.39	11.9	clear	NO	76
0900	0.24	0.88	0.08	5.52	6.06	27.95	0.482	1.33	11.0	clear	NO	84
0903	0.24	1.12	0.08	5.52	6.05	28.10	0.483	1.30	9.6	clear	NO	89
<u>Well purged 2/1/08</u>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil Smith</u>	SAMPLING INITIATED AT: <u>0905</u>	SAMPLING ENDED AT: <u>0910</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>14</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>AW-1</u>	<u>3</u>	<u>CG</u>	<u>400ml</u>	<u>HCL</u>	<u>-</u>	<u>6.2</u>	<u>Salicylic Acid ~ RFPF</u>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevese FL
WELL NO: RW-2	SAMPLE ID: RW-2 DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 4	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 5.03	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 8:55	PURGING ENDED AT: 9:25	TOTAL VOLUME PURGED (gallons): 7.0

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
9:00	0.5	0.5	0.1	5.05	5.66	27.19	0.420	1.04	7.64	clear	none	131
9:05	1.0	1.0	0.1	5.05	5.68	27.58	0.415	0.80	3.02	"	"	119
9:10	1.5	1.5	0.1	5.05	5.70	28.10	0.413	0.63	2.37	"	"	106
9:15	2.0	2.0	0.1	5.05	5.69	27.87	0.415	0.58	0.64	"	"	86
9:20	2.5	2.5	0.1	5.05	5.69	27.83	0.414	0.55	1.95	"	"	76
9:25	3.0	3.0	0.1	5.05	5.69	27.92	0.411	0.54	1.31	"	"	70

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / Arcadis	SAMPLER(S) SIGNATURE(S): <i>[Signature]</i>	SAMPLING INITIATED AT: 9:26	SAMPLING ENDED AT: 9:30
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 4100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
RW-2	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC</u>	SITE LOCATION: <u>Talleast</u>
WELL NO: <u>MW-3</u>	SAMPLE ID: <u>MW-3</u>
DATE: <u>1-25-08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>23.5</u> feet to <u>25</u> feet	STATIC DEPTH TO WATER (feet): <u>4.4</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= 0 gallons + 60026 gallons/foot X 31.5 feet + 125 gallons = gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>28.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>26</u>	PURGING INITIATED AT: <u>13:19</u>	PURGING ENDED AT: <u>13:48</u>	TOTAL VOLUME PURGED (gallons): <u>28</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TD WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
13:24	1.5	1.5	0.1	4.1	7.27	24.49	.057	5.07	77.4	clear	none
13:27	1.3	1.8	0.1	4.09	7.33	24.03	.052	4.88	90	brown	none
13:30	1.3	1.1	0.1	4.10	7.42	23.99	.052	5.25	109.1	brown	none
13:33	1.3	1.4	0.1	4.11	7.47	23.89	.051	4.70	128.0	brown	none
13:36	1.3	1.7	0.1	4.10	7.53	23.46	.049	4.55	128.0	brown	none
13:39	1.3	2.0	0.1	4.11	7.53	23.65	.049	4.22	150.0	brown	none
13:42	1.3	2.3	0.1	4.10	7.38	23.25	.048	2.90	152.0	brown	none
13:45	1.3	2.5	0.1	4.10	7.32	23.26	.048	2.89	152.0	brown	none
13:48	1.3	2.8	0.1	4.10	7.35	23.25	.048	2.88	153.0	brown	none

95
91
86
82
81
83
100
101
102

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Darrin Johnson</u>	SAMPLER(S) SIGNATURES: <u>Darrin Johnson</u>	SAMPLING INITIATED AT: <u>13:49</u>	SAMPLING ENDED AT: <u>13:54</u>
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PUMP OR TUBING DEPTH IN WELL (feet): <u>28.5</u> <u>MDS 9/1/08</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>400</u> <u>MDS 9/1/08</u>	TUBING MATERIAL CODE: <u>PE</u>
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FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: Y <u>N</u> FILTER SIZE: µm	DUPLICATE: Y <u>N</u>
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SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-3	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: This well needs to be re developed

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallahassee, FL
WELL NO: MW-4	SAMPLE ID: MW-4 DATE: 125-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 4 feet to 19 feet	STATIC DEPTH TO WATER (feet): 4.07 / 5.23	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 19 feet) + 0.25 gallons = 0.25 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 12	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 12	PURGING INITIATED AT: 1310	PURGING ENDED AT: 1343	TOTAL VOLUME PURGED (gallons): 3.3

TIME	VOLUME PURGED (gallons)	CUMULATIVE VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1328	1.8	1.8	0.1	5.23	5.91	25.55	0.107	4.06	115	red brown	none noted
1331	0.3	2.1		5.32	5.64	25.50	0.129	2.33	116		
1334		2.4		5.43	5.68	25.67	0.133	2.38	119		
1337		2.7		5.88	6.06	25.76	0.149	2.46	122		
1340		3.0		5.33	6.09	25.73	0.151	2.32	119		
1343		3.3		5.33	6.08	25.72	0.154	2.32	116		

ORP

146
112
95
72
62
60

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amy Coak / Arcadis	SAMPLER(S) SIGNATURES: <i>Amy Coak</i>	SAMPLING INITIATED AT: 1344	SAMPLING ENDED AT: 1345
PUMP OR TUBING DEPTH IN WELL (feet): 12	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 0.45 µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-4	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallevase, FL	
WELL NO: MW-5	SAMPLE ID: MW-5	DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 5.5	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 10 feet + 0.25 gallons = 0.276 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 11:01	PURGING ENDED AT: 11:25	TOTAL VOLUME PURGED (gallons): 2.1
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
11:05	0.1	0.1	0.1	5.5	5.64	22.28	0.390	9.42	49	slt turb	none	-12
11:10	0.5	0.6	0.1	5.5	6.18	24.13	0.467	0.57	0	clear	none	-53
11:15	0.5	1.1	0.1	5.5	6.21	24.18	0.428	0.58	0	"	"	-57
11:20	0.5	1.6	0.1	5.5	6.22	24.11	0.417	0.61	0	"	"	-57
11:25	0.5	2.1	0.1	5.5	6.22	24.39	0.413	0.64	0	"	"	-57

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes (ARCADIS)	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 11:26	SAMPLING ENDED AT: 11:30
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 500 (MDS 4/11/08)	TUBING MATERIAL CODE: PE	(MDS 4/11/08)
FIELD DECONTAMINATION: <i>[Signature]</i>	FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/> FILTER SIZE: µm	DUPLICATE: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-5	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU, optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: TALLEVAST
WELL NO: mw-6	SAMPLE ID: mw6
DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/2	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 5.47	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 25 feet) + .25 gallons = 0.32 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 7	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 7	PURGING INITIATED AT: 1101	PURGING ENDED AT: 1114	TOTAL VOLUME PURGED (gallons): 2.73

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1106	1.05	1.05	0.21	5.60	6.53	23.9	272	0.71	1.36	clear	NONE
1108	.42	1.47	0.21	5.60	6.46	24.0	277	0.67	1.53	clear	NONE
1110	.42	1.89	0.21	5.60	6.40	24.0	280	0.65	1.38	clear	NONE
1112	.42	2.31	0.21	5.60	6.39	23.9	281	0.66	1.88	clear	NONE
1114	.42	2.73	0.21	5.60	6.38	24.0	283	0.65	0.32	clear	NONE

ORP
44
38
83
34
32

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1115	SAMPLING ENDED AT: 1118
PUMP OR TUBING DEPTH IN WELL (feet): 7	SAMPLE PUMP FLOW RATE (mL per minute): ~800	TUBING MATERIAL CODE: PE	MDX 3/11/08
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y Filtration Equipment Type: <input checked="" type="checkbox"/>	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-6	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Talleast
WELL NO: MW-75	SAMPLE ID: MW-75
DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (Inches): 2	TUBING DIAMETER (Inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 4.08	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + 1.0026 gallons/foot X 12 feet + 1.25 gallons = 1.29 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 10	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 8	PURGING INITIATED AT: 10:31	PURGING ENDED AT: 10:48	TOTAL VOLUME PURGED (gallons): 1.7

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
10:39	0.8	0.8	0.1	4.13	6.20	21.88	1.25	1.41	36.6	amber	none
10:42	0.3	1.1	0.1	4.15	6.23	21.87	1.29	1.39	30.6	clear	none
10:45	0.3	1.4	0.1	4.15	6.24	21.88	1.31	1.28	29.8	clear	none
10:48	0.3	1.7	0.1	4.15	6.25	21.98	1.31	1.15	19.8	clear	none

2
3
3
4

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: Darrin Johnson	SAMPLING INITIATED AT: 10:49	SAMPLING ENDED AT: 10:54
PUMP OR TUBING DEPTH IN WELL (feet): 8 (MDS 4/11/08)	SAMPLE PUMP FLOW RATE (mL per minute): <100 (MDS 4/11/08)	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-75	3	CG	40ml	HCl	—	—	8260	RFPP

MDS
4/11/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
 Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-7D	SAMPLE ID: MW-7D DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 3.85	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)

= **0** gallons + **1.0026** gallons/foot X **22** feet + **.25** gallons = **0.3** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 20	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 11:01	PURGING ENDED AT: 11:11	TOTAL VOLUME PURGED (gallons): 1.0 MDS 4/11/08
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
11:05	0.4	0.4	0.1	4.08	4.87	23.06	1.47	1.03	5.01	clear	none	297
11:08	0.3	0.7	0.1	4.09	4.88	23.32	1.47	0.91	3.14	↓	↓	302
11:11	0.3	1.0	0.1	4.08	4.90	23.61	1.46	0.88	3.02	↓	↓	289

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson		SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>		SAMPLING INITIATED AT: 11:12	SAMPLING ENDED AT: 11:17
PUMP OR TUBING DEPTH IN WELL (feet): 17.5 MDS 4/11/08	SAMPLE PUMP FLOW RATE (mL per minute): 100 MDS 4/11/08		TUBING MATERIAL CODE: PE		
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N		

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-7D	3	CG	40ml	HCl	—	—	8260	RFPP MDS 4/11/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallvest
WELL NO: MW-8S	SAMPLE ID: MW-8S
DATE: 1-25-08	

PURGING DATA

WELL DIAMETER (Inches): 2	TUBING DIAMETER (Inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 3.7	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (1.0026 gallons/foot X 12 feet) + .25 gallons = 0.28 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 10	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 8	PURGING INITIATED AT: 12:58	PURGING ENDED AT: 13:11	TOTAL VOLUME PURGED (gallons): 1.3

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
13:05	.7	.7	0.1	3.88	6.20	23.05	1.87	1.69	0.99	clear	none
13:08	.3	1.0	0.1	3.88	6.15	23.08	1.85	1.55	0.85	clear	none
13:11	.3	1.3	0.1	3.88	6.14	23.02	1.85	1.50	0.85	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: Darrin Johnson	SAMPLING INITIATED AT: 13:12	SAMPLING ENDED AT: 13:17
PUMP OR TUBING DEPTH IN WELL (feet): 8	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-8S	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallewas, FL
WELL NO: MW-8D	DATE: 2-1-08
SAMPLE ID: MW-8D	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 4.45	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 = (**20** feet - **4.45** feet) X **0.25** gallons/foot = **3.55** gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 = **0** gallons + (**0.0026** gallons/foot X **20** feet) + **0.25** gallons = **0.25** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 0837	PURGING ENDED AT: 0903	TOTAL VOLUME PURGED (gallons): 2.9
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle) (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0854	2.0	2.0	0.1	4.78	4.87	24.34	1.74	1.52	0.109	clear w/ sm. flecks	none
0857	0.3	2.3	0.1	4.82	4.87	24.25	2.00	1.12	1.18	sm. flecks	
0900	0.3	2.6	0.1				1.27				
0900	0.3	2.6	0.1	4.83	4.81	24.25	1.28	1.05	2.01		
0903	0.3	2.9	0.1	4.84	4.81	24.26	1.28	1.04	0.72		

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal/Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amy Cook / Arcadis	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 0905	SAMPLING ENDED AT: 0909
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N	FILTER SIZE: µm	DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-8D	3	CG	40mL	HCl			8260	RFPP

REMARKS: **0835** soil transducer accidentally **near** pump for minute before measure static dth. **Replace transducer 0908**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Flow rate in 1 hour faster to purge. total vol. removed ≈ 3.35 gall
 Replace tubing back down well 0907

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-9S	SAMPLE ID: MW-9S DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 4.27	PURGE PUMP, TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (10 feet - 4.27 feet) X 0.16 gallons/foot = 0.92 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 33 feet) + 0.25 gallons = 0.25 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 7	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 7	PURGING INITIATED AT: 1348	PURGING ENDED AT: 1402	TOTAL VOLUME PURGED (gallons): 1.82

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circled or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8 1356	1.04	1.04	0.13	4.36	6.27	25.1	256	0.52	3.58	clear	NONE
1358	0.26	1.30	0.13	4.36	6.25	25.2	254	0.50	2.26	clear	NONE
1400	0.26	1.56	0.13	4.36	6.27	25.2	261	0.51	1.43	clear	NONE
1402	0.26	1.82	0.13	4.36	6.27	25.3	262	0.51	0.74	clear	NONE

ORP
-37
-39
-44
-48

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STALL ARCADIS		SAMPLER(S) SIGNATURES: <i>Bridget Stall</i>		SAMPLING INITIATED AT: 1403	SAMPLING ENDED AT: 1408
PUMP OR TUBING DEPTH IN WELL (feet): 7		SAMPLE PUMP FLOW RATE (mL per minute): ~430		TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm		DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-9S	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: JALLEVAST	SITE LOCATION:
WELL NO: mw-9D	SAMPLE ID: mw-9D DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 4.00	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 33 feet) + .25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 1335	PURGING ENDED AT: 1354	TOTAL VOLUME PURGED (gallons):

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1345	1.10	1.10	0.11	4.45	4.9	26.1	0.184	6.4	1.72	clear	none
1347	0.22	1.32	0.11	4.45	4.9	26.1	0.183	7.0	2.32	clear	none
1349	0.22	1.54	0.11	4.45	4.9	26.1	0.201	6.4	1.60	clear	none
1351	0.22	1.76	0.11	4.45	4.8	26.1	0.200	7.0	2.18	clear	none

ORP

207
183
175
159

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1356	SAMPLING ENDED AT: 1400
PUMP OR TUBING DEPTH IN WELL (feet): 17.5	SAMPLE PUMP FLOW RATE (mL per minute): ~416	TUBING MATERIAL CODE: AG	MS 4/11/08 T
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-9D	3	CG	40mL	HCL	40mL	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.02

SITE NAME: Tallerast		SITE LOCATION: Tallerast, FL	
WELL NO: MW-10	SAMPLE ID: MW-10	DATE: 1/31/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 5.40	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 35 feet + 0.25 gallons = 34 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 1006	PURGING ENDED AT: 1019	TOTAL VOLUME PURGED (gallons): 1.95
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
7 1013	1.05	1.05	0.15	5.70	5.63	28.6	0.418	0.84	51.4	cloudy	none	131
1015	0.3	1.35	0.15	5.70	5.50	28.6	0.417	0.79	34.9	cloudy	none	132
1017	0.3	1.65	0.15	↓	5.59	28.7	0.421	0.74	15.1	cloudy	none	135
1019	0.3	1.95	0.15	↓	5.59	28.6	0.411	0.72	9.99	clear	none	136
										etc		

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1020	SAMPLING ENDED AT: 1030
PUMP OR TUBING DEPTH IN WELL (feet): 17.5	SAMPLE PUMP FLOW RATE (mL per minute): ~560	TUBING MATERIAL CODE: T	MPS 4/11/08
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	FILTER SIZE: _____ µm	DUPLICATE: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-10	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **sampled with transducer**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: MW-11	DATE: 1/21/08
SAMPLE ID: MW-11	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 5.29	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 20 feet) + 0.25 gallons = .302 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1030	PURGING ENDED AT: 1059	TOTAL VOLUME PURGED (gallons): 2.9
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1044	1.4	1.4	0.1	5.40	4.34	25.04	1.61	1.77	2.32	None	None noted	231
1047	0.3	1.7	↓	5.38	4.34	25.00	1.64	1.44	1.07	↓	↓	215
1050	↓	2.0	↓	5.38	4.34	25.08	1.67	0.95	0.94	↓	↓	213
1053	↓	2.3	↓	5.38	4.34	25.00	1.69	0.82	1.26	↓	↓	208
1056	↓	2.6	↓	5.38	4.34	24.90	1.69	0.81	1.18	↓	↓	207
1059	↓	2.9	↓	5.38	4.35	24.99	1.67	0.79	1.09	↓	↓	202
							MS/cm					

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amie Coats Arcadis	SAMPLER(S) SIGNATURES: <i>Amie Coats</i>	SAMPLING INITIATED AT: 1100	SAMPLING ENDED AT: 1105
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100 ^{MPS 4/11/08}	TUBING MATERIAL CODE: T ^{MPS 4/11/08}	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: _____ µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-11	3	CG	40mL	HCl	—	—	8260	LFPP ^{MPS 4/11/08}

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After-Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

MDS 9/11/08

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-12	SAMPLE ID: MW-12	DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL (feet): 15 to 20	STATIC DEPTH TO WATER (feet): 4.92	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (15 feet - 4.92 feet) X 0.0026 gallons/foot = 23.22 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 20 feet) + 0.25 gallons = 0.077 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 10:02	PURGING ENDED AT: 10:30	TOTAL VOLUME PURGED (gallons): 3

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	GRP
1005	0.5	0.5	0.1	5.05	4.88	28.90	0.349	1.17	56	turbid	none	42
1010	1.0	1.0	0.1	5.05	4.80	28.97	0.347	0.79	17	"	"	42
1015	1.5	1.5	0.1	5.05	4.74	28.96	0.351	0.58	7	"	"	35
1020	2.0	2.0	0.1	5.05	4.72	29.08	0.362	0.53	5	"	"	30
1025	2.5	2.5	0.1	5.05	4.67	29.27	0.384	0.49	4	"	"	23
1030	3.0	3.0	0.1	5.05	4.66	29.06	0.390	0.48	3	"	"	22

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1031	SAMPLING ENDED AT: 1035
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 500	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-12	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-13S	SAMPLE ID: MW-13S DATE: 1-29-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 4.70	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0260 gallons/foot X 10 feet) + 0.25 gallons = 0.28 gallons				

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1230	0.6	0.6		4.88	6.69	24.73	0.317	1.62	4.47	None	None	-85
1233		0.9		4.88	6.62	24.73	0.315	1.89	2.58			-112
1236		1.2		4.88	6.58	24.78	0.314	1.66	2.15			-122
1239				4.88	6.57	24.80	0.314	1.52	2.01			-126
1242				4.88	6.55	24.73	0.315	1.42	2.17			-13
1245				4.88	6.54	24.81	0.314	1.39	1.73			-134
1248				4.88	6.53	24.74	0.316	1.33	1.45			-136
1251				4.88	6.53	24.79	0.334	1.30	1.37			-140
1254				4.88	6.53	24.60	0.316	1.26	1.20			-145
1257					6.54	24.76	0.316	1.24	1.47			-144

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Anny Coats / Arcadis	SAMPLER(S) SIGNATURES: <i>Anny Coats</i>	SAMPLING INITIATED AT: 1302	SAMPLING ENDED AT: 1305
PUMP OR TUBING DEPTH IN WELL (feet): 7	SAMPLE/PUMP FLOW RATE (mL per minute): < 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-13S	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **1220 pull transducer** time **1300** → **6.54** pH **24.77** temp **0.318** cond **1.25** DO **1.50** turb

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

1314 Replace transducer

MW-13D

DEP-SOP-001/01
Form FD 9000-24

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility MDS 4/11/08 SITE LOCATION: Tallcrest
 WELL NO: ~~MW-14D~~ SAMPLE ID: ~~MW-14D~~ MW-13D DATE: 1-29-08

PURGING DATA

WELL DIAMETER (inches): 2 TUBING DIAMETER (inches): 1/4 WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet STATIC DEPTH TO WATER (feet): 4.91 PURGE PUMP TYPE OR BAILER: PP
 WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 = (15 - 4.91) feet X gallons/foot = gallons
 EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = 0 gallons + (1.0026 gallons/foot X 22 feet) + 1.25 gallons = 0.3 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 20 FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5 PURGING INITIATED AT: 12:25 PURGING ENDED AT: 12:38 TOTAL VOLUME PURGED (gallons): 1.3

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
12:29	0.4	0.4	0.1	5.3	4.82	26.09	0.069	2.57	2.40	clear	none
12:32	0.3	0.7	0.1	5.18	4.76	26.26	0.070	1.76	1.81	clear	none
12:35	0.3	1.0	0.1	5.18	4.71	26.33	0.071	1.63	1.78	clear	none
12:38	0.3	1.3	0.1	5.18	4.72	26.36	0.071	1.59	1.62	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson SAMPLER(S) SIGNATURES: [Signature] SAMPLING INITIATED AT: 12:39 SAMPLING ENDED AT: 12:44
 PUMP OR TUBING DEPTH IN WELL (feet): 17.5 MDS 4/11/08 SAMPLE PUMP FLOW RATE (mL per minute): 2100 MDS 4/11/08 TUBING MATERIAL CODE: PE
 FIELD DECONTAMINATION: Y N FIELD-FILTERED: Y N FILTER SIZE: _____ µm DUPLICATE: Y N
 Filtration Equipment Type: _____

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-13D	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-14S	SAMPLE ID: MW-14S DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 386	PURGE PUMP TYPE OR BAILER: geopump (PP)								
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (10.0 feet - 3-86 feet) X 0.16 gallons/foot = 0.98 gallons												
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 13 feet) + 0.25 gallons = 0.28 gallons												
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 7	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 7	PURGING INITIATED AT: 0949	PURGING ENDED AT: 1007	TOTAL VOLUME PURGED (gallons): 2.16								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm) (off)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	SRP
9 0958	1.08	1.08	0.12	3.99	6.5	23.8	0.222	1.1	18.4	clear	NONE	-32
3 1001	0.36	1.44	0.12	3.99	6.5	23.8	0.220	1.0	18.4	clear	NONE	-43
3 1004	0.36	1.80	0.12	3.99	6.5	23.9	0.217	0.9	14.3	clear	NONE	-59
3 1007	0.36	2.16	0.12	3.99	6.5	24.0	0.216	0.8	13.9	clear	NONE	-70
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016												

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STALL/ARGAIS			SAMPLER(S) SIGNATURES: Bridget Stall			SAMPLING INITIATED AT: 1008		SAMPLING ENDED AT: 1012	
PUMP OR TUBING DEPTH IN WELL (feet): 7			SAMPLE PUMP FLOW RATE (mL per minute): ~440			TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: Y (N)			FIELD-FILTERED: Y (N) FILTER SIZE: _____ µm			DUPLICATE: Y (N)			
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH			
MW-14S	3	CG	40ml	HCl	-	-	8260	RFPP	
REMARKS:									
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)									
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump									
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)									

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 7)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-14D	SAMPLE ID: MW-14D DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 3.96	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 33 feet) + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 0950	PURGING ENDED AT: 1004	TOTAL VOLUME PURGED (gallons): 1.82

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0958	1.04	1.04	0.13	4.13	5.24	26.1	0.087	0.66	4.13	clear	NONE	56
1000	0.26	1.30	0.13	4.14	5.21	26.2	0.087	0.63	3.07	clear	NONE	51
1002	0.26	1.56	0.13	4.14	5.20	26.2	0.087	0.60	2.20	clear	NONE	44
1004	0.26	1.82	0.13	4.14	5.18	26.2	0.087	0.58	2.07	clear	NONE	42

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1007	SAMPLING ENDED AT: 1009
PUMP OR TUBING DEPTH IN WELL (feet): 17.5	SAMPLE PUMP FLOW RATE (mL per minute): 1492	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-14D	3	CG	40mL	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALKEVAST	SITE LOCATION: Sarasota FL
WELL NO: MW-155	SAMPLE ID: MW-155
DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 4.36	PURGE PUMP TYPE OR BAILER: Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (10 feet - 4.36 feet) X 0.16 gallons/foot = 0.90 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 13 feet) + 1.25 gallons = 0.28 gallons				
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INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 7	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 7	PURGING INITIATED AT: 0853	PURGING ENDED AT: 0911	TOTAL VOLUME PURGED (gallons): 1.98
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9 0902	0.99	0.99	0.11	4.50	6.52	23.6	1.394	0.71	2.03	Clear	NONE
3 0905	0.33	1.32	0.11	4.51	6.54	23.6	0.401	0.70	1.59	Clear	NO
3 0908	0.33	1.65	0.11	4.51	6.54	23.6	0.401	0.70	3.02	Clear	NO
3 0911	0.33	1.98	0.11	4.51	6.55	23.6	0.410	0.70	2.01	Clear	NO

ORP
-74
-93
-97
-105

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 0912	SAMPLING ENDED AT: 0918
PUMP OR TUBING DEPTH IN WELL (feet): 7	SAMPLE PUMP FLOW RATE (mL per minute): ~416	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-155	3	CG	40mL	HCL	40mL	-	P260	RPPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump

EQUIPMENT CODES: RPPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: TALLEVAST, FL.
WELL NO: mw-15D	SAMPLE ID: mw-15D DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 449	PURGE PUMP TYPE OR BAILER: Geo Pump
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable) = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = **0** gallons + (**0.0026** gallons/foot X **343** feet) + **.25** gallons = **0.91** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 0848	PURGING ENDED AT: 0904	TOTAL VOLUME PURGED (gallons): 1.76
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0858	1.10	1.10	0.11	5.19	5.0	24.2	ms/cm 0.213	0.6	2.19	clear	NO
0900	0.22	1.32	0.11	5.19	5.0	24.3	0.211	0.6	1.24	clear	NO
0902	0.22	1.54	0.11	5.19	5.0	24.3	0.211	0.6	0.67	clear	NO
0904	0.22	1.76	0.11	5.19	5.0	24.3	0.211	0.6	0.80	clear	NO

ORP
74
55
45
35

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 906	SAMPLING ENDED AT: 0910
PUMP OR TUBING DEPTH IN WELL (feet): 17.5 ^{MDS} 4711/02	SAMPLE PUMP FLOW RATE (mL per minute): 416	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-15D	3	CG	40mL	HCL	40mL	-	8260	RFPP

REMARKS: **1.08**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-16S	SAMPLE ID: MW-16S	DATE: 1.29.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 7 feet 0 feet	STATIC DEPTH TO WATER (feet): 2.9	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (7 feet - 2.9 feet) X 0.0026 gallons/foot = 0.026 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 10 feet) + 0.25 gallons = 0.276 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 8	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 8	PURGING INITIATED AT: 15:15	PURGING ENDED AT: 16:00	TOTAL VOLUME PURGED (gallons): 4.4

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
15:20	0.4	0.4	0.1	2.95	5.78	21.14	0.226	0.32	0	Clear	None	-41
15:25	0.5	0.9	0.1	2.99	5.20	21.18	0.276	0.28	0	"	"	-71
15:30	0.5	1.4	0.1	3.00	5.16	21.19	0.340	0.29	0	"	"	-72
15:35	0.5	1.9	0.1	3.00	5.06	21.18	0.936	0.28	0	"	"	-74
15:40	0.5	2.4	0.1	3.00	5.06	21.20	0.999	0.31	0	"	"	-72
15:45	0.5	2.9	0.1	3.00	5.00	21.24	0.589	0.41	0	"	"	-77
15:50	0.5	3.4	0.1	3.00	4.98	21.25	0.242	0.41	0	"	"	-76
15:55	0.5	3.9	0.1	3.00	4.98	21.25	0.213	0.41	0	"	"	-77
16:00	0.5	4.4	0.1	3.00	4.98	21.20	0.227	0.41	0	"	"	-78

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 16:01	SAMPLING ENDED AT: 16:05
PUMP OR TUBING DEPTH IN WELL (feet): 8	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	FILTER SIZE: 0.45 µm	DUPLICATE: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-16S	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellems	SITE LOCATION: Sarasota, FL
WELL NO: MW-16D	SAMPLE ID: MW-16D DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: (pp)geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = 10.0026 ²⁵ feet - feet X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (0.0026 gallons/foot X 25 feet) + 0.25 gallons = 0.215 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~17.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~17.5	PURGING INITIATED AT: 1513	PURGING ENDED AT: 1527	TOTAL VOLUME PURGED (gallons): 1.26

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle % or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1518	0.45	0.45	0.09	3.45	5.47	22.99	374	0.86	34.2	clear	yes	60
1521	0.27	0.72	0.09	3.45	5.45	22.97	375	0.85	35.6	clear	yes	59
1524	0.27	0.99	0.09	3.45	5.41	23.00	377	0.82	37.2	clear	yes	37
1527	0.27	1.26	0.09	3.45	5.40	23.02	377	0.82	35.6	clear	yes	33
<i>Not back 1/29/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Nail Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Nail Smith</i>	SAMPLING INITIATED AT: 1530	SAMPLING ENDED AT: 1535
PUMP OR TUBING DEPTH IN WELL (feet): ~17.5	SAMPLE PUMP FLOW RATE (mL per minute): 6100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-16D	3	CG	40ml	HCL	—	7.2	ppbO3 + ppbO2 in RFP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

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Form FD 9000-24 GROUNDWATER SAMPLING LOG

SITE NAME: TAIlevast		SITE LOCATION: Sarasota FL		MDS 4/11/08	
WELL NO: mw-17S		SAMPLE ID: mw-17S		DATE: 1-23-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 4.41	PURGE PUMP TYPE OR BAILER: PP 600 Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 13 feet) + 0.25 gallons = 0.28 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 7	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 7	PURGING INITIATED AT: 1437	PURGING ENDED AT: 1449	TOTAL VOLUME PURGED (gallons): 1.92
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1443	0.96	0.96	0.16	4.58	6.34	23.7	0.334	1.23	3.99	clear	ylb
1445	0.32	1.28	0.16	4.58	6.34	23.7	0.319	1.14	5.65	clear	ylb
1447	0.32	1.6	0.16	4.58	6.34	23.8	0.314	1.03	6.38	clear	ylb
1449	0.32	1.92	0.16	4.58	6.34	23.8	0.302	0.98	6.61	clear	ylb

ORP
-45
-57
-68
-73

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCAAS		SAMPLER(S) SIGNATURES: Jennifer King		SAMPLING INITIATED AT: 1451	SAMPLING ENDED AT: 1454
PUMP OR TUBING DEPTH IN WELL (feet): 7 MDS 4/11/08		SAMPLE PUMP FLOW RATE (mL per minute): 600		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm		DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-17S	3	CG	40mL	HCL	40mL	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: Sarasota FL MDC 4/11/08
WELL NO: mw-170	SAMPLE ID: mw-170 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 23 feet) + .25 gallons = 0.31 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 1434	PURGING ENDED AT: 1448	TOTAL VOLUME PURGED (gallons): 1.68

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm³)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8 1442	0.96	0.96	0.12	4.90	4.9	25.4	.236	0.6	4.13	clear	yes
2 1444	0.24	1.2	0.12	4.92	4.8	25.4	.228	0.5	3.55	clear	yes
2 1446	0.24	1.44	0.12	4.92	4.8	25.4	.227	0.5	3.12	clear	yes
2 1448	0.24	1.68	0.12	4.92	4.8	25.4	.225	0.5	2.77	clear	yes

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1449	SAMPLING ENDED AT: 1455
PUMP OR TUBING DEPTH IN WELL (feet): 17.5	SAMPLE PUMP FLOW RATE (mL per minute): ~454	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: Y <input checked="" type="radio"/> N <input type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N <input type="radio"/> Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-170	3	CG	40mL	HEL	40mL	-	8260	RF PP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-18S	SAMPLE ID: MW-18S DATE: 1-28-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 4 feet to 10 feet	STATIC DEPTH TO WATER (feet): 3.09	PURGE PUMP TYPE OR BAILER: PP MDS 4/11/08
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (10 feet - 3.09 feet) X 0.16 gallons/foot = 1.11 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 7 MDS 4/11/08	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 7 MDS 4/11/08	PURGING INITIATED AT: 1022	PURGING ENDED AT: 1045	TOTAL VOLUME PURGED (gallons): 2.3
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1036	1.4	1.4	0.1	3.24	5.16	22.19	0.091	1.49	33.6	NONE	None noted
1039	0.3	1.7	↓	3.24	5.16	21.95	0.088	1.52	29.4	↓	↓
1042	0.3	2.0	↓	3.24	5.14	22.03	0.091	1.54	30.0	↓	↓
1045	0.3	2.3	↓	3.24	5.13	22.15	0.089	1.50	30.6	↓	↓
µS/cm											

36
33
29
23

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amc Coats Acadis	SAMPLER(S) SIGNATURES: <i>Amc Coats</i>	SAMPLING INITIATED AT: 1046	SAMPLING ENDED AT: 1050
PUMP OR TUBING DEPTH IN WELL (feet): 7'	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/11/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 1 µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-18S	3	CG	40mL	HCl	—	—	8260	RFPP

MDS 4/11/08

REMARKS: **1022 - remove transducer 1022 basin pump 1056 finish pumping 1058 replace**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

total volume removed: 11 L (graduated bucket)



GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Talleast
WELL NO: MW-18D	SAMPLE ID: MW-18D DATE: 1-28-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 3.14	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (20 feet - 3.14 feet) X .16 gallons/foot = 2.69 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 10:12	PURGING ENDED AT: 10:32	TOTAL VOLUME PURGED (gallons): 2.0					
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1023	1.1	1.1	0.1	3.39	4.98	23.93	0.163	6.21	2.35	clear	none
1026	.3	1.4	0.1	↓	4.98	24.04	0.166	6.15	2.54	↓	↓
1029	.3	1.7	0.1	↓	4.98	23.90	0.169	6.99	2.86	↓	↓
1032	.3	2.0	0.1	↓	4.98	24.03	0.172	6.04	1.73	↓	↓

-66
-70
-74
-78

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 1033	SAMPLING ENDED AT: 1035
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100 (MDS 2/1/08)	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-18D	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

3.09

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST	
WELL NO: mw-19	SAMPLE ID: mw-19	DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH (feet): 27.5 to 29.5	STATIC DEPTH TO WATER (feet): 19.23	PURGE PUMP TYPE OR BAILER: PP Bess Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 300.5 feet) + .25 gallons = 1.03 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 287.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 287.5	PURGING INITIATED AT: 1520	PURGING ENDED AT: 1634	TOTAL VOLUME PURGED (gallons): 3.36

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1628	3.12	3.12	.04	22.54	7.9	24.4	.566	0.4	35.1	clear	yo
1630	.08	3.20	.04	22.57	7.9	24.4	.564	0.4	37.3	clear	yo
1632	.08	3.28	.04	22.59	7.9	24.4	.562	0.4	28.9	clear	yo
1634	.08	3.36	.04	22.60	7.9	24.6	.561	0.4	26.5	clear	yo

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1636	SAMPLING ENDED AT: 1643
PUMP OR TUBING DEPTH IN WELL (feet): 287.5	SAMPLE PUMP FLOW RATE (mL per minute): ~150	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: 0 N	FIELD-FILTERED: Y Filtration Equipment Type: 0	FILTER SIZE: _____ µm	DUPLICATE: Y 0

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-19	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS: **Sampled with transducer missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: Tallevast, FL
WELL NO: mw-20	SAMPLE ID: mw-20 DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35 feet to 40 feet	STATIC DEPTH TO WATER (feet): 5.22	PURGE PUMP TYPE OR BAILER: Geo. Pump
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable) = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = **0** gallons + (**0.0026** gallons/foot X **43** feet) + **1.25** gallons = **.36** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 37.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 37.5	PURGING INITIATED AT: 1147	PURGING ENDED AT: 1203	TOTAL VOLUME PURGED (gallons): 1.76
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1157	1.10	1.10	0.11	5.59	5.40	24.2	.223	0.64	0.41	Clear	yes
1159	.22	1.32	0.11	5.59	5.40	24.2	.223	0.58	3.53	clear	yes
1201	.22	1.54	0.11	5.59	5.42	24.2	.228	0.60	1.64	Clear	yes
1203	.22	1.76	0.11	5.59	5.42	24.2	.231	0.55	2.12	Clear	yes

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAAL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Staal</i>	SAMPLING INITIATED AT: 1205	SAMPLING ENDED AT: 1211
PUMP OR TUBING DEPTH IN WELL (feet): 37.5	SAMPLE PUMP FLOW RATE (mL per minute): ~420	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-20	3	CG	40mL	HCL	40mL	—	8260	RPPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RPPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.89

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST, RI	
WELL NO: mw-21	SAMPLE ID: mw-21	DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 135 feet to 145 feet	STATIC DEPTH TO WATER (feet): 178	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 148 feet) + .25 gallons = 0.63 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 140	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 140	PURGING INITIATED AT: 1051	PURGING ENDED AT: 1113	TOTAL VOLUME PURGED (gallons): 2.73
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
15 1106	1.95	1.95	0.13	21.10	8.93	24.5	0.488	0.51	3.37	clear	yes
1108	0.24	2.21	0.13	21.27	8.89	24.5	0.493	0.50	2.90	clear	yes
1110	0.26	2.47	0.13	21.39	8.83	24.5	0.496	0.50	2.46	clear	yes
1112	0.26	2.73	0.13	21.49	8.77	24.5	0.497	0.49	2.59	clear	yes

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1115	SAMPLING ENDED AT: 1127
PUMP OR TUBING DEPTH IN WELL (feet): 140	SAMPLE PUMP FLOW RATE (mL per minute): 490	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-21	3	CG	40ml	HCl	40 ml	-	8260	RF PP
Dup-14	3	CG	40ml	HCl	40 ml	-	8260	RF PP

REMARKS: **Well NOT Labeled with RED TAG / F labeled concrete PAD with Paint Pen.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-22	SAMPLE ID: MW-22 DATE: 1/25/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 277 feet to 277 feet	STATIC DEPTH TO WATER (feet): 16.85	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 330 feet + 0.25 gallons = 1.11 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 287	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 287	PURGING INITIATED AT: 0915	PURGING ENDED AT: 1112	TOTAL VOLUME PURGED (gallons): 3.51

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or M/S/cm)	DISSOLVED OXYGEN (circ mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	CRP
1106	3.33	3.33	0.03	27.12	9.5	23.3	1709	2.5	132	cloudy	yes	-322
1108	.06	3.39	0.03	27.16	9.4	23.2	1717	2.6	140	cloudy	yes	-308
1110	.06	3.45	0.03	27.18	9.4	23.1	1727	2.5	142	cloudy	yes	-290
1112	.06	3.51	0.03	27.19	9.3	23.0	1740	2.5	145	cloudy	yes	-304

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1114	SAMPLING ENDED AT: 1121
PUMP OR TUBING DEPTH IN WELL (feet): 287	SAMPLE PUMP FLOW RATE (mL per minute): ~114	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-22	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **tubing needs changing; multiple kinks causing water blockage,**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

234

SITE NAME: Tallevast		SITE LOCATION: Tallevast, FL	
WELL NO: MW-23	SAMPLE ID: MW-23	DATE: 1/25/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 152 feet to 172 feet	STATIC DEPTH TO WATER (feet): 16.54	PURGE PUMP TYPE OR BAILER: PP	MDS 4/11/02
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable					
= (feet - feet) X gallons/foot = gallons					

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)					
= 0 gallons + (0.0026 gallons/foot X 205 feet) + 0.25 gallons = 0.78 gallons					

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 162	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 162	PURGING INITIATED AT: 1004	PURGING ENDED AT: 1030	TOTAL VOLUME PURGED (gallons): 2.58
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78

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1022	2.34	2.34	0.03	18.83	7.33	23.5	511	.47	1.65	cloudy	yes	-258
1024	0.06	2.40	0.03	18.79	7.32	23.5	512	.54	2.63	clear	yes	-257
1026	0.06	2.46	0.03	18.79	7.41	23.5	511	.49	1.62	clear	yes	-263
1028	0.06	2.52	0.03	18.79	7.42	23.3	511	.49	1.21	clear	yes	-263
1030	0.06	2.58	0.03	18.79	7.42	23.4	511	.47	1.14	clear	yes	-264

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Bridget Stahl / Archon	SAMPLER(S) SIGNATURES: <i>[Signatures]</i>	SAMPLING INITIATED AT: 1032	SAMPLING ENDED AT: 1041
PUMP OR TUBING DEPTH IN WELL (feet): 162	SAMPLE PUMP FLOW RATE (mL per minute): 114	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> <i>[initials]</i>

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-23	3	CG	40ml	HCl	—	—	8260	RFPP
DWP-11	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **Tubing should be changed; lots of leaks - lots of sediment in well; larger tubing recommended.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.05

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST, FL	
WELL NO: mw-24	SAMPLE ID: mw-24	DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 30.5 feet to 35.5 feet	STATIC DEPTH TO WATER (feet): 5.32	PURGE PUMP TYPE OR BAILER: PP 6ro Pump								
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable												
= (feet - feet) X gallons/foot = gallons												
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)												
= 0 gallons + (0.0026 gallons/foot X 38.5 feet) + .25 gallons = 0.35 gallons												
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33	PURGING INITIATED AT: 0859	PURGING ENDED AT: 0912	TOTAL VOLUME PURGED (gallons): 2.21								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle right) or % saturation	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
0906	1.19	1.19	0.17	6.04	5.4	24.2	1450	0.5	6.30	clear	yes	-11
0908	0.34	1.53	0.17	6.04	5.4	24.2	1447	0.5	4.85	clear	yes	-14
0910	0.34	1.87	0.17	6.04	5.4	24.2	1442	0.5	2.81	clear	yes	-20
0912	0.34	2.21	0.17	6.04	5.4	24.2	1439	0.5	1.93	clear	yes	-24
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016												

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BUDGET STAHL/ARCADIS			SAMPLER(S) SIGNATURES: <i>Budget Stahl</i>			SAMPLING INITIATED AT: 0913		SAMPLING ENDED AT: 0916		
PUMP OR TUBING DEPTH IN WELL (feet): 33			SAMPLE PUMP FLOW RATE (mL per minute): 2640			TUBING MATERIAL CODE:				
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N			FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm			DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N				
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH				
mw-24	3	CG	40mL	HCl	40mL	-	8260		RFPP	
REMARKS:										
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)										
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)										

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALKEVAST	SITE LOCATION: Sarasota FL	MDS 4/11/08
WELL NO: MW-25	SAMPLE ID: MW-25	DATE: 4-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 36.4 feet to 43.4 feet	STATIC DEPTH TO WATER (feet): 3.80	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 46.4 feet) + .25 gallons = 0.37 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 40	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 40	PURGING INITIATED AT: 1528	PURGING ENDED AT: 1543	TOTAL VOLUME PURGED (gallons): 1.95
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1537	1.17	1.17	0.13	4.88	5.35	26.4	.689	0.53	3.51	CLEAR	NONE
1539	0.26	1.43	0.13	4.84	5.32	26.3	.694	0.51	3.48	CLEAR	NONE
1541	0.26	1.69	0.13	4.81	5.31	26.4	.696	0.49	1.90	CLEAR	NONE
1543	0.26	1.95	0.13	4.81	5.31	26.4	.694	0.49	1.78	CLEAR	NONE

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / AREA DIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1545	SAMPLING ENDED AT: 1550
PUMP OR TUBING DEPTH IN WELL (feet): 40	SAMPLE PUMP FLOW RATE (mL per minute): 492	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: Y <input checked="" type="radio"/> N <input type="radio"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N <input type="radio"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-25	3	CG	40mL	HCL	40mL	-	8260	REPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/14/08	SITE LOCATION: Sarasota FL MDS 4/14/08
WELL NO: MW-26	SAMPLE ID: MW-26 DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 21.5 feet to 26.5 feet	STATIC DEPTH TO WATER (feet): 3.77	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)	= 0 gallons + (0.0026 gallons/foot X 26.5 feet) + 0.25 gallons = 1.72 gallons
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INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 23.5 MDS 4/14/08	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 23.5 MDS 4/14/08	PURGING INITIATED AT: 1440	PURGING ENDED AT: 1457	TOTAL VOLUME PURGED (gallons): 1.7 MDS 4/14/08
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1445	0.5	0.5	0.1	4.35	6.8	23.2	63	5.3	4.18	Clear	SuRv
1448	0.3	0.8	↓	4.35	6.7	23.1	63	3.8	2.85		
1451	0.3	1.1		4.38	6.7	23.3	63	3.6	2.73		
1454	0.3	1.4		4.42	6.7	23.2	63	3.4	2.73		
1457	0.3	1.7		4.45	6.7	23.1	63	3.6	2.70		
1500	Sampled @ 1500										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: K. ABCY	SAMPLER(S) SIGNATURES: K. ABCY / ARCADIS	SAMPLING INITIATED AT: 1500	SAMPLING ENDED AT: 1503
PUMP OR TUBING DEPTH IN WELL (feet): 23.5	SAMPLE PUMP FLOW RATE (mL per minute): 1100 MDS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 1 µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-26	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **Initial purge clear, noticeable sulfur odor.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tellevast	SITE LOCATION: Sarasota, FL
WELL NO: MW-27	SAMPLE ID: MW-27
DATE: 1/28/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 30 feet to 35 feet	STATIC DEPTH TO WATER (feet): 3.05	PURGE PUMP TYPE OR BAILER: (pp) geo pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 38 feet) + 0.50 gallons = (0.60) gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~32.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~32.5	PURGING INITIATED AT: 1000	PURGING ENDED AT: 1030	TOTAL VOLUME PURGED (gallons): 3.00
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle method or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1006	0.60	0.60	0.10	3.50	6.22	21.82	681	6.75	3.79	clear	NO	ORP 290
1009	0.30	0.90	0.10	3.54	6.25	22.10	674	6.38	2.44	clear	NO	287
1012	0.30	1.20	0.10	3.54	6.25	22.24	675	6.13	2.33	clear	NO	285
1015	0.30	1.50	0.10	3.54	6.24	22.40	683	4.56	2.32	clear	NO	280
1018	0.30	1.80	0.10	3.54	6.23	22.51	684	4.32	1.89	clear	NO	133
1021	0.30	2.10	0.10	3.54	6.23	22.44	683	3.24	1.89	clear	NO	102
1024	0.30	2.40	0.10	3.54	6.25	22.38	641	3.14	2.28	clear	NO	11
1027	0.30	2.70	0.10	3.54	6.29	22.45	637	3.16	1.80	clear	NO	-40
1030	0.30	3.00	0.10	3.54	6.35	22.41	629	3.12	1.80	clear	NO	-85

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1033	SAMPLING ENDED AT: 1040
PUMP OR TUBING DEPTH IN WELL (feet): ~32.5	SAMPLE PUMP FLOW RATE (mL per minute): ~100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y Filtration Equipment Type: X	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> X

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-27	3	CG	40ml	HCL	—	~2	8260B+8260SK	RFPP

REMARKS: **ToC Not flush with manhole cover**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

9000

DEP-SOP-001/01
Form FD 9000-24

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellevast</u>		SITE LOCATION: <u>Sarasota, FL</u>	
WELL NO: <u>MW-28</u>	SAMPLE ID: <u>MW-28</u>	DATE: <u>1/28/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>25</u> feet to <u>30</u> feet	STATIC DEPTH TO WATER (feet): <u>3.13</u>	PURGE PUMP TYPE OR BAILER: <u>(PP) gas pump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (<u> </u> feet - <u> </u> feet) X <u> </u> gallons/foot = <u> </u> gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = <u> </u> gallons + (<u>0.0026</u> gallons/foot X <u>36</u> feet) + <u>0.50</u> gallons = <u>0.59</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>228</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>228</u>	PURGING INITIATED AT: <u>1225</u>	PURGING ENDED AT: <u>1242</u>	TOTAL VOLUME PURGED (gallons): <u>1.36</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>8 1233</u>	<u>0.64</u>	<u>0.64</u>	<u>0.08</u>	<u>3.74</u>	<u>5.12</u>	<u>24.66</u>	<u>424</u>	<u>0.59</u>	<u>26.8</u>	<u>clear</u>	<u>yes</u>
<u>1236</u>	<u>0.24</u>	<u>0.88</u>	<u>0.08</u>	<u>3.78</u>	<u>5.18</u>	<u>24.65</u>	<u>449</u>	<u>0.47</u>	<u>7.62</u>	<u>clear</u>	<u>yes</u>
<u>1239</u>	<u>0.24</u>	<u>1.12</u>	<u>0.08</u>	<u>3.78</u>	<u>5.17</u>	<u>24.85</u>	<u>451</u>	<u>0.42</u>	<u>5.45</u>	<u>clear</u>	<u>yes</u>
<u>1242</u>	<u>0.24</u>	<u>1.36</u>	<u>0.08</u>	<u>3.81</u>	<u>5.21</u>	<u>24.86</u>	<u>452</u>	<u>0.45</u>	<u>2.28</u>	<u>clear</u>	<u>yes</u>

ORP
29
22
21
15

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil J. J. / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil J. J.</u>	SAMPLING INITIATED AT: <u>1242</u>	SAMPLING ENDED AT: <u>1255</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>228</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2.6</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N Filtration Equipment Type: <u> </u>	FILTER SIZE: <u> </u> µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-28</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u> </u>	<u>2.2</u>	<u>8260B+(260)11- RFPP</u>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)



GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-29	SAMPLE ID: MW-29 DATE: 1-28-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 25 feet to 30 feet	STATIC DEPTH TO WATER (feet): 2.99	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (30 feet - 2.99 feet) X .16 gallons/foot = 4.32 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (.0026 gallons/foot X 32 feet) + .25 gallons = .33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 30	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 27	PURGING INITIATED AT: 10:45	PURGING ENDED AT: 11:10	TOTAL VOLUME PURGED (gallons): 2.5

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
10:49	.4	.4	0.1	3.88	5.41	24.63	0.434	3.89	9.89	clear	none	66
10:52	.3	.7	0.1	3.94	5.44	24.80	0.448	3.22	10.42			
10:55	.3	1.0	0.1	4.0	5.53	24.75	0.654	2.13	10.53			5
10:58	.3	1.3	0.1	4.02	5.78	24.88	0.463	1.37	4.50			58
11:01	.3	1.6	0.1	4.04	5.96	24.90	0.505	0.91	3.98			94
11:04	.3	1.9	0.1	3.93	6.04	25.06	0.528	0.91	2.55			
11:07	.3	2.2	0.1	3.93	6.07	25.06	0.535	0.91	2.03			113
11:10	.3	2.5	0.1	3.94	6.10	25.19	0.540	0.90	2.06	↓	↓	116

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Dorin Johnson	SAMPLER(S) SIGNATURES: <i>Dorin Johnson</i>	SAMPLING INITIATED AT: 11:11	SAMPLING ENDED AT: 11:16
PUMP OR TUBING DEPTH IN WELL (feet): 27	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: pe	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-29	3	CG	40mL	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Seasota, FL
WELL NO: MW-30	SAMPLE ID: MW-30
DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 25.5 feet to 28.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0024 gallons/foot X 31.5 feet) + 0.5 gallons 0.58 gallons 0.13 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26	PURGING INITIATED AT: 1548	PURGING ENDED AT: 1612	TOTAL VOLUME PURGED (gallons): 2.9

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmho/5cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1600	1.2	1.2	0.1	8.78	6.74	25.64	509	0.73	10.65	clear	yes
1603	0.3	1.5	0.1	9.84	6.75	25.57	475	0.72	4.69	Yellow	yes
1606	0.3	1.8	0.1	10.40	6.74	25.59	472	0.59	4.24	"	yes
1609	0.3	2.1	0.1	10.85	6.73	25.60	468	0.50	4.22	"	yes
1612	0.3	2.4	0.1	11.5	6.72	25.66	474	0.47	4.80	"	yes

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW TH / AREADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1614	SAMPLING ENDED AT: 1620
PUMP OR TUBING DEPTH IN WELL (feet): 26	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTRATION EQUIPMENT TYPE: N	FILTER SIZE: µm
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-30	3	CG	40ml	HCl	40ml	—	82607+82608+sim	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-31	SAMPLE ID: MW-31 DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 275 feet to 295 feet	STATIC DEPTH TO WATER (feet): 16.96	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 298 feet) + .25 gallons = 0.788 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 285	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 285	PURGING INITIATED AT: 1215	PURGING ENDED AT: 1300	TOTAL VOLUME PURGED (gallons): 3.60

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
3 ⁹ 1254	3.12	3.12	0.08	23.97	7.6	24.1	0.666	0.4	4.65	clear	NO	-184
1256	0.16	3.28	0.08	24.09	7.7	24.2	0.667	0.4	3.94	clear	NO	-186
1258	0.16	3.44	0.08	24.25	7.7	24.2	0.672	0.4	3.66	clear	NO	-187
1300	0.16	3.60	0.08	24.35	7.7	24.3	0.669	0.4	4.11	clear	NO	-188

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1301	SAMPLING ENDED AT: 1305
PUMP OR TUBING DEPTH IN WELL (feet): 285	SAMPLE PUMP FLOW RATE (mL per minute): ~290	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-31	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC</u>	SITE LOCATION: <u>Tallevast</u>
WELL NO: <u>MW-32</u>	SAMPLE ID: <u>MW-32</u>
DATE: <u>1-24-08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>29.5</u> feet to <u>29.5</u> feet	STATIC DEPTH TO WATER (feet): <u>4.61</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + <u>6.0026</u> gallons/foot X <u>29.5</u> feet + <u>0.25</u> gallons = <u>0.33</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>29.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>27</u>	PURGING INITIATED AT: <u>9:35</u>	PURGING ENDED AT: <u>10:06</u>	TOTAL VOLUME PURGED (gallons): <u>3.1</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>9:57</u>	<u>2.2</u>	<u>2.2</u>	<u>0.1</u>	<u>6.93</u>	<u>5.60</u>	<u>30.25</u>	<u>0.755</u>	<u>1.35</u>	<u>1.86</u>	<u>clear</u>	<u>none</u>
<u>10:00</u>	<u>0.3</u>	<u>2.5</u>	<u>0.1</u>	<u>6.85</u>	<u>5.61</u>	<u>30.18</u>	<u>0.759</u>	<u>1.08</u>	<u>3.19</u>	<u>clear</u>	<u>none</u>
<u>10:03</u>	<u>0.3</u>	<u>2.8</u>	<u>0.1</u>	<u>6.85</u>	<u>5.61</u>	<u>30.24</u>	<u>0.753</u>	<u>1.06</u>	<u>1.86</u>	<u>clear</u>	<u>none</u>
<u>10:06</u>	<u>0.3</u>	<u>3.1</u>	<u>0.1</u>	<u>6.84</u>	<u>5.61</u>	<u>30.25</u>	<u>0.759</u>	<u>1.12</u>	<u>1.68</u>	<u>clear</u>	<u>none</u>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Darrin Johnson</u>	SAMPLER(S) SIGNATURES: <u>Darrin Johnson</u>	SAMPLING INITIATED AT: <u>10:07</u>	SAMPLING ENDED AT: <u>10:12</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>27</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	
DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-32</u>	<u>3</u>	<u>CG</u>	<u>90ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast	SITE LOCATION: Sarasota, FL
WELL NO: MW-33	SAMPLE ID: MW-33
DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 43.5 feet) + 0.5 gallons = 0.61 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 0945	PURGING ENDED AT: 1013	TOTAL VOLUME PURGED (gallons): 2.24

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1001	1.28	1.28	0.08	9.64	6.50	28.78	1257	1.26	2.60	clear	NONE
1004	0.24	1.52	0.08	9.66	6.53	28.93	1262	0.79	1.60	clear	NONE
1007	0.24	1.76	0.08	9.65	6.56	28.91	1267	0.83	1.05	clear	NONE
1010	0.24	2.0	0.08	9.66	6.60	28.96	1274	0.85	1.80	clear	NONE
1013	0.24	2.24	0.08	9.67	6.61	28.98	1275	0.72	0.86	clear	NONE

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: JHSW / ARCAOIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1014	SAMPLING ENDED AT: 1020
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-33	3	CG	40ml	HCl	40ml	—	82008+826051A-RFPP	

REMARKS: **Pizzomake removed @ 0932 put back @ 1024**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABL Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-34	SAMPLE ID: MW-34
DATE: 1.30.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet): 18.05	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 155 feet) + 0.25 gallons = 0.65 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150	PURGING INITIATED AT: 7:48	PURGING ENDED AT: 8:15	TOTAL VOLUME PURGED (gallons): 2.7

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
7:55	0.7	0.7	0.1	18.5	7.27	23.9	0.578	0.4	4	clear	none	-292
8:00	0.5	1.2	0.1	18.6	7.32	24.0	0.572	0.36	12	"	"	-303
8:05	0.5	1.7	0.1	18.5	7.36	23.8	0.570	0.34	0	"	"	-301
8:10	0.5	2.2	0.1	18.4	7.38	23.7	0.569	0.34	2	"	"	-312
8:15	0.5	2.7	0.1	18.4	7.38	23.8	0.569	0.36	0	"	"	-313
None												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 8:16	SAMPLING ENDED AT: 8:20
PUMP OR TUBING DEPTH IN WELL (feet): 150 MDS 4/14/08	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-34	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellest</u>	SITE LOCATION: <u>Jarvis, FL</u>
WELL NO: <u>MW-35</u>	SAMPLE ID: <u>MW-35</u> DATE: <u>1/30/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>25</u> feet to <u>30</u> feet	STATIC DEPTH TO WATER (feet): <u>2.81</u>	PURGE PUMP TYPE OR BAILER: <u>(pp) geom pump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<u>0.026</u> gallons/foot X <u>32</u> feet) + <u>0.25</u> gallons = <u>0.33</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~27</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~27</u>	PURGING INITIATED AT: <u>0755</u>	PURGING ENDED AT: <u>0808</u>	TOTAL VOLUME PURGED (gallons): <u>1.82</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle <u>OD</u> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
0759	0.56	0.56	0.14	<u>3.59</u>	5.38	23.76	1560	0.69	3.31	clear	yes	-20
0802	0.42	0.98	0.14	3.59	5.38	23.76	1550	0.63	2.14	clear	yes	-21
0805	0.42	1.40	0.14	3.59	5.37	23.78	1554	0.63	1.01	clear	yes	-24
0808	0.42	1.82	0.14	3.61	5.37	23.82	1.53	0.62	1.48	clear	yes	-17
<i>Discontinued</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Ned Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Ned Smith</u>	SAMPLING INITIATED AT: <u>0808</u>	SAMPLING ENDED AT: <u>0820</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~27</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>~100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> N FILTER SIZE: _____ µm	DUPLICATE: Y <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-35	3	CG	400mL	HCL	~	~2	8260B+260mL RFP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.02

SITE NAME: Tallevast		SITE LOCATION: Tallevast, FL	
WELL NO: MW-36	SAMPLE ID: MW-36	DATE: 1/31/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 5.63	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0006 gallons/foot X 35 feet) + 0.25 gallons = .34 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1056	PURGING ENDED AT: 1112	TOTAL VOLUME PURGED (gallons): 1.80
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9 1106	1.08	1.08	0.12	7.41	5.34	28.8	395	.80	4.00	clear	NO
1108	.24	1.32	0.12	7.47	5.35	28.9	397	.77	2.52	clear	NO
1110	.24	1.56	0.12	7.50	5.34	28.8	403	.72	2.35	clear	NO
1112	.24	1.80	0.12	7.52	5.35	28.9	406	.70	2.00	clear	NO
			0.12								

ORP
61
59
58
57

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / Arcadis	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1113	SAMPLING ENDED AT: 1120
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): 450	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> NO FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-36	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.09

SITE NAME: Tallerast		SITE LOCATION: Tallerast, FL	
WELL NO: MW-37	SAMPLE ID: MW-37	DATE: 1/31/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet): 10.18	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 43.5 feet) + 0.25 gallons = 36 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 1004	PURGING ENDED AT: 1030	TOTAL VOLUME PURGED (gallons): 1.76
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1024	1.10	1.10	0.11	10.45	6.8	27.1	1.23	0.5	6.61	clear	NONE
1026	.22	1.32	0.11	10.47	6.8	27.1	1.20	0.5	8.27	clear	NONE
1028	.22	1.54	0.11	10.48	6.8	27.1	1.18	0.4	5.65	clear	NONE
1030	.22	1.76	0.11	10.48	6.8	27.1	1.16	0.4	5.94	clear	NONE

REF
-81
-83
-86
-87

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STALL/ARCADIS		SAMPLER(S) SIGNATURES: <i>Bridget Stall</i>		SAMPLING INITIATED AT: 1031	SAMPLING ENDED AT: 1038
PUMP OR TUBING DEPTH IN WELL (feet): 38		SAMPLE PUMP FLOW RATE (mL per minute): 430		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> N FILTER SIZE: _____ µm		DUPLICATE: <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-37	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **sampled w/ transducer in well, missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast		SITE LOCATION: Sarasota, FL	
WELL NO: MW-38	SAMPLE ID: MW-38	DATE: 2-1-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (28 - 23) feet X 0.75 gallons/foot = 3.75 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 31 feet) + 0.25 gallons = .33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 0916	PURGING ENDED AT: 0934	TOTAL VOLUME PURGED (gallons): 1.98

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0922	0.66	0.66	0.11	5.93	5.25	29.74	836	1.60	23.5	stained	NO
0925	0.33	0.99	0.11	5.99	5.24	29.99	865	2.94	28	clear	NO
0928	0.53	1.32	0.11	6.02	5.16	30.07	866	1.44	8.45	clear	NO
0931	0.33	1.65	0.11	6.04	5.14	30.10	866	1.52	7.52	clear	NO
0934	0.33	1.98	0.11	6.05	5.13	30.06	865	1.55	5.47	clear	NO

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: TH/SW/ DCAAS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 0936	SAMPLING ENDED AT: 0940
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): ±100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-38	3	CA	40ml	HCL	40ML	—	8260B 8260S1M RFPP	

REMARKS: **Lead sulfate removed @ 0913 Replaced @ 0945**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Seasata, FL
WELL NO: MW-39	SAMPLE ID: MW-39 DATE: 2-1-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 43.5 feet) + 0.25 gallons = 0.36 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 0955	PURGING ENDED AT: 1013 ^{NDX} 4/14/08	TOTAL VOLUME PURGED (gallons): 2.16
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle fig/ or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1001	0.72	0.72	0.12	13.22	10.92	29.94	526	0.94	27.9	clear	NO
1004	0.36	1.08	0.12	13.95	10.93	29.92	518	0.90	21.7	clear	NO
1007	0.36	1.44	0.12	14.47	10.88	29.80	488	0.82	19.0	clear	NO
1010	0.36	1.80	0.12	14.67	10.85	29.71	472	0.77	18.4	clear	NO
1013	0.36	2.16	0.12	14.87	10.70	29.59	445	0.69	15.1	clear	NO
1016	0.36	2.52	0.12							clear	NO

NDX
4/14/08

ORP
-200
-229
-229
-231
-225

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SWITH / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1014	SAMPLING ENDED AT: 1016
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): ≤100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-39	3	CG	40ml	HCl	40ml	—	82600 82651m RFPP	

REMARKS: **Transducer removed @ 0947 Replaced @ 1037**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; -O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleyast		SITE LOCATION: Seasata, FL	
WELL NO: MW-40	SAMPLE ID: MW-40	DATE: 1-31-08	

PURGING DATA


WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0024) gallons/foot X 31 feet + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1129	PURGING ENDED AT: 1144	TOTAL VOLUME PURGED (gallons): 1.80

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1135	0.72	0.72	0.12	5.69	5.23	30.52	854	1.20	2.43	clear	NONE
1138	0.36	1.08	0.12	5.70	5.23	30.57	858	0.98	1.85	clear	NONE
1141	0.36	1.44	0.12	5.69	5.21	30.55	861	0.87	1.80	clear	NONE
1144	0.36	1.80	0.12	5.68	5.20	30.57	857	0.89	2.41	clear	NONE

ORP
-152
-161
-167
-171

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Sw, Tgt / ARCADIS	SAMPLER(S) SIGNATURES: 	SAMPLING INITIATED AT: 1145	SAMPLING ENDED AT: 1150
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): <100 ^(NPL 7/19/08)	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N Filtration Equipment Type: <input checked="" type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW 40	3	CG	40ml	HCl	40ml	—	82605-182605im	RFP

REMARKS: **pizometer removed @ 1125 Replaced @ 1152**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Seasata, FL
WELL NO: MW-41	SAMPLE ID: MW-41
DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026**) gallons/foot X **43.5** feet + **0.5** gallons = **0.6** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 1042	PURGING ENDED AT: 1107
TOTAL VOLUME PURGED (gallons): 2.0			

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1058	1.28	1.28	0.08	9.89	6.74	29.74	1070	1.00	4.19	clear	NONE
1101	0.24	1.52	0.08	9.89	6.73	29.74	1105	0.75	1.50	clear	NONE
1104	0.24	1.76	0.08	9.89	6.72	29.68	1121	0.71	1.79	clear	NONE
1107	0.24	2.0	0.08	9.89	6.72	29.72	1131	0.68	1.24	clear	NONE

OEP
-145
-148
-150
-138

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SWITH / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1108	SAMPLING ENDED AT: 1111
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): <100 (MDS 9/14/02)	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-41	3	CG	40ml	HCl	40ml	—	8260 + 8200 SIM RFP	

REMARKS: **piezometer pulled @ 1032 Replaced @ 1116**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
- pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Sarasota, FL
WELL NO: MW-42	SAMPLE ID: MW-42 DATE: 2-1-08

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 540	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 31 feet) + 0.25 gallons = .33 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 10:33	PURGING ENDED AT: 10:50	TOTAL VOLUME PURGED (gallons): 1.7
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle m or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
10:38	.5	.5	0.1	7.53	5.06	30.36	870	1.69	1.49	clear	none
10:41	.3	.8	0.1	7.51	5.04	30.40	877	1.34	2.68	clear	none
10:44	.3	1.1	0.1	7.51	5.04	30.18	888	1.12	1.74	clear	none
10:47	.3	1.4	0.1	7.51	5.04	30.11	881	1.15	.93	clear	none
10:50	.3	1.7	0.1	6.78	5.03	30.41	882	1.16	1.03	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: S. J. H. / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:51	SAMPLING ENDED AT: 10:56
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): 1100 (MPS 9/14/08)	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-42	3	CG	40ml	HCl	40ml	—	8260B 8260S1m RFPP	

REMARKS: **pulled transducer @ 1030 Replaced at 1102**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast		SITE LOCATION: Sarasota, FL	
WELL NO: MW-43	SAMPLE ID: MW-43	DATE: 2-1-08	

MDS 4/14/08

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 355 feet to 405 feet	STATIC DEPTH TO WATER (feet): 10.03	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

MDS 4/14/08

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.026 gallons/foot X 43.5 feet) + 0.25 gallons = 0.3 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 10:35	PURGING ENDED AT: 10:57	TOTAL VOLUME PURGED (gallons): 2.2
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MDS 4/14/08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
10:42	.7	.7	0.1	10.25	6.25	28.95	1.53	1.15	1.81	clear	none
10:45	.3	1.0	0.1	10.26	6.29	28.95	1.46	1.00	1.03	clear	none
10:48	.3	1.3	0.1	10.25	6.3	28.93	1.45	.99	.93	clear	none
10:51	.3	1.6	0.1	10.25	6.34	28.96	1.40	0.63	1.04	clear	none
10:54	.3	1.9	0.1	10.25	6.41	28.92	1.35	1.47	1.0	clear	none
10:57	.3	2.2	0.1	10.25	6.42	28.93	1.33	0.45	1.03	clear	none

134
146
154
757

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: ARCAOIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:58	SAMPLING ENDED AT: 11:03
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): 2100 (MDS 4/14/08)	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-43	3	CG	40ml	HCl	40ml	—	82005226 SM	RFPP
Dup-31	3	CG	40ml	HCl	—	—	8260	RFPP

MDS 4/14/08

REMARKS: **pulled transducer @ 1032 / re-inserted @ 10:59 / Dup 31 taken**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-44	SAMPLE ID: MW-44
DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 142 feet to 152 feet	STATIC DEPTH TO WATER (feet): 18.9	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (152 feet - 18.9 feet) X .16 gallons/foot = 148.9 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (10026 gallons/foot X 155 feet) + .25 gallons = .65 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 152	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 148	PURGING INITIATED AT: 12:39	PURGING ENDED AT: 13:12	TOTAL VOLUME PURGED (gallons): 3.4

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
12:49	1.0	1.0	0.1	23.8	7.52	25.01	537	1.62	29.4	clear	none
12:52	1.3	1.3	0.1	24.69	7.52	24.99	535	1.66	618.0	cloudy	none
12:55	1.3	1.6	0.1	24.83	7.51	24.89	535	1.71	533.0	cloudy	none
12:58	1.3	1.9	0.1	25.13	7.51	24.97	535	1.71	458.0		
13:01	1.3	2.2	0.1	25.79	7.5	24.95	537	1.68	392.0		
13:04	1.3	2.5	0.1	25.80	7.5	24.97	539	1.65	278.0		
13:07	1.3	2.5	0.1	25.8	7.49	25.02	540	1.64	225.0		
13:10	1.3	3.1	0.1	25.83	7.49	24.99	541	1.62	219.0		
13:13	1.3	3.4	0.1	25.88	7.5	24.99	540	1.62	218.5		

-256
-268
-272
-273
-279
-280
-286
-281
-282

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darin Johnson	SAMPLER(S) SIGNATURES: <i>Darin Johnson</i>	SAMPLING INITIATED AT: 13:14	SAMPLING ENDED AT: 13:19
PUMP OR TUBING DEPTH IN WELL (feet): 148 ^{MDS} 4/19/08	SAMPLE PUMP FLOW RATE (mL per minute): 0.1	TUBING MATERIAL CODE: PT	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> N	FILTRATION EQUIPMENT TYPE: <input type="checkbox"/> µm	DUPLICATE: <input checked="" type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-44	3	CG	40mL	HCl	-	-	8260	RFPP

(MDS 4/19/08)

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Talleast
WELL NO: MW-45	SAMPLE ID: MW-45
DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 150 feet to 160 feet	STATIC DEPTH TO WATER (feet): 19.04	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (1,0026 gallons/foot X 153 feet) + 125 gallons = 0.65 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 13:10	PURGING ENDED AT: 13:47	TOTAL VOLUME PURGED (gallons): 3.6					
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
13:40	3.0	3.0	0.1	26.15	7.94	26.25	1670	1.32	257	cloudy	none
13:43	0.3	3.3	0.1	26.40	7.94	26.38	1669	1.25	255	cloudy	none
13:47	0.3	3.6	0.1	26.68	7.95	26.38	1672	1.22	249	cloudy	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson		SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>		SAMPLING INITIATED AT: 13:48	SAMPLING ENDED AT: 13:53
PUMP OR TUBING DEPTH IN WELL (feet):		SAMPLE PUMP FLOW RATE (mL per minute): < 100		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N		FILTER SIZE: _____ µm	
Filtration Equipment Type:		DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-45	3	CG	40ml	HCl	—	—	8260	RF PP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: MW-46	SAMPLE ID: MW-46 DATE: 1.31.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet): 15.1	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.6026 gallons/foot X 360 feet + 0.25 gallons = 1.03 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 280	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 280	PURGING INITIATED AT: 9:33	PURGING ENDED AT: 10:05	TOTAL VOLUME PURGED (gallons): 3.2
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
9:45	1.2	1.20	0.1	16.7	7.77	23.6	2.88	0.47	130	slt. turbid	none	ORP -309
9:50	0.5	1.70	0.1	16.7	7.65	23.68	1.46	0.47	50	"	"	-309
9:55	0.5	2.2	0.1	17.5	7.63	23.7	0.628	0.41	6	"	"	-315
10:00	0.5	2.7	0.1	17.5	7.62	23.8	0.620	0.39	2	"	"	-316
10:05	0.5	3.2	0.1	17.7	7.61	23.9	0.619	0.38	2	"	"	-316

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:06	SAMPLING ENDED AT: 10:10
PUMP OR TUBING DEPTH IN WELL (feet): 280	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE: NDS F/14/08	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-46	3	CG	40ml	HCl	—	—	8260	RFPP
Dup-27	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **collect Dup-27 here**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast	SITE LOCATION: Sarasota, FL
WELL NO: MW-47	SAMPLE ID: MW-47 DATE: 1-30-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 22 feet to 27 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 30 feet) + 0.5 gallons = 0.58 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 24.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 24.5	PURGING INITIATED AT: 1631	PURGING ENDED AT: 1700	TOTAL VOLUME PURGED (gallons): 2.03

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1648	1.19	1.19	0.07	5.99	6.68	25.34	807	0.89	11.7	clear	yes	-168
1651	0.21	1.4	0.07	6.00	6.68	25.32	815	0.72	7.44	clear	yes	-171
1654	0.21	1.61	0.07	6.01	6.68	25.31	822	0.65	5.87	clear	yes	-186
1657	0.21	1.82	0.07	6.01	6.69	25.31	822	0.63	5.58	clear	yes	-202
1700	0.21	2.03	0.07	6.02	6.72	25.33	819	0.60	2.37	clear	yes	-219

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW, TH / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1701 TH	SAMPLING ENDED AT: 1705 TH
PUMP OR TUBING DEPTH IN WELL (feet): 30 24.5	SAMPLE PUMP FLOW RATE (mL per minute): £100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-47	3	CG	40ml	HCl	40ml	---	8260B+8260SSIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallecra	SITE LOCATION: Seasata, FL
WELL NO: Mw-48	SAMPLE ID: Mw-48 DATE: 1/30/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 38.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 43.5 feet + 0.5 gallons = 0.61 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 36	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 36	PURGING INITIATED AT: 11:35	PURGING ENDED AT: 11:58	TOTAL VOLUME PURGED (gallons): 2.07

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
11:49	1.26	1.26	0.09	10.92	6.61	24.58	1201	0.67	2.58	clear	NO
11:52	0.27	1.53	0.09	10.93	6.62	24.55	1203	0.55	2.47	clear	NO
11:55	0.27	1.80	0.09	10.91	6.60	24.57	1204	0.51	2.42	clear	NO
11:58	0.27	2.07	0.09	10.92	6.60	24.54	1205	0.48	2.36	clear	NO

ORP
-163
-179.4
-193.7
-197.5

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: S. Wells, T. Hollingsworth / Arcadis	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1200	SAMPLING ENDED AT: 1206
PUMP OR TUBING DEPTH IN WELL (feet): 36	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
Mw-48	3	CG	40ml	HCl	40ml	—	8208+8208sim RFPF	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tallevast	SITE LOCATION: Sarasota FL	MDS 4/14/08
WELL NO: MW-49	SAMPLE ID: MW-49	DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 146 feet to 156 feet	STATIC DEPTH TO WATER (feet): 16.45	PURGE PUMP TYPE OR BAILER: Geopump (PP)
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable) = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = gallons + **0.0026** gallons/foot X **159** feet + **.25** gallons = **0.66** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 151	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 151	PURGING INITIATED AT: 1046	PURGING ENDED AT: 1136	TOTAL VOLUME PURGED (gallons): 2.50
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1126	2.00	2.00	0.05	27.10	7.3	24.7	592	0.4	190	cloudy	yes	-253
1128	.10	2.10	0.05	27.16	7.3	24.7	593	0.4	97.9	slightly cloudy	yes	-254
1130	.10	2.20	0.05	27.21	7.3	24.7	593	0.4	67.8	slightly cloudy	yes	-256
1132	.10	2.30	0.05	27.25	7.4	24.7	592	0.4	52.9	slightly cloudy	yes	-257
1134	.10	2.40	0.05	27.28	7.4	24.7	593	0.4	50.1	slightly cloudy	yes	-259
1136	.10	2.50	0.05	27.31	7.4	24.7	593	0.4	48.8	slightly cloudy	yes	-261

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1138	SAMPLING ENDED AT: 1141
PUMP OR TUBING DEPTH IN WELL (feet): 151	SAMPLE PUMP FLOW RATE (mL per minute): ~190	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-49	3	CG	40mL	HCL	40mL	—	8260	RPPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

EB 17.15

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 4/14/08		SITE LOCATION: <u>Sarasota FL</u> MDS 4/14/08	
WELL NO: <u>MW-50</u>	SAMPLE ID: <u>MW-50</u>	DATE: <u>1/20/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>0.25</u>	WELL SCREEN INTERVAL DEPTH: <u>275</u> feet to <u>255</u> feet	STATIC DEPTH TO WATER (feet): <u>15.61</u>	PURGE PUMP TYPE OR BAILER: <u>Peristaltic (2)</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= 0 gallons + (0.0026 gallons/foot X 255 feet) + 0.25 gallons = 0.91 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>11:35</u>	PURGING ENDED AT: <u>13:10</u>	TOTAL VOLUME PURGED (gallons): <u>77</u>
--	--	------------------------------------	--------------------------------	--

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1140	0.5	0.5	0.1	17.29	7.66	25.33	0.683	1.62	828	gray	-	-282
1145	0.5	1.0	0.1	17.69	7.68	25.21	0.697	1.62	60.5	gray	-	-297
1155	1.0	2.0	0.1	17.82	7.70	25.95	0.696	0.91	7999	gray	-	-339
1200	0.5	2.5	0.1	18.00	7.68	25.99	0.696	0.90	155	gray	-	-347
1205	0.5	3.0	0.1	18.18	7.68	25.38	0.697	1.05	76.0	gray	-	-334
1210	0.5	3.5	0.1	18.22	7.67	25.31	0.694	1.75	88.3	gray	-	-319
1215	0.5	4.0	0.1	18.25	7.65	25.67	0.696	0.96	812	gray	-	-343
1225	1.0	5.0	0.1	18.27	7.65	25.96	0.696	0.91	1559	gray	-	-362
1230	0.5	5.5	0.1	18.27	7.64	25.82	0.696	0.91	1318	gray	-	-364
1235	0.5	6.0	0.1	18.27	7.64	25.78	0.698	0.99	1058	gray	-	-361

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith</u>	SAMPLER(S) SIGNATURES: <u>Nathan Smith</u>	SAMPLING INITIATED AT: <u>1313</u>	SAMPLING ENDED AT: <u>1318</u>
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PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u>	TUBING MATERIAL CODE:
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N MDS 4/14/08	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N MDS 4/14/08	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-50	3	CG	40ml	HCL	—	—	8266	RFPP

REMARKS: * Ran 2 pumps in series.

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

page 2 of 2

SITE NAME:		SITE LOCATION:	
WELL NO: <i>MW-50</i>	SAMPLE ID: <i>MW-50</i>	DATE: <i>1/20/08</i>	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:								
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (<i>See Page 1</i> feet - <i>See Page 1</i> feet) X _____ gallons/foot = _____ gallons												
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = _____ gallons + (_____ gallons/foot X _____ feet) + _____ gallons = _____ gallons												
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT:	PURGING ENDED AT:	TOTAL VOLUME PURGED (gallons):								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<i>1240</i>	<i>0.5</i>	<i>6.5</i>	<i>0.1</i>	<i>18.27</i>	<i>7.63</i>	<i>25.95</i>	<i>0.698</i>	<i>0.92</i>	<i>903</i>	<i>95y</i>	—	<i>-363</i>
<i>1245</i>	<i>0.5</i>	<i>7.0</i>	<i>0.1</i>	<i>18.27</i>	<i>7.63</i>	<i>25.45</i>	<i>0.699</i>	<i>1.07</i>	<i>774</i>	<i>91y</i>	—	<i>-361</i>
<i>1250</i>	<i>0.25</i>	<i>7.25</i>	<i>0.05</i>	<i>18.27</i>	<i>7.63</i>	<i>25.67</i>	<i>0.699</i>	<i>0.93</i>	<i>145</i>	<i>light gray</i>	—	<i>-366</i>
<i>1255</i>	<i>0.25</i>	<i>7.5</i>	<i>0.05</i>	<i>18.27</i>	<i>7.63</i>	<i>25.82</i>	<i>0.699</i>	<i>0.99</i>	<i>153</i>	<i>clearing up</i>	—	<i>-350</i>
<i>1258</i>		<i>7.7</i>	<i>0.05</i>	<i>18.27</i>	<i>7.63</i>	<i>25.88</i>	<i>0.699</i>	<i>0.96</i>	<i>161</i>	" "	—	<i>-367</i>
<i>1301</i>				<i>18.27</i>	<i>7.63</i>	<i>25.69</i>	<i>0.698</i>	<i>0.93</i>	<i>146</i>	" "	—	<i>-368</i>
<i>1304</i>				<i>18.27</i>	<i>7.63</i>	<i>25.83</i>	<i>0.700</i>	<i>0.91</i>	<i>168</i>	" "	—	<i>-378</i>
<i>1307</i>				<i>18.27</i>	<i>7.63</i>	<i>25.82</i>	<i>0.700</i>	<i>0.91</i>	<i>160</i>	" "	—	<i>-371</i>
<i>1310</i>				<i>18.27</i>	<i>7.63</i>	<i>25.63</i>	<i>0.698</i>	<i>0.91</i>	<i>165</i>	" "	—	<i>-371</i>
<i>1313</i>	→ <i>Sampled at 1313</i> ←											
WELL CAPACITY (Gallons Per Foot): <i>0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88</i>												
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): <i>1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016</i>												

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION:		SAMPLER(S) SIGNATURES:		SAMPLING INITIATED AT:	SAMPLING ENDED AT:		
PUMP OR TUBING DEPTH IN WELL (feet):		SAMPLE PUMP FLOW RATE (mL per minute):		TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: Y N		FIELD-FILTERED: Y N		FILTER SIZE: _____ µm	DUPLICATE: Y N		
Filtration Equipment Type:							
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION		INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)		

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/14/08 SITE LOCATION: Sarasota FL MDS 4/14/08
WELL NO: MW-51 SAMPLE ID: MW-51 DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches): 2" TUBING DIAMETER (inches): 1/4" WELL SCREEN INTERVAL DEPTH: 261.6 feet to 271.6 feet STATIC DEPTH TO WATER (feet): 12.77 PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)
= 0 gallons + (0.0026 gallons/foot X 271.6 feet) + 0.25 gallons = 0.96 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): _____ FINAL PUMP OR TUBING DEPTH IN WELL (feet): _____ PURGING INITIATED AT: 4:05 / 3:20 PURGING ENDED AT: 4:30 TOTAL VOLUME PURGED (gallons): 2.2

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
4:10	0.5	0.5	0.5	17.64	7.26	22.36	0.301	2.63	36.2	clear	-	-159
4:15	0.5	1.0	0.1	10.67	7.27	22.17	0.302	2.22	23.7	clear	-	-167
4:18	0.3	1.3	0.1	19.51	7.17	22.16	0.302	2.17	14.6	clear	-	-168
4:21	0.3	1.6	0.1	20.49	7.16	21.90	0.302	2.11	15.1	clear	-	-167
4:24	0.3	1.9	0.1	20.99	7.15	22.02	0.306	2.13	16.7	clear	-	-166
4:27	0.7	2.2	0.1	22.10	7.16	21.60	0.313	2.03	8.45	clear	-	-180
4:30	sampled @ 10:36											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN P SWIRE SAMPLER(S) SIGNATURES: Nathan Swire SAMPLING INITIATED AT: 4:30 SAMPLING ENDED AT: 4:40

PUMP OR TUBING DEPTH IN WELL (feet): _____ SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/08 TUBING MATERIAL CODE: _____

FIELD DECONTAMINATION: YES MDS 4/14/08 FIELD-FILTERED: YES MDS 4/14/08 FILTER SIZE: _____ µm DUPLICATE: YES N

SAMPLE CONTAINER SPECIFICATION: _____ SAMPLE PRESERVATION: _____ INTENDED ANALYSIS AND/OR METHOD: _____ SAMPLING EQUIPMENT CODE: _____

SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
MW-51	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: _____

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: MW-52	SAMPLE ID: MW-52
DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 147 feet to 157 feet	STATIC DEPTH TO WATER (feet): 15.53	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 157 feet) + 0.25 gallons = .66 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 11:06	PURGING ENDED AT: 11:37	TOTAL VOLUME PURGED (gallons): 3.1

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle, mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1116	1.0	1.0	0.1	19.48	9.55	23.75	0.277	0.60	6.53	grey	no odor
1119	0.3	1.3		19.72	9.40	23.71	0.287	0.57	7.37		
1122		1.6		19.88	9.26	23.74	0.297	0.55	8.12		
1125		1.9		19.98	9.16	23.74	0.305	0.54	8.07		
1128		2.2		20.06	9.08	23.76	0.313	0.52	8.78		
1131		2.5		20.30	8.79	23.65	0.346	0.45	11.29		
1134		2.8		20.40	8.74	23.79	0.350	0.43	12.20		
1137		3.1		20.48	8.68	23.68	0.359	0.41	14.60		

-250
-254
-259
-262
-269
-270
-272
-273

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: MDS 9/14/08 IARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1138	SAMPLING ENDED AT: 1140
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 5100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N Filtration Equipment Type:	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW.52	3	CG	40mL	HCl	—	—	B260	RFPP

MDS
9/14/08

REMARKS: **transducer removed @ 11:04 Pumping stopped 11:45 Transducer replaced**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

2 gallons removed

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG



SITE NAME: <u>Telleme</u>		SITE LOCATION: <u>Sarasota FL</u>	
WELL NO: <u>MW-53</u>		SAMPLE ID: <u>MW-53</u>	
DATE: <u>1/28/08</u>			

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>141</u> feet to <u>151</u> feet	STATIC DEPTH TO WATER (feet): <u>16.19</u>	PURGE PUMP TYPE OR BAILER: <u>(APP) gas pump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X <u>160</u> feet) + 0.50 gallons = <u>0.91</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>-146</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>-146</u>	PURGING INITIATED AT: <u>1145</u>	PURGING ENDED AT: <u>1206</u>	TOTAL VOLUME PURGED (gallons): <u>1.89</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or S/cm)	DISSOLVED OXYGEN (circle <u>DO</u> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1157	1.08	1.08	0.09	21.33	7.46	24.69	620	0.45	42.4	clear	yes
1200	0.27	1.35	0.09	22.31	7.45	24.81	632	0.42	79.4	clear	yes
1203	0.27	1.62	0.09	22.72	7.45	24.81	631	0.41	79.0	clear	yes
1206	0.27	1.89	0.09	22.77	7.45	24.69	635	0.42	80.5	clear	yes
<i>Neel Sen 1/28/08</i>											

*ORP
-246
-252
-253
-253*

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neel Sen / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neel Sen</u>	SAMPLING INITIATED AT: <u>1210</u>	SAMPLING ENDED AT: <u>1215</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~146</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> <u>N</u> <input type="checkbox"/> Filtration Equipment Type: _____	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> <u>AS</u> <input checked="" type="checkbox"/> <u>N</u>

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-53	3	CG	40ml	HCC	-	~2	<u>DO + ORP + pH</u>	<u>RFPP</u>

REMARKS: DUP-17

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellepsen</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>MW-54</u>	SAMPLE ID: <u>MW-54</u> DATE: <u>1/28/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>145</u> feet to <u>155</u> feet	STATIC DEPTH TO WATER (feet): <u>15.15</u>	PURGE PUMP TYPE OR BAILER: <u>pp gear pump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X <u>160</u> feet) + 0.50 gallons = <u>(1.00)</u> <u>0.21</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~150</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~150</u>	PURGING INITIATED AT: <u>1050</u>	PURGING ENDED AT: <u>1108</u>	TOTAL VOLUME PURGED (gallons): <u>2.16</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1059	1.08	1.08	0.12	17.36	7.52	22.93	659	0.55	3.30	clear	yes
10102	0.36	1.44	0.12	17.44	7.52	22.87	659	0.54	2.42	clear	yes
1105	0.36	1.80	0.12	17.46	7.52	22.97	660	0.48	2.96	clear	yes
1108	0.36	2.16	0.12	17.49	7.52	22.94	659	0.49	2.35	clear	yes
<i>Next day 1/28/08</i>											

28P
-248
-257
-257
-266

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: <u>1109</u>	SAMPLING ENDED AT: <u>1115</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~150</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>~100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> <u>N</u>	FIELD-FILTERED: <u>Y</u> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-54</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCL</u>	<u>-</u>	<u>7.2</u>	<u>826B-1026JW</u>	<u>RFP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.83

SITE NAME: Tallevast	SITE LOCATION: Tallevast FL
WELL NO: MW-55	SAMPLE ID: MW-55 DATE: 1/25/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 127 feet to 137 feet	STATIC DEPTH TO WATER (feet): 16.80	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 140 feet) + 0.25 gallons = 0.61 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 132	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 132	PURGING INITIATED AT: 0816	PURGING ENDED AT: 0839	TOTAL VOLUME PURGED (gallons): 2.53

17

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0833	1.87	1.87	0.11	20.00	10.24	23.5	.737	.49	7.68	clear	yes
0835	0.22	2.09	0.11	20.10	9.69	23.6	.400	.47	5.36	clear	yes
0837	0.22	2.31	0.11	20.10	9.57	23.6	.396	.46	5.17	clear	yes
0839	0.22	2.53	0.11	20.10	9.42	23.6	.395	.47	4.86	clear	yes

ORP
-324
-323
-323
-323

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Denrick King/ALCAA	SAMPLER(S) SIGNATURES: <i>Denrick King</i>	SAMPLING INITIATED AT: 0842	SAMPLING ENDED AT: 0855
PUMP OR TUBING DEPTH IN WELL (feet): 132	SAMPLE PUMP FLOW RATE (mL per minute): 416	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-55	3	CG	40ml	HCl	-	-	8260	RFPP
DUP-1a	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

1.98

SITE NAME: TALLENAST	SITE LOCATION: TALLENAST, P1
WELL NO: mw-56	SAMPLE ID: mw-56
DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet): 15.43	PURGE PUMP TYPE OR BAILER: PP Peristaltic Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 158 feet) + .25 gallons = .66 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150	PURGING INITIATED AT: 1525	PURGING ENDED AT: 1550	TOTAL VOLUME PURGED (gallons): 2.75
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1544	2.09	2.09	.11	18.04	7.75	24.4	.258	0.57	25.0	clear	yes -243
1546	.22	2.31	.11	18.05	7.76	24.4	.251	0.56	19.9	clear	yes -247
1548	.22	2.53	.11	18.10	7.76	24.4	.255	0.55	18.9	clear	yes -249
1550	.22	2.75	.11	18.10	7.77	24.4	.258	0.55	18.1	clear	yes -252

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Bridget STALL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stall</i>	SAMPLING INITIATED AT: 1551	SAMPLING ENDED AT: 1557
PUMP OR TUBING DEPTH IN WELL (feet): 150	SAMPLE PUMP FLOW RATE (mL per minute): 430	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw56	3	CG	40ml	Hel	40ml	-	8260	RPPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RPPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-57	SAMPLE ID: MW-57 DATE: 1/30/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 136 feet to 146 feet	STATIC DEPTH TO WATER (feet): 1843	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 149 feet) + 0.25 gallons = 0.64 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 141	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 141	PURGING INITIATED AT: 0857	PURGING ENDED AT: 0921	TOTAL VOLUME PURGED (gallons): 2.64

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mgS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
0915	1.98	1.98	0.11	20.31	10.3	25.0	.448	0.6	14.3	Clear	NONE	-255
0917	0.22	2.20	0.11	20.35	10.2	25.0	.445	0.7	12.2	Clear	NONE	-256
0919	0.22	2.42	0.11	20.37	10.2	25.0	.445	0.7	12.2	Clear	NONE	-259
0921	0.22	2.64	0.11	20.38	10.1	25.0	.441	0.6	9.49	Clear	NONE	-258

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 0922	SAMPLING ENDED AT: 0930
PUMP OR TUBING DEPTH IN WELL (feet): 141	SAMPLE PUMP FLOW RATE (mL per minute): ~420	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	
SAMPLE CONTAINER SPECIFICATION		SAMPLE PRESERVATION	

SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
MW-57	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **missing 2 bottles**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

1.95

GROUNDWATER SAMPLING LOG

SITE NAME: TALEUAST		SITE LOCATION: TALEUAST, FL	
WELL NO: mw-58		SAMPLE ID: mw-58	DATE: 1-30-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 140 feet to 150 feet	STATIC DEPTH TO WATER (feet): 1930	PURGE PUMP TYPE OR BAILER: PP 600 Pump									
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons													
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 153 feet) + 0.25 gallons = 0.40 gallons													
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 145	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 145	PURGING INITIATED AT: 1402	PURGING ENDED AT: 1423	TOTAL VOLUME PURGED (gallons): 2.73									
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle <input checked="" type="radio"/> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP	
1417	1.95	1.95	0.13	22.25	7.5	25.3	.517	0.4	16.0	clear	yes		-201
1419	0.26	2.21	0.13	22.26	7.5	25.2	.515	0.4	15.1	clear	yes		-201
1421	0.26	2.47	0.13	22.28	7.5	25.2	.514	0.4	14.5	clear	yes		-201
1423	0.26	2.73	0.13	22.30	7.5	25.2	.517	0.4	8.70	clear	yes		-203
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016													

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STANL/ARCADIS			SAMPLER(S) SIGNATURES: <i>Bridget Stanl</i>			SAMPLING INITIATED AT: 1424		SAMPLING ENDED AT: 1436		
PUMP OR TUBING DEPTH IN WELL (feet): 145			SAMPLE PUMP FLOW RATE (mL per minute): 490			TUBING MATERIAL CODE:				
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N			FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N			FILTER SIZE: _____ µm		DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N		
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH				
mw-58	3	CG	40ml	HCl	40ml	-	8260		RPPP	
DUP23	3	CG	40ml	HCl	40ml	-	8260		RPPP	
REMARKS: missing 1 bolt										
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)										
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump EQUIPMENT CODES: RPPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)										

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.95

SITE NAME: Tallevast		SITE LOCATION: Tallevast, FL	
WELL NO: MW-59	SAMPLE ID: MW-59	DATE: 1/28/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 140 feet to 150 feet	STATIC DEPTH TO WATER (feet): 168.5	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 153 feet + 0.25 gallons = 0.65 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 145	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 145	PURGING INITIATED AT: 1211	PURGING ENDED AT: 1232	TOTAL VOLUME PURGED (gallons): 2.73
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15

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1226	1.95	1.95	0.13	21.01	8.29	24.6	0.40	0.51	56.0	cloudy	ys
1228	0.26	2.21	0.13	21.19	8.24	24.7	0.476	0.50	51.1	cloudy	ys
1230	0.26	2.47	0.13	21.26	8.15	24.6	0.470	0.50	46.3	cloudy	ys
1232	0.26	2.73	0.13	21.35	8.07	24.4	0.482	0.49	39.6	"	ys

MDS 4/14/08

CRP

-315
-311
-310
-310

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / Arcadis	SAMPLER(S) SIGNATURES: Jennifer King	SAMPLING INITIATED AT: 1235	SAMPLING ENDED AT: 1240
PUMP OR TUBING DEPTH IN WELL (feet): 145	SAMPLE PUMP FLOW RATE (mL per minute): ~490	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-59	3	CG	40 ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: TALLEVAST, F1
WELL NO: mw-60	SAMPLE ID: mw-60 DATE: 1-28-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet): 18.08	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 158 feet) + .25 gallons = 0.66 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150	PURGING INITIATED AT: 1600	PURGING ENDED AT: 1622	TOTAL VOLUME PURGED (gallons): 2.86

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1616	2.08	2.08	0.13	21.35	7.27	24.1	527	0.49	2.71	clear	yes
1618	0.26	2.34	0.13	21.45	7.32	24.2	527	0.47	3.06	clear	yes
1620	0.26	2.60	0.13	21.55	7.33	24.2	527	0.48	3.61	clear	yes
1622	0.26	2.86	0.13	21.64	7.34	24.2	528	0.47	4.54	clear	yes

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS			SAMPLER(S) SIGNATURES: <i>Jennifer King</i>			SAMPLING INITIATED AT: 1624		SAMPLING ENDED AT: 1627	
PUMP OR TUBING DEPTH IN WELL (feet): 150			SAMPLE PUMP FLOW RATE (mL per minute): 490			TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <input type="radio"/> Y <input checked="" type="radio"/> N			FIELD-FILTERED: <input type="radio"/> Y <input checked="" type="radio"/> N FILTER SIZE: _____ µm			DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-60	3	CG	40ml	HCl	40mL	-	8260	RFPP

REMARKS: **missing 1 bolt**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> ^{MDS 4/14/08}	SITE LOCATION: <u>Sarasota FL</u> ^{MDS 4/14/08}
WELL NO: <u>MW-61</u>	SAMPLE ID: <u>MW-61</u> DATE: <u>1/30/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>135</u> feet to <u>145</u> feet	STATIC DEPTH TO WATER (feet): <u>173.5</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (<u>145</u> feet - <u>173.5</u> feet) X <u>0.0026</u> gallons/foot = <u>0.63</u> gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>145</u> feet) + <u>.25</u> gallons = <u>0.63</u> gallons ^{MDS 4/14/08}				

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
<u>0935</u>	<u>0.5</u>	<u>0.5</u>	<u>0.1</u>	<u>19.40</u>	<u>7.4</u>	<u>25.5</u>	<u>.55</u>	<u>4.8</u>	<u>8.63</u>	<u>Clear</u>	<u>none</u>	<u>160</u>
<u>0938</u>				<u>20.22</u>	<u>7.3</u>	<u>25.7</u>	<u>.59</u>	<u>3.3</u>	<u>5.92</u>			<u>138</u>
<u>0941</u>				<u>21.59</u>	<u>7.4</u>	<u>25.7</u>	<u>.59</u>	<u>3.3</u>	<u>4.14</u>			<u>131</u>
<u>0944</u>				<u>22.82</u>	<u>7.4</u>	<u>25.8</u>	<u>.60</u>	<u>3.1</u>	<u>11.9</u>			<u>97</u>
<u>0947</u>				<u>24.01</u>	<u>9.3</u>	<u>25.8</u>	<u>.49</u>	<u>2.9</u>	<u>12.6</u>			<u>72</u>
<u>0950</u>				<u>24.37</u>	<u>9.2</u>	<u>25.8</u>	<u>.48</u>	<u>2.9</u>	<u>10.51</u>			<u>72</u>
<u>0953</u>				<u>24.69</u>	<u>9.2</u>	<u>25.8</u>	<u>.47</u>	<u>2.9</u>	<u>9.68</u>	<u>V</u>	<u>✓</u>	<u>69</u>
<u>0956</u>	→ <u>Sampled @ 0956</u> ←											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Rosey Cornwell / Arcadis</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>0956</u>	SAMPLING ENDED AT: <u>1000</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>190</u> ^{MDS 4/14/08}	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u> ^{MDS 4/14/08}	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N ^{MDS 4/14/08}	FIELD-FILTERED: <input checked="" type="checkbox"/> Y ^{MDS 4/14/08}	DUPLICATE: <input checked="" type="checkbox"/> Y ^{MDS 4/14/08} <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-61</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellesant	SITE LOCATION: Sarasota, FL
WELL NO: MW-62	SAMPLE ID: MW-62 DATE: 1/31/08

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: 18 feet to 23 feet	STATIC DEPTH TO WATER (feet): 3.76	PURGE PUMP TYPE OR BAILER: (PP) geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 25 feet) + 0.25 gallons = 0.315 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~21	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~21	PURGING INITIATED AT: 1005	PURGING ENDED AT: 1018	TOTAL VOLUME PURGED (gallons): 1.30

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle µg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1009	0.40	0.40	0.10	3.89	6.02	23.52	0.572	1.10	18.1	clear	yes	-111
1012	0.30	0.70	0.10	3.91	5.97	23.65	0.551	0.89	6.4	clear	yes	-109
1015	0.30	1.00	0.10	3.91	5.90	23.77	0.541	0.73	5.9	clear	yes	-106
1018	0.30	1.30	0.10	3.91	5.90	23.79	0.527	0.74	3.6	clear	yes	-106
<i>Handwritten signature and date 1/31/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Butcher ARCADIS			SAMPLER(S) SIGNATURES: <i>Neil Butcher</i>			SAMPLING INITIATED AT: 1020	SAMPLING ENDED AT: 1025
PUMP OR TUBING DEPTH IN WELL (feet): ~21			SAMPLE PUMP FLOW RATE (mL per minute): ~100			TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N			FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: _____ µm			DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-62	3	CG	40mL	HCL	-	~2	RFPP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellepsst	SITE LOCATION: Jarvis, FL
WELL NO: MW-63	SAMPLE ID: MW-63 DATE: 1/31/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 25 feet to 30 feet	STATIC DEPTH TO WATER (feet): 3.19	PURGE PUMP TYPE OR BAILER: (pp) geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet -- feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 32 feet) + 0.25 gallons = 0.33 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~27	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~27	PURGING INITIATED AT: 0935	PURGING ENDED AT: 0948	TOTAL VOLUME PURGED (gallons): 1.30
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle NO or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0939	0.40	0.40	0.10	3.45	7.15	23.35	0.759	3.85	5.21	clear	yes	ORP -222
0942	0.30	0.70	0.10	3.45	7.16	23.35	0.793	3.86	3.34	clear	yes	-214
0945	2.30	1.00	0.10	3.45	7.16	23.46	0.798	3.88	9.25	clear	yes	-210
0948	0.30	1.30	0.10	3.45	6.98	23.49	0.810	3.84	4.1	clear	yes	-203
<i>Next 1/31/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ned Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Ned Smith</i>	SAMPLING INITIATED AT: 0950	SAMPLING ENDED AT: 0955
PUMP OR TUBING DEPTH IN WELL (feet): ~27	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y / N	FIELD-FILTERED: Y FILTER SIZE: µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW63	3	CG	40m	HCL	—	7.2	200B of 20m RFP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-64	SAMPLE ID: MW-64
DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1.4	WELL SCREEN INTERVAL DEPTH: 25 feet to 30 feet	STATIC DEPTH TO WATER (feet): 2.94	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (30 feet - 2.94 feet) X .16 gallons/foot = 4.32 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (.0026 gallons/foot X 33 feet) + .25 gallons = .34 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 30	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 27.5	PURGING INITIATED AT: 11:37	PURGING ENDED AT: 11:59	TOTAL VOLUME PURGED (gallons): 2.2
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
11:47	1.0	1.0	0.1	3.65	4.89	24.93	1208	3.8	5.04	clear	none
11:50	0.3	1.3	0.1	3.65	4.98	25.04	1199	3.51	5.20		
11:53	0.3	1.6	0.1	3.60	5.09	24.95	1202	3.37	5.22		
11:56	0.3	1.9	0.1	3.58	5.10	24.89	1200	3.34	5.21		
11:59	0.3	2.2	0.1	3.56	5.11	24.86	1199	3.33	5.20	✓	✓

MDS
4/14/08

91
40
13
13

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 12:00	SAMPLING ENDED AT: 12:05
PUMP OR TUBING DEPTH IN WELL (feet): 27.5	SAMPLE PUMP FLOW RATE (mL per minute): 5100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	
DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-64	3	CG	40ml	HCl	—	—	8260	RFPP

MDS
4/14/08

REMARKS: **need new T-plug**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 40% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast MDS 4/14/08	SITE LOCATION: Sarasota FL MDS 4/14/08
WELL NO: MW-65	SAMPLE ID: MW-65 DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 19 feet to 24 feet	STATIC DEPTH TO WATER (feet): 3.51	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= 0 gallons + (0.0026 gallons/foot X **24** feet) + **.25** gallons = **0.31** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 22	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 22	PURGING INITIATED AT: 0955	PURGING ENDED AT: 1009	TOTAL VOLUME PURGED (gallons): 1.4
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1000	0.5	0.5	6.1	4.75	6.2	18.2	50	4.9	15.7	Clear	none
1003	0.3	0.8	↓	6.10	6.5	19.3	48	4.1	17.2	↓	↓
1006	0.3	1.1	↓	7.11	6.5	19.1	48	4.1	15.1	↓	↓
1009	0.3	1.4	↓	7.15	6.5	19.2	47	4.1	11.6	↓	↓
1012	→ Sampled @ 1012 ←										

MDS
4/14/08
MDS
4/14/08
ORP
234
816
206
155

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KLC / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1012	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 22	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: 1 µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-65	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS: ***initial purge clear no odor.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

-needs 2 Bolts

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: Sarasota FL MMS 4/19/08
WELL NO: mW-66	SAMPLE ID: mW-66 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 19.5 feet to 23.5 feet	STATIC DEPTH TO WATER (feet): 3.68	PURGE PUMP TYPE OR BAILER: PP Get Pump
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026** gallons/foot X **25.5** feet) + **.25** gallons = **0.32** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 21	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 21	PURGING INITIATED AT: 1354	PURGING ENDED AT: 1408	TOTAL VOLUME PURGED (gallons): 1.68
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1402	0.96	0.96	0.12	4.21	5.66	25.5	.304	0.74	15.8	Clear	NONE	-50
1404	.24	1.20	0.12	4.24	5.63	25.5	.303	0.7	5.94	Clear	NONE	-56
1406	.24	1.44	0.12	4.25	5.63	25.5	.305	0.68	1.78	clear	NONE	-59
1408	.24	1.68	0.12	4.25	5.63	25.5	.303	0.66	3.96	clear	NONE	-63

ORP

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STALL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stall</i>	SAMPLING INITIATED AT: 1409	SAMPLING ENDED AT: 1415
PUMP OR TUBING DEPTH IN WELL (feet): 21	SAMPLE PUMP FLOW RATE (mL per minute): ~454	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mW-66	3	CG	40mL	HCL	40mL	-	8260	RFPP
DWP 2	3	CG	40mL	HCL	40mL	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump

EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.05

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST, FL	
WELL NO: mw-67		SAMPLE ID: mw-67	
DATE: 1-28-08			

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 24 feet to 29 feet	STATIC DEPTH TO WATER (feet): 5.19	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0024 gallons/foot X 37 feet) + .25 gallons = 0.35 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26.5	PURGING INITIATED AT: 1636	PURGING ENDED AT: 1648	TOTAL VOLUME PURGED (gallons): 2.3

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1642	1.14	1.14	0.19	6.65	5.71	25.7	0.398	1.51	13.0	clear	none
1644	0.38	1.52	0.19	6.72	5.71	25.8	0.432	1.10	9.54	clear	none
1646	0.38	1.90	0.19	6.78	5.71	25.8	.462	0.74	10.13	clear	none
1648	0.38	2.28	0.19	6.79	5.71	25.8	.470	0.63	14.9	clear	none

ORP
28
-10
-32
-41

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1650	SAMPLING ENDED AT: 1657
PUMP OR TUBING DEPTH IN WELL (feet): 26.5	SAMPLE PUMP FLOW RATE (mL per minute): ~700	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y Filtration Equipment Type:	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-67	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-68	SAMPLE ID: MW-68 DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet): 10.57	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (40.5 feet - 10.57 feet) X 0.16 gallons/foot = 4.79 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + 0.0026 gallons/foot X 43.5 feet + 0.25 gallons = 0.36 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 1247	PURGING ENDED AT: 1340	TOTAL VOLUME PURGED (gallons): 9.01
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29
8
8
8

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle ng/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1316	4.93	4.93	0.17	11.29	7.23	25.1	1561	0.46	4.10	Clear	NONE
1324	6.36	6.29	0.17	11.30	7.27	25.1	1563	0.46	0.52	Clear	NONE
1332	6.36	7.65	0.17	11.30	7.29	25.0	1564	0.44	0.29	Clear	NONE
1340	6.36	9.01	0.17	11.30	7.29	25.0	1568	0.44	0.26	Clear	NONE

ORP
-210
-215
-220
-222

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/ALCOA'S	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1342	SAMPLING ENDED AT: 1350
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): 640	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-68	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **missing 1 bolt**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast	MDS 4/14/08	SITE LOCATION: Sarasota FL	MDS 4/14/08
WELL NO: MW-69	SAMPLE ID: MW-69		DATE: 1/30/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 4.32	PURGE PUMP TYPE OR BAILER: Peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY = (28 feet - 4.32 feet) X 0.25 gallons/foot = 5.92 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME = 0 gallons + (0.0026 gallons/foot X 26 feet) + 0.25 gallons = 0.3 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26	PURGING INITIATED AT: 9:31	PURGING ENDED AT: 11:02	TOTAL VOLUME PURGED (gallons): 3.2

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1039	0.6	0.6	0.1	5.92	5.64	26.92	0.573	1.24	4.43	clear	-	-157
1041	0.3	0.9	0.1	5.92	5.64	26.93	0.545	1.13	4.16	clear	-	-159
1043	0.3	1.4	0.1	5.92	5.64	26.95	0.521	1.10	4.50	clear	-	-164
1047	0.3	1.7	0.1	5.92	5.65	26.96	0.529	1.05	4.33	" "	-	-174
1050	0.3	2.0	0.1	5.91	5.65	27.00	0.579	1.01	3.67	" "	-	-182
1053	0.3	2.3	0.1	5.91	5.65	26.96	0.508	1.03	4.34	" "	-	-186
1056	0.3	2.6	0.1	5.91	5.64	26.93	0.584	1.00	4.66	" "	-	-190
1059	0.3	2.9	0.1	5.91	5.62	27.06	0.584	1.05	4.24	" "	-	-194
1102	0.3	3.2	0.1	5.91	5.61	27.03	0.581	1.01	4.20	" "	-	-197

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN SMITH	SAMPLER/SIGNATURE: <i>Nathan Smith</i>	SAMPLING INITIATED AT: 11:02	SAMPLING ENDED AT: 11:08
PUMP OR TUBING DEPTH IN WELL (feet): 26	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y	FIELD-FILTERED: <input checked="" type="checkbox"/> Y	DUPLICATE: <input checked="" type="checkbox"/> Y	<input checked="" type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-69	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **Outer casing full of water. Had to bail it out.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVA87		SITE LOCATION: TALLEVA87 P1	
WELL NO: MW-70	SAMPLE ID: MW-70	DATE: 1-30-07	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 29 feet	STATIC DEPTH TO WATER (feet): 5.85	PURGE PUMP TYPE OR BAILER: 600 Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 37 feet) + .25 gallons = .35 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26	PURGING INITIATED AT: 1232	PURGING ENDED AT: 1250	TOTAL VOLUME PURGED (gallons): 2.21 2.89
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
7 1240	1.19	1.19	0.17	9.35	5.86	25.0	1609	56	20.9	clear	none	-68
1242	.34	1.53	0.17	9.62	5.79	25.1	1709	53	17.9	clear	none	-69
1244	.34	1.87	0.17	9.78	5.73	25.1	1783	52	14.9	clear	none	-70
1246	.34	2.21	0.17	9.91	5.68	25.1	1849	51	11.9	clear	none	-71
1248	.34	2.55	.17	9.95	5.65	25.1	1876	50	10.68	clear	none	-72
1250	.34	2.89	.17	10.00	5.63	25.2	1880	50	9.67	clear	none	-74
JK 1252	.34	3.23	.17							clear	none	

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1251	SAMPLING ENDED AT: 1257
PUMP OR TUBING DEPTH IN WELL (feet): 26	SAMPLE PUMP FLOW RATE (mL per minute): 643	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-70	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS: **missing 2 bolts; ADDED 5' teflon tubing; well cap destroyed**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.02

SITE NAME: Tallerast		SITE LOCATION: Tallerast	
WELL NO: mw-71	SAMPLE ID: mw-71	DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 24 feet to 29 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 35 feet) + .25 gallons = .34 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26.5	PURGING INITIATED AT: 1321	PURGING ENDED AT: 1333	TOTAL VOLUME PURGED (gallons): 2.28
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1327	1.14	1.14	.19	9.21	5.51	27.0	.346	.70	9.10	clear	none
1329	.38	1.52	.19	9.13	5.53	27.0	.375	.65	5.52	clear	none
1331	.38	1.90	.19	9.40	5.53	27.0	.383	.63	7.30	clear	none
1333	.38	2.28	.19	9.59	5.54	27.0	.389	.61	8.10	clear	none

ORP
-34
-40
-43
-47

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King-Jacobs	SAMPLER(S) SIGNATURES: Jennifer King	SAMPLING INITIATED AT: 1334	SAMPLING ENDED AT: 1340
PUMP OR TUBING DEPTH IN WELL (feet): 26.5	SAMPLE PUMP FLOW RATE (mL per minute): ~700	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: 0 N	FIELD-FILTERED: Y 0 N FILTER SIZE: _____ µm	DUPLICATE: Y 0 N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-71	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS: **missing 2 bolts, sampled w/ transducer; ADDED 35' teflon tubing.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-72	SAMPLE ID: MW-72 DATE: 1/30/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23.5 feet to 28.5 feet	STATIC DEPTH TO WATER (feet): 4.56	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 31.5 feet) + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26	PURGING INITIATED AT: 1006	PURGING ENDED AT: 1022	TOTAL VOLUME PURGED (gallons): 1.76

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1016	1.10	1.10	0.11	5.03	5.5	28.8	1.21	0.5	11.5	clear	yes	-73
1018	0.22	1.32	0.11	5.04	5.5	28.8	1.22	0.5	11.6	clear	yes	-76
1020	0.22	1.54	0.11	5.06	5.5	28.9	1.22	0.4	11.1	clear	yes	-81
1022	0.22	1.76	0.11	5.08	5.5	28.9	1.21	0.4	10.68	clear	yes	-84

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Bridget Stahl / AECADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1023	SAMPLING ENDED AT: 1027
PUMP OR TUBING DEPTH IN WELL (feet): 26	SAMPLE PUMP FLOW RATE (mL per minute): 416	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-72	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **Sampled w/ transducer in well ; missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-73		SAMPLE ID: MW-73	
		DATE: 1.25.08	

PURGING DATA *MDA 4/14/08*

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 22 feet to 27 feet	STATIC DEPTH TO WATER (feet): 2.7	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 14:19		PURGING ENDED AT: 1450		TOTAL VOLUME PURGED (gallons): 3.1			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1420	0.1	0.1	0.1	2.7	6.00	24.04	1.26	0.80	212	Clear	None
1425	0.5	0.6	0.1	2.7	4.75	24.2	1.27	0.40	41	ll	ll
1430	0.5	1.1	0.1	2.7	4.94	24.3	1.32	0.37	50	ll	ll
1435	0.5	1.6	0.11	2.7	5.03	24.5	1.33	0.34	10	ll	ll
1440	0.5	2.1	0.11	2.7	5.09	24.4	1.33	0.33	5	ll	ll
1445	0.5	2.6	0.1	2.7	5.10	24.4	1.32	0.32	3	ll	ll
1450	0.5	3.1	0.1	2.7	5.11	24.4	1.32	0.31	2	ll	ll

MDA 4/14/08
ORP
56
88
45
24
12
9
5

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / Arcadis		SAMPLER(S) SIGNATURES: <i>[Signature]</i>		SAMPLING INITIATED AT: 1451		SAMPLING ENDED AT: 1455	
PUMP OR TUBING DEPTH IN WELL (feet):		SAMPLE PUMP FLOW RATE (mL per minute): <100 (9/14/08)		TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FILTER SIZE: 0.45 µm		DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-73	3	CG	40ml	HCl	—	—	8260	RFPP

MDA 4/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-74	SAMPLE ID: MW-74 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 5.5 feet to 32.5 feet	STATIC DEPTH TO WATER (feet): 2.46	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 35.5 feet) + 0.25 gallons = 0.34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 30	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 30	PURGING INITIATED AT: 1600	PURGING ENDED AT: 1615	TOTAL VOLUME PURGED (gallons): 1.80

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1609	1.08	1.08	0.12	3.62	6.49	24.1	468	0.50	0.84	clear	NONE	-136
1611	0.24	1.32	0.12	3.66	6.66	24.1	468	0.49	1.26	clear	NONE	-137
1613	0.24	1.56	0.12	3.68	6.66	24.1	467	0.48	0.26	clear	NONE	-139
1615	0.24	1.80	0.12	3.71	6.65	24.1	465	0.47	0.10	clear	NONE	-139

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STANTARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stantarcadis</i>	SAMPLING INITIATED AT: 1616	SAMPLING ENDED AT: 1622
PUMP OR TUBING DEPTH IN WELL (feet): 30	SAMPLE PUMP FLOW RATE (mL per minute): 2460	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-74	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-75	SAMPLE ID: MW-75	DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 49.5 feet to 44.5 feet	STATIC DEPTH TO WATER (feet): 5.5	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (49.5 feet - 5.5 feet) X 0.026 gallons/foot = 0.96 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 44.5 feet) + 0.25 gallons = 0.37 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 41	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 41	PURGING INITIATED AT: 9:12	PURGING ENDED AT: 9:40	TOTAL VOLUME PURGED (gallons): 2.0
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	GRP
9:20	0.8	0.8	0.1	5.5	5.59	23.8	0.468	0.76	28.6	clear	none	-1
9:25	0.5	1.3	0.1	5.51	5.45	23.9	0.458	0.53	18.8	"	"	3
9:30	0.5	1.8	0.1	5.5	5.49	23.7	0.459	0.42	17.3	"	"	5
9:35	0.5	2.3	0.1	5.5	5.51	23.7	0.456	0.41	13.8	"	"	9
9:40	0.5	2.8	0.1	5.5	5.62	23.7	0.450	0.38	14	"	"	2

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Mican Forbes ARCADIS		SAMPLER(S) SIGNATURES: <i>[Signature]</i>		SAMPLING INITIATED AT: 9:41	SAMPLING ENDED AT: 9:45
PUMP OR TUBING DEPTH IN WELL (feet): 41	MDS 4/19/08	SAMPLE PUMP FLOW RATE (mL per minute): <100	MDS 9/16/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: <input type="checkbox"/> µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW 75	3	CG	40ml	ACI			8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast		SITE LOCATION: Tallevast FL	
WELL NO: MW-76	SAMPLE ID: MW-76	DATE: 1/24/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 4.76	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 31 feet + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1423	PURGING ENDED AT: 1438	TOTAL VOLUME PURGED (gallons): 1.65

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or MS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	CRP
9 1432	0.99	0.99	0.11	6.39	5.3	26.4	.361	0.4	16.7	CLEAR	yes	30
2 1434	0.22	1.21	0.11	6.42	5.4	26.4	.319	0.4	19.9	CLEAR	yes	20
2 1436	0.22	1.43	0.11	6.44	5.4	26.4	.372	0.4	15.6	CLEAR	yes	9
2 1438	0.22	1.65	0.11	6.46	5.4	26.4	.377	0.4	11.3	CLEAR	yes	3

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1439	SAMPLING ENDED AT: 1445
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): ~410	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-76	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-77	SAMPLE ID: MW-77 DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 36 feet to 41 feet	STATIC DEPTH TO WATER (feet): 5.85	PURGE PUMP TYPE OR BAILER: Geo pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 64 feet) + 0.25 gallons = 0.42 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	PURGING INITIATED AT: 1055	PURGING ENDED AT: 1110	TOTAL VOLUME PURGED (gallons): 2.90

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or M/S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1104	1.26	1.26	0.14	8.75	6.42	24.3	1645	1.39	6.50	Clear	NONE-79
1106	0.28	1.54	0.14	8.95	6.43	24.3	1648	1.03	4.77	Clear	NONE-95
1108	0.28	1.82	0.14	9.03	6.44	24.3	1648	0.81	3.11	Clear	NONE-108
1110	0.28	2.10	0.14	9.07	6.45	24.3	1649	0.82	1.88	Clear	NONE-108

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1112	SAMPLING ENDED AT: 1122
PUMP OR TUBING DEPTH IN WELL (feet): 38.5	SAMPLE PUMP FLOW RATE (mL per minute): ~490	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N MDS 2/14/08	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-77	3	CG	40ml	HCl	—	—	8260	RFPP
DVP-8	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-78	SAMPLE ID: MW-78 DATE: 1-28-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 36 feet to 41 feet	STATIC DEPTH TO WATER (feet): 6.49	PURGE PUMP TYPE OR BAILER: PP								
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (41 feet - 6.49 feet) X 0.16 gallons/foot = 5.5 gallons												
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons												
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1246	PURGING ENDED AT: 1339	TOTAL VOLUME PURGED (gallons): 10								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1326	5.5	5.5	0.3*	7.33	6.54	25.85	0.742	1.33	2.93	none - crystal	none noted	-140
1331	1.5	7.0	↓	7.28	6.54	25.89	0.740	1.00	0.87	crystal	↓	-142
1336	1.5	8.5	↓	7.26	6.52	25.93	0.729	0.94	1.17	clear	↓	-142
1339	1.5	10.0	✓	7.26	6.52	25.89	0.727	0.93	1.17	↓	↓	-142
							mS/cm					
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016												

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ann Coats / Arcadis			SAMPLER(S) SIGNATURES: <i>[Signature]</i>			SAMPLING INITIATED AT: 1340	SAMPLING ENDED AT: 1345	
PUMP OR TUBING DEPTH IN WELL (feet):			SAMPLE PUMP FLOW RATE (mL per minute): <100			TUBING MATERIAL CODE: PE		
FIELD DECONTAMINATION: (Y) N			FIELD-FILTERED: Y (N) FILTER SIZE: 0.7µm			DUPLICATE: Y (N)		
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-78	3	CG	40mL	HCl	—	—	8260	RF PP
REMARKS: One bolt broken in hole								
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)								
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)								

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

*Purge rate was higher before readings. purged 5.5 gallons before 1st reading.

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tallerast	SITE LOCATION: Sarasota FL MDS 4/19/08
WELL NO: mw-79	SAMPLE ID: mw-79 DATE: 1-25-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 36 feet to 41 feet	STATIC DEPTH TO WATER (feet): 9.10	PURGE PUMP TYPE OR BAILER: Geo Pump								
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons												
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 44 feet) + 125 gallons = 0.36 gallons												
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	PURGING INITIATED AT: 1520	PURGING ENDED AT: 1526	TOTAL VOLUME PURGED (gallons): 1.76								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1520	1.10	1.10	0.11	10.16	6.4	25.7	701	0.5	3.46	Clear	NONE	ORP -81
1522	0.22	1.32	0.11	10.18	6.4	25.5	706	0.6	0.99	Clear	NONE	-89
1524	0.22	1.54	0.11	10.21	6.4	25.4	707	0.6	3.97	Clear	NONE	-89
1526	0.22	1.76	0.11	10.21	6.4	25.2	706	0.6	3.37	Clear	NONE	-91
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016												

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BUDGET STAHL / ARCADIS			SAMPLER(S) SIGNATURES: <i>Budget Stahl</i>			SAMPLING INITIATED AT: 1527		SAMPLING ENDED AT: 1534	
PUMP OR TUBING DEPTH IN WELL (feet): 38.5			SAMPLE PUMP FLOW RATE (mL per minute): ~416			TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N			FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm			DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH			
mw-79	3	CG	40mL	HCL	40mL	-	8260		RFPP
REMARKS:									
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)									
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump									
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)									

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

1-08

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-80	SAMPLE ID: MW-80
DATE: 1/30/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 36 feet to 41 feet	STATIC DEPTH TO WATER (feet): 8.77	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 44 feet) + 0.25 gallons = 0.36 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	PURGING INITIATED AT: 1002	PURGING ENDED AT: 1015	TOTAL VOLUME PURGED (gallons): 2.08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1009	1.12	1.12	0.16	9.42	6.88	28.6	1.27	0.64	1.09	Clear	NONE	-135
1011	0.32	1.44	0.16	9.45	6.86	28.7	1.23	0.61	1.19	Clear	NONE	-145
1013	0.32	1.76	0.16	9.46	6.86	28.7	1.22	0.61	0.19	Clear	NONE	-147
1015	0.32	2.08	0.16	9.45	6.86	28.7	1.22	0.60	0.16	Clear	NONE	-149

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/AREADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1016	SAMPLING ENDED AT: 1026
PUMP OR TUBING DEPTH IN WELL (feet): 38.5	SAMPLE PUMP FLOW RATE (mL per minute): ~610	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTRATION EQUIPMENT TYPE: <input checked="" type="radio"/> <input type="radio"/>	FILTER SIZE: _____ µm
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-80	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **sampled with transducer in well; missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA, FL
WELL NO: NW-81	SAMPLE ID: MW-81 DATE: 1/30/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 36 feet to 41 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 46 feet) + 0.5 gallons = 0.62 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	PURGING INITIATED AT: 1033	PURGING ENDED AT: 1054	TOTAL VOLUME PURGED (gallons): 2.31

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1045	1.32	1.32	0.11	11.01	7.11	26.38	865	1.04	0.74	clear yellow	YES	-255
1048	0.33	1.65	0.11	11.02	7.13	26.45	961	0.60	0.11	clear	YES	-272
1051	0.33	1.98	0.11	11.02	7.14	26.42	980	0.48	0.50	clear	YES	-276
1054	0.33	2.31	0.11	11.03	7.13	26.46	987	0.44	0.44	clear	YES	-277

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: S. Wells / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1058	SAMPLING ENDED AT: 1106
PUMP OR TUBING DEPTH IN WELL (feet): 38.5	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
NW-81	3	CG	40mL	HCl	40mL	—	B260B+B260BSM	RFPP

REMARKS: **TRANSOMER REMOVED FROM 0957 TO**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallewa st	SITE LOCATION: Tallewa st FL
WELL NO: MW-82	SAMPLE ID: MW-82 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 37 feet to 42 feet	STATIC DEPTH TO WATER (feet): 4.18	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 45 feet + 0.25 gallons = 0.37 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 39.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 39.5	PURGING INITIATED AT: 1427	PURGING ENDED AT: 1443	TOTAL VOLUME PURGED (gallons): 1.92

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle) (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1437	1.20	1.20	0.12	5.50	7.06	24.4	.602	0.58	1.79	Clear	NONE -174
1439	.24	1.44	0.12	5.51	7.08	24.4	.598	0.57	2.82	Clear	NONE -173
1441	.24	1.68	0.12	5.53	7.09	24.4	.594	0.56	5.01	Clear	NONE -173
1443	.24	1.92	0.12	5.55	7.10	24.4	.591	0.54	3.60	Clear	NONE -173

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BUDGET STATE/ARCADIS	SAMPLER(S) SIGNATURES: <i>Budget State</i>	SAMPLING INITIATED AT: 1444	SAMPLING ENDED AT: 1450
PUMP OR TUBING DEPTH IN WELL (feet): 39.5	SAMPLE PUMP FLOW RATE (mL per minute): ~450	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
NW-82	3	CG	40ml	Hel	—	—	8260	RFPP

REMARKS: **missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Former ABC Facility</i>	SITE LOCATION: <i>Tallahassee FL</i>
WELL NO: <i>MW-83</i>	SAMPLE ID: <i>MW-83</i> DATE: <i>1.24.08</i>

PURGING DATA

WELL DIAMETER (inches): <i>2</i>	TUBING DIAMETER (inches): <i>1/4</i>	WELL SCREEN INTERVAL DEPTH: <i>102</i> feet to <i>112</i> feet	STATIC DEPTH TO WATER (feet): <i>13.6</i>	PURGE PUMP TYPE OR BAILER: <i>PP</i>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <i>1446</i>	PURGING ENDED AT: <i>1520</i>	TOTAL VOLUME PURGED (gallons): <i>3.0</i> MDS 4/14/08
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<i>1450</i>	<i>0.5</i>	<i>0.5</i>	<i>0.1</i>	<i>13.6</i>	<i>7.18</i>	<i>23.8</i>	<i>0.365</i>	<i>0.94</i>	<i>12</i>	<i>clear</i>	<i>none</i>	<i>ORP -195</i>
<i>1455</i>	<i>0.5</i>	<i>1.0</i>	<i>0.1</i>	<i>13.6</i>	<i>7.25</i>	<i>23.9</i>	<i>0.370</i>	<i>0.42</i>	<i>13</i>	<i>"</i>	<i>"</i>	<i>-227</i>
<i>1510</i>	<i>0.5</i>	<i>2.0</i>	<i>0.1</i>	<i>13.6</i>	<i>7.28</i>	<i>23.81</i>	<i>0.369</i>	<i>0.32</i>	<i>30</i>	<i>"</i>	<i>"</i>	<i>-271</i>
<i>1515</i>	<i>0.5</i>	<i>2.5</i>	<i>0.1</i>	<i>13.6</i>	<i>7.28</i>	<i>23.82</i>	<i>0.369</i>	<i>0.32</i>	<i>34</i>	<i>"</i>	<i>"</i>	<i>-275</i>
<i>1520</i>	<i>0.5</i>	<i>3.0</i>	<i>0.1</i>	<i>13.6</i>	<i>7.28</i>	<i>23.89</i>	<i>0.369</i>	<i>0.30</i>	<i>38</i>	<i>"</i>	<i>"</i>	<i>-279</i>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 6.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Micah Forbes / ARCADIS</i>	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: <i>1521</i>	SAMPLING ENDED AT: <i>1525</i>
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PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <i><100</i> MDS 4/14/08	TUBING MATERIAL CODE:
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: <i>1</i> µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>MW83</i>	<i>3</i>	<i>CG</i>	<i>40ml</i>	<i>HCl</i>	<i>-</i>	<i>-</i>	<i>8260</i>	<i>RFPP</i> MDS 4/14/08

REMARKS: *missed two readings on the phone*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 9/14/08	SITE LOCATION: Sarasota FL MDS 9/14/08
WELL NO: MW-84	SAMPLE ID: MW-84 DATE: 1.25.08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 35.5 feet to 40.5 feet	STATIC DEPTH TO WATER (feet): 9.38	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) ← 0 feet - feet X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 40 feet) + 0.25 gallons = 0.35 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1432	PURGING ENDED AT: 1454	TOTAL VOLUME PURGED (gallons): 2.2

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1442	1.0	1.0	0.1	9.75	6.73	27.21	1.28	1.49	0.07	NONE	none noted
1445	0.3	1.3	↓	9.75	6.68	27.22	1.20	1.25	0.09		
1448	1	1.6	↓	9.75	6.65	27.24	1.25	1.16	0.07		
1451	1	1.9	↓	9.75	6.65	27.26	1.25	1.25 ^{1.12}	0.07		
1454	1	2.2	↓	9.75	6.65	27.20	1.25	1.09	0.05		
							MScm				

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ann Peaty Arcadis	SAMPLER(S) SIGNATURES: <i>Ann Peaty</i>	SAMPLING INITIATED AT: 1455	SAMPLING ENDED AT: 1500
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): < 100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 0.45 µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-84	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

1.2

SITE NAME: Tallevast	SITE LOCATION: Tallevast
WELL NO: MW-85	SAMPLE ID: MW-85 DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 50 feet to 55 feet	STATIC DEPTH TO WATER (feet): 544	PURGE PUMP TYPE OR BAILER: Geopump (PP)
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 58 feet) + 0.25 gallons = 0.40 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 52.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 52.5	PURGING INITIATED AT: 1525	PURGING ENDED AT: 1541	TOTAL VOLUME PURGED (gallons): 1.92							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1535	1.20	1.20	0.12	6.40	7.2	25.8	0.783	0.5	1.52	clear	yes
1537	0.24	1.44	0.12	6.40	7.2	25.8	0.796	0.4	0.77	clear	yes
1539	0.24	1.68	0.12	6.40	7.2	25.8	0.804	0.4	1.01	clear	yes NO
1541	0.24	1.92	0.12	6.40	7.2	25.8	0.814	0.4	1.01	clear	yes NO

ORP
-84
-92
-98
-101

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / Arcadis	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1544	SAMPLING ENDED AT: 1452
PUMP OR TUBING DEPTH IN WELL (feet): 52.5	SAMPLE PUMP FLOW RATE (mL per minute): ~454	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: Y <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-85	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

TB-14-1130

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAH	SITE LOCATION: TALLEVAH, FL
WELL NO: MW-11086	SAMPLE ID: MW-11086 DATE: 1-28-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 10.95	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (28 feet - 10.95 feet) X 0.16 gallons/foot = 2.73 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 36 feet) + .25 gallons = .34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1004	PURGING ENDED AT: 1018	TOTAL VOLUME PURGED (gallons): 1.82

8

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle right) or % saturation	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1012	1.04	1.04	0.13	13.55	7.06	24.1	0.562	0.56	3.60	Clear	yes
1014	0.26	1.30	0.13	13.73	7.16	24.1	0.553	0.54	3.29	Clear	yes
1016	0.26	1.56	0.13	13.81	7.18	24.2	0.549	0.54	3.52	Clear	yes
1018	0.26	1.82	0.13	13.89	7.17	24.1	0.547	0.53	2.87	Clear	yes

ORP
-209
-207
-205
-204

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT)/AFFILIATION: Jennifer King/ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1020	SAMPLING ENDED AT: 1028
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): 490	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW110	3	CG	40ml	Hel	40ml	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-87	SAMPLE ID: MW-87 DATE: 1-30-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 0.36 feet to 41 feet	STATIC DEPTH TO WATER (feet): 4.8	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.026 gallons/foot X 41 feet) + 0.25 gallons = 0.35 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38.5	PURGING INITIATED AT: 8:32	PURGING ENDED AT: 8:50	TOTAL VOLUME PURGED (gallons): 1.9

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
8:35	0.4	0.4	0.1	5.3	6.88	23.8	1.33	1.10	0	clear	none	-183
8:40	0.5	0.9	0.1	5.4	6.70	23.9	1.34	0.38	0	"	"	-188
8:45	0.5	0.4 0.9	0.1	5.4	6.70	24.01	1.35	0.35	0	"	"	-189
8:50	0.5	0.4 1.4	0.1	5.4	6.70	24.07	1.36	0.33	0	"	"	-190
		1.9										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS		SAMPLER(S) SIGNATURES: <i>[Signature]</i>		SAMPLING INITIATED AT: 8:51	SAMPLING ENDED AT: 8:55
PUMP OR TUBING DEPTH IN WELL (feet): 38.5 MRS 9/14/08		SAMPLE PUMP FLOW RATE (mL per minute): 5100 MDS 9/14/08		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm		DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-87	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.44

SITE NAME: TALEVAST		SITE LOCATION: TALEVAST FI	
WELL NO: mw-88	SAMPLE ID: mw-88	DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 76 feet to 86 feet	STATIC DEPTH TO WATER (feet): 12.38	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 89 feet) + .25 gallons = .48 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 81	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 81	PURGING INITIATED AT: 1454	PURGING ENDED AT: 1515	TOTAL VOLUME PURGED (gallons): 2.10
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	OCF
1509	1.50	1.50	0.16	19.74	8.42	24.7	.124	1.77	14.6	Clear	NONE	-131
1511	0.20	1.70	0.10	20.08	8.40	24.7	.130	1.60	19.8	Clear	NONE	-137
1513	0.20	1.90	0.10	20.51	8.38	24.7	.136	1.06	17.2	Clear	NONE	-147
1515	0.20	2.10	0.10	20.84	8.37	24.7	.141	1.18	19.3	Clear	NONE	-151

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAN / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stan</i>	SAMPLING INITIATED AT: 1516	SAMPLING ENDED AT: 1519
PUMP OR TUBING DEPTH IN WELL (feet): 81	SAMPLE PUMP FLOW RATE (mL per minute): 390	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-88	3	CG	40ml	HEI	—	—	P260	RFPP

REMARKS: **MISSING 2 BOLTS**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: Sarasota FL MDS 4/19/08
WELL NO: MW-89	SAMPLE ID: MW-89 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 27 feet to 32 feet	STATIC DEPTH TO WATER (feet): 4.45	PURGE PUMP TYPE OR BAILER: PP Geo Pump
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable) = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = **0** gallons + (**0.0026** gallons/foot X **35** feet) + **.25** gallons = **0.34** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 29.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 29.5 MDS 4/19/08	PURGING INITIATED AT: 1357	PURGING ENDED AT: 1412	TOTAL VOLUME PURGED (gallons): 1.92
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1405	1.12	1.12	0.14	6.25	6.8	25.2	0.673	0.5	2.56	clear	yes
1407	.28	1.40	0.14	6.25	6.8	25.1	0.674	0.8	3.70	clear	yes
1409	.28	1.68	0.14	6.20	6.8	25.1	0.673	0.9	2.68	clear	yes
1411	.28	1.92	0.14	6.20	6.8	25.1	0.669	0.9	2.85	clear	yes

ORP
-132
-134
-136
-136

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1414	SAMPLING ENDED AT: 1417
PUMP OR TUBING DEPTH IN WELL (feet): 29.5 MDS 4/19/08	SAMPLE PUMP FLOW RATE (mL per minute): 530	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-89	3	CG	40mL	HCL	40mL	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tallevast	SITE LOCATION: Sarasota FL MDS 2/14/08
WELL NO: MW-90	SAMPLE ID: MW-90 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 25.5 feet to 30.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: Geopump (PP)
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 33.5 feet) + 0.25 gallons = 0.34 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 28	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 28	PURGING INITIATED AT: 0937	PURGING ENDED AT: 0953	TOTAL VOLUME PURGED (gallons): 1.80
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm) (µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
0946	1.08	1.08	0.12	4.97	6.7	24.5	.613	1.1	0.40	Clear	NONE	-64
0948	.24	1.32	0.12	5.02	6.7	24.5	.612	0.7	0.56	Clear	NONE	-73
0950	.24	1.56	0.12	5.04	6.7	24.5	.612	0.6	0.52	CLEAR	NONE	-78
0952	.24	1.80	0.12	5.05	6.7	24.5	.613	0.6	1.33	CLEAR	NONE	-80

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 0955	SAMPLING ENDED AT: 1002
PUMP OR TUBING DEPTH IN WELL (feet): 28	SAMPLE PUMP FLOW RATE (mL per minute): ~455	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input type="radio"/> N <input checked="" type="radio"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-90	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump

EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Form ABC Facility</u>		SITE LOCATION: <u>Tallahassee, FL</u>	
WELL NO: <u>NW-91</u>	SAMPLE ID: <u>MW-91</u>	DATE: <u>1-25-05</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>4"</u>	WELL SCREEN INTERVAL DEPTH: <u>37.5</u> feet to <u>37.9</u> feet	STATIC DEPTH TO WATER (feet): <u>7.75</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (<u>37.5</u> feet - <u>7.75</u> feet) X <u>.16</u> gallons/foot = <u>4.76</u> gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>37</u> feet) + <u>0.25</u> gallons = <u>0.35</u> gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>37.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>35</u>	PURGING INITIATED AT: <u>9:12</u>	PURGING ENDED AT: <u>9:29</u>	TOTAL VOLUME PURGED (gallons): <u>1.7</u>							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (μ mhos/cm or μ S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:23	1.1	1.1	0.1	12.95	6.52	23.74	0.669	0.88	3.11	clear	none
9:26	0.3	1.4	0.1	13.83	6.50	23.65	0.667	1.0	3.91	↓	↓
9:29	0.3	1.7	0.1	14.17	6.49	23.82	0.603	1.01	3.15	↓	↓
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.008; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Ami Coats / Arcadis</u>			SAMPLER(S) SIGNATURES: <u>Ami Coats</u>			SAMPLING INITIATED AT: <u>9:30</u>			SAMPLING ENDED AT: <u>9:35</u>		
PUMP OR TUBING DEPTH IN WELL (feet): <u>35'</u>			SAMPLE PUMP FLOW RATE (mL per minute): <u>< 100</u>			TUBING MATERIAL CODE: <u>pe</u>					
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N			FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N			FILTER SIZE: <u>0.1</u> μ m			DUPLICATE: <input checked="" type="radio"/> <input type="radio"/> N		
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH					
<u>MW-91</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>---</u>	<u>---</u>	<u>8260</u>		<u>RFPP</u>		
REMARKS: <u>transducer reinserted 9:57</u> <u>transducer removed @ 9:07</u> / <u>7.75 liter removed total</u>											
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)											
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump											
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)											

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: \pm 0.2 units Temperature: \pm 0.2 °C Specific Conductance: \pm 5% Dissolved Oxygen: all readings \leq 20% saturation (see Table FS 2200-2); optionally, \pm 0.2 mg/L or \pm 10% (whichever is greater) Turbidity: all readings \leq 20 NTU; optionally \pm 5 NTU or \pm 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: MW-92	SAMPLE ID: MW-92 DATE: 1-31-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 32.5 feet to 37.5 feet	STATIC DEPTH TO WATER (feet): 11.9	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 37.5 feet) + 0.25 gallons = 0.34 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 35	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 35	PURGING INITIATED AT: 8:58	PURGING ENDED AT: 9:20	TOTAL VOLUME PURGED (gallons): 2.2
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:05	0.7	0.7	0.1	12.6	7.16	23.57	0.54	0.65	0	clear	none
9:10	0.5	1.2	0.1	12.75	7.25	23.8	0.53	0.47	0	"	"
9:15	0.5	1.7	0.1	12.75	7.28	23.8	0.54	0.44	0	"	"
9:20	0.5	2.2	0.1	12.75	7.28	23.8	0.53	0.41	0	"	"

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT:	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 35	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-92	3	CG	40ml	HCl	—	—	8260	R.FPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

25-25

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 4/14/08		SITE LOCATION: <u>Sarasota FL</u> MDS 4/14/08	
WELL NO: <u>MW-93</u>	SAMPLE ID: <u>MW-93</u>	DATE: <u>1/28/08</u>	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>32.5</u> feet to <u>37.5</u> feet	STATIC DEPTH TO WATER (feet): <u>12.46</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u> MDS 4/14/08
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (<u>0.0026</u> gallons/foot X <u>37</u> feet) + <u>0.25</u> gallons = <u>0.34</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>37</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>37</u>	PURGING INITIATED AT: <u>1035</u>	PURGING ENDED AT: <u>1050</u>	TOTAL VOLUME PURGED (gallons): <u>2.5</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
10:38	0.3	0.3	0.1	13.62	6.95	24.52	0.777	1.53	light yellow	light brown	-	-131
10:41	0.3	0.6	0.1	13.99	7.05	24.47	0.776	1.21	1.74	clear	-	-181
10:44	0.3	0.9	0.1	14.23	7.09	24.47	0.776	1.17	4.80	"	-	-210
10:47	0.3	1.2	0.1	14.35	7.11	24.46	0.780	1.15	1.22	"	-	-223
10:50	0.3	1.5	0.1	14.45	7.13	24.45	0.778	1.15	1.58	"	-	-238
MDS 4/14/08												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>NATHAN SMITH</u>	SAMPLER(S) SIGNATURES: <u>Nathan P. Smith</u>	SAMPLING INITIATED AT: <u>1050</u>	SAMPLING ENDED AT: <u>1055</u>
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PUMP OR TUBING DEPTH IN WELL (feet): <u>37</u> MDS 4/14/08	SAMPLE PUMP FLOW RATE (mL per minute): <u>L100</u> MDS 4/14/08	TUBING MATERIAL CODE:
FIELD DECONTAMINATION: <u>Y</u> MDS 4/14/08	FIELD-FILTERED: <u>Y</u> MDS 4/14/08	DUPLICATE: <u>Y</u> MDS 4/14/08

SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-93	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Sarasota FL	2/14/08 MBS
WELL NO: MW-94	SAMPLE ID: MW-94	DATE: 1/22/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 24.5 feet to 29.5 feet	STATIC DEPTH TO WATER (feet): 3.17	PURGE PUMP TYPE OR BAILER: geopump (PP)
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

= gallons + (0.0026 gallons/foot X 29 feet) + 0.25 gallons = 0.33 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 27	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 27	PURGING INITIATED AT: 1607	PURGING ENDED AT: 1623	TOTAL VOLUME PURGED (gallons): 1.80
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle <input checked="" type="checkbox"/> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1616	1.08	1.08	0.12	3.50	6.0	23.4	.476	.7	6.04	Clear	NONE	-42
1618	.24	1.32	0.12	3.54	6.0	23.4	.464	.7	2.03	Clear	NONE	-48
1620	.24	1.56	0.12	3.55	6.0	23.4	.470	.7	3.38	Clear	NONE	-52
1622	.24	1.80	0.12	3.55	6.0	23.4	.470	.7	2.79	Clear	NONE	-58

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Bridget Stahl TARCADIS	SAMPLER(S) SIGNATURES: Bridget Stahl	SAMPLING INITIATED AT: 1628	SAMPLING ENDED AT: 1630
PUMP OR TUBING DEPTH IN WELL (feet): 27	SAMPLE PUMP FLOW RATE (mL per minute): ~455	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input checked="" type="radio"/> N	FIELD-FILTERED: Y <input checked="" type="radio"/> N	FILTRATION EQUIPMENT TYPE: _____	DUPLICATE: Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-94	3	CG	40mL	HCL	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tallevast	SITE LOCATION: Sarasota FL MDS 4/14/08
WELL NO: MW-95	SAMPLE ID: MW-95 DATE: 1/22/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 2.96	PURGE PUMP TYPE OR BAILER: Geopump (PP)
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 27.5 feet) + 0.25 gallons = 0.32 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1522	PURGING ENDED AT: 1538	TOTAL VOLUME PURGED (gallons): 1.6

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1532	1.0	1.0	0.1	3.20	6.2	23.6	429	0.7	3.54	Clear	NONE
1534	0.2	1.2	0.1	3.21	6.2	23.6	426	0.5	1.81	clear	NONE
1536	0.2	1.4	0.1	3.21	6.1	23.6	424	0.5	0.50	Clear	NONE
1538	0.2	1.6	0.1	3.21	6.1	23.6	424	0.4	2.05	Clear	NONE

ORP
-35
-43
-51
-57

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1540	SAMPLING ENDED AT: 1547
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): ~380	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>	
SAMPLE CONTAINER SPECIFICATION		SAMPLE PRESERVATION	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME
PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD
SAMPLING EQUIPMENT CODE			
MW-95	3	CG	40mL
			HCL
			-
			-
			8260
			RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tallahassee	SITE LOCATION: Sarasota, FL
WELL NO: MW-96	SAMPLE ID: MW-96 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 196 feet to 206 feet	STATIC DEPTH TO WATER (feet): 14.5	PURGE PUMP TYPE OR BAILER: pp/perist pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 210 feet) + 0.25 gallons = 0.25 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~201	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~201	PURGING INITIATED AT: 1035	PURGING ENDED AT: 1105	TOTAL VOLUME PURGED (gallons): 2.24

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mol or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1039	0.32	0.32	0.08	18.22	7.37	22.76	422	0.43	24.9	clear	yes	0.28
1045	0.48	0.64	0.08	18.45	7.24	22.77	313	2.32	22.3	clear	yes	-267
1050	0.40	1.04	0.08	19.09	7.18	22.32	709	0.82	23.7	clear	yes	-277
1055	0.40	1.44	0.08	19.86	7.21	22.36	309	0.55	12.7	clear	yes	-287
1100	0.40	1.84	0.08	20.61	7.20	22.38	308	0.46	6.94	clear	yes	-294
1105	0.40	2.24	0.08	21.41	7.20	22.41	308	0.41	5.34	clear	yes	-297
<i>ended 1/29/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 6.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1108	SAMPLING ENDED AT: 1115
PUMP OR TUBING DEPTH IN WELL (feet): ~201	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-96	3	CG	400mL	HCL	-	~2	8260B + 8260J in RFP	

REMARKS: **Initial purge - Blue color then cleared - Placed new 2" well cap & Dolphin lock on well** *needs bolts for cover*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC Facility</u>	SITE LOCATION: <u>Tallahassee, FL</u>
WELL NO: <u>MW-97</u>	SAMPLE ID: <u>MW-97</u> DATE: <u>1-30-08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>208.6</u> feet to <u>226.6</u> feet	STATIC DEPTH TO WATER (feet): <u>4.8</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>226.6</u> feet) + <u>0.25</u> gallons = <u>0.84</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>220</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>220</u>	PURGING INITIATED AT: <u>11:58</u>	PURGING ENDED AT: <u>12:30</u>	TOTAL VOLUME PURGED (gallons): <u>3.4</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
12:05	0.9	0.9	0.1	19.4	7.60	23.09	0.625	0.32	23	H. gray	none
12:10	0.5	1.4	0.1	19.4	7.59	23.6	0.604	0.31	24	clear	none
12:15	0.5	1.9	0.1	21.9	7.44	23.7	0.267	0.29	28	"	none
12:20	0.5	2.4	0.1	23.6	7.26	23.7	0.173	0.30	19	"	"
12:25	0.5	2.9	0.1	24.4	7.28	24.0	0.167	0.31	8	"	"
12:30	0.5	3.4	0.1	25.5	7.23	24.06	0.185	0.30	0	"	"

ORP
-309
-309
-295
-274
-265
-267

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nicah Forbes / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>12:31</u>	SAMPLING ENDED AT: <u>12:35</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>220</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>1100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTRATION EQUIPMENT TYPE: <u>L100</u>	FILTER SIZE: <u>0.45µm</u>
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-97	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: lid cap for well is damaged

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-98	SAMPLE ID: MW-98	DATE: 1.25.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 33 feet to 38 feet	STATIC DEPTH TO WATER (feet): 4.3	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 13:45	PURGING ENDED AT: 14:05	TOTAL VOLUME PURGED (gallons): 2.0
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1350	0.5	0.5	0.1	4.9	6.46	24.19	1.28	1.15	6.2	clear	none
1355	0.5	1.0	0.1	4.9	6.47	24.2	1.28	0.66	4	"	"
1400	0.5	1.5	0.1	4.9	6.48	24.3	1.28	0.57	4.52	"	"
1405	0.5	2.0	0.1	6.1	6.97	24.4	1.29	0.50	4	"	"

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 14:06	SAMPLING ENDED AT: 14:10
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N	FILTER SIZE: µm	DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW98	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEYST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-99	SAMPLE ID: MW-99
DATE: 1/25/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 71 feet to 81 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 86 feet) + 0.5 gallons = 0.72 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 76	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 76	PURGING INITIATED AT: 1235	PURGING ENDED AT: 1255	TOTAL VOLUME PURGED (gallons): 1.40

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1246	0.77	0.77	0.07	14.62	7.42	26.05	416	0.55	11.9	clear	yes	-257
1249	0.21	0.98	0.07	15.10	7.42	26.05	416	0.52	12.6	clear	yes	-258
1252	0.21	1.19	0.07	15.34	7.42	25.92	417	0.46	13.4	clear	yes	-259
1255	0.21	1.40	0.07	15.72	7.41	26.03	416	0.46	15.1	clear	yes	-261

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NS/SW/ ARCADIS	SAMPLER(S) SIGNATURES: <i>Sa Wlu</i>	SAMPLING INITIATED AT: 1257	SAMPLING ENDED AT: 1303
PUMP OR TUBING DEPTH IN WELL (feet): 76	SAMPLE PUMP FLOW RATE (mL per minute): ≤100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-99	3	CG	40mL	HCl	40mL	—	8260B+8260BSM	RFPP
REP-10	3	CG	40mL	HCl	40mL	—	8260B+8260BSM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallahassee</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>MW-100</u>	SAMPLE ID: <u>MW-100</u> DATE: <u>1/25/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>24.5</u> feet to <u>29.5</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH -- STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)	=	gallons + (<u>00.0020</u> gallons/foot X <u>33</u> feet) + <u>0.50</u> gallons = <u>0.58</u> gallons
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INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~26</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~26</u>	PURGING INITIATED AT: <u>1240</u>	PURGING ENDED AT: <u>1255</u>	TOTAL VOLUME PURGED (gallons): <u>3.15</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1243	0.63	0.63	0.21	6.29	5.99	27.46	339	1.04	6.0	Clear	Yes
1246	0.63	1.26	0.21	6.31	6.00	27.57	344	0.82	4.14	Clear	Yes
1249	0.63	1.89	0.21	6.31	6.06	27.71	356	0.76	3.91	Clear	Yes
1252	0.63	2.52	0.21	6.31	6.12	27.70	364	0.82	3.14	Clear	Yes
1255	0.63	3.15	0.21	6.31	6.14	27.76	364	0.83	2.59	Clear	Yes

-79
-86
-91
-98
-100

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>J. Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>1256</u>	SAMPLING ENDED AT: <u>1300</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~26</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>~100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y N</u>	FIELD-FILTERED: <u>Y</u> FILTER SIZE: <u> </u> µm	DUPLICATE: <u>Y X</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-100</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u>-</u>	<u>~2</u>	<u>8260B+8260S1M RFP</u>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1,2

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST FL	
WELL NO: mw-101	SAMPLE ID: mw-101	DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 57.7 feet to 57.7 feet	STATIC DEPTH TO WATER (feet): 60.7	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 60.7 feet) + .25 gallons = 0.40 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 55	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 55	PURGING INITIATED AT: 0818	PURGING ENDED AT: 0832	TOTAL VOLUME PURGED (gallons): 2.38

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0826	1.36	1.36	0.17	8.10	7.5	23.9	.420	0.6	20.4	Clear	yes
0828	0.34	1.70	0.17	8.13	7.5	24.0	.420	0.6	3.75	clear	yes
0830	0.34	2.04	0.17	8.18	7.5	24.0	.419	0.5	3.98	clear	yes
0832	0.34	2.38	0.17	8.21	7.4	24.0	.419	0.5	2.29	clear	yes

ORP
-133
-138
-144
-148

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL/ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 0833	SAMPLING ENDED AT: 0845
PUMP OR TUBING DEPTH IN WELL (feet): 55	SAMPLE PUMP FLOW RATE (mL per minute): ~643	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-101	3	CG	40ml	HCl	40ml	-	8260	RFPP

REMARKS: **missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-102	SAMPLE ID: MW-102 DATE: 1.25.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 89.8 feet to 99.8 feet	STATIC DEPTH TO WATER (feet): 12.9	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026** gallons/foot X feet) + **0.25** gallons = gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 9:53	PURGING ENDED AT: 10:25	TOTAL VOLUME PURGED (gallons): 3.1
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:55	0.1	0.1	0.1	12.9	7.54	24.17	0.528	3.41	260	clear	none
10:00	0.5	0.6	0.1	12.9	7.2	24.4	0.618	0.75	87	"	"
10:05	0.5	1.1	0.1	12.9	7.19	24.4	0.644	0.59	63	"	"
10:10	0.5	1.6	0.1	12.9	7.19	24.4	0.661	0.51	32	"	"
10:15	0.5	2.1	0.1	12.9	7.18	24.5	0.667	0.47	18	"	"
10:20	0.5	2.6	0.1	12.9	7.19	24.5	0.673	0.45	11	"	"
10:25	0.5	3.1	0.1	12.9	7.19	24.6	0.675	0.41	7	"	"

ORP
-198
-234
-244
-250
-253
-257
-260

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 6.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:26	SAMPLING ENDED AT: 10:30
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): < 100 ^{MPS} _{9/14/08}	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N FILTER SIZE: µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW102	3	CG	40ml	HCT	—	—	826B	RFPP

MPS
9/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallahassee, FL
WELL NO: MW-103	SAMPLE ID: MW-103 DATE: 1.25.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL (feet): 25 to 30	STATIC DEPTH TO WATER (feet): 4.3	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (05 feet - 0 feet) X _____ gallons/foot = _____ gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.026 gallons/foot X _____ feet) + 0.25 gallons = _____ gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 10:40	PURGING ENDED AT: 11:10	TOTAL VOLUME PURGED (gallons): 2.8

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	GRP
10:45	0.3	0.3	0.1	4.3	6.18	23.9	1.28	2.6	27	sl turbid	none	-37
10:50	0.5	0.8	0.1	4.5	5.98	24.2	1.33	1.36	38	"	"	-128
10:55	0.5	1.3	0.1	4.5	6.07	24.3	1.36	0.70	46	"	"	-150
11:00	0.5	1.8	0.1	4.5	6.12	24.4	1.35	0.58	31	"	"	-151
11:05	0.5	2.3	0.1	4.5	6.16	24.6	1.36	0.48	35	"	"	-156
11:10	0.5	2.8	0.1	4.5	6.19	24.6	1.36	0.43	37	"	"	-159

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT)/AFFILIATION: Nicah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 11:11	SAMPLING ENDED AT: 11:15
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/08	TUBING MATERIAL CODE: PE MPS 4/14/08	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: _____ µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-103	3	CG	40ML	HCl	—	—	8260	RFPP

REMARKS: **Missing tubing, only ~ ten feet deep, used disposable poly. for tubing**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee FL
WELL NO: MW-104	SAMPLE ID: MW-104 DATE: 1.28.08

PURGING DATA

WELL DIAMETER (Inches): 2	TUBING DIAMETER (Inches): 1/4	WELL SCREEN INTERVAL DEPTH: 25.9 feet to 30.4 feet	STATIC DEPTH TO WATER (feet): 3.15	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 10:43	PURGING ENDED AT: 11:10	TOTAL VOLUME PURGED (gallons): 2.7 MDS 4/14/08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
10:50	0.7	0.7	0.1	3.15	5.35	23.6	0.9	0.58	2.1	clear	none	ORP -57
10:55	0.5	1.2	0.1	3.15	5.39	23.6	1.0	0.49	5.5	"	"	-61
11:00	0.5	1.7	0.1	3.15	5.32	23.8	2.61	0.36	2.5	"	"	-60
11:05	0.5	2.2	0.1	3.15	5.29	23.8	*	0.35	2.16	"	"	-62
11:10	0.5	2.7	0.1	3.15	5.29	23.9	*	0.34	1.94	"	"	-64

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 11:11	SAMPLING ENDED AT: 11:15
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-104	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **conductivity meter readings very strange, may be malfunctioning**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

EB-21 @ 1650

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-105	SAMPLE ID: MW-105	DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 41.8 feet to 46.8 feet	STATIC DEPTH TO WATER (feet): 4.7	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 9:30	PURGING ENDED AT: 10:05	TOTAL VOLUME PURGED (gallons): 2.8
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (umhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:35	0.4	0.4	0.1	4.7	6.89	22.8	0.478	0.79	9.94	clear	none
9:40	0.4	0.8	0.1	4.7	6.91	22.9	0.741	0.43	4.9	"	"
9:45	0.4	1.2	0.1	4.9	6.99	23.0	0.9	0.48	2.2	"	"
9:50	0.4	1.6	0.1	4.9	7.10	23.0	0.9	0.37	0	"	"
9:55	0.4	2.0	0.1	5.05	7.11	23.1	2.2	0.36	0	"	"
10:00	0.4	2.4	0.1	5.05	7.12	23.0	3.8	0.36	0	"	"
10:05	0.4	2.8	0.1	5.05	7.16	23.0	*	0.32	0	"	"

MDS 2/14/08

ORP

-178
-199
-205
-217
-225
-228
-232

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes ARADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:06	SAMPLING ENDED AT: 10:10
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: µm	DUPLICATE: Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-105	3	CG	40ml	Hel	—	—	8260	RFPP

MDS 2/14/08

REMARKS: **conductivity meter readings very strange. The meter may be malfunctioning.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALEFAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-106	SAMPLE ID: MW-106 DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 40 feet to 45 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 50 feet) + 0.5 gallons = 0.63 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 42.5		FINAL PUMP OR TUBING DEPTH IN WELL (feet): 42.5		PURGING INITIATED AT: 0953		PURGING ENDED AT: 1010		TOTAL VOLUME PURGED (gallons): 1.36			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circ mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1001	0.64	0.64	0.08	7.40	6.71	24.70	293	1.21	5.79	clear	no
1004	0.24	0.88	0.08	7.40	6.82	24.68	334	0.83	4.38	clear	no
1007	0.24	1.12	0.08	7.41	6.89	24.67	349	0.79	2.86	clear	no
1010	0.24	1.36	0.08	7.40	6.99	24.68	339	1.08	2.08	clear	no

0.02
2
-39
-80

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/NS / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1012	SAMPLING ENDED AT: 1018
PUMP OR TUBING DEPTH IN WELL (feet): 42.5	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
NW-106	3	CG	40mL	HCl	40mL	—	8260+82605m	RFP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 4/19/08	SITE LOCATION: <u>Sarasota FL</u> MDS 4/19/08
WELL NO: <u>MW-107</u>	SAMPLE ID: <u>MW-107</u> DATE: <u>1/24/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>21</u> feet to <u>26</u> feet	STATIC DEPTH TO WATER (feet): <u>3.13</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<u>0.0026</u> gallons/foot X <u>26</u> feet) + <u>0.25</u> gallons = <u>0.32</u> gallons				

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1355	0.5	0.5	0.1	3.77	6.6	22.7	2.94	4.3	6.88	Clear	Cows
1400	0.5	1.0	↓	3.83	6.6	23.4	2.94	3.5	4.06	↓	↓
1405	0.5	1.5	↓	3.89	6.6	23.5	2.94	3.1	2.67	↓	↓
1410	0.5	2.0	↓	3.90	6.6	23.8	2.94	3.0	2.00	↓	↓
1415	0.5	2.5	↓	3.93	6.6	24.2	2.95	2.9	1.57	↓	↓
1420	→ <u>Sampled @ 1420</u> ←										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>IARLADES</u>	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: <u>1440</u>	SAMPLING ENDED AT: <u>1450</u>
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u> MDS 4/19/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-107	3	CG	40ml	HCL	—	—	8260	RFPP

REMARKS: * Agricultural odors overwhelming GW odor, initial purge.

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast MDS 9/19/08	SITE LOCATION: Sarasota FL MDS 9/19/08
WELL NO: MW-108	SAMPLE ID: MW-108 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 3.58	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= \emptyset gallons + (0.0026 gallons/foot X 28 feet) + .25 gallons = 0.32 gallons MDS 9/19/08				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25	PURGING INITIATED AT: 1512	PURGING ENDED AT: 1527	TOTAL VOLUME PURGED (gallons): 1.4
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1515	.5	.5	0.1	4.25	5.9	26.7	.55	3.5	8.63	clear	none	145
1518	.3	.8	0.1	4.25	5.8	26.7	.54	5.2	4.37	↓	↓	115
1521	.3	1.1	0.1	4.25	5.8	26.7	.54	5.0	2.27	↓	↓	110
1524	.3	1.4	0.1		5.8	26.8	.54	4.9	1.93	↓	↓	108
1527	→ Sampled @ 1527 ←											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KASU CORNWELL AREADIS MDS 9/19/08	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1527	SAMPLING ENDED AT: 1532
PUMP OR TUBING DEPTH IN WELL (feet): 25	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-108	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: initial purge clear w/ no noticeable odor.

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

*needs 2 Bolts.

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEUDST		SITE LOCATION: TALLEN AVE, FL	
WELL NO: mw-109	SAMPLE ID: mw-109	DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/2	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 4.35	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 36 feet) + .25 gallons = 0.34 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1510	PURGING ENDED AT: 1524	TOTAL VOLUME PURGED (gallons): 1.82

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mL or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8 1518	1.04	1.04	0.13	7.26	6.50	25.1	0.95	0.85	22.3	clear	NONE -39
1520	0.26	1.30	0.13	7.96	6.50	25.1	0.93	0.90	16.8	clear	NONE -49
1522	0.26	1.56	0.13	8.37	6.49	25.2	0.91	0.74	13.5	clear	NONE -60
1524	0.26	1.82	0.13	8.69	6.50	25.2	0.92	0.90	8.14	clear	NONE -71

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1525	SAMPLING ENDED AT: 1528
PUMP OR TUBING DEPTH IN WELL (feet): 25.5	SAMPLE PUMP FLOW RATE (mL per minute): 490	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N Filtration Equipment Type:	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-109	3	CG	40ml	HCl	40ml	-	8260	RFPP

REMARKS: **missing 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Tallevast, FL
WELL NO: MW-110	SAMPLE ID: MW-110 DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet): 5.11	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (28 feet - 5.11 feet) X 0.16 gallons/foot = 3.66 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 31 feet) + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 25.5	PURGING INITIATED AT: 1053	PURGING ENDED AT: 1107	TOTAL VOLUME PURGED (gallons): 1.54

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or M/S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1101	0.88	0.88	0.11	5.69	6.5	23.6	.721	0.7	7.20	clear	NONE	-73
1103	0.22	1.10	0.11	5.81	6.5	23.6	.720	0.6	4.87	clear	NONE	-78
1105	0.22	1.32	0.11	5.91	6.5	23.6	.717	0.6	5.86	clear	NONE	-80
1107	0.22	1.54	0.11	5.99	6.5	23.7	.721	0.6	3.77	clear	NONE	-81

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1108	SAMPLING ENDED AT: 1112
PUMP OR TUBING DEPTH IN WELL (feet): 25.5 ^{MPS} 4/14/08	SAMPLE PUMP FLOW RATE (mL per minute): ~ 430	TUBING MATERIAL CODE: Teflon	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-110	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **2nd purge - New Tubing placed in well; 2 bottles missing.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	MDS 4/14/08	SITE LOCATION: Sarasota FL	MDS 4/14/08
WELL NO: MW-111	SAMPLE ID: MW-111	DATE: 1/28/08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 23 feet to 28 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= \emptyset gallons + (0.0036 gallons/foot X 28 feet) + 0.25 gallons = 0.32 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26	PURGING INITIATED AT: 16:33	PURGING ENDED AT: 16:54	TOTAL VOLUME PURGED (gallons): 2.0
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1638	0.5	0.5	0.1	3.90	6.2	24.4	65	4.8	5.37	Clear	Slight
1641	0.3	0.8	0.1	3.83	6.3	24.4	59	5.1	2.63		
1644	0.3	1.1	0.1	3.81	6.3	24.4	60	5.1	2.20		
1647	0.3	1.4	0.1	3.91	6.3	24.5	64	5.3	1.32		
1650	0.3	1.7	0.1	3.90	6.3	24.5	66	5.5	1.06		
1653	0.3	2.0	0.1		6.3	24.6	66	5.6	1.05		
→ Sampled @ 1656 ←											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KUC / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1656	SAMPLING ENDED AT: 1700
PUMP OR TUBING DEPTH IN WELL (feet): 26	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y	Filtration Equipment Type: MDS 4/14/08	DUPLICATE: <input checked="" type="checkbox"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-111	3	CG	40ml	HLL	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Need 1 bolt.

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tallewas	SITE LOCATION: Sarasota, FL
WELL NO: MW-112	DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 76 feet to 86 feet	STATIC DEPTH TO WATER (feet): -	PURGE PUMP TYPE OR BAILER: PP Geopac
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable
 = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = gallons + **0.0026** gallons/foot X **94** feet + **0.50** gallons = **0.74** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 281	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 281	PURGING INITIATED AT: 1020	PURGING ENDED AT: 1036	TOTAL VOLUME PURGED (gallons): 1.76
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle OL or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
7 1027	0.77	0.77	0.11	25.34	7.70	24.96	434	0.43	1.76	clear	yes
1030	0.33	1.10	0.11	26.59	7.88	24.76	414	0.39	1.58	clear	yes
1033	0.33	1.43	0.11	27.25	7.90	24.65	412	0.36	1.69	clear	yes
1036	0.33	1.76	0.11	28.14	7.93	24.60	406	0.38	1.64	clear	yes

ORL
-219
-213
-218
-223

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: Neil Smith	SAMPLING INITIATED AT: 1038	SAMPLING ENDED AT: 1045
PUMP OR TUBING DEPTH IN WELL (feet): 281	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100 mL	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y Filtration Equipment Type: _____	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-112	2	CG	40mL	HCC	-	7.2	5260B + 8260S1m RFPP	

REMARKS: **Well pressurized**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	MDS 4/14/08	SITE LOCATION: Sarasota FL	MDS 4/14/08
WELL NO: MW-113	SAMPLE ID: MW-113	DATE: 1/28/08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 37 feet to 42 feet	STATIC DEPTH TO WATER (feet): 11.31	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= _____ feet - _____ feet X _____ gallons/foot = _____ gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 10.002 gallons/foot X 42 feet + 0.25 gallons = _____ gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1503	PURGING ENDED AT: 1610	TOTAL VOLUME PURGED (gallons):

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
15 1352	0.5	0.5	0.1	12.75	7.3	24.6	28	3.4	3.58	Clear	Slight	167
15 1355				13.10	7.2	24.6	29	3.1	2.26			166
15 1358				13.45	7.2	24.6	30	3.0	2.82			166
16 1401				13.47	7.1	24.6	50	2.9	3.45			105
16 1404				13.33	7.1	24.5	53	2.8	3.31			92
16 1407				13.35	7.1	24.5	55	2.6	4.24			60
16 1410					7.1	24.4	55	2.6	2.57			
16 1413	→ Sampled @ 1413 ←											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Kasey Cornwell / ABBL	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1613	SAMPLING ENDED AT: 1618
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N Filtration Equipment Type: MDS 4/14/08	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	MDS 4/14/08

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-113	3	CG	90ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

A needs 2 bolts.

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee FL
WELL NO: MW-114	SAMPLE ID: MW-114 DATE: 1-24-08

PURGING DATA

WELL DIAMETER (Inches): 2	TUBING DIAMETER (Inches): 1/4	WELL SCREEN INTERVAL DEPTH: 35 feet to 70 feet	STATIC DEPTH TO WATER (feet): 245	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 13:53	PURGING ENDED AT: 14:15	TOTAL VOLUME PURGED (gallons): 1.7
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1400	0.2	0.2	0.11	2.7	6.37	23.82	0.250	1.75	0	clear none	
1405	0.3	0.7	0.1	2.7	6.36	23.79	0.247	0.56	0	clear none	
1410	0.5	1.2	0.1	2.7	6.39	24.34	0.244	0.45	0	" "	
1415	0.3	1.7	0.1	2.7	6.42	24.36	0.245	0.39	0	" "	

MWS
4/19/08
ORP
-103
-129
-134
-137

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Michal Forbes / Arcadis	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1416	SAMPLING ENDED AT: 1420
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> FILTER SIZE: µm	DUPLICATE: Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW 114	3	CG	40ml	HCl	—	—	8260	RFPP

MWS
4/19/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-115	SAMPLE ID: MW-115
DATE: 1-29-05	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 20 feet to 25 feet	STATIC DEPTH TO WATER (feet): 6.74	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (.0026 gallons/foot X 28 feet) + .25 gallons = 32 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 9:36	PURGING ENDED AT: 4:14/2005	TOTAL VOLUME PURGED (gallons): 2.4

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle) (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
951	1.5	1.5	0.1	8.12	6.59	25.41	0.578	0.74	3.69	None	None
954	0.3	1.8	↓	8.11	6.59	25.54	0.579	1.08	3.13	↓	↓
957	↓	2.1	↓	8.12	6.59	25.62	0.583	1.15	1.62	↓	↓
1000	↓	2.4	↓	8.13	6.56	25.61	0.584	1.16	1.31	↓	↓

MDS
4/14/05
ORP
-122
-126
-128
-130

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 1001	SAMPLING ENDED AT: 1005
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/05	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTRATION EQUIPMENT TYPE: <input checked="" type="radio"/> N	FILTER SIZE: _____ µm
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-115	3	CG	40mL	HCl	—	—	8260	RFPP

MDS
4/14/05

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallewaast</u>		SITE LOCATION: <u>Sarasota FL</u>	
WELL NO: <u>MW-116</u>	SAMPLE ID: <u>MW-116</u>	DATE: <u>1/25/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>0.25</u>	WELL SCREEN INTERVAL DEPTH: <u>21</u> feet to <u>26</u> feet	STATIC DEPTH TO WATER (feet): <u>1.76</u>	PURGE PUMP TYPE OR BAILER: <u>peristaltic</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (<u>26</u> feet - <u>24.24</u> feet) X <u>0.163</u> gallons/foot = <u>0.29</u> gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = _____ gallons + (_____ gallons/foot X <u>MD5 4/14/08</u>) + _____ gallons = _____ gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: <u>1350</u>		PURGING ENDED AT: <u>1420</u>		TOTAL VOLUME PURGED (gallons): <u>3.5</u>				
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	DRP
<u>1350</u>	<u>0.5</u>	<u>0.5</u>	<u>0.1</u>	<u>3.35</u>	<u>6.9</u>	<u>25.0</u>	<u>0.430</u>	<u>4.1</u>	<u>2.10</u>	<u>clear</u>	<u>—</u>	<u>127</u>
<u>1355</u>	<u>1.0</u>	<u>1.0</u>	<u>0.1</u>	<u>3.59</u>	<u>6.9</u>	<u>25.2</u>	<u>0.440</u>	<u>3.2</u>	<u>1.40</u>	<u>clear</u>	<u>—</u>	<u>121</u>
<u>1400</u>	<u>0.5</u>	<u>1.5</u>	<u>0.1</u>	<u>3.64</u>	<u>6.9</u>	<u>25.1</u>	<u>*70</u>	<u>5.0</u>	<u>1.32</u>	<u>clear</u>	<u>—</u>	<u>109</u>
<u>1405</u>	<u>0.5</u>	<u>2.0</u>	<u>0.1</u>	<u>3.67</u>	<u>6.3</u>	<u>25.2</u>	<u>69</u>	<u>3.2</u>	<u>0.75</u>	<u>clear</u>	<u>—</u>	<u>827</u>
<u>1410</u>	<u>0.5</u>	<u>2.5</u>	<u>0.1</u>	<u>3.69</u>	<u>6.3</u>	<u>25.2</u>	<u>66</u>	<u>3.1</u>	<u>0.70</u>	<u>clear</u>	<u>—</u>	<u>61</u>
<u>1415</u>	<u>0.5</u>	<u>3.0</u>	<u>0.1</u>	<u>3.69</u>	<u>6.3</u>	<u>25.2</u>	<u>65</u>	<u>3.0</u>	<u>0.61</u>	<u>"</u>	<u>—</u>	<u>37</u>
<u>1420</u>	<u>0.5</u>	<u>3.5</u>	<u>0.1</u>	<u>3.69</u>	<u>6.3</u>	<u>25.3</u>	<u>63</u>	<u>3.0</u>	<u>0.50</u>	<u>"</u>	<u>—</u>	<u>27</u>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith</u>	SAMPLER(S) SIGNATURES: <u>Nathan Smith</u>	SAMPLING INITIATED AT: <u>1420</u>	SAMPLING ENDED AT: <u>1430</u>
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u>< 100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u>	FIELD-FILTERED: <u>Y</u>	DUPLICATE: <u>Y</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-116</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFP</u>

REMARKS: * re-calibrated Horiba @ 1401 (after recording readings). Time sampled = 1420

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: **Former ABC** SITE LOCATION: **Talleast**
WELL NO: **MW-117** SAMPLE ID: **MW-117** DATE: **1/25/07**

PURGING DATA

WELL DIAMETER (inches): **2** TUBING DIAMETER (inches): **1/8** WELL SCREEN INTERVAL: **4.2** STATIC DEPTH: **6.60** PURGE PUMP TYPE OR BAILER: **PP**
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
only fill out if applicable

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): **41.8** FINAL PUMP OR TUBING DEPTH IN WELL (feet): **38.5** PURGING INITIATED AT: **8:45** PURGING ENDED AT: **8:57** TOTAL VOLUME PURGED (gallons): **1.1**

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8:48	0.3	0.3	0.1	8.72	7.22	24.37	0.530	0.78	5.68	clear	none
8:51	0.3	0.6	0.1	9.8	7.24	24.41	0.528	0.65	1.86	clear	none
8:54	0.3	0.9	0.1	11.00	7.25	24.49	0.526	0.55	1.35	clear	none
8:57	0.3	1.1	0.1	11.41	7.25	24.50	0.525	0.53	1.76	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: **Darrin Johnson** SAMPLER(S) SIGNATURES: **Darrin Johnson** SAMPLING INITIATED AT: **8:58** SAMPLING ENDED AT: **9:03**
PUMP OR TUBING DEPTH IN WELL (feet): **38.5** SAMPLE PUMP FLOW RATE (mL per minute): **<100** TUBING MATERIAL CODE: **T**
FIELD DECONTAMINATION: **Y** N FIELD-FILTERED: **Y** N FILTER SIZE: **0** µm DUPLICATE: **Y** N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-117	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 4/14/08	SITE LOCATION: <u>Sarasota FL</u> MDS 4/14/08
WELL NO: <u>MW-118</u>	SAMPLE ID: <u>MW-118</u> DATE: <u>1/25/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/8</u>	WELL SCREEN INTERVAL DEPTH: <u>20</u> feet to <u>40</u> feet	STATIC DEPTH TO WATER (feet): <u>5.95</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>24.07</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>21.57</u>	PURGING INITIATED AT: <u>8:16</u>	PURGING ENDED AT: <u>8:34</u>	TOTAL VOLUME PURGED (gallons): <u>1.8</u>

TIME	VOLUME PURGED (gallons)	CUMUL VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8:25	0.9	0.9	0.1	6.57	7.10	24.85	1.01	0.53	10.22	black	none
8:28	0.3	1.2	0.1	6.60	6.73	24.96	1.01	0.29	7.11	clear	none
8:31	0.3	1.5	0.1	6.61	6.67	24.92	1.01	0.24	3.56		
8:34	0.3	1.8	0.1	6.62	6.68	24.97	1.01	0.23	3.64		

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Darrin Johnson</u>	SAMPLER(S) SIGNATURES: <u>Darrin Johnson</u>	SAMPLING INITIATED AT: <u>8:35</u>	SAMPLING ENDED AT: <u>8:40</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>21.57</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u> MDS 4/14/08	TUBING MATERIAL CODE: <u>T</u>	
FIELD DECONTAMINATION: <u>(Y) N</u>	FIELD-FILTERED: <u>(Y) N</u> FILTER SIZE: _____ µm	DUPLICATE: <u>Y (N)</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-118</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCL</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC</u>	SITE LOCATION: <u>Tallevast</u>
WELL NO: <u>MW-119</u>	SAMPLE ID: <u>MW-119</u>
DATE: <u>1-25-05</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>31</u> feet to <u>35.55</u> feet	STATIC DEPTH TO WATER (feet): <u>9.33</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable
 = (35.55 feet - 9.33 feet) X 0.026 gallons/foot = 0.35 gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = 0 gallons + 10026 gallons/foot X 37.5 feet + 125 gallons = 0.35 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>35.55</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>32.75</u>	PURGING INITIATED AT: <u>7:50</u>	PURGING ENDED AT: <u>8:00</u>	TOTAL VOLUME PURGED (gallons): <u>1.0</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>7:54</u>	<u>0.4</u>	<u>0.4</u>	<u>0.1</u>	<u>13.02</u>	<u>11.65</u>	<u>23.78</u>	<u>1.62</u>	<u>1.70</u>	<u>3.68</u>	<u>clear</u>	<u>none</u>
<u>7:57</u>	<u>0.3</u>	<u>0.7</u>	<u>0.1</u>	<u>14.84</u>	<u>11.68</u>	<u>23.78</u>	<u>1.67</u>	<u>1.84</u>	<u>3.81</u>	<u>clear</u>	<u>none</u>
<u>8:00</u>	<u>0.3</u>	<u>1.0</u>	<u>0.1</u>	<u>16.75</u>	<u>11.68</u>	<u>23.85</u>	<u>1.69</u>	<u>1.89</u>	<u>5.03</u>	<u>clear</u>	<u>none</u>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Darrin Johnson</u>	SAMPLER(S) SIGNATURES: <u>Darrin Johnson</u>	SAMPLING INITIATED AT: <u>8:01</u>	SAMPLING ENDED AT: <u>8:06</u>
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PUMP OR TUBING DEPTH IN WELL (feet): <u>32.75</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>L100</u>	TUBING MATERIAL CODE: <u>T</u>
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTRATION EQUIPMENT TYPE: _____

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-119</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS: Will check cal solution due to high pH readings

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-120	SAMPLE ID: MW120 DATE: 1-25-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 20 feet to 25 feet	STATIC DEPTH TO WATER (feet): 2.85	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - 9/14/08 feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0076 gallons/foot X 27 feet + .25 gallons = 0.32 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 25.3	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 22.5	PURGING INITIATED AT: 7:30	PURGING ENDED AT: 7:45	TOTAL VOLUME PURGED (gallons): 1.5

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
7:36	0.6	0.6	0.1	4.20	6.58	23.72	0.791	0.51	1.15	clear	none
7:39	0.3	0.9	0.1	3.61	6.62	23.08	0.796	0.38	0.97		
7:42	0.3	1.2	0.1	3.61	6.65	23.03	0.798	0.51	1.0		
7:45	0.3	1.5	0.1	3.61	6.68	23.03	0.800	0.49	0.93		

-87
-99
-104
-112

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 846	SAMPLING ENDED AT: 850
PUMP OR TUBING DEPTH IN WELL (feet): 22.5	SAMPLE PUMP FLOW RATE (mL per minute): 200	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: (Y) N	FIELD-FILTERED: (Y) N FILTER SIZE: _____ µm	DUPLICATE: Y (N)	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-120	3	CG	40ml	HCl	—	—	8260	RFPP

MDS
9/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast MDS 4/14/08	SITE LOCATION: Sarasota FL MDS 4/14/08
WELL NO: MW-121	SAMPLE ID: MW-121 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/8	WELL SCREEN INTERVAL DEPTH: 18 feet to 23 feet	STATIC DEPTH TO WATER (feet): 8.77	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 20.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 20.5	PURGING INITIATED AT: 1600	PURGING ENDED AT: 1612	TOTAL VOLUME PURGED (gallons): 1.02 MDS 4/14/08
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1603	0.3	0.3	0.1	9.94	7.01	26.53	0.1600	1.21	4.93	Clear	
1606	0.3	0.6	0.1	9.94	7.01	26.65	0.1602	1.06	2.08	Clear	
1609	0.3	0.9	0.1	10.30	6.98	26.65	0.1609	0.93	2.22		
1612	0.3	1.2	0.1	10.39	6.95	26.60	0.1606	0.93	1.64		
1615	→ Sampled @ 1615 ←										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: IARCADIS	SAMPLER(S) SIGNATURES: MDS 4/14/08	SAMPLING INITIATED AT: 1615	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 20.5	SAMPLE PUMP FLOW RATE (mL per minute): 2100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 1 µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-121	3	CG	90ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 4/14/08		SITE LOCATION: <u>Sarasota FL</u> MDS 4/14/08	
WELL NO: <u>MW-122</u>	SAMPLE ID: <u>MW-122</u>	DATE: <u>1/25/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>21</u> feet to <u>20</u> feet	STATIC DEPTH TO WATER (feet): <u>6.45</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (<u> </u> feet - <u> </u> feet) X <u> </u> gallons/foot = <u> </u> gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (0.0026 gallons/foot X <u>20'</u> feet) + 0.25 gallons = <u>0.32</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>23</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>23</u>	PURGING INITIATED AT: <u>0900</u>	PURGING ENDED AT: <u>0925</u>	TOTAL VOLUME PURGED (gallons): <u>2.0</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0905	0.5	0.5	0.1	7.05	6.80	20.07	0.906	1.54	1.78	Clear	None
0910	0.5	1.0		7.21	6.92	21.51	0.918	1.15	0.64		
0915	0.5	1.5		7.31	6.92	21.57	0.917	1.09	0.47		
0920	0.5	2.0		7.39	6.92	21.61	0.914	1.06	0.55		
0925	→ Sampled @ 0925 ←										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasey Connell / ARCADIS</u>		SAMPLER(S) SIGNATURES: <u>[Signature]</u>		SAMPLING INITIATED AT: <u>0925</u>	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): <u>23</u>		SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u> MDS 4/14/08		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: <u> </u> µm		DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-122	3	CG	40ml	HCl	-	-	3260	RFPP

REMARKS: Initial Purge clear, no noticeable odor.

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-123	SAMPLE ID: MW-123
DATE: 2-01-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 375 feet to 395 feet	STATIC DEPTH TO WATER (feet): 16.81	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (.0026 gallons/foot X 398 feet) + .25 gallons = 1.28 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 395	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 375	PURGING INITIATED AT: 0839	PURGING ENDED AT: 9:15	TOTAL VOLUME PURGED (gallons): 3.6							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:09	3.0	3.0	0.1	17.79	7.41	24.83	1.20	.28	1.11	clear	none
9:12	.3	3.3	0.1	16.73	7.43	24.85	1.19	.24	1.13	clear	none
9:15	.3	3.6	0.1	16.74	7.43	24.81	1.20	.24	.92	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darvin Johnson	SAMPLER(S) SIGNATURES: <i>Darvin Johnson</i>	SAMPLING INITIATED AT: 9:16	SAMPLING ENDED AT: 9:21
PUMP OR TUBING DEPTH IN WELL (feet): 375	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	FILTRATION EQUIPMENT TYPE: _____	FILTER SIZE: _____ µm
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-123	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **transducer removed @ 0836 / transducer re-inserted @ 0930**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-124	SAMPLE ID: MW-124 DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 127 feet to 137 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 142 feet) + 0.5 gallons = 0.87 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 129.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 129.5	PURGING INITIATED AT: 0900	PURGING ENDED AT: 0925	TOTAL VOLUME PURGED (gallons): 1.75

TIME	VOLUME PURGED (gallons)	CUMUL VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0913	0.91	0.91	0.07	17.02	7.28	25.63	1240	3.13	15.1	clear	no
0917	0.28	1.19	0.07	17.15	7.29	25.79	1240	3.21	16.5	clear	yes
0921	0.28	1.47	0.07	17.32	7.29	25.75	1250	3.24	14.3	clear	yes
0925	0.28	1.75	0.07	17.40	7.29	25.79	1240	3.23	16.1	clear	yes

ORP
-218
-219
-220
-220

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NS/SW/ ARCADIS	SAMPLER(S) SIGNATURES: <i>Steve Will</i>	SAMPLING INITIATED AT: 0926	SAMPLING ENDED AT: 0930
PUMP OR TUBING DEPTH IN WELL (feet): 129.5	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: PE/T	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-124	3	CG	40mL	HCl	40mL	—	8260 + 8260SIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Talleast
WELL NO: MW-125	SAMPLE ID: MW-125
DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 30 feet to 35 feet	STATIC DEPTH TO WATER (feet): 7.94	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (1,0026 gallons/foot X 38 feet) + .25 gallons = .35 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 35	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 32.5	PURGING INITIATED AT: 10:50	PURGING ENDED AT: 11:08	TOTAL VOLUME PURGED (gallons): 1.8

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
10:56	.6	.6	0.1	17.57	9.14	25.79	222	6.91	19.3	clear	none
10:59	1.3	.9	0.1	18.15	9.13	25.83	221	6.50	19.4	clear	none
11:02	1.3	1.2	0.1	20.76	9.36	25.57	226	6.11	11.40	clear	none
11:05	.3	1.5	0.1	20.56	9.27	25.63	232	6.03	8.37	clear	none
11:08	.3	1.8	0.1	21.07	9.33	25.63	230	5.96	8.56	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson			SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>			SAMPLING INITIATED AT: 11:09		SAMPLING ENDED AT: 11:14		
PUMP OR TUBING DEPTH IN WELL (feet): 32.5 MDS 4/14/08			SAMPLE PUMP FLOW RATE (mL per minute): 4100 MDS 4/14/08			TUBING MATERIAL CODE: PE				
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N			FIELD-FILTERED: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N			FILTER SIZE: _____ µm		DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N		

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-125	3	CG	40ml	HCL	—	—	2250	RFPP

REMARKS: **new teflon tubing / needs 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Telleavast	SITE LOCATION: Surasota, FL
WELL NO: MW-126	SAMPLE ID: MW-126
DATE: 1/24/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 27 feet to 32 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: pp geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0024 gallons/foot X 35 feet) + 0.50 gallons = 0.591 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 230	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 230	PURGING INITIATED AT: 0945	PURGING ENDED AT: 1000	TOTAL VOLUME PURGED (gallons): 3.15
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle µg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
3 0948	0.63	0.63	0.21	14.00	6.40	25.50	604	2.19	43.8	clear	yes	-79
0951	0.63	1.26	0.21	14.53	6.46	25.41	612	1.95	24.6	clear	yes	-93
0954	0.63	1.89	0.21	15.01	6.51	25.40	622	1.52	17.1	clear	yes	-102
0957	0.63	2.52	0.21	15.94	6.53	25.36	634	1.24	16.4	clear	yes	-106
1000	0.63	3.15	0.21	16.46	6.54	25.33	639	1.15	13.0	clear	yes	-108
<i>End (Sed) 1/24/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1000	SAMPLING ENDED AT: 1010
PUMP OR TUBING DEPTH IN WELL (feet): 230	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y FILTER SIZE: 0.1µm	DUPLICATE: Y X	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-126	3	CG	40ml	HCL	-	6.2	8260B+8260SM RFPF	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.63

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST, FL	
WELL NO: mw-127	SAMPLE ID: mw-127	DATE: 1-30-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 100 feet to 116 feet	STATIC DEPTH TO WATER (feet): 15.78	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 113 feet) + 0.25 gallons = 0.54 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 105	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 105	PURGING INITIATED AT: 1146	PURGING ENDED AT: 1209	TOTAL VOLUME PURGED (gallons): 2.30
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle NO or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1203	1.70	1.70	0.1	16.01	6.9	26.9	1.14	0.9	5.49	clear	NONE
1205	0.2	1.90	0.1	16.03	6.9	26.9	1.14	0.8	7.97	clear	NONE
1207	0.2	2.10	0.1	16.03	6.9	26.9	1.14	0.8	6.40	clear	NONE
1209	0.2	2.30	0.1	16.03	7.0	27.0	1.14	0.8	6.39	clear	NONE

ORP
-97
-99
-100
-102

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1210	SAMPLING ENDED AT: 1221
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PUMP OR TUBING DEPTH IN WELL (feet): 105	SAMPLE PUMP FLOW RATE (mL per minute): 1360	TUBING MATERIAL CODE:
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FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N
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SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-127	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS: **Sampled w/ transfer or in well; 2 bolts missing.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.94

SITE NAME: <u>Tallevast</u>		SITE LOCATION: <u>Tallevast, FL</u>	
WELL NO: <u>MW-128</u>	SAMPLE ID: <u>MW-128</u>	DATE: <u>1/31/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>140</u> feet to <u>150</u> feet	STATIC DEPTH TO WATER (feet): <u>19.43</u>	PURGE PUMP TYPE OR BAILER: <u>Geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>153</u> feet) + <u>0.25</u> gallons = <u>0.65</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>145</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>145</u>	PURGING INITIATED AT: <u>1135</u>	PURGING ENDED AT: <u>1202</u>	TOTAL VOLUME PURGED (gallons): <u>260</u>
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20

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle <u>NO</u> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1155	2.00	2.00	0.1	25.12	11.60	28.3	2.14	0.66	20.6	Clear	NO
1157	0.2	2.20	0.1	25.18	11.60	28.3	2.12	0.66	22.8	Clear	NO
1159	0.2	2.40	0.1	25.25	11.60	28.0	2.11	0.66	22.7	Clear	NO
1202	0.2	2.60	0.1	25.45	11.58	28.2	2.04	0.63	22.6	Clear	NO

ORP
-283
-285
-286
-290

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Jennifer King / ARONIS</u>	SAMPLER(S) SIGNATURES: <u>Jennifer King</u>	SAMPLING INITIATED AT: <u>1205</u>	SAMPLING ENDED AT: <u>1220</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>145</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>370</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-28	3	CG	40ml	HCl	—	—	8260	RFPP
DUP-28	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: sampled with transducer, missing 2 bolts

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

165

SITE NAME: <u>Tallevast</u>		SITE LOCATION: <u>Tallevast</u>	
WELL NO: <u>mw-129</u>	SAMPLE ID: <u>mw-129</u>	DATE: <u>1-30-08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/2</u>	WELL SCREEN INTERVAL DEPTH: <u>103</u> feet to <u>113</u> feet	STATIC DEPTH TO WATER (feet): <u>116.25</u>	PURGE PUMP TYPE OR BAILER: <u>PP Geo Pump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>116</u> feet) + <u>.25</u> gallons = <u>.55</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>108</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>108</u>	PURGING INITIATED AT: <u>1400</u>	PURGING ENDED AT: <u>1423</u>	TOTAL VOLUME PURGED (gallons): <u>2.30</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
17 1407	1.70	1.70	.10	17.70	7.57	25.4	.921	.49	12.6	clear	yes
1419	0.2	1.90	.10	17.71	7.38	25.4	.932	.48	13.3	↓	yes
1421	0.2	2.10	.10	↓	7.30	25.3	.936	.48	13.9	↓	yes
1423	0.2	2.30	.10	↓	7.25	25.3	.940	.47	13.9	↓	yes

ORP
-223
-220
-220
-220

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Jennifer King/ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Jennifer King</u>	SAMPLING INITIATED AT: <u>1425</u>	SAMPLING ENDED AT: <u>1431</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>108</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>~370</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>mw-129</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.62

SITE NAME: Tallevast		SITE LOCATION: Tallevast, FL	
WELL NO: MW-130	SAMPLE ID: MW-130	DATE: 1/30/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 100 feet to 110 feet	STATIC DEPTH TO WATER (feet): 15.26	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 113 feet) + .25 gallons = .54 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 105	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 105	PURGING INITIATED AT: 0853	PURGING ENDED AT: 0911	TOTAL VOLUME PURGED (gallons): 2.55
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or m/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
905	1.65	1.65	0.15	16.59	7.19	25.3	654	.84	.62	clear	yes
907	0.30	1.95	0.15	16.40	7.21	25.4	659	.81	.58	clear	yes
909	0.30	2.25	0.15	16.61	7.21	25.3	660	.79	.60	clear	yes
911	0.30	2.55	0.15	16.62	7.22	25.4	661	.78	.25	clear	yes

RFPP
-223
-227
-229
-231

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 6.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / M&C's	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 913	SAMPLING ENDED AT: 922
PUMP OR TUBING DEPTH IN WELL (feet): 105	SAMPLE PUMP FLOW RATE (mL per minute): 567	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-130	CG ↔ 3		40ml	HCl	—	—	8260	RFPP

REMARKS: **missing 2 bolts; sampled w/ transducer in well**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellerat	SITE LOCATION: Sarasota, FL
WELL NO: MW-131	SAMPLE ID: MW-131 DATE: 1/31/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 100 feet to 110 feet	STATIC DEPTH TO WATER (feet): 1438	PURGE PUMP TYPE OR BAILER: (pp) gear pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
= gallons + (0.0026 gallons/foot X 112 feet) + 0.25 gallons = 0.54 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 105		FINAL PUMP OR TUBING DEPTH IN WELL (feet): 105		PURGING INITIATED AT: 0855		PURGING ENDED AT: 0911		TOTAL VOLUME PURGED (gallons): 1.50			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or S/cm)	DISSOLVED OXYGEN (circle method or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0900	0.60	0.60	0.10	15.56	7.28	23.49	0.684	0.62	15.0	clear	yes
0903	0.30	0.90	0.10	15.56	7.27	23.50	0.684	0.67	13.4	clear	yes
0908	0.30	1.20	0.10	15.60	7.25	23.53	0.683	0.60	10.24	clear	yes
0911	0.30	1.50	0.10	15.60	7.25	23.52	0.683	0.61	8.1	clear	yes

ORP
-302
-304
-309
-310

Neil Smith 1/31/08

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS		SAMPLER(S) SIGNATURES: <i>Neil Smith</i>		SAMPLING INITIATED AT: 0915	SAMPLING ENDED AT: 0920
PUMP OR TUBING DEPTH IN WELL (feet): 105		SAMPLE PUMP FLOW RATE (mL per minute): 100		TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: X N		FIELD-FILTERED: Y <input checked="" type="checkbox"/> FILTER SIZE: _____ µm		DUPLICATE: Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-131	3	CG	40ML	HCL	2 2	2	82603 + 82605in RFPP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

1.65

14
72

Form FD 9000-24 GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: Sarasota FL	MOS 4/19/08
WELL NO: mw-132	SAMPLE ID: mw-132	DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 101 feet to 111 feet	STATIC DEPTH TO WATER (feet): 16.10	PURGE PUMP TYPE OR BAILER: 6 to Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 114 feet) + .25 gallons = 0.55 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 106	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 106	PURGING INITIATED AT: 1508	PURGING ENDED AT: 1528	TOTAL VOLUME PURGED (gallons): 2.40
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle 10/1 or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1522	1.68	1.68	0.12	16.85	7.30	24.9	0.565	2.0	1.89	clear	yes
1524	0.24	1.92	0.12	16.70	7.30	24.8	0.562	0.5	1.34	clear	yes
1526	0.24	2.16	0.12	16.74	7.3	24.9	0.560	0.5	1.06	clear	yes
1528	0.24	2.40	0.12	16.76	7.2	24.8	0.565	0.4	1.28	clear	yes

ORP
-242
-255
-265
-279

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1530	SAMPLING ENDED AT: 1540
PUMP OR TUBING DEPTH IN WELL (feet): 106	SAMPLE PUMP FLOW RATE (mL per minute): 454	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> <input checked="" type="checkbox"/> JK	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-132	3	CG	40mL	HCL	40mL	-	8260	RFPP
DUP 2	3	CG	40mL	HCL	40mL	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

EB16 @ 1625

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallevast	
WELL NO: MW-133	SAMPLE ID: MW-133	DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 100 feet to 110 feet	STATIC DEPTH TO WATER (feet): 14.88	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 108 feet) + 0.25 gallons = 0.53 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 9:12	PURGING ENDED AT: 9:31	TOTAL VOLUME PURGED (gallons): 1.9
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:25	1.3	1.3	0.1	15.82	7.38	23.57	553	0.97	6.69	clear	none
9:28	1.3	1.6	0.1	15.85	7.36	23.54	561	0.94	5.47	↓	↓
9:31	1.3	1.9	0.1	15.85	7.34	23.65	565	0.92	5.30	↓	↓

-280
-283
-286

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson / Arcadis	SAMPLER(S) SIGNATURE(S): <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 9:32	SAMPLING ENDED AT: 9:37
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTRATION SIZE: 100 µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-133	CG	40mL						
MW-133	3	CG	40mL	HCl			8260	RFPP
Dup-16	3	CG	40mL	HCl			8260	RFPP

yes
1/2/08

REMARKS: **transducer reinserted @ 10:00**
transducer removed @ 9:07
4.5 liters removed, total
DUP 16-012808 taken

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Sarasota, FL
WELL NO: MW-134	SAMPLE ID: MW-134 DATE: 2-1-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 103 feet to 113 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 116 feet + 0.25 gallons = 0.55 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 108	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 108	PURGING INITIATED AT: 1109	PURGING ENDED AT: 1128	PURGED VOLUME (gallons): 2.09

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1119	1.10	1.10	6.11	17.38	7.50	27.43	650	1.21	12.2	Clear	NOYS-341
1122	0.33	1.43	0.11	17.25	7.43	27.34	649	1.27	10.54	Clear	NOYS-347
1125	0.33	1.76	0.11	17.36	7.38	27.42	649	1.23	7.38	Clear	NOYS-347
1128	0.33	2.09	0.11	17.37	7.36	27.36	646	1.31	7.64	Clear	NOYS-323

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 6.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/TH / ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1129	SAMPLING ENDED AT: 1132
PUMP OR TUBING DEPTH IN WELL (feet): 108	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-134	3	CG	40ml	HCl	40ml	—	82608 82608m RFPP	

REMARKS: **Transducer removed @ 1040 Replaced @ 1131**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: 110-135 Talleast	SITE LOCATION: Sarasota FL
WELL NO: MW-135	DATE: 1/28/08

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: 94 feet to 104 feet	STATIC DEPTH TO WATER (feet): 15.43	PURGE PUMP TYPE: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 104 feet) + $.25$ gallons = 0.50 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 99	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 99	PURGING INITIATED AT: 1035	PURGING ENDED AT: 1055	TOTAL VOLUME PURGED (gallons): 2.0
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1037	0.2	0.2	0.1	15.75	7.2	23.2	81	4.1	8.35	Clear	Slight
1040	0.3	0.5		15.79	7.2	23.3	81	3.6	50.0	Cloudy	0
1043	0.3	0.8		15.79	7.1	23.8	82	3.2	46.7		
1046	0.3	1.1		15.79	7.0	23.8	82	2.9	21.2		
1049	0.3	1.4		15.80	7.0	23.8	82	2.8	21.4		
1052	0.3	1.7		15.80	7.0	23.8	82	2.7	21.7		
1055	0.3	2.0		15.80	7.0	23.8	82	2.7	21.4		
1100	→ Sampled @ 1100 ←										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KLC / ARCA ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1100	SAMPLING ENDED AT: 1105
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PUMP OR TUBING DEPTH IN WELL (feet): 99	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N DUP 15

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-135	3	CG	40ml	HCL	-	-	8260	RFPP
Dup-15	3	CG	40ml	HCL	-	-	8260	RFPP

REMARKS: **Initial purge clear, slight sulfur odor.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* Transducer in well

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-136	SAMPLE ID: MW-136 DATE: 1.29.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 8.5 feet to 108.5 feet	STATIC DEPTH TO WATER (feet): 14.5	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 108.5 feet) + 0.25 gallons = 0.53 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 108.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 108.5	PURGING INITIATED AT: 8:15	PURGING ENDED AT: 8:45	TOTAL VOLUME PURGED (gallons): 2.00

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8:25	1.0	1.0	6.1	16.3	7.32	22.9	0.76	0.63	0	Clear	none
8:30	0.25	1.25	0.05	16.05	7.38	22.5	0.84	0.59	0	"	"
8:35	0.25	1.50	0.05	16.9	7.35	22.7	0.90	0.45	0	"	"
8:40	0.25	1.75	0.05	15.8	7.37	22.8	0.90	0.51	0	"	"
8:45	0.25	2.00	0.05	15.7	7.37	22.8	0.9	0.50	0	"	"

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

ORP

-253
-253
-260
-261
-263

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS			SAMPLER(S) SIGNATURES: <i>[Signature]</i>			SAMPLING INITIATED AT: 8:46	SAMPLING ENDED AT: 8:50
PUMP OR TUBING DEPTH IN WELL (feet): 108.5 MOS 4/14/08			SAMPLE PUMP FLOW RATE (mL per minute): 5100 MOS 4/14/08			TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N			FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm			DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-136	3	CG	40ml	He1	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

MOS 4/14/08

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ±0.2-units Temperature: ±0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ±0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellevar	SITE LOCATION: Sarasota, FL
WELL NO: MW-137	SAMPLE ID: MW-137 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 17 feet to 22 feet	STATIC DEPTH TO WATER (feet): 2.62	PURGE PUMP TYPE OR BAILER: (PP) geopump
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~19	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~19	PURGING INITIATED AT: 0855	PURGING ENDED AT: 0909	TOTAL VOLUME PURGED (gallons): 1.40
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle <input checked="" type="checkbox"/> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0900	0.50	0.50	0.10	3.86	6.94	22.31	825	1.32	2.78	clear	yes	ORP -141
0903	0.30	0.80	0.10	4.04	6.84	22.85	794	1.29	2.05	clear	yes	-145
0906	0.30	1.20	0.10	4.06	6.85	22.92	796	1.16	1.80	clear	yes	-148
0909	0.30	1.40	0.10	4.06	6.84	22.99	808	1.15	2.10	clear	yes	-148
<i>End Pump 1/29/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Fisher / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Fisher</i>	SAMPLING INITIATED AT: 0910	SAMPLING ENDED AT: 0915
PUMP OR TUBING DEPTH IN WELL (feet): ~19	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: _____ µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-137	3	CG	4 each	HCl	—	~2	826BPEPESM RFPP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellerast	SITE LOCATION: Sarasota, FL
WELL NO: MW-138	SAMPLE ID: MW-138 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 38 feet to 43 feet	STATIC DEPTH TO WATER (feet): 1009	PURGE PUMP TYPE OR BAILER: (pp) geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 47 feet) + 0.25 gallons = 0.272 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~40 1/2	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~40 1/2	PURGING INITIATED AT: 0810	PURGING ENDED AT: 0835	TOTAL VOLUME PURGED (gallons): 3.00

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
0814	0.48	0.48	0.12	12.74	6.90	22.25	1140	1.81	52.0	clear	yes	-142
0817	0.36	0.84	0.12	13.05	6.95	22.36	1140	1.64	23.1	clear	yes	-149
0820	0.36	1.20	0.12	13.52	6.98	22.44	1130	1.56	18.6	clear	yes	-152
0823	0.36	1.56	0.12	13.61	6.99	22.53	1130	1.45	11.0	clear	yes	-155
0826	0.36	1.92	0.12	13.62	7.08	22.70	1120	1.15	6.70	clear	yes	-169
0829	0.36	2.28	0.12	13.63	7.11	22.70	1120	1.08	4.2	clear	yes	-175
0832	0.36	2.64	0.12	13.75	7.13	22.79	1120	1.05	4.04	clear	yes	-181
0835	0.36	3.00	0.12	13.76	7.16	22.86	1110	0.90	3.30	clear	yes	-187

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 0838	SAMPLING ENDED AT: 0845
PUMP OR TUBING DEPTH IN WELL (feet): ~40 1/2	SAMPLE PUMP FLOW RATE (mL per minute): £100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-138	3	CG	40mL	HCL	-	~2	8260B+8260H RFP	

REMARKS: **Need one (1) extra bail for manhole cover**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Telleary+</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>MW-139</u>	SAMPLE ID: <u>MW-139</u> DATE: <u>1/29/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>147</u> feet to <u>157</u> feet	STATIC DEPTH TO WATER (feet): <u>149.2</u>	PURGE PUMP TYPE OR BAILER: <u>(pp) geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME <u>(0.70)</u> (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X <u>163</u> feet) + 0.25 gallons = <u>0.67</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~152</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~152</u>	PURGING INITIATED AT: <u>0940</u>	PURGING ENDED AT: <u>0958</u>	TOTAL VOLUME PURGED (gallons): <u>1.60</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<u>0947</u>	<u>0.20</u>	<u>0.20</u>	<u>0.10</u>	<u>6.92</u>	<u>7.47</u>	<u>23.31</u>	<u>715</u>	<u>1.47</u>	<u>3.09</u>	<u>clear</u>	<u>yes</u>	<u>-23c</u>
<u>0952-0945</u>	<u>0.30</u>	<u>1.00</u>	<u>0.10</u>	<u>6.92</u>	<u>7.48</u>	<u>23.35</u>	<u>715</u>	<u>1.29</u>	<u>1.96</u>	<u>clear</u>	<u>yes</u>	<u>-240</u>
<u>0955-0948</u>	<u>0.30</u>	<u>1.30</u>	<u>0.10</u>	<u>6.92</u>	<u>7.47</u>	<u>23.41</u>	<u>715</u>	<u>1.24</u>	<u>2.31</u>	<u>clear</u>	<u>yes</u>	<u>-243</u>
<u>0958-0944</u>	<u>0.30</u>	<u>1.60</u>	<u>0.10</u>	<u>6.92</u>	<u>7.47</u>	<u>23.37</u>	<u>715</u>	<u>2.15</u>	<u>1.45</u>	<u>clear</u>	<u>yes</u>	<u>-245</u>
<u>stopped 1/25/08</u>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil Smith</u>	SAMPLING INITIATED AT: <u>1000</u>	SAMPLING ENDED AT: <u>1005</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~152</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>500</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>X</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-139</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u>-</u>	<u>~2</u>	<u>82608+82608M RFPP</u>	

REMARKS: Placed new tubing in well to 152' - purge initial - color "blue" then clear - Need BOTTLES for cover

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-140	SAMPLE ID: MW-140 DATE: 1.29.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet): 13.8	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (300 feet - 13.8 feet) X 0.16 gallons/foot = 45.8 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 280 feet) + 0.25 gallons = 0.978 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 280	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 280	PURGING INITIATED AT: 9:15	PURGING ENDED AT: 10:05	TOTAL VOLUME PURGED (gallons): 4.5

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
9:25	1.0	1.0	0.1	15.45	8.00	21.8	0.74	1.4	35	turbid	none	-260
9:30	0.5	1.5	0.1	15.8	7.77	21.9	0.74	1.66	33	"	"	-258
9:35	0.5	2.0	0.1	15.8	7.67	21.8	0.74	1.67	156	"	"	-253
9:40	0.5	2.5	0.1	15.8	7.67	22.4	0.74	1.54	75	"	"	-253
9:45	0.5	3.0	0.1	16.4	7.69	22.3	0.74	1.48	74	"	"	-258
9:50	0.5	3.5	0.1	16.4	7.71	22.4	0.74	1.43	39	"	"	-256
9:55	0.5	4.0	0.1	16.6	7.75	22.7	0.751	1.30	21	"	"	-262
10:00	0.25	4.25	0.05	16.6	7.78	22.4	0.743	1.21	20	"	"	-257
10:05	0.25	4.50	0.05	16.6	7.78	22.7	0.741	1.20	17	"	"	-269

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 10:06	SAMPLING ENDED AT: 10:10
PUMP OR TUBING DEPTH IN WELL (feet): 280	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/>	FIELD-FILTERED: Y (N) FILTER SIZE: _____ µm	DUPLICATE: Y (N)	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW140	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallahassee</u>	SITE LOCATION: <u>Jamiatay Fl</u>
WELL NO: <u>MW-141</u>	SAMPLE ID: <u>MW-141</u> DATE: <u>1/29/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>15</u> feet to <u>20</u> feet	STATIC DEPTH TO WATER (feet): <u>265</u>	PURGE PUMP TYPE OR BAILER: <u>(pp) geopump</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= gallons + (0.0026 gallons/foot X 35 feet) + 0.25 gallons = 0.34 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~17.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~17.5</u>	PURGING INITIATED AT: <u>1235</u>	PURGING ENDED AT: <u>1244</u>	TOTAL VOLUME PURGED (gallons): <u>2.10</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle % or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1238	0.42	0.42	0.14	4.12	7.10	24.19	727	0.80	4.31	clear	yes
1241	0.42	0.84	0.14	4.19	7.08	24.12	727	0.54	2.13	clear	yes
1244	0.42	2.10	0.14	4.20	7.06	24.17	720	0.44	1.53	clear	yes
<i>new pump 1/29/08</i>											

*OKP
-202
-206*

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil Smith</u>	SAMPLING INITIATED AT: <u>1245</u>	SAMPLING ENDED AT: <u>1250</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~17.5</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>£100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>X</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-141</u>	<u>3</u>	<u>CG</u>	<u>4one</u>	<u>HCL</u>	<u>-</u>	<u>£2</u>	<u>8260B+8260B in RFPP</u>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tellevast	SITE LOCATION: Sarasota, FL
WELL NO: MW-142	SAMPLE ID: MW-142 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 4/4	WELL SCREEN INTERVAL DEPTH: 26 feet to 31 feet	STATIC DEPTH TO WATER (feet): 2.55	PURGE PUMP TYPE OR BAILER: PP peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.026 gallons/foot X 40 feet) + 0.25 gallons = 0.35 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~28	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~28	PURGING INITIATED AT: 1255	PURGING ENDED AT: 1308	TOTAL VOLUME PURGED (gallons): 1.56
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
4 1259	0.48	0.48	0.12	3.02	6.14	23.78	641	0.47	3.42	clear	yes	ORP -149
1302	0.36	0.84	0.12	3.04	6.08	23.77	625	0.40	4.99	clear	yes	-150
1305	0.36	1.20	0.12	3.04	6.07	23.75	621	0.36	4.38	clear	yes	-152
1308	0.36	1.56	0.12	3.06	6.07	23.75	623	0.34	4.10	clear	yes	-152
Hand Sample / 1/29/08												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1309	SAMPLING ENDED AT: 1315
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PUMP OR TUBING DEPTH IN WELL (feet): ~28	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE
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FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y Filtration Equipment Type: N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y N
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SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-142	3	CG	4ane	HCC	-	6.2	8200B peristaltic RFPP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-143	SAMPLE ID: MW-143 DATE: 1.29.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 96 feet to 106 feet	STATIC DEPTH TO WATER (feet): 13.4	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 106 feet + 0.25 gallons = 0.52 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 101	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 101	PURGING INITIATED AT: 12:44	PURGING ENDED AT: 13:10	TOTAL VOLUME PURGED (gallons): 1.75
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
12:50	0.5	0.5	0.1	13.4	7.24	23.9	0.55	0.58	22	clear	none
12:55	0.5	1.0	0.1	13.4	7.22	23.9	0.52	0.92	0	"	"
13:00	0.25	1.25	0.05	13.45	7.21	23.9	0.52	0.33	0	"	"
13:05	0.25	1.50	0.05	13.4	7.20	23.9	0.52	0.31	0	"	"
13:10	0.25	1.75	0.05	13.4	7.20	23.9	0.52	0.32	0	"	"

QPP
-246
-257
-262
-265
-265

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 13:11	SAMPLING ENDED AT: 13:15
PUMP OR TUBING DEPTH IN WELL (feet): 101	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: N	FIELD-FILTERED: N FILTER SIZE: µm	DUPLICATE: Y (N)	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-143	3	CG	40ml	HCl	-	-	8260	RFPP

MDS 2/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Tellefont</i>	SITE LOCATION: <i>Sarasota, FL</i>
WELL NO: <i>MW-144</i>	SAMPLE ID: <i>MW-144</i> DATE: <i>1/29/08</i>

PURGING DATA

WELL DIAMETER (inches): <i>2</i>	TUBING DIAMETER (inches): <i>1/4</i>	WELL SCREEN INTERVAL DEPTH: <i>140</i> feet to <i>150</i> feet	STATIC DEPTH TO WATER (feet): <i>14.82</i>	PURGE PUMP TYPE OR BAILER: <i>(PP) gear pump</i>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)

= gallons + (*0.0026* gallons/foot X *160* feet) + *0.25* gallons = *0.66* gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~145</i>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <i>145</i>	PURGING INITIATED AT: <i>1118</i>	PURGING ENDED AT: <i>1213</i>	TOTAL VOLUME PURGED (gallons): <i>3.30</i>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<i>12</i> 1130	<i>0.72</i>	<i>0.72</i>	<i>0.06</i>	<i>18.36</i>	<i>7.47</i>	<i>23.75</i>	<i>685</i>	<i>1.09</i>	<i>56.8</i>	<i>clear</i>	<i>yes</i>	<i>-283</i>
1136	0.36	1.08	0.06	18.81	7.44	23.72	688	0.92	114	clear	yes	-285
1142	0.36	1.44	0.06	19.03	7.42	23.76	692	0.86	101.7	clear	yes	-288
1148	0.36	1.80	0.06	19.33	7.40	23.84	694	0.80	70.6	clear	yes	-289
1153	0.30	2.10	0.06	19.56	7.37	23.82	698	0.77	44.2	clear	yes	-291
1158	0.30	2.40	0.06	19.60	7.37	23.84	698	0.74	28.8	clear	yes	-292
1203	0.30	2.70	0.06	19.65	7.36	23.86	699	0.70	20.0	clear	yes	-293
1208	0.30	3.00	0.06	19.71	7.34	23.87	700	0.67	13.8	clear	yes	-295
1213	0.30	3.30	0.06	19.74	7.35	23.86	700	0.65	10.37	clear	yes	-299

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Neil Smith ARCADIS</i>	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: <i>1214</i>	SAMPLING ENDED AT: <i>1220</i>
PUMP OR TUBING DEPTH IN WELL (feet): <i>~145</i>	SAMPLE PUMP FLOW RATE (mL per minute): <i>5100</i>	TUBING MATERIAL CODE: <i>PE</i>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>MW-144</i>	<i>3</i>	<i>CG</i>	<i>40mL</i>	<i>HCL</i>	<i>-</i>	<i>< 2</i>	<i>8260B + 8260SM RFPP</i>	

REMARKS: *Placed new dedicated tubing in well (NS) = Neil Smith*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahesse, FL
WELL NO: MW-145	SAMPLE ID: MW-145
DATE: 1.29.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet): 13.18	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 280 feet) + 0.25 gallons = 0.978 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 280	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 280	PURGING INITIATED AT: 10:50	PURGING ENDED AT: 11:25	TOTAL VOLUME PURGED (gallons): 3
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
11:00	1.0	1.0	0.1	15.3	7.54	23.9	0.640	0.38	16	clear	none	ORP -292
11:05	0.5	1.5	0.1	15.3	7.52	23.97	0.641	0.35	21	"	"	-299
11:10	0.5	2.0	0.1	16.05	7.51	23.9	0.640	0.34	9	"	"	-301
11:15	0.5	2.5	0.1	16.05	7.50	23.9	0.641	0.35	13	"	"	-302
11:20	0.25	2.75	0.05	16.25	7.50	24.0	0.642	0.34	5	"	"	-306
11:25	0.25	3.00	0.05	16.15	7.51	23.9	0.643	0.37	7	"	"	-309

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 11:26	SAMPLING ENDED AT: 11:30
PUMP OR TUBING DEPTH IN WELL (feet): 280 MDS 4/14/08	SAMPLE PUMP FLOW RATE (mL per minute): 2100 MDS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-145	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **Missing j-plug**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.05

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST	
WELL NO: mw-146		SAMPLE ID: mw-146	
DATE: 1-29-08			

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 19.5 feet to 24.5 feet	STATIC DEPTH TO WATER (feet): 2.95	PURGE PUMP TYPE OR BAILER: PP MDS 4/14/08
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 37.5 feet) + .25 gallons = .35 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 22	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 22	PURGING INITIATED AT: 1128	PURGING ENDED AT: 1141	TOTAL VOLUME PURGED (gallons): 2.21
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle right or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1135	1.19	1.19	0.17	3.50	6.6	24.8	.558	1.5	19.0	clear	N
1137	.34	1.53	0.17	3.69	6.5	24.7	.557	0.6	14.9	clear	N
1139	.34	1.87	0.17	3.80	6.4	24.8	.565	0.5	4.0	clear	N
1141	.34	2.21	0.17	3.80	6.4	24.8	.571	0.5	3.0	clear	N

ORP
-18
-10
-19
-33

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ACCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1143	SAMPLING ENDED AT: 1149
PUMP OR TUBING DEPTH IN WELL (feet): 22	SAMPLE PUMP FLOW RATE (mL per minute): ~640	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-146	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.05

SITE NAME: Tallevast		SITE LOCATION: Tallevast FL	
WELL NO: mw-147	SAMPLE ID: mw-147	DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 29.5 feet to 34.5 feet	STATIC DEPTH TO WATER (feet): 2.91	PURGE PUMP TYPE OR BAILER: PP	MOS 4/14/08
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)					

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)					
= 0 gallons + (0.0024 gallons/foot X 37.5 feet) + .25 gallons = 0.35 gallons					

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 32	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 32	PURGING INITIATED AT: 1038	PURGING ENDED AT: 1051	TOTAL VOLUME PURGED (gallons): 2.21
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1045	1.19	1.19	0.17	4.19	7.2	24.4	.511	1.6	0.05	clear	yno
1047	0.34	1.53	0.17	4.25	7.3	24.4	.507	1.5	0.46	clear	yno
1049	0.34	1.87	0.17	4.29	7.3	24.4	.502	0.5	0.11	clear	yno
1051	0.34	2.21	0.17	4.31	7.3	24.4	.508	0.5	0.25	clear	yno

0PP
-174
-186
-194
-203

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: Jennifer King	SAMPLING INITIATED AT: 1053	SAMPLING ENDED AT: 1059
PUMP OR TUBING DEPTH IN WELL (feet): 32	SAMPLE PUMP FLOW RATE (mL per minute): 640	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-147	3	CG	40ml	He1	40ml	-	8260	RFPF

REMARKS: **missing bolt**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.59

SITE NAME: Tallevast		SITE LOCATION: Tallevast Pl	
WELL NO: mw-148	SAMPLE ID: mw-148	DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 95.5 set to 105.5 feet	STATIC DEPTH TO WATER (feet): 13.52	PURGE PUMP TYPE OR BAILER: PP bell Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 108.5 feet) + .25 gallons = 0.53 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 100.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 100.5	PURGING INITIATED AT: 1042	PURGING ENDED AT: 1102	TOTAL VOLUME PURGED (gallons): 2.40
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or m/cm)	DISSOLVED OXYGEN (circle fig) or % saturation	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1056	1.68	1.68	0.12	13.75	7.20	24.5	.564	0.50	9.07	clear	yes
1058	0.24	1.92	0.12	13.75	7.21	24.6	.564	0.50	7.15	clear	yes
1100	0.24	2.16	0.12	13.75	7.22	24.6	.562	0.49	6.76	clear	yes
1102	0.24	2.40	0.12	13.75	7.22	24.5	.563	0.48	4.25	clear	yes

0.94
-260
-265
-270
-274

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1103	SAMPLING ENDED AT: 1108
PUMP OR TUBING DEPTH IN WELL (feet): 100.5	SAMPLE PUMP FLOW RATE (mL per minute): ~440	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: (N) N	FIELD-FILTERED: (N) Y	FILTER SIZE: _____ µm	DUPLICATE: Y (N)

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID / CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-148	3	CG	10mL	He1	-	-	8260	RFPP

REMARKS: **MISSING 2 bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.98

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST FL 30	
WELL NO: mw-149	SAMPLE ID: mw-149	DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet): 154	PURGE PUMP TYPE OR BAILER: PP 600 Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0024 gallons/foot X 158 feet) + .25 gallons = .46 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150	PURGING INITIATED AT: 1518	PURGING ENDED AT: 1540	TOTAL VOLUME PURGED (gallons): 2.86
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle 0/1 or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1534	2.08	2.08	.13	23.28	7.27	25.0	.568	0.45	4.56	clear	yes
1536	.26	2.34	.13	23.63	7.27	25.0	.568	0.45	4.56	clear	yes
1538	.26	2.60	.13	23.97	7.28	25.1	.568	0.45	4.80	clear	yes
1540	.26	2.86	.13	24.31	7.28	25.1	.568	0.45	4.49	clear	yes

RFPP
-217
-217
-219
-221

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1541	SAMPLING ENDED AT: 1549
PUMP OR TUBING DEPTH IN WELL (feet): 150	SAMPLE PUMP FLOW RATE (mL per minute): ~490	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-149	3	CG	40ml	HCl	40ml	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Tallevest</i>		SITE LOCATION: <i>Tallevest, FL</i>	
WELL NO: <i>MW-150</i>	SAMPLE ID: <i>MW-150</i>	DATE: <i>1/29/08</i>	

PURGING DATA

WELL DIAMETER (inches) <i>2</i>	TUBING DIAMETER (inches) <i>4</i>	WELL SCREEN INTERVAL DEPTH: <i>285</i> feet to <i>305</i> feet	STATIC DEPTH TO WATER (feet): <i>1393</i>	PURGE PUMP TYPE OR BAILER: <i>Geopump</i>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = <i>0</i> gallons + (<i>0.0026</i> gallons/foot X <i>338</i> feet) + <i>0.25</i> gallons = <i>1.13</i> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <i>285</i>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <i>285</i>	PURGING INITIATED AT: <i>1142</i>	PURGING ENDED AT: <i>1231</i>	TOTAL VOLUME PURGED (gallons): <i>3.92</i>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	<i>ORP</i>
<i>1225</i>	<i>3.44</i>	<i>3.44</i>	<i>0.08</i>	<i>22.26</i>	<i>7.67</i>	<i>24.1</i>	<i>1612</i>	<i>0.47</i>	<i>4.83</i>	<i>clear</i>	<i>yes</i>	<i>-259</i>
<i>1227</i>	<i>0.16</i>	<i>3.60</i>	<i>0.08</i>	<i>22.50</i>	<i>7.69</i>	<i>24.1</i>	<i>1611</i>	<i>0.47</i>	<i>3.86</i>	<i>clear</i>	<i>yes</i>	<i>-261</i>
<i>1229</i>	<i>0.16</i>	<i>3.76</i>	<i>0.08</i>	<i>22.58</i>	<i>7.69</i>	<i>24.1</i>	<i>1613</i>	<i>0.48</i>	<i>2.28</i>	<i>clear</i>	<i>yes</i>	<i>-261</i>
<i>1231</i>	<i>0.16</i>	<i>3.92</i>	<i>0.08</i>	<i>22.64</i>	<i>7.69</i>	<i>24.1</i>	<i>1613</i>	<i>0.47</i>	<i>2.86</i>	<i>clear</i>	<i>yes</i>	<i>-261</i>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>BUDGET STAFF / ARCADIS</i>	SAMPLER(S) SIGNATURES: <i>Budget Staff</i>	SAMPLING INITIATED AT: <i>1233</i>	SAMPLING ENDED AT: <i>1240</i>
PUMP OR TUBING DEPTH IN WELL (feet): <i>285</i>	SAMPLE PUMP FLOW RATE (mL per minute): <i>~320</i>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>MW-150</i>	<i>3</i>	<i>CG</i>	<i>40ml</i>	<i>HCl</i>	<i>—</i>	<i>—</i>	<i>8260</i>	<i>RFPP</i>

REMARKS: *missing 2 bolts*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevest MDS 4/19/08	SITE LOCATION: Garasota FL MDS 4/19/08
WELL NO: MW-151	DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 17 feet to 22 feet	STATIC DEPTH TO WATER (feet): 3.95	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 15 gallons + (0.0026 gallons/foot X 22 feet) + .25 gallons = .31 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 20	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 20	PURGING INITIATED AT: 1028	PURGING ENDED AT: 1050	TOTAL VOLUME PURGED (gallons): 0.6

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1033			0.1	4.99	6.1	26.2	0.785	3.7	27.4	Clear	none
1036			0.1	5.08	6.1	26.2	0.784	3.4	16.6		
1039				5.10	6.0	26.2	0.781	2.9	14.9		
1042				5.10	6.0	26.2	0.777	2.7	10.86		
1045				5.11	6.0	26.3	0.778	2.6	6.29		
1048				5.10	6.0	26.3	0.775	2.5	7.41		
1050 → Sampled @ 1150 ←											

58
60
64
64
65
64

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KLC/ARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1050	SAMPLING ENDED AT: 1155
PUMP OR TUBING DEPTH IN WELL (feet): 20	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/19/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-151	3	CG	40ml	HCL	—	—	8260	RFPP

MDS
4/19/08

REMARKS: **Initial purge: clear no distinct odor.**

MATERIAL CODES: JAG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talhoust.</u>	SITE LOCATION: <u>Sarasota FL</u>
WELL NO: <u>MW-152</u>	SAMPLE ID: <u>MW-152</u> DATE: <u>1/23/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>0.25</u>	WELL SCREEN INTERVAL DEPTH: <u>37.5</u> feet to <u>42.5</u> feet	STATIC DEPTH TO WATER (feet): <u>7.21</u>	PURGE PUMP TYPE OR BAILER: <u>GeoPump (Peristaltic)</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable $= (42.5 \text{ feet} - 7.21 \text{ feet}) \times 0.163 \text{ gallons/foot} = 5.75 \text{ gallons}$				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) $= \text{gallons} + (\text{gallons/foot} \times \text{feet}) + \text{gallons} = \text{gallons}$				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>40</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>40</u>	PURGING INITIATED AT: <u>10:10</u>	PURGING ENDED AT: <u>10:44</u>	TOTAL VOLUME PURGED (gallons): <u>30</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
10:15	0.3	0.3	0.1	7.22	7.40	27.22	0.587	5.07	1.48	clear	clear	190
10:18	0.3	0.6	0.1	7.99	7.38	26.55	0.625	3.90	1.18	"	"	175
10:21		0.9	0.1	8.59	7.38	26.47	0.625	3.77	1.23	"	"	167
10:24		1.2	0.1	9.47	7.37	26.44	0.624	3.69	1.78	"	"	75
10:27		1.5	0.1	10.15	7.36	26.42	0.628	3.60	1.55	"	"	-61
10:30		1.8	0.1	10.62	7.36	26.44	0.631	3.55	1.41	"	"	-84
10:33		2.1	0.1	11.05	7.35	26.52	0.631	3.40	1.39	"	"	-113
10:36		2.4	0.1	11.60	7.35	26.53	0.639	3.18	1.90	"	"	-131
10:39		2.7	0.1	12.14	7.36	26.44	0.639	3.08	1.56	"	"	-134
10:42		3.0	0.1	12.69	7.36	26.28	0.650	3.00	0.99	"	"	-146

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Nathan Smith</u>	SAMPLING INITIATED AT: <u>10:44</u>	SAMPLING ENDED AT: <u>10:48</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>40</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>N</u>	FIELD-FILTERED: <u>N</u>	DUPLICATE: Y <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-152	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

needs 2 bolts.

DEP-SOP-001/01
Form FD 9000-24

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 9/14/08	SITE LOCATION: <u>Sarasota FL</u> MDS 9/14/08
WELL NO: <u>MW-153</u>	SAMPLE ID: <u>MW-153</u> DATE: <u>1/24/08</u>

WELL DIAMETER (inches):	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>9.7</u> feet to <u>10.7</u> feet	STATIC DEPTH TO WATER (feet): <u>10.70</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME	<u>0.53</u> MDS 9/14/08
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>10.7</u> feet) + <u>0.25</u> gallons = <u>0.028</u> gallons	

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>103</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>103</u>	PURGING INITIATED AT: <u>1005</u>	PURGING ENDED AT: <u>1031</u>	TOTAL VOLUME PURGED (gallons):
---	---	-----------------------------------	-------------------------------	--------------------------------

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1008				10.37	7.13	25.57	0.558	1.47	143	GRAY	Strong
1011				10.43	7.16	25.65	0.561	1.12	60.5	Cloudy	
1014				10.69	7.18	25.70	0.569	1.04	37.0	Clear	
1017				10.72	7.18	25.77	0.570	1.01	27.90		
1020				10.75	7.18	25.80	0.573	0.99	20.4		
1023				10.78	7.18	25.79	0.580	0.98	24.4		
1026				10.78	7.18	25.77	0.591	0.98	14.5		
1029				10.78	7.18	25.80	0.593	0.97	10.79		
1031				10.78	7.18	25.73	0.583	0.97	9.48		

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasey Cornwall / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>1035</u>	SAMPLING ENDED AT: <u>1045</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>103</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u> MDS 9/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: <u>1</u> µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-153</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS: * Strong odor, initial purge dark gray let Purge 3 minutes before hooking up to Hombic's

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* @ time of first reading light Gray incolor.

GROUNDWATER SAMPLING LOG

11

SITE NAME: Talleast ^{MOS} 4/14/08		SITE LOCATION: Sarasota FL ^{MOS} 4/14/08	
WELL NO: MW-154	SAMPLE ID: MW-154 ^{MOS} 4/14/08	DATE: 1/23/08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 144.5 feet to 154.5 feet	STATIC DEPTH TO WATER (feet): 11.25	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= \emptyset gallons + (0.0021 gallons/foot X **154.5** feet) + 0.25 gallons = **.65** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150'	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150'	PURGING INITIATED AT: 936	PURGING ENDED AT: 956 ^{MOS} 4/14/08	TOTAL VOLUME PURGED (gallons):
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939
942
950
953
956
1000

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1039				11.97	7.4	24.9	0.433	3.3	2.16	Clear	Slight*
1042				12.34	7.4	25.3	0.427	2.8	1.56	Clear	
1050				12.82	7.4	25.4	0.427	2.7	1.27		
1053				13.03	7.4	25.3	0.428	2.6	1.03		
1056					7.4	25.5	0.424	2.5	1.03		
→ SAMPLED @ 1000 ←											

159
32
8
-9
-3

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Kasei / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1000	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 155	SAMPLE PUMP FLOW RATE (mL per minute): <100 ^{MOS} 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-154	3	CG	40ml	HCl	—	—	8260	RFPP

^{MOS} **4/14/08**

REMARKS: **Slight rotten egg/sulfur smell @ initial purge**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

3.785 L = 1 Gallon.

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u>	<u>MDS 9/19/08</u>	SITE LOCATION: <u>Sarasota FL</u>	<u>MDS 9/19/08</u>
WELL NO: <u>MW-155</u>	SAMPLE ID: <u>MW-155</u>	DATE: <u>1/24/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>0.25</u>	WELL SCREEN INTERVAL DEPTH: <u>284</u> feet to <u>304</u> feet	STATIC DEPTH TO WATER (feet): <u>9.93</u>	PURGE PUMP TYPE OR BAILER: <u>Peristaltic</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (<u>304</u> feet - <u>9.93</u> feet) X <u>47.0163</u> gallons/foot = <u>47.9</u> gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>10:15</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>10:15</u>	PURGING ENDED AT: <u>11:25</u>	TOTAL VOLUME PURGED (gallons): <u>5.5</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
10:25	0.5	1.0	0.1	11.62	7.4	22.6	1.05	9.1	84.0	slightly gray	-
10:35		1.5	0.1	14.01	7.6	24.3	0.763	3.7	1060	blue gray	-
10:45		2.0	0.1	15.25	-	-	-	8.7	330	light gray	-
10:55		2.5	0.1	16.31	7.6	24.1	1.00	3.2	114	clear	-
11:00		3.0	0.1	16.75	7.6	23.9	1.00	3.0	123	clear	-
11:05		3.5	0.1	17.20	7.6	23.8	1.00	2.9	40.3	clear	-
11:10		4.0	0.1	17.36	7.6	23.9	1.00	2.9	24.7	clear	-
11:15		4.5	0.1	17.58	7.6	23.8	0.99	2.8	20.6	clear	-
11:20		5.0	0.1	17.74	7.6	23.8	0.99	2.7	15.0	clear	-
11:25		5.5	0.1	17.86	7.6	23.9	0.99	2.7	10.88	clear	-

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith</u>	SAMPLER(S) SIGNATURES: <u>Nathan Smith</u>	SAMPLING INITIATED AT: <u>11:25</u>	SAMPLING ENDED AT: <u>11:30</u> *
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u> <u>MDS 9/19/08</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>(Y)</u> N	FIELD-FILTERED: Y <u>(N)</u> FILTER SIZE: _____ µm	DUPLICATE: Y <u>(N)</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-155	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS: * Time Sampled @ 11:30

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* Turning blue gray. Removed flow thru cell until turbidity decreases.
** clearing up.
No tubing in well. We replaced tubing (teflon). Needs 2 cover bolts.
Revision Date: February 1, 2004

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallerast
WELL NO: MW-156	SAMPLE ID: MW-156
DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 1.84	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.026 gallons/foot X 20 feet) + 0.25 gallons = 0.30 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 20	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5	PURGING INITIATED AT: 13:58	PURGING ENDED AT: 14:23	TOTAL VOLUME PURGED (gallons): 2.5

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1408	1.0	1.0	0.1	5.08	6.19	23.89	0.689	0.30	13.2			-131
1411	0.3	1.3		5.14	6.17	24.08	0.650	0.30	6.16 12.0			-135
1414	0.3	1.6		5.22	6.15	24.09	0.628	0.28	10.93	AC		-143
1417	0.3	1.9		5.34	6.11	24.24	0.573	0.20	9.42			-145
1420	0.3	2.2		5.34	6.11	24.24	0.573	0.20	7.96			-146
1423	0.3	2.5		5.36	6.11	24.25	0.573	0.20	7.84			-147
							MS/cm					

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amey Coats / Accords	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1424	SAMPLING ENDED AT: 1425
PUMP OR TUBING DEPTH IN WELL (feet): 17.5	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: Y	FIELD-FILTERED: Y	FILTRATION EQUIPMENT TYPE: N	DUPLICATE: Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-156	3	CG	40mL	HCl	—	—	8260	PP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC</u>		SITE LOCATION: <u>Talleast</u>	
WELL NO: <u>MW-157</u>	SAMPLE ID: <u>MW-157</u>	DATE: <u>1-24-08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>33</u> feet to <u>38</u> feet	STATIC DEPTH TO WATER (feet): <u>9.82</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0020</u> gallons/foot X <u>38</u> feet) + <u>0.35</u> gallons = <u>0.35</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>38</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>35.5</u>	PURGING INITIATED AT: <u>13:57</u>	PURGING ENDED AT: <u>14:16</u>	PURGED (gallons): <u>2.0</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
14:07	1.0	1.0	0.1	16.1	7.29	25.05	0.775	1.43	5.02	amber	none
14:10	0.3	1.3	0.1	18.18	7.21	25.14	0.753	1.05	7.88	clear	none
14:13	0.3	1.7	0.1	18.63	7.19	25.21	0.747	0.99	5.88	clear	none
14:16	0.3	2.0	0.1	17.55	7.18	25.32	0.743	0.95	5.80	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Darrin Johnson Arrows</u>	SAMPLER(S) SIGNATURES: <u>Darrin Johnson</u>	SAMPLING INITIATED AT: <u>14:17</u>	SAMPLING ENDED AT: <u>14:22</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>35.5</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>400</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> N	FILTRATION EQUIPMENT TYPE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> <u>Y</u> <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-157</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

REMARKS: Dep 7-012408 no dupe

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 4/14/08	SITE LOCATION: <u>Sarasota FL</u> MDS 4/14/08
WELL NO: <u>MW-158</u>	SAMPLE ID: <u>MW-158</u> DATE: <u>1/24/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>100</u> feet to <u>110</u> feet	STATIC DEPTH TO WATER (feet): <u>12.82</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1450 1448	<u>0.5</u>	<u>0.5</u>	<u>0.1</u>	<u>13.14</u>	<u>7.6</u>	<u>24.5</u>	<u>0.515</u>	<u>3.5</u>	<u>11.4</u>	<u>Clear</u>	
				<u>13.13</u>	<u>7.6</u>	<u>24.8</u>	<u>0.510</u>	<u>2.8</u>	<u>21.1</u>		
1500 1445				<u>13.18</u>	<u>7.6</u>	<u>24.8</u>	<u>0.503</u>	<u>2.5</u>	<u>16.8</u>		
				<u>13.14</u>	<u>7.6</u>	<u>24.7</u>	<u>0.502</u>	<u>2.5</u>	<u>8.06</u>		
				<u>13.14</u>	<u>7.5</u>	<u>24.7</u>	<u>0.502</u>	<u>2.4</u>	<u>4.40</u>		
				<u>13.11</u>	<u>7.5</u>	<u>24.7</u>	<u>0.501</u>	<u>2.3</u>	<u>3.10</u>		
					<u>7.5</u>	<u>24.7</u>	<u>0.51</u>	<u>2.3</u>	<u>2.31</u>		
				<u>Sampled @ 1525</u>							

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>ARCADIS</u> MDS 4/14/08	SAMPLER(S) SIGNATURES:
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u>400</u> MDS 4/14/08
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm
Duplication Equipment Type: _____	
SAMPLING INITIATED AT: <u>1525</u>	SAMPLING ENDED AT: <u>1535</u>
TUBING MATERIAL CODE: _____	
DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-158</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-159	SAMPLE ID: MW-159 DATE: 1-24-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 140 feet to 150 feet	STATIC DEPTH TO WATER (feet): 13.34	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 60026 gallons/foot X 153 feet + .25 gallons = 0.65 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 145	PURGING INITIATED AT: 14:48	PURGING ENDED AT: 15:09	TOTAL VOLUME PURGED (gallons): 2.1
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
15:03	1.5	1.5	0.1	19.12	8.38	24.96	0.639	0.93	3.38	clear	none
15:06	0.3	1.8	0.1	19.18	8.52	24.87	0.639	1.02	3.30	clear	none
15:09	0.3	2.1	0.1	19.30	8.55	24.91	0.638	0.92	3.35	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURE: Darrin Johnson	SAMPLING INITIATED AT: 15:10	SAMPLING ENDED AT: 15:15
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PUMP OR TUBING DEPTH IN WELL (feet): 145	SAMPLE PUMP FLOW RATE (mL per minute): 5100	TUBING MATERIAL CODE:
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	Filtration Equipment Type: FILTER SIZE: µm
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: Y <input checked="" type="radio"/> N

SAMPLE ID CODE	SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-159	3	CG	90ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallevast, FL
WELL NO: MW-160	SAMPLE ID: MW-160
DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/8	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet): 1253	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1448	PURGING ENDED AT: 1	TOTAL VOLUME PURGED (gallons):
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle) (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1504				15.50	7.48	24.55	0.801	0.47	2.72	NONE	NONE
1507				15.90	7.52	24.44	0.798	0.45	2.08		
1510				16.40	7.54	24.53	0.792	0.42	1.54	↓	↓
1513				16.56	7.54	24.40	0.793	0.41	1.31	↓	↓

ORP
-277
-278
-283
-284

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: MDS 2/15/08 IARCADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1514	SAMPLING ENDED AT: 1515
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 2100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-160	3	CG	40mL	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/19/08	SITE LOCATION: Sarasota FL MDS 4/19/08
WELL NO: MW-161	SAMPLE ID: MW-161 DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 381 feet to 40 feet	STATIC DEPTH TO WATER (feet): 11.85	PURGE PUMP TYPE OR BAILER: Geopump (Peristaltic)
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (401 feet - 11.85 feet) X 63.4 gallons/foot = 63.4 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 3		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 15:05	PURGING ENDED AT: 15:45	TOTAL VOLUME PURGED (gallons): 4.0						
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
15:20	1.5	1.5	0.1	12.05	8.39	24.82	1.16	4.24	23.7	clear	—	029
15:25	0.5	2.0	0.1	12.10	8.52	24.54	1.16	1.96	26.8	"	—	-298
15:30	0.5	2.5	0.1	12.11	8.65	23.92	1.16	1.56	26.3	"	—	-300
15:35	0.5	3.0	0.1	12.11	8.71	24.10	1.16	1.31	22.4	"	—	-307
15:40	0.5	3.5	0.1	12.11	8.71	24.07	1.16	1.33	17.4	"	—	-311
15:45	0.5	4.0	0.1	12.11	8.63	23.94	1.17	1.29	13.5	"	—	-315
MDS 4/19/08												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN SMITH / MCDONALD'S		SAMPLER(S) SIGNATURES: <i>Nathan Smith</i>		SAMPLING INITIATED AT: 15:45	SAMPLING ENDED AT: 15:54 *
PUMP OR TUBING DEPTH IN WELL (feet):		SAMPLE PUMP FLOW RATE (mL per minute): 4100 MDS 4/19/08		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: <input type="text"/> µm		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-161	3	CG	90ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* Time sample @ 1554

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallahassee</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>MW-162</u>	SAMPLE ID: <u>MW-162</u> DATE: <u>1/30/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>4</u>	WELL SCREEN INTERVAL DEPTH: <u>11</u> feet to <u>16</u> feet	STATIC DEPTH TO WATER (feet): <u>2.47</u>	PURGE PUMP TYPE OR BAILER: <u>PPS geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + <u>0.0026</u> gallons/foot X <u>30</u> feet + <u>0.25</u> gallons = <u>0.328</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>5/3.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~13.5</u>	PURGING INITIATED AT: <u>1150</u>	PURGING ENDED AT: <u>1215</u>	TOTAL VOLUME PURGED (gallons): <u>1.68</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/l)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<u>1157</u>	<u>0.42</u>	<u>0.42</u>	<u>0.06</u>	<u>3.10</u>	<u>6.60</u>	<u>23.34</u>	<u>0.649</u>	<u>2.53</u>	<u>9.52</u>	<u>clear</u>	<u>yes</u>	<u>-70</u>
<u>1200</u>	<u>0.18</u>	<u>0.60</u>	<u>0.06</u>	<u>3.15</u>	<u>6.46</u>	<u>23.35</u>	<u>0.755</u>	<u>2.04</u>	<u>7.69</u>	<u>clear</u>	<u>yes</u>	<u>-84</u>
<u>1203</u>	<u>0.18</u>	<u>0.78</u>	<u>0.06</u>	<u>3.20</u>	<u>6.44</u>	<u>23.48</u>	<u>0.772</u>	<u>1.99</u>	<u>6.40</u>	<u>clear</u>	<u>yes</u>	<u>-82</u>
<u>1206</u>	<u>0.18</u>	<u>0.96</u>	<u>0.06</u>	<u>3.21</u>	<u>6.41</u>	<u>23.43</u>	<u>0.631</u>	<u>1.94</u>	<u>3.1</u>	<u>clear</u>	<u>yes</u>	<u>-81</u>
<u>1209</u>	<u>0.18</u>	<u>1.14</u>	<u>0.06</u>	<u>3.20</u>	<u>6.37</u>	<u>23.18</u>	<u>0.803</u>	<u>0.95</u>	<u>1.6</u>	<u>clear</u>	<u>yes</u>	<u>-79</u>
<u>1212</u>	<u>0.18</u>	<u>1.32</u>	<u>0.06</u>	<u>3.20</u>	<u>6.35</u>	<u>23.34</u>	<u>0.819</u>	<u>0.82</u>	<u>1.9</u>	<u>clear</u>	<u>yes</u>	<u>-77</u>
<u>1215</u>	<u>0.18</u>	<u>1.50</u>	<u>0.06</u>	<u>3.20</u>	<u>6.34</u>	<u>23.46</u>	<u>0.831</u>	<u>0.81</u>	<u>2.24</u>	<u>clear</u>	<u>yes</u>	<u>-75</u>
<u>1218</u>	<u>0.18</u>	<u>1.68</u>	<u>0.06</u>	<u>3.21</u>	<u>6.34</u>	<u>23.47</u>	<u>0.834</u>	<u>0.72</u>	<u>1.76</u>	<u>clear</u>	<u>yes</u>	<u>-71</u>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCAAS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>1220</u>	SAMPLING ENDED AT: <u>1225</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~13.5</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-162</u>	<u>3</u>	<u>CG</u>	<u>42ml</u>	<u>HCL</u>	<u>-</u>	<u>4.2</u>	<u>8260B/garson RFP</u>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC Facility</u>	SITE LOCATION: <u>Talleras, FL</u>
WELL NO: <u>MW-163</u>	SAMPLE ID: <u>MW-163</u> DATE: <u>1.30.08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>3</u> feet to <u>35</u> feet	STATIC DEPTH TO WATER (feet): <u>9.8</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + <u>0.0020</u> gallons/foot X <u>35</u> feet + <u>0.25</u> gallons = <u>0.34</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>32</u>		FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>32</u>		PURGING INITIATED AT: <u>14:12</u>		PURGING ENDED AT: <u>14:35</u>		TOTAL VOLUME PURGED (gallons): <u>2.3</u>			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>14:20</u>	<u>0.8</u>	<u>0.8</u>	<u>0.1</u>	<u>10.4</u>	<u>7.10</u>	<u>23.33</u>	<u>0.673</u>	<u>0.58</u>	<u>0</u>	<u>clear</u>	<u>none</u>
<u>14:25</u>	<u>0.5</u>	<u>1.3</u>	<u>0.1</u>	<u>10.9</u>	<u>6.91</u>	<u>23.17</u>	<u>0.669</u>	<u>0.32</u>	<u>10</u>	<u>"</u>	<u>"</u>
<u>14:30</u>	<u>0.5</u>	<u>1.8</u>	<u>0.1</u>	<u>10.9</u>	<u>6.91</u>	<u>23.15</u>	<u>0.662</u>	<u>0.30</u>	<u>9</u>	<u>"</u>	<u>"</u>
<u>14:35</u>	<u>0.5</u>	<u>2.3</u>	<u>0.1</u>	<u>10.9</u>	<u>6.92</u>	<u>23.14</u>	<u>0.669</u>	<u>0.30</u>	<u>7</u>	<u>"</u>	<u>"</u>

ORP
-206
-242
-248
-252

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Micah Forbes / ARCADIS</u>		SAMPLER(S) SIGNATURES: <u>[Signature]</u>		SAMPLING INITIATED AT: <u>14:36</u>	SAMPLING ENDED AT: <u>14:40</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>32</u> <u>MOS 4/14/08</u>		SAMPLE PUMP FLOW RATE (mL per minute): <u>100</u> <u>MOS 4/14/08</u>		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u> <u>N</u>		FIELD-FILTERED: <u>Y</u> <u>N</u> FILTER SIZE: <u>0.1</u> µm		DUPLICATE: <u>Y</u> <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW163</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>2260</u>	<u>RFPP</u> <u>MOS 4/14/08</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC Facility</u>	SITE LOCATION: <u>Tallevast, FL</u>
WELL NO: <u>MW-164</u>	SAMPLE ID: <u>MW-164</u> DATE: <u>1-30-08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>92</u> feet to <u>102</u> feet	STATIC DEPTH TO WATER (feet): <u>16.3</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + <u>0.0026</u> gallons/foot X <u>102</u> feet + <u>0.25</u> gallons = <u>0.51</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>98</u>		FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>98</u>		PURGING INITIATED AT: <u>11:08</u>		PURGING ENDED AT: <u>11:35</u>		TOTAL VOLUME PURGED (gallons): <u>2.7</u>			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>11:15</u>	<u>0.7</u>	<u>0.7</u>	<u>0.1</u>	<u>16.6</u>	<u>7.21</u>	<u>23.09</u>	<u>0.695</u>	<u>0.39</u>	<u>6</u>	<u>clear</u>	<u>none</u>
<u>11:20</u>	<u>0.5</u>	<u>1.2</u>	<u>0.1</u>	<u>16.6</u>	<u>7.22</u>	<u>23.15</u>	<u>0.692</u>	<u>0.34</u>	<u>10</u>	<u>"</u>	<u>"</u>
<u>11:25</u>	<u>0.5</u>	<u>1.7</u>	<u>0.1</u>	<u>16.6</u>	<u>7.20</u>	<u>23.08</u>	<u>0.696</u>	<u>0.33</u>	<u>4</u>	<u>"</u>	<u>"</u>
<u>11:30</u>	<u>0.5</u>	<u>2.2</u>	<u>0.1</u>	<u>16.6</u>	<u>7.21</u>	<u>23.07</u>	<u>0.690</u>	<u>0.32</u>	<u>4</u>	<u>"</u>	<u>"</u>
<u>11:35</u>	<u>0.5</u>	<u>2.7</u>	<u>0.1</u>	<u>16.6</u>	<u>7.19</u>	<u>23.05</u>	<u>0.689</u>	<u>0.33</u>	<u>7</u>	<u>"</u>	<u>"</u>

ORP
-295
-300
-302
-302
-303

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Micah Forbes / ARCADIS</u>		SAMPLER(S) SIGNATURES: <u>[Signature]</u>		SAMPLING INITIATED AT: <u>11:36</u>	SAMPLING ENDED AT: <u>11:40</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>98</u>		SAMPLE PUMP FLOW RATE (mL per minute): <u>1100</u>		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u> <u>N</u>		FIELD-FILTERED: <u>Y</u> <u>N</u>		FILTER SIZE: <u>1</u> µm	
SAMPLE CONTAINER SPECIFICATION		SAMPLE PRESERVATION		DUPLICATE: <u>Y</u>	

SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
<u>MW164</u>	<u>3</u>	<u>CG</u>	<u>90ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

MDS
A/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u>	SITE LOCATION: <u>Swasoty FL</u>
WELL NO: <u>MW-165</u>	SAMPLE ID: <u>MW-165</u> DATE: <u>1/20/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>142</u> feet to <u>152</u> feet	STATIC DEPTH TO WATER (feet): <u>14.70</u>	PURGE PUMP TYPE: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<u>0.0026</u> gallons/foot X <u>180</u> feet) + <u>0.25</u> gallons = <u>0.66</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~147</u>		FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~147</u>		PURGING INITIATED AT: <u>1115</u>		PURGING ENDED AT: <u>1134</u>		TOTAL VOLUME PURGED (gallons): <u>1.33</u>			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle µg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>1125</u>	<u>0.70</u>	<u>0.70</u>	<u>0.07</u>	<u>21.51</u>	<u>7.43</u>	<u>23.80</u>	<u>0.767</u>	<u>0.39</u>	<u>2.26</u>	<u>clear</u>	<u>yes</u>
<u>1128</u>	<u>0.21</u>	<u>0.91</u>	<u>0.07</u>	<u>23.33</u>	<u>7.43</u>	<u>23.88</u>	<u>0.764</u>	<u>0.38</u>	<u>1.80</u>	<u>clear</u>	<u>yes</u>
<u>1131</u>	<u>0.21</u>	<u>1.12</u>	<u>0.07</u>	<u>24.14</u>	<u>7.43</u>	<u>23.97</u>	<u>0.765</u>	<u>0.39</u>	<u>0.64</u>	<u>clear</u>	<u>yes</u>
<u>1134</u>	<u>0.21</u>	<u>1.33</u>	<u>0.07</u>	<u>24.35</u>	<u>7.43</u>	<u>23.95</u>	<u>0.766</u>	<u>0.39</u>	<u>0.81</u>	<u>clear</u>	<u>yes</u>

ORP
-231
-237
-240
-240

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>		SAMPLER(S) SIGNATURE(S): <u>Neil Smith</u>		SAMPLING INITIATED AT: <u>1135</u>	SAMPLING ENDED AT: <u>1140</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~147</u>		SAMPLE PUMP FLOW RATE (mL per minute): <u>2100ml</u>		TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>⓪</u> N		FIELD-FILTERED: <u>X</u> <u>FF</u> FILTER SIZE: _____ µm		DUPLICATE: Y <u>X</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-165</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCL</u>	<u>-</u>	<u>~2</u>	<u>8260B+8260S</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallahassee</u>	SITE LOCATION: <u>Swasoh, FL</u>
WELL NO: <u>MW-166</u>	SAMPLE ID: <u>MW-166</u> DATE: <u>1/30/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>281</u> feet to <u>301</u> feet	STATIC DEPTH TO WATER (feet): <u>1420</u>	PURGE PUMP TYPE OR BAILER: <u>(pp) geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<u>0.001</u> gallons/foot X <u>295</u> feet) + <u>0.25</u> gallons = <u>1.00</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~281</u>		FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~281</u>		PURGING INITIATED AT: <u>1400</u>		PURGING ENDED AT: <u>1439</u>		TOTAL VOLUME PURGED (gallons): <u>1.50</u>			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/c or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>1425</u>	<u>1.00</u>	<u>1.00</u>	<u>0.04</u>	<u>15.31</u>	<u>8.16</u>	<u>23.69</u>	<u>0.702</u>	<u>0.42</u>	<u>17.6</u>	<u>clear</u>	<u>yes</u>
<u>1432</u>	<u>0.25</u>	<u>1.25</u>	<u>0.04</u>	<u>15.32</u>	<u>8.05</u>	<u>23.66</u>	<u>0.711</u>	<u>0.38</u>	<u>15</u>	<u>clear</u>	<u>yes</u>
<u>1439</u>	<u>0.25</u>	<u>1.50</u>	<u>0.04</u>	<u>15.32</u>	<u>8.02</u>	<u>23.65</u>	<u>0.712</u>	<u>0.37</u>	<u>15.4</u>	<u>clear</u>	<u>yes</u>

25

ORP
-277
-288
-296

Neil Smith

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: <u>1439</u>	SAMPLING ENDED AT: <u>1445</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~281</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>≤ 100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> <u>N</u>	FIELD-FILTERED: <u>Y</u> <u>N</u> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-166</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCL</u>	<u>—</u>	<u>~2</u>	<u>22603626036</u>	<u>RFPP</u>
<u>DUP-227!!</u>								

REMARKS: Ants!

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: STA	SITE LOCATION: Sarasota, FL
WELL NO: MW-167	DATE: 1/23/08
SAMPLE ID: MW-167	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 20 feet to 25 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP geom
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 30 feet) + 0.50 gallons = 0.578 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~22.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~22.5	PURGING INITIATED AT: 1540	PURGING ENDED AT: 1555	TOTAL VOLUME PURGED (gallons): 1.50
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle read or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1546	0.60	0.60	0.10	5.60	4.95	29.00	352	4.26	9.68	clear	NO	53
1549	0.30	0.90	0.10	5.60	4.92	29.01	349	5.12	7.50	clear	NO	52
1552	0.30	1.20	0.10	5.60	4.90	29.10	349	4.89	6.4	clear	NO	50
1555	0.30	1.50	0.10	5.60	4.93	29.12	349	5.00	4.1	clear	NO	50
<i>Handwritten scribble</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1556	SAMPLING ENDED AT: 1600
PUMP OR TUBING DEPTH IN WELL (feet): ~22.5	SAMPLE PUMP FLOW RATE (mL per minute): ~100ml	TUBING MATERIAL CODE: PG	
FIELD DECONTAMINATION: X N	FIELD-FILTERED: Y N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-167	3	CG	40ml	HCL	-	~2	SRG08 FSM	RFP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-168	SAMPLE ID: MW-168 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 42 feet to 47 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 52 feet) + 0.5 gallons = 0.64 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 44.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 44.5	PURGING INITIATED AT: 1619	PURGING ENDED AT: 1639	TOTAL VOLUME PURGED (gallons): 1.60								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circled mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	CRP
1627	0.64	0.64	0.08	6.55	6.51	27.61	493	6.71	17.3	clear	no	-133
1630	0.24	0.88	0.08	6.55	6.52	27.62	493	2.02	14.4	clear	no	-132
1633	0.24	1.12	0.08	6.55	6.55	27.73	494	6.72	12.7	clear	no	-131
1636	0.24	1.36	0.08	6.55	6.56	27.81	495	6.30	12.3	clear	no	-131
1639	0.24	1.60	0.08	6.55	6.57	27.91	497	6.20	11.2	clear	no	-133

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NS/SW A RCADIS			SAMPLER(S) SIGNATURES: <i>[Signature]</i>			SAMPLING INITIATED AT: 1642		SAMPLING ENDED AT: 1647			
PUMP OR TUBING DEPTH IN WELL (feet): 44.5			SAMPLE PUMP FLOW RATE (mL per minute): 100			TUBING MATERIAL CODE: PE					
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N			FIELD-FILTERED: Y <input checked="" type="checkbox"/> N <input type="checkbox"/> FILTER SIZE: _____ µm			DUPLICATE: Y <input checked="" type="checkbox"/> N <input type="checkbox"/>					
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH					
MW-168	3	CG	40mL	HCl	40mL	-	8260+8260SIM		RFPP		

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALUEAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-169	SAMPLE ID: MW-169 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 108 feet to 110 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026** gallons/foot X **123** feet) + **0.5** gallons = **0.82** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 113	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 113	PURGING INITIATED AT: 1448	PURGING ENDED AT: 1506	TOTAL VOLUME PURGED (gallons): 1.8
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circled mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1457	0.9	0.9	0.1	25.05	11.30	27.65	1640	0.60	1.69	clear	no	-278
1500	0.3	1.2	0.1	25.73	11.33	27.67	1640	0.89	1.39	clear	no	-262
1503	0.3	1.5	0.1	26.41	11.34	27.69	1640	1.34	0.95	clear	no	-245
1506	0.3	1.8	0.1	27.03	11.35	27.57	1640	1.39	0.77	clear	no	-240
1510												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/NS / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1510	SAMPLING ENDED AT: 1515
PUMP OR TUBING DEPTH IN WELL (feet): 113	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-169	3	CG	40mL	HCl	40mL	—	B260+826051M	RFPP

REMARKS: *** PRESSURE IN WELL - BEWARE OPENING CAP!**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellerus STA</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>MW-170</u>	SAMPLE ID: <u>MW-170</u> DATE: <u>1/23/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>280</u> feet to <u>300</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>(pp) geopump</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= gallons + (0.0026 gallons/foot X 291 feet) + 0.50 gallons = 1.25 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~290</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~290</u>	PURGING INITIATED AT: <u>1615</u>	PURGING ENDED AT: <u>1649</u>	TOTAL VOLUME PURGED (gallons): <u>1.70</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/L)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1640	1.25	1.25	0.05	22.74	7.68	27.68	1080	0.30	35	clear	yes
1643	0.15	1.40	0.05	22.84	7.67	27.64	1060	0.29	13.1	clear	yes
1646	0.15	1.55	0.05	22.86	7.67	27.65	1070	0.29	9.37	clear	yes
1649	0.15	1.70	0.05	22.96	7.67	27.66	1070	0.28	8.1	clear	yes

NOS
4/17/08
ORP
-231
-238
-239
-238

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil Smith</u>	SAMPLING INITIATED AT: <u>1655</u>	SAMPLING ENDED AT: <u>1705</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~290</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>~100ml</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="checkbox"/> X	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-170	3	CG	40ml	HCL	-	~2	8260B1SIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast	SITE LOCATION: Sarasota FL
WELL NO: HW-171	SAMPLE ID: MW-171
DATE: 1/23/08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 35 feet to 40 feet	STATIC DEPTH TO WATER (feet): 7.93	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY				
= (40 feet - 7.93 feet) X 0.0026 gallons/foot = 0.35 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME				
= 0.35 gallons + (0.0026 gallons/foot X 40') + 0.25 gallons = 0.35 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 37	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 37	PURGING INITIATED AT: 1412	PURGING ENDED AT: 1433	TOTAL VOLUME PURGED (gallons): 2.0
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1414	0.2	0.2	.1	5.45	7.03	27.07	1.12	3.57	1.88	Clear	none
1417	0.3	0.5	.1	9.41	6.97	27.14	1.02	1.22	1.22		
1420	0.3	0.8	.1	10.45	6.95	27.12	1.01	0.99	0.41		
1423	0.3	1.1	.1	11.83	6.95	27.20	1.02	0.92	0.73		
1426	0.3	1.4	.1	12.92	6.93	27.12	1.01	1.23	0.35		
1429	0.3	1.7	.1	13.51	6.93	27.13	1.03	1.13	0.25		
1431	0.3	2.0	.1	13.87	6.93	27.05	1.04	1.06	0.71		

WELL CAPACITY (Gallons Per Foot): 0.15" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: IARLADIS	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1433	SAMPLING ENDED AT:
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PUMP OR TUBING DEPTH IN WELL (feet): 37	SAMPLE PUMP FLOW RATE (mL per minute): 500	TUBING MATERIAL CODE:
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-171	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: MW-172 Tallavast ^{MDS 9/14/08} SITE LOCATION: Sarasota FL ^{MDS 4/14/08}
 WELL NO: MW-172 SAMPLE ID: MW-172 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2" TUBING DIAMETER (inches): 1/4" WELL SCREEN INTERVAL DEPTH: 100 feet to 10 feet STATIC DEPTH TO WATER (feet): 1230 PURGE PUMP TYPE OR BAILER: PP
 WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable: $\emptyset = 110$ feet - \emptyset feet X \emptyset gallons/foot = \emptyset gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = \emptyset gallons + $(0.0026 \text{ gallons/foot} \times 110 \text{ feet}) + .25$ gallons = \emptyset gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 105 FINAL PUMP OR TUBING DEPTH IN WELL (feet): 105 PURGING INITIATED AT: 1444 PURGING ENDED AT: 1509 TOTAL VOLUME PURGED (gallons): 110

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1449				13.30	7.07	25.87	0.684	1.08	1230	cloudy	Slight
1451				13.44	7.06	25.79	0.684	0.99	37.4		
1454				13.62	7.06	25.53	0.684	0.90	13.7	Clear	
1457				13.70	7.06	25.61	0.683	0.88	7.26		
1500				13.73	7.06	25.34	0.680	0.88	8.61		
1503				13.75	7.06	25.34	0.680	0.86	2.25		
1506				13.80	7.06	25.67	0.677	0.83	1.85		
1509				13.79	7.06	25.71	0.678	0.84	1.50		
1510	→ Sampled at 1510 ←										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KUCINAS SAMPLER(S) SIGNATURES: [Signature] SAMPLING INITIATED AT: 1510 SAMPLING ENDED AT: 1515
 PUMP OR TUBING DEPTH IN WELL (feet): 105 SAMPLE PUMP FLOW RATE (mL per minute): 400 ^{MDS 9/14/08} TUBING MATERIAL CODE:
 FIELD DECONTAMINATION: (Y) N FIELD-FILTERED: (Y) N FILTER SIZE: — µm DUPLICATE: (Y) (N)

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-172	3	CG	90ml	HCl	—	—	8260	RFPP

REMARKS:
 MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	SITE LOCATION: Sarasota FL
WELL NO: MW-173	SAMPLE ID: MW-173
DATE: 1/23/08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 142 feet to 152 feet	STATIC DEPTH TO WATER (feet): 12.24	PURGE PUMP TYPE OR BAILER:
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY				
= (142 feet - 12.24 feet) X 0.0026 gallons/foot = 2.9 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME				
= 0 gallons + (0.0026 gallons/foot X 152 feet) + 0.25 gallons = 0.4 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 147	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 147	PURGING INITIATED AT: 1520	PURGING ENDED AT: 1549	TOTAL VOLUME PURGED (gallons): 2.9
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1530	1.0	1.0								Gray	none
1534	0.4	1.4	0.1	15.32	7.30	25.94	0.583	1.12	170	Cloudy	
1537	0.3	1.7		15.87	7.28	25.76	0.581	0.99	42.2		
1540	0.3	2.0		16.16	7.27	25.68	0.580	0.97	10.59	Clear	
1543	0.3	2.3		16.32	7.27	25.73	0.578	0.96	5.22		
1546	0.3	2.6		16.46	7.27	25.83	0.578	0.93	3.28		
1549	0.3	2.9			7.26	25.27	0.576	0.97	2.55		
1552	Sampled @ 1552										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KLC / NPS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1552	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 147	SAMPLE PUMP FLOW RATE (mL per minute): 4100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: 1 µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-173	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **initial purge clear/went silty gray very quickly. no noticeable odor**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 9/19/08	SITE LOCATION: <u>Sarasota FL</u> MDS 9/19/08
WELL NO: <u>MW-174</u>	SAMPLE ID: <u>MW-174</u> DATE: <u>1/25/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>225</u> feet to <u>295</u> feet	STATIC DEPTH TO WATER (feet): <u>10.78</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
(only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>285</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>285</u>	PURGING INITIATED AT: <u>0805</u>	PURGING ENDED AT: <u>0840</u>	TOTAL VOLUME PURGED (gallons):
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0810				12.97	7.80	21.78	0.720	1.32	1.25	Clear	Slight
0815				13.99	7.69	21.88	0.721	1.18	37.1	Gray	
0820				15.18	7.68	21.99	0.724	1.16	13.0	↓	
0825				16.01	7.69	21.01	0.722	1.12	3.40	Clear	
0830				16.75	7.69	22.51	0.719	1.09	3.51	↓	
0835					7.69	21.78	0.722	1.11	1.40	↓	
0840	→ Sampled @ 0840 →										

012P

327
333
337
341
343
343

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasey Cornwall ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Kasey Cornwall</u>	SAMPLING INITIATED AT: <u>0840</u>	SAMPLING ENDED AT: <u>0850</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>285</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>5100</u> MDS 9/19/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N MDS 9/19/08	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-174	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: Initial purge clear after 10 minutes light gray slight sulfur odor

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* Needs 2 bolts

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-175	SAMPLE ID: MW-175 DATE: 1-29-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 98.3 feet to 168.3 feet	STATIC DEPTH TO WATER (feet): 191.5	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 108 feet) + .25 gallons = 0.53 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1054	PURGING ENDED AT: 1124	PURGING VOLUME PURGED (gallons): 30
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1112	1.8	1.8	0.1	19.19	7.04	24.49	0.1089	1.38	1.38	none	none noted
1115	0.3	2.1	↓ MDS 4/14/08	19.16	7.06	24.38	0.1089	1.25	0.63		
1118	0.3	2.4		19.14	7.07	24.36	0.1085	1.09	0.38		
1121	0.3	2.7		19.14	7.06	24.35	0.1089	1.05	0.71		
1124	0.3	3.0		19.14	7.07	24.39	0.1085	0.99	0.17		

MDS
4/14/08
MDS
4/14/08
 -250
 -260
 -263
 -264
 -266

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amie Coats / Arcadis	SAMPLER(S) SIGNATURES: <i>Amie Coats</i>	SAMPLING INITIATED AT: 1126	SAMPLING ENDED AT: 1130
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 1100	TUBING MATERIAL CODE: TE	Teflon
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-175	3	CG	40mL	HCl	—	—	8260	RFPP
DUP-20012908	CG	CG	40mL	HCl	—	—	8260	RFPP

REMARKS: *** New dedicated teflon tubing as of today / lost about 15 nuts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; ? = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

down well

DUP-20012908

*** nuts in bottom of well, may be tubing too**

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <u>Former ABC Facility</u>	SITE LOCATION: <u>Tallevast, FL</u>
WELL NO: <u>MW-176</u>	SAMPLE ID: <u>MW-176</u> DATE: <u>1-29-08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>150</u> feet to <u>160</u> feet	STATIC DEPTH TO WATER (feet): <u>2083</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>1100</u> feet) + <u>0.25</u> gallons = <u>0.67</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>900</u>	PURGING ENDED AT: <u>918</u>	TOTAL VOLUME PURGED (gallons): <u>1.8</u>							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>912</u>	<u>1.2</u>	<u>1.2</u>	<u>0.1</u>	<u>24.10</u>	<u>7.29</u>	<u>23.18</u>	<u>0.615</u>	<u>0.60</u>	<u>2.53</u>	<u>none</u>	<u>none noted</u>
<u>915</u>	<u>0.3</u>	<u>1.5</u>	<u>↓</u>	<u>24.38</u>	<u>7.32</u>	<u>23.30</u>	<u>0.614</u>	<u>0.62</u>	<u>3.93</u>	<u>crystal</u>	<u>↓</u>
<u>918</u>	<u>0.3</u>	<u>1.8</u>	<u>↓</u>	<u>24.56</u>	<u>7.33</u>	<u>23.25</u>	<u>0.614</u>	<u>0.62</u>	<u>4.21</u>	<u>clear</u>	<u>↓</u>

ORP
-235
-238
-237

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Ann Coats / Arcadis</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>919</u>	SAMPLING ENDED AT: <u>920</u>
PUMP/OR TUBING DEPTH IN WELL (feet): <u>155</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u> MDS 4/14/08	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N FILTER SIZE: <u>1µm</u>	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-176</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

MDS
4/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Talleast
WELL NO: MW-177	SAMPLE ID: MW-177
DATE: 1-29-05	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 285 feet to 305 feet	STATIC DEPTH TO WATER (feet): 18.19	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)

= **0** gallons + (**.0026** gallons/foot X **310** feet) + **.25** gallons = **1.0** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 305	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 285	PURGING INITIATED AT: 9:05	PURGING ENDED AT: 9:21	TOTAL VOLUME PURGED (gallons): 2.2
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
9:15	1.0	1.0	0.1	19.38	8.24	22.56	1609	1.49	25.8	clear	none
9:18	0.3	1.3	0.1	19.40	8.19	22.77	1600	1.14	25.8	clear	none
9:21	0.3	1.6	0.1	19.41	8.17	22.94	1597	1.09	14.4	clear	none
9:24	0.3	1.9	0.1	19.40	8.15	22.97	1596	1.11	9.49	clear	none
9:27	0.3	2.2	0.1	19.40	8.13	23.08	1597	1.08	7.62	clear	none

MDS 4/14/05

261
265
270
272
276

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 9:28	SAMPLING ENDED AT: 9:33
PUMP OR TUBING DEPTH IN WELL (feet): 285	SAMPLE PUMP FLOW RATE (mL per minute): <100 (MDS 4/14/05)	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N Filtration Equipment Type: _____	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-177	3	CG	40ml	HCl	—	—	8260	RFPP

MDS 4/14/05

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/14/08	SITE LOCATION: Sarasota FL MDS 4/14/08
WELL NO: MW-178	SAMPLE ID: MW-178 DATE: 1/25/08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 3.5 feet to 36.5 feet	STATIC DEPTH TO WATER (feet): 13.20	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X 3 gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (3.002 gallons/foot X 36.5 feet) + 3.25 gallons = 0.39 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33	PURGING INITIATED AT: 1055	PURGING ENDED AT: 1115	PURGE VOLUME PURGED (gallons):

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (umhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1100				14.22	6.97	24.85	1.00	1.38	4.01	Clear	None
1105				14.72	6.96	24.84	0.99	1.13	2.93	Clear	
1110				14.95	6.96	24.99	0.99	1.08	3.07	↓	↓
1115				15.15	6.95	24.80	0.99	1.04	4.34	↓	↓

ORP
-114
-187
-237
-263

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Kasey Cornwell / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1115	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 33	SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-178	3	CG	40ml	HCL	—	—	8260	RFPP

MDS 4/14/08

REMARKS: **Initial Purge is clear w/ no noticeable odor.**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u>		SITE LOCATION: <u>Sarasota FL</u>	
WELL NO: <u>MW-179</u>	SAMPLE ID: <u>MW-179</u>	DATE: <u>1/25/09</u>	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>103</u> feet to <u>93</u> feet	STATIC DEPTH TO WATER (feet): <u>14.48</u>	PURGE PUMP TYPE OR BAILER: <u>Peristaltic</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY = (<u>103</u> feet - <u>14.48</u> feet) X <u>0.163</u> gallons/foot = <u>14.9</u> gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME = _____ gallons + (_____ gallons/foot X _____ feet) + _____ gallons = _____ gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: <u>11:00</u>	PURGING ENDED AT: <u>11:25</u>	TOTAL VOLUME PURGED (gallons): <u>2.5</u>					
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
11:05	0.5	0.5	0.1	14.65	7.0	23.5	0.635	3.0	1.34	clear	odor
11:10	0.5	1.0	0.1	14.75	7.3	24.4	0.644	2.3	1.45	"	"
11:15	0.5	1.5	0.1	14.75	7.3	24.6	0.655	2.3	1.03	"	"
11:20	0.5	2.0	0.1	14.75	7.3	24.7	0.653	2.4	0.99	"	"
11:25	0.5	2.5	0.1	14.75	7.3	24.8	0.651	2.3	0.98	"	"

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith</u>	SAMPLER(S) SIGNATURES: <u>Nathan Smith</u>	SAMPLING INITIATED AT: <u>11:25</u>	SAMPLING ENDED AT: <u>11:35</u>
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-179	3	CG	40ml	HCl	—	—	8260	RFPP
Dup-13	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: Samples taken @ 11:25; Dup-13 collected at this location

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleuas</u>	SITE LOCATION: <u>Sarasota FL</u>
WELL NO: <u>MW-180</u>	SAMPLE ID: <u>MW-180</u>
	DATE: <u>1/25/68</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/8</u>	WELL SCREEN INTERVAL DEPTH: <u>145</u> feet to <u>150.3</u> feet	STATIC DEPTH TO WATER (feet): <u>157.7</u>	PURGE PUMP TYPE OR BAILER: <u>AP</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>1345</u>	PURGING ENDED AT: <u>1410</u>	TOTAL VOLUME PURGED (gallons):
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>1350</u>				<u>17.48</u>	<u>7.27</u>	<u>25.44</u>	<u>0.712</u>	<u>1.26</u>	<u>0.64</u>	<u>Clear</u>	<u>none</u>
<u>1355</u>				<u>17.89</u>	<u>7.23</u>	<u>25.41</u>	<u>0.708</u>	<u>1.11</u>	<u>0.82</u>	<u>↓</u>	<u>↓</u>
<u>1400</u>				<u>18.48</u>	<u>7.22</u>	<u>25.43</u>	<u>0.707</u>	<u>1.02</u>	<u>0.88</u>	<u>↓</u>	<u>↓</u>
<u>1405</u>				<u>18.88</u>	<u>7.21</u>	<u>25.49</u>	<u>0.706</u>	<u>0.99</u>	<u>1.24</u>	<u>↓</u>	<u>↓</u>
<u>1410</u>				<u>→ Sample @ 1410 ←</u>							

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>MDS 9/14/08</u>	SAMPLING INITIATED AT: <u>1410</u>	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u>5100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: <u>µm</u>	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-180</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>✓</u>	<u>—</u>	<u>8260</u>	<u>RFP</u>

REMARKS: → initial purge clear, no noticeable odor.

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 9/14/08	SITE LOCATION: Sarasota MDS 9/14/08
WELL NO: MW-181	SAMPLE ID: MW-181 DATE: 1/25/09

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/8	WELL SCREEN INTERVAL DEPTH: 27.5 feet to 29.5 feet	STATIC DEPTH TO WATER (feet): 1340	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 1440		PURGING ENDED AT: 1505		TOTAL VOLUME PURGED (gallons):			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1445				15.26	8.71	24.87	0.578	1.80	0.73	clear	none
1450				18.21	8.70	25.04	0.557	1.09	0.63		
1455				18.93	8.57	25.07	0.566	1.03	0.91		
1500				19.59	8.45	24.97	0.591	1.00	0.91		
1505				19.99	8.42	24.88	0.612	1.01	0.99		

-351
-382
-383
-384
-384

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KLC / ARCADIS		SAMPLER(S) SIGNATURES: KLC MDS 9/14/08		SAMPLING INITIATED AT: 1510		SAMPLING ENDED AT:	
PUMP OR TUBING DEPTH IN WELL (feet):		SAMPLE PUMP FLOW RATE (mL per minute): 2100 MDS 9/14/08		TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 1 µm		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-181	3	CG	40ml	HCl	—	—	8260	LFPP

MDS
9/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: STA	SITE LOCATION: Sarasota FL
WELL NO: MW-182	SAMPLE ID: MW-182 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 164 feet to 174 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: app geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 178 feet) + 0.50 gallons = ⁽¹⁰⁰⁾ 0.96 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~169	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~169	PURGING INITIATED AT: 1450	PURGING ENDED AT: 1506	TOTAL VOLUME PURGED (gallons): 1.44
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TIME	VOLUME PURGED (gallons)	CUMUL VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	<i>ORP</i>
1502	1.08	1.08	0.09	18.18	7.40	28.03	634	0.31	13.5	clear	yes	-254
1503	0.18	1.26	0.09	18.18	7.39	27.97	635	0.30	13.0	clear	yes	-259
1506	0.18	1.44	0.09	18.18	7.38	27.97	641	0.29	8.69	clear	yes	
<i>NS 1/23/08</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: N. Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>N. Smith</i>	SAMPLING INITIATED AT: 1507	SAMPLING ENDED AT: 1515
PUMP OR TUBING DEPTH IN WELL (feet): ~169	SAMPLE PUMP FLOW RATE (mL per minute): <100mL	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-182	3	CG	40mL	HCL	-	<2	8260B + Sim	RFPP
<i>NS 1/23/08</i>								

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEAST	SITE LOCATION: SARASOTA, FL
WELL NO: NW-183	SAMPLE ID: NW-183 DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 30 feet) + 0.5 gallons = 0.56 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5		FINAL PUMP OR TUBING DEPTH IN WELL (feet): 17.5		PURGING INITIATED AT: 1444		PURGING ENDED AT: 1511		TOTAL VOLUME PURGED (gallons): 1.50			MPS 4/14/08	
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORD
1456	0.60	0.60	0.05	3.36	7.17	24.83	417	3.95	188	brown/clear	no	18
1459	0.15	0.90	0.05	3.40	7.10	25.16	421	2.79	182	brown/clear	no	-21
1502	0.15	1.05	0.05	3.43	7.02	25.51	425	2.39	168	brown/clear	no	-46
1505	0.15	1.20	0.05	3.50	6.93	25.44	429	1.81	130	brown/clear	no	-75
1508	0.15	1.35	0.05	3.67	6.92	25.46	430	1.76	133	brown/clear	no	-78
1511	0.15	1.50	0.05	3.48	6.91	25.38	432	1.70	127	brown/clear	no	-80

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/NS / ARCADIS		SAMPLER(S) SIGNATURES: <i>Stu Well</i>		SAMPLING INITIATED AT: 1513	SAMPLING ENDED AT: 1519
PUMP OR TUBING DEPTH IN WELL (feet): 17.5		SAMPLE PUMP FLOW RATE (mL per minute): 5100		TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> FILTER SIZE: _____ µm		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
NW-183	3	CG	40mL	HCl	40mL	—	8260+8260SIM	RFP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallahassee</u>	SITE LOCATION: <u>Jacksonville, FL</u>
WELL NO: <u>MW-184</u>	SAMPLE ID: <u>MW-184</u> DATE: <u>1/24/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>28</u> feet to <u>33</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>PP geopump</u>
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

= gallons + (0.0026 gallons/foot X 38 feet) + 0.50 gallons = 0.38 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~31</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~31</u>	PURGING INITIATED AT: <u>1445</u>	PURGING ENDED AT: <u>1502</u>	TOTAL VOLUME PURGED (gallons): <u>1.80</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
<u>1450</u>	<u>0.60</u>	<u>0.60</u>	<u>0.12</u>	<u>10.92</u>	<u>7.69</u>	<u>25.71</u>	<u>254</u>	<u>0.81</u>	<u>39.4</u>	<u>clear</u>	<u>yes</u>	<u>-133</u>
<u>1452</u>	<u>0.24</u>	<u>0.84</u>	<u>0.12</u>	<u>10.94</u>	<u>7.71</u>	<u>25.73</u>	<u>258</u>	<u>0.64</u>	<u>32.9</u>	<u>clear</u>	<u>yes</u>	<u>-151</u>
<u>1454</u>	<u>0.24</u>	<u>0.96</u>	<u>0.12</u>	<u>11.01</u>	<u>7.73</u>	<u>25.74</u>	<u>317</u>	<u>0.47</u>	<u>23.0</u>	<u>clear</u>	<u>yes</u>	<u>-168</u>
<u>1456</u>	<u>0.24</u>	<u>1.08</u>	<u>2.12</u>	<u>11.01</u>	<u>7.56</u>	<u>25.65</u>	<u>384</u>	<u>0.38</u>	<u>13.6</u>	<u>clear</u>	<u>yes</u>	<u>-172</u>
<u>1458</u>	<u>0.24</u>	<u>1.32</u>	<u>0.12</u>	<u>11.02</u>	<u>7.60</u>	<u>25.63</u>	<u>420</u>	<u>0.32</u>	<u>12.4</u>	<u>clear</u>	<u>yes</u>	<u>-174</u>
<u>1500</u>	<u>0.24</u>	<u>1.56</u>	<u>0.12</u>	<u>11.02</u>	<u>7.57</u>	<u>25.69</u>	<u>430</u>	<u>0.37</u>	<u>13.3</u>	<u>clear</u>	<u>yes</u>	<u>-174</u>
<u>1502</u>	<u>0.24</u>	<u>1.80</u>	<u>0.12</u>	<u>11.03</u>	<u>7.51</u>	<u>25.75</u>	<u>460</u>	<u>0.31</u>	<u>10.52</u>	<u>clear</u>	<u>yes</u>	<u>-177</u>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Neil Smith</u>	SAMPLING INITIATED AT: <u>1503</u>	SAMPLING ENDED AT: <u>1510</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~31</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>≤ 100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> X FILTER SIZE: <u> </u> µm	DUPLICATE: <u>Y</u> X	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-184</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u>-</u>	<u>< 2</u>	<u>P260B + P260JL</u>	<u>RFPF</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 4/14/08	SITE LOCATION: <u>Sarasota FL</u> MDS 4/14/08
WELL NO: <u>MW-158 185</u>	SAMPLE ID: <u>MW-158 185</u> DATE: <u>4/22/08 1/24/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>15</u> feet to <u>15</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>100</u> feet) + <u>0.5</u> gallons = <u>0.76</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>90</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>90</u>	PURGING INITIATED AT: <u>1580</u>	PURGING ENDED AT: <u>1553</u>	TOTAL VOLUME PURGED (gallons): <u>1.61</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle) (mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	<u>ORP</u>
<u>1501</u>	<u>0.77</u>	<u>0.77</u>	<u>0.07</u>	<u>13.25</u>	<u>7.56</u>	<u>24.96</u>	<u>694</u>	<u>2.15</u>	<u>1.04</u>	<u>clear</u>	<u>yes</u>	<u>-171</u>
<u>15044</u>	<u>0.21</u>	<u>0.98</u>	<u>0.07</u>	<u>13.28</u>	<u>7.55</u>	<u>24.99</u>	<u>696</u>	<u>2.30</u>	<u>0.86</u>	<u>clear</u>	<u>yes</u>	<u>-190</u>
<u>150047</u>	<u>0.21</u>	<u>1.19</u>	<u>0.07</u>	<u>13.28</u>	<u>7.53</u>	<u>25.11</u>	<u>700</u>	<u>0.85</u>	<u>0.94</u>	<u>clear</u>	<u>yes</u>	<u>-193</u>
<u>1550</u> <u>1600</u>	<u>0.21</u>	<u>1.40</u>	<u>0.07</u>	<u>13.28</u>	<u>7.52</u>	<u>25.12</u>	<u>701</u>	<u>0.75</u>	<u>0.71</u>	<u>clear</u>	<u>yes</u>	<u>-196</u>
<u>1553</u>	<u>0.21</u>	<u>1.61</u>	<u>0.07</u>	<u>13.29</u>	<u>7.51</u>	<u>25.00</u>	<u>706</u>	<u>0.64</u>	<u>0.46</u>	<u>clear</u>	<u>yes</u>	<u>-195</u>

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Fl.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>SW/NS/ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>1558</u>	SAMPLING ENDED AT: <u>1600</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>90</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>≤ 100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>⊖</u> N	FIELD-FILTERED: Y <u>⊖</u> FILTER SIZE: _____ µm	DUPLICATE: Y <u>⊖</u> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-185</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u>40mL</u>	<u>—</u>	<u>8260+8260 SIM</u>	<u>RFP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST		SITE LOCATION: SARASOTA, FL	
WELL NO: MW-186	SAMPLE ID: MW-157	DATE: 1/22/2008	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 0.2 feet to 2.3 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: HP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (38 feet - 1.60 feet) X _____ gallons/foot = _____ gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 165 feet) + 0.5 gallons = 0.93 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 155	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 155	PURGING INITIATED AT: 1608	PURGING ENDED AT: 1633	TOTAL VOLUME PURGED (gallons): 1.82
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1620	0.98	0.98	0.07	15.33	7.54	25.13	668	0.30	10.72	clear	no	-217
1625	0.28	1.26	0.07	15.51	7.44	25.09	678	0.29	4.36	clear	no	-218
1630	0.28	1.54	0.07	15.62	7.36	25.05	688	0.52	11.7	clear	no	-202
1633	0.28	1.82	0.07	15.68	7.28	25.02	696	0.30	3.80	clear	no	-216

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SN/NS/ARLADIS	SAMPLER(S) SIGNATURES: <i>Stu WM</i>	SAMPLING INITIATED AT: 1635	SAMPLING ENDED AT: 1640
PUMP OR TUBING DEPTH IN WELL (feet): 155	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-186	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellicoast</u>	SITE LOCATION: <u>Parasota, FL</u>
WELL NO: <u>MW-187</u>	SAMPLE ID: <u>MW-187</u>
DATE: <u>1/22/2008</u>	

PURGING DATA

WELL DIAMETER (inches): <u>4</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL OPEN INTERVAL DEPTH: <u>15</u> feet to <u>20</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)
= (1750 feet - _____ feet) X _____ gallons/foot = _____ gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME 1.30
(only fill out if applicable)
= _____ gallons + 0.026 gallons/foot X 295 feet + 0.50 gallons = 1.267 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 175 FINAL PUMP OR TUBING DEPTH IN WELL (feet): 290 PURGING INITIATED AT: 1530 PURGING ENDED AT: 1559 TOTAL VOLUME PURGED (gallons): 2.32

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle % or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
17 1547	1.36	1.36	0.08	18.74	11.85	25.09	3130	0.31	30.2	clear	yes
1550	0.24	1.60	0.08	19.13	11.87	25.02	3070	0.31	3.62	clear	yes
1553	0.24	1.84	0.08	20.33	11.88	24.76	3050	0.33	2.24	clear	yes
1556	0.24	2.08	0.08	20.97	11.88	24.56	3030	0.34	0.63	clear	yes
1558	0.24	2.32	0.08	21.41	11.88	24.50	3020	0.36	0.59	clear	yes
<i>Not Sampled 1/24/08</i>											

ORP
-289
-289
-289
-289
-281

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>SW/NS/ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Naloff</u>	SAMPLING INITIATED AT: <u>1600</u>	SAMPLING ENDED AT: <u>1610</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>290</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>≤ 100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-187	3	CG	40mL	ACL	-	7.2	8260BFSIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 40% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast FL
WELL NO: MW-188	SAMPLE ID: MW-188 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 12 feet to 17 feet	STATIC DEPTH TO WATER (feet): 3.7	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + .25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 10:20		PURGING ENDED AT: 10:55		TOTAL VOLUME PURGED (gallons): 3.5			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1035	2.0	2.0	0.25	4.41	6.29	24.69	0.705	0.41	10	clear	no odor
1040	2.5	2.5	0.50	4.42	6.33	24.66	0.706	0.38	17	"	"
1045	2.8	2.8	0.1	4.41	6.37	24.67	0.709	0.37	7	"	"
1050	3.0	3.0	0.1	4.41	6.37	24.66	0.706	0.37	3	"	"
1055	3.5	3.5	0.1	4.41	6.39	24.66	0.705	0.36	4	"	"

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micha Forbes	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 10:56	SAMPLING ENDED AT: 11:00
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100 MOS	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	
SAMPLE CONTAINER SPECIFICATION		SAMPLE PRESERVATION	

SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
MW-188	3	CG	40ml	Hcl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump

EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast, FL
WELL NO: MW-189	SAMPLE ID: MW-189 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 28 feet to 33 feet	STATIC DEPTH TO WATER (feet): 8.05	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1018	PURGING ENDED AT: 1049	TOTAL VOLUME PURGED (gallons): 3.1
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1040	2.2	2.2	0.1	9.49	7.15	24.36	0.712	0.25	0.0	none	none noted	-190
1043	0.3	2.5	↓	9.53	7.10	24.34	0.709	0.22	0.0	↓	↓	-196
1046	0.3	2.8	↓	9.56	7.14	24.34	0.709	0.20	0.0	↓	↓	-202
1049	0.3	3.1	↓	9.57	7.12	24.36	0.708	0.20	0.0	↓	↓	-205
					7.139							
							<i>mS/cm</i>					

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amy Coats / Arcadis	SAMPLER(S) SIGNATURES: <i>Amy Coats</i>	SAMPLING INITIATED AT: 1050	SAMPLING ENDED AT: 1055
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-189	3	CG	40mL	HCl	—	—	8260	RFPP
DUP-3012308	3	CG	40mL	HCl	—	—	8260	RFPP

REMARKS: **DUP - 3012308**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-190	SAMPLE ID: MW-190 DATE: 1.25.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 90 feet to 100 feet	STATIC DEPTH TO WATER (feet): 12.09	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 8:25		PURGING ENDED AT: 8:50		TOTAL VOLUME PURGED (gallons): 2.5			
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8:35	0.5	1.0	0.1	13.05	7.22	23.32	0.475	0.58	0	clear	none
8:40	0.5	1.5	0.1	13.05	7.24	23.15	0.47	0.47	0	"	"
8:45	0.5	2.0	0.1	13.05	7.24	23.12	0.471	0.46	0	"	"
8:50	0.5	2.5	0.1	13.05	7.33	23.49	0.469	0.44	0	"	"

ORP
-263
-272
-272
-275

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 8:51	SAMPLING ENDED AT: 8:55
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 2100 MAS 4/14/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-190	3	CG	40ml	HCl	—	—	8260	RF PP

MAS
4/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallerost, FL
WELL NO: MW-191	SAMPLE ID: MW-191
DATE: 1.23.08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 146 feet to 156 feet	STATIC DEPTH TO WATER (feet): 12.06	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1118	PURGING ENDED AT: 1145	TOTAL VOLUME PURGED (gallons): 2.5
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1125	1.0	1.0	0.1	15.91	7.56	25.62	0.441	0.42	29	clear	none	-281
1130	1.5	1.5	0.1	16.05	7.61	25.25	0.439	0.38	64	clear	none	-297
1135	2.0	2.0	0.1	16.05	7.61	25.62	0.435	0.35	60	clear	none	-302
1140	2.3	2.3	0.1	16.18	7.61	24.46	0.435	0.35	40	"	"	-304
1145	2.5	2.5	0.1	16.18	7.62	25.33	0.435	0.34	52	"	"	-307
							ms/cm					

MDS
4/14/08
ORP
-281
-297
-302
-304
-307

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Nicah Forbes / Arcadis	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1146	SAMPLING ENDED AT: 11:50
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-191	3	CG	40ml	HCl	—	—	8260	RFPP

MDS
4/14/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-192		SAMPLE ID: MW-192	DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet): 8.68	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet): 1138		PURGING INITIATED AT: 1120 # AC	PURGING ENDED AT: 1159	TOTAL VOLUME PURGED (gallons): 2.1					
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1147	0.9	0.9	0.1	9.55	11.49	24.89	1.82	0.42	7.14		
1150	0.3	1.2	↓	9.54	11.49	24.97	1.80	0.39	7.80		
1153	0.3	1.5	↓	12.99	11.50	24.69	1.81	0.36	7.27		
1150	0.3	1.8	↓	14.17	11.50	24.69	1.82	0.35	7.31		
1159	0.3	2.1	↓	16.84	11.51	24.75	1.81	0.34	7.80		
							µS/cm				

ORD
-376
-379
-379
-379
-379

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Am Coats / Arcadis		SAMPLER(S) SIGNATURES: <i>[Signature]</i>		SAMPLING INITIATED AT: 1200	SAMPLING ENDED AT: 1205
PUMP OR TUBING DEPTH IN WELL (feet):		SAMPLE PUMP FLOW RATE (mL per minute): <100		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y N		FIELD-FILTERED: Y N FILTER SIZE: µm		DUPLICATE: Y N	
Filtration Equipment Type:					

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-192	3	CG	40ml	HCl	-	-	2260	RFPP

MDS 2-19-08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: <u>Telleport</u>	SITE LOCATION: <u>Subsata, FL</u>
WELL NO: <u>MW-193</u>	DATE: <u>1/24/08</u>
SAMPLE ID: <u>MW-193</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>105</u> feet to <u>165</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>(PP) geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (<u>0.0026</u> gallons/foot X <u>98</u> feet) + <u>0.5</u> gallons = <u>0.75</u> gallons				
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INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~95</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~95</u>	PURGING INITIATED AT: <u>1223</u>	PURGING ENDED AT: <u>1242</u>	TOTAL VOLUME PURGED (gallons): <u>1.52</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1233	0.80	0.80	0.08	12.24	7.56	23.71	614	0.43	0.61	clear	no
1236	0.24	1.04	0.08	12.27	7.56	23.73	613	0.53	0.57	clear	no
1239	0.24	1.28	0.08	12.27	7.57	23.75	612	0.52	0.25	clear	no
1242	0.24	1.52	0.08	12.27	7.57	23.88	611	0.39	0.65	clear	no

4/14/08
MPS

CRD
-226
-232
-238
-224

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>N. Smith ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>1244</u>	SAMPLING ENDED AT: <u>1248</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~95</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>600</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>X</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <input checked="" type="checkbox"/>	
FILTRATION EQUIPMENT TYPE: _____			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-193</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCL</u>	<u>-</u>	<u>~2</u>	<u>826-B + 826-SM</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: <u>Tellegast</u>	SITE LOCATION: <u>Sarasota, FL</u>
WELL NO: <u>MW-194</u>	SAMPLE ID: <u>MW-194</u> DATE: <u>1/24/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>4/4</u>	WELL SCREEN INTERVAL DEPTH: <u>145</u> feet to <u>155</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>pp (geopump)</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<u>0.0026</u> gallons/foot X <u>158</u> feet) + <u>0.58</u> gallons = <u>0.91</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~150</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~150</u>	PURGING INITIATED AT: <u>1223</u>	PURGING ENDED AT: <u>1237</u>	TOTAL VOLUME PURGED (gallons): <u>1.82</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle <u>pp</u> or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	<u>ORP</u>
1231	1.04	1.04	0.13	15.15	7.62	24.02	650	0.33	11.3	clear	yes	-234
1234	0.39	1.43	0.13	15.20	7.61	24.02	651	0.33	15.0	clear	yes	-232
1237	0.39	1.82	0.13	16.31	7.61	24.07	648	0.30	13.0	clear	yes	-245
<i>Not used</i>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>S. Wells / AREAS</u>	SAMPLER(S) SIGNATURES: <u>Neil Reed</u>	SAMPLING INITIATED AT: <u>1238</u>	SAMPLING ENDED AT: <u>1245</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~150</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>≤ 100 mL</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> <u>N</u>	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-194</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u>-</u>	<u>~2</u>	<u>8260B + 8260S12</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-195	SAMPLE ID: MW-195 DATE: 1/24/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 4	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026** gallons/foot X **305** feet) + **0.5** gallons = **0.793** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 290	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 290	PURGING INITIATED AT: 1128	PURGING ENDED AT: 1208	TOTAL VOLUME PURGED (gallons): 2.40
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circled or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1150	1.32	1.32	0.06	17.11	8.31	23.85	928	5.37	6.65	clear	yes	-258
1156	0.36	1.68	0.06	18.41	8.31	23.87	919	5.47	6.28	clear	yes	-245
1202	0.36	2.04	0.06	19.08	8.32	23.87	927	5.36	7.67	clear	yes	-245
1208	0.36	2.40	0.06	19.88	8.31	23.90	928	5.70	10.87	clear	yes	-244

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT)/ AFFILIATION: NS/SW	SAMPLER(S) SIGNATURES: <i>St...</i>	SAMPLING INITIATED AT: 1210	SAMPLING ENDED AT: 1214
PUMP OR TUBING DEPTH IN WELL (feet): 290	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-195	3	CG	40mL	HCl	40mL	—	8260+8260 514	RFPP
DUP-9	3	CG	40mL	HCl	40mL	—	8260+8260 514	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> ^{MDS} <u>4/19/08</u>	SITE LOCATION: <u>Sarasota</u> ^{MDS} <u>4/19/08</u>
WELL NO: <u>MW-196</u>	SAMPLE ID: <u>MW-196</u> DATE: <u>4/30/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>90</u> feet to <u>100</u> feet	STATIC DEPTH TO WATER (feet): <u>16.8</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (<u> </u> feet - <u> </u> feet) X <u> </u> gallons/foot = <u> </u> gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>Ø</u> gallons + <u>0.0026</u> gallons/foot X <u>100</u> feet + <u>.25</u> gallons = <u>0.5</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>95</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>95</u>	PURGING INITIATED AT: <u>1433</u>	PURGING ENDED AT: <u>1452</u>	TOTAL VOLUME PURGED (gallons): <u>1.2</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>1442</u>	<u>0.12</u>	<u>.22</u>	<u>.08</u>	<u>16.76</u>	<u>6.7</u>	<u>25.1</u>	<u>.890</u>	<u>4.4</u>	<u>0.46</u>	<u>Clear</u>	<u>None</u>
<u>1445</u>	<u>.24</u>	<u>.96</u>	<u>↓</u>	<u>16.80</u>	<u>6.7</u>	<u>25.0</u>	<u>.890</u>	<u>3.4</u>	<u>0.93</u>	<u>↓</u>	<u>↓</u>
<u>1449</u>	<u>.24</u>	<u>1.2</u>	<u>↓</u>	<u>16.80</u>	<u>6.7</u>	<u>25.1</u>	<u>.89</u>	<u>3.2</u>	<u>0.28</u>	<u>↓</u>	<u>↓</u>
<u>1452</u>	<u>.24</u>	<u>1.44</u>	<u>.08</u>	<u>16.80</u>	<u>6.7</u>	<u>25.1</u>	<u>.890</u>	<u>3.3</u>	<u>0.32</u>	<u>↓</u>	<u>↓</u>
<u>1455</u>	<u>→ Sampled @ 1455 ←</u>										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

initial pump silt run clear at **SAMPLING DATA** *time of 1st reading*

SAMPLED BY (PRINT) / AFFILIATION: <u>Natesmith / Arcadis</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>1455</u>	SAMPLING ENDED AT: <u>1500</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>95</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>400</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> N	FILTRATION EQUIPMENT TYPE: <u>[Signature]</u>	FILTER SIZE: <u> </u> µm

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-196</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

200ml/min = 0.5 gal

GROUNDWATER SAMPLING LOG 1.2 Gallons remaining

SITE NAME: Tallewas t MDS 4/15/08 SITE LOCATION: Sarasota FL MDS 4/15/08
 WELL NO: MW-197 SAMPLE ID: MW-197 DATE: 1/30/08

MDS 4/15/08 **PURGING DATA**
 WELL DIAMETER (inches): _____ TUBING DIAMETER (inches): 1/4 WELL SCREEN INTERVAL DEPTH: 106 feet to 116 feet STATIC DEPTH TO WATER (feet): 21.12 PURGE PUMP TYPE OR BAILER: PP
 WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 = (_____ feet - _____ feet) X _____ gallons/foot = _____ gallons
 EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = 0 gallons + (0.0026 gallons/foot X 116 feet) + .25 gallons = .55 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 111 FINAL PUMP OR TUBING DEPTH IN WELL (feet): 111 PURGING INITIATED AT: 1327 PURGING ENDED AT: 1413 TOTAL VOLUME PURGED (gallons): 3.25

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1332	.5	.5	0.1	24.85	7.36	26.12	0.636	1.27	0.41	clear	none
1335	.3	.8	0.1	25.61	7.38	26.09	0.631	1.17	2.01		
1338	.3	1.10	0.1	26.00	7.40	26.15	0.631	1.08	1.65		
1359				29.16	7.33		0.702	1.1			
1404	1.7	2.8	200ml/min	28.4	7.40	26.40	0.630	3.00	2.01	clear	none
1407	.15	2.95	0.05	29.55	7.35	26.25	0.661	1.24	4.59		
1410	.15	3.10	0.05	29.69	7.35	26.29	0.651	1.14	4.96		
1413	.15	3.25	0.05	29.81	7.35	26.32	0.640	1.07	3.19		
1415	Sampled @ 1415										

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

* Stopped taking readings, waited for drawdown to stabilize

SAMPLED BY (PRINT) AFFILIATION: Nate Smith / ARCADIS SAMPLER(S) SIGNATURES: _____
 PUMP OR TUBING DEPTH IN WELL (feet): 111 SAMPLE PUMP FLOW RATE (mL per minute): <100 MDS 4/15/08 TUBING MATERIAL CODE: _____
 FIELD DECONTAMINATION: (P) N FIELD-FILTERED: Y (N) FILTER SIZE: _____ µm DUPLICATE: Y (N)
 Filtration Equipment Type: _____

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-197	3	CG	40ml	HCL	-	-	8260	RFPP

REMARKS:
 MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

1345: - timed off pump, pulled up tubing to check length
 350 resumed pumping

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-198	SAMPLE ID: MW-198	DATE: 1-28-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 10 feet to 16 feet	STATIC DEPTH TO WATER (feet): 3.0	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (16 feet - 3.0 feet) X 0.0026 gallons/foot = 0.29 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + 0.0026 gallons/foot X 16 feet + 0.25 gallons = 0.29 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 16.14	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 14	PURGING INITIATED AT: 14:15	PURGING ENDED AT: 14:50	TOTAL VOLUME PURGED (gallons): 2.25

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
14:20	0.5	0.5	0.1	4.95	6.30	23.4	0.828	0.73	280	clear	none	79
14:25	0.5	1.0	0.1	6.60	6.19	24.1	0.854	0.56	240	sl. turbid	none	37
14:30	0.25	1.25	0.05	6.17	6.17	24.8	1.00	0.47	140	"	"	15
14:35	0.25	1.50	0.05	6.1	6.2	25.1	1.16	0.41	89	clear	none	41
14:40	0.25	1.75	0.05	5.8	6.24	25.1	1.00	0.38	60	"	"	53
14:45	0.25	2.00	0.05	5.75	6.29	25.1	0.95	0.39	45	"	"	58
14:50	0.25	2.25	0.05	5.75	6.31	25.1	0.94	0.39	53	"	"	63

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Michal Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 14:51	SAMPLING ENDED AT: 14:55
PUMP OR TUBING DEPTH IN WELL (feet): 14 ^{MDS 9/15/08}	SAMPLE PUMP FLOW RATE (mL per minute): 5100 ^{MDS 9/15/08}	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N Filtration Equipment Type:	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-198	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-199	SAMPLE ID: MW-199 DATE: 1.28.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 20 feet to 35 feet	STATIC DEPTH TO WATER (feet): 8.85	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X 35 feet + 0.25 gallons = 0.341 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33	PURGING INITIATED AT: 3:22	PURGING ENDED AT: 15:45	TOTAL VOLUME PURGED (gallons): 1.34

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
15:25	0.34	0.34	0.1	9.7	11.3	25.6	1.43	0.60	17	clear	none
15:30	0.25	0.59	0.05	9.8	11.29	25.39	1.44	0.41	4	"	"
15:35	0.25	0.84	"	9.7	11.24	25.32	1.43	0.35	1	"	"
15:40	0.25	1.09	"	9.7	11.13	25.32	1.31	0.32	2	"	"
15:45	0.25	1.34	"	9.7	11.10	25.30	1.27	0.32	1	"	"

CRP
-272
-283
-288
-290
-290

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 15:46	SAMPLING ENDED AT: 15:50
PUMP OR TUBING DEPTH IN WELL (feet): 33	SAMPLE PUMP FLOW RATE (mL per minute): 100 mL/min	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-199	3	CG	40ml	HCl	—	—	8260	RFPP

MCS 4/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

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DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Tellevast	SITE LOCATION: Green Street
WELL NO: MW-200	SAMPLE ID: MW-200
DATE: 1/28/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 90 feet to 100 feet	STATIC DEPTH TO WATER (feet): 9.50	PURGE PUMP TYPE OR BAILER: (pp) geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
= gallons + (0.022 gallons/foot X 105 feet) + 0.25 gallons = 0.52 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~95	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~95	PURGING INITIATED AT: 1420	PURGING ENDED AT: 1431	TOTAL VOLUME PURGED (gallons): 1.21
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1425	0.55	0.55	0.11	9.84	11.24	24.91	1230	0.78	8.15	clear	yes
1428	0.33	0.88	0.11	9.84	11.24	25.08	1230	0.74	7.24	clear	yes
1431	0.33	1.21	0.11	9.84	11.25	25.03	1230	0.72	4.49	clear	yes
<i>Hold Stand 1/28/08</i>											

ORP
-259
-261
-261

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1432	SAMPLING ENDED AT: 1440
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PUMP OR TUBING DEPTH IN WELL (feet): ~95	SAMPLE PUMP FLOW RATE (mL per minute): 500	TUBING MATERIAL CODE: PE
FIELD DECONTAMINATION: 0 N	FIELD-FILTERED: Y <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="checkbox"/>

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-200	3	CG	40ml	HCL	-	~2	8260B + 8260JW RFP	

REMARKS: **Need new bolts for manhole covers**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3):
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Telleast	SITE LOCATION: Jacksonville, FL
WELL NO: MW-201	SAMPLE ID: MW-201
DATE: 1/28/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 150 feet to 160 feet	STATIC DEPTH TO WATER (feet): 142.7	PURGE PUMP TYPE OR BAILER: (PP) gear pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 165 feet) + 0.25 gallons = 0.674 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~155	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~155	PURGING INITIATED AT: 1515	PURGING ENDED AT: 1542	TOTAL VOLUME PURGED (gallons): 2.80

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1521	0.70	0.70	0.10	12.84	7.85	24.99	1060	1.93	127	clear	yes	-187
1524	0.30	1.00	0.10	13.34	7.68	25.06	1060	1.78	62.4	clear	yes	-189
1527	0.30	1.30	0.10	13.62	7.64	25.06	1060	1.66	50.9	clear	yes	-188
1530	0.30	1.60	0.10	14.56	7.55	24.84	1050	1.50	43.0	clear	yes	-187
1533	0.30	1.90	0.10	15.15	7.53	24.99	1050	1.39	87.8	clear	yes	-189
1539	0.60	2.50	0.10	15.71	7.53	25.06	1050	1.23	85.6	clear	yes	-191
1542	0.30	2.80	0.10	16.15	7.53	25.12	1050	1.13	86.1	clear	yes	-192

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1543	SAMPLING ENDED AT: 1550
PUMP OR TUBING DEPTH IN WELL (feet): ~155	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> MS	FIELD-FILTERED: <input checked="" type="checkbox"/> X	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> X

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-201	3	CG	4am	HCL	-	~2	82608 + 82608 SM RFPP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellus st</u>	SITE LOCATION: <u>Saratoga Pl</u>
WELL NO: <u>MW-202</u>	SAMPLE ID: <u>MW-202</u> DATE: <u>1/2 2/1/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>280</u> feet to <u>300</u> feet	STATIC DEPTH TO WATER (feet): <u>8.87</u>	PURGE PUMP TYPE OR BAILER: <u>pp geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0024 gallons/foot X <u>290</u> feet) + 0.25 gallons = <u>1.00</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~280</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~280</u>	PURGING INITIATED AT: <u>0745</u>	PURGING ENDED AT: <u>0805</u>	TOTAL VOLUME PURGED (gallons): <u>1.90</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0755	1.00	1.00	0.10	12.15	7.48	24.20	0.856	2.87	31.4	clear	yes
0758	0.30	1.30	0.10	12.39	7.48	24.31	0.853	2.73	17.6	clear	yes
0801	0.30	1.60	0.10	12.80	7.50	24.33	0.856	2.60	12.1	clear	yes
0804	0.30	1.90	0.10	12.99	7.50	24.35	0.867	2.50	11.5	clear	yes
<i>ended here 8/8/08</i>											

ORP
-164
-165
-168
-162

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Neil Davis / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>0806</u>	SAMPLING ENDED AT: <u>0810</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>280</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>6.60</u>	TUBING MATERIAL CODE: <u>R</u>	
FIELD DECONTAMINATION: <u>N</u>	FIELD-FILTERED: <u>Y</u> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-202	3	CG	40mL	HCL	-	6.2	22005/22006/22007 RFPF	

REMARKS: Placed new 'dedicated' tubing in well

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: Tellicoast - STA	SITE LOCATION: Pinusota, FL
WELL NO: MW-204	SAMPLE ID: MW-204 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 11 feet to 16 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR: BAILER: PP (geopump)
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 25 feet) + 0.50 gallons = 0.566 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~13.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 13.5	PURGING INITIATED AT: 1020	PURGING ENDED AT: 1034	TOTAL VOLUME PURGED (gallons): 1.68
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
5 1025	0.60	0.6	0.12	5.06	6.67	24.54	650	0.70	17.1	clear	yes
1028	0.36	0.96	0.12	5.07	6.66	24.69	736	0.49	12.0	clear	yes
1031	0.36	1.32	0.12	5.07	6.65	24.72	758	0.47	9.4	clear	yes
1034	0.36	1.68	0.12	5.07	6.65	24.72	753	0.39	8.2	clear	yes
<i>Analysis 1/23/08</i>											

ORP:
-145
-50
-149
-148

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Smith / ARCADIS	SAMPLER(S) SIGNATURES: <i>Neil Smith</i>	SAMPLING INITIATED AT: 1035	SAMPLING ENDED AT: 1040
PUMP OR TUBING DEPTH IN WELL (feet): ~13.5	SAMPLE PUMP FLOW RATE (mL per minute): ±100ml	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> NO FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="radio"/> N	
Filtration Equipment Type: _____			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-204	3	CG	40ml	HCL	-	<2	P200B+SM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-205	SAMPLE ID: MW-205 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 30 feet to 35 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 40 feet) + 0.5 gallons = 0.6 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 32.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 32.5	PURGING INITIATED AT: 1018	PURGING ENDED AT: 1037	TOTAL VOLUME PURGED (gallons): 1.14
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1028	0.60	0.60	0.06	11.92	7.23	25.05	729	3.95	2.27	clear	no
1031	0.18	0.78	0.06	12.29	7.27	25.13	727	4.12	2.31	clear	no
1034	0.18	0.96	0.06	12.69	7.28	25.14	729	4.15	2.46	clear	no
1037	0.18	1.14	0.06	13.05	7.31	25.10	712	4.13	2.87	clear	no

**MS
4/13/08
APP**

192
192
192
184

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SWINS ARCADIS	SAMPLER(S) SIGNATURES: 	SAMPLING INITIATED AT: 1039	SAMPLING ENDED AT: 1044
PUMP OR TUBING DEPTH IN WELL (feet): 32.5	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-205	3	CG	40mL	HCl	40mL	—	8260 + 8260 SIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES:** 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Tellevert</i>		SITE LOCATION: <i>Bradley, FL</i>	
WELL NO: <i>MW-206</i>		SAMPLE ID: <i>MW-206</i>	
DATE: <i>1/29/08</i>			

PURGING DATA

WELL DIAMETER (inches): <i>2</i>	TUBING DIAMETER (inches): <i>4/4</i>	WELL SCREEN INTERVAL DEPTH: <i>90</i> feet to <i>100</i> feet	STATIC DEPTH TO WATER (feet): <i>1098</i>	PURGE PUMP TYPE OR BAILER: <i>APP Geopon</i>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<i>0.0026</i> gallons/foot X <i>105</i> feet) + <i>0.25</i> gallons = <i>0.52</i> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~95</i>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~95</i>	PURGING INITIATED AT: <i>1345</i>	PURGING ENDED AT: <i>1355</i>	TOTAL VOLUME PURGED (gallons): <i>1.60</i>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<i>1349</i>	<i>0.64</i>	<i>0.64</i>	<i>0.16</i>	<i>20.49</i>	<i>7.32</i>	<i>24.70</i>	<i>769</i>	<i>0.74</i>	<i>14.8</i>	<i>clear</i>	<i>yes</i>
<i>1352</i>	<i>0.48</i>	<i>1.12</i>	<i>0.16</i>	<i>20.49</i>	<i>7.31</i>	<i>24.73</i>	<i>770</i>	<i>0.72</i>	<i>12.1</i>	<i>clear</i>	<i>yes</i>
<i>1355</i>	<i>0.48</i>	<i>1.60</i>	<i>0.16</i>	<i>20.51</i>	<i>7.31</i>	<i>24.75</i>	<i>770</i>	<i>0.71</i>	<i>10.9</i>	<i>clear</i>	<i>yes</i>
<i>no further 1/29/08</i>											

ORP
-260
-261
-262

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Neil Smith / ARCADIS</i>		SAMPLER(S) SIGNATURES: <i>Neil Smith</i>		SAMPLING INITIATED AT: <i>1359</i>	SAMPLING ENDED AT: <i>1405</i>
PUMP OR TUBING DEPTH IN WELL (feet): <i>~95</i>		SAMPLE PUMP FLOW RATE (mL per minute): <i>200</i>		TUBING MATERIAL CODE: <i>PE</i>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm		DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>MW-206</i>	<i>3</i>	<i>CG</i>	<i>42ml</i>	<i>HCL</i>	<i>120 ml</i>	<i>12</i>	<i>260B+820514 RPPP</i>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RPPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

165 + D

SITE NAME: Former ABC Facility		SITE LOCATION: Tallahassee, FL	
WELL NO: MW-207	SAMPLE ID: MW-207	DATE: 1.29.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 281 feet to 301 feet	STATIC DEPTH TO WATER (feet): 9.5	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 281 feet) + 0.25 gallons = 0.98 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 281	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 281	PURGING INITIATED AT: 13:40	PURGING ENDED AT: 14:10	TOTAL VOLUME PURGED (gallons): 2.0							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
13:50	1.0	1.0	0.1	13.3	11.45	24.95	1.57	1.05	1	clear	none
13:55	0.25	1.25	0.05	14.4	11.49	25.13	1.57	0.90	0	"	"
14:00	0.25	1.50	0.05	14.4	11.49	24.9	1.58	0.91	0	"	"
14:05	0.25	1.75	0.05	14.4	11.48	25.02	1.57	0.87	0	"	"
14:10	0.25	2.00	0.05	14.4	11.47	25.01	1.57	0.83	0	"	"

ORP
-224
-223
-232
-233
-284

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 14:11	SAMPLING ENDED AT: 14:15
PUMP OR TUBING DEPTH IN WELL (feet): 281 (MDS 9/15/08)	SAMPLE PUMP FLOW RATE (mL per minute): 100 (MDS 9/15/08)	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: _____ µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-207	3	CG	40ml	HCl	—	—	8260	RFPP

MDS 9/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast, FL
WELL NO: MW-208	SAMPLE ID: MW-208 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 16 feet to 21 feet	STATIC DEPTH TO WATER (feet): 3.30	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

= (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026** gallons/foot X feet) + **0.25** gallons = gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1418	PURGING ENDED AT: 1438	TOTAL VOLUME PURGED (gallons): 20
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1429	1.1	1.1	0.1	3.67	7.05	24.87	1.79	0.33	2.15	none	none
1432	0.3	1.4	↓	3.68	7.04	24.99	1.78	0.30	3.34	↓	↓
1435	0.3	1.7	↓	3.68	7.03	25.05	1.75	0.26	1.91	↓	↓
1438	0.3	2.0	↓	3.68	7.03	25.15	1.74	0.31	3.44	↓	↓
							µS/cm				

ORP
-192
-196
-197
-196

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ami Cook / Arcadis	SAMPLER(S) SIGNATURES: Ami Cook	SAMPLING INITIATED AT: 1439	SAMPLING ENDED AT: 1440
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 1100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-208	3	CG	40mL	HCl	—	—	8260	RFPP

MPS 9/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. **STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)**
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallerost, FL
WELL NO: MW-209	SAMPLE ID: MW-209 DATE: 1.23.08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 41 feet to 46 feet	STATIC DEPTH TO WATER (feet): 7.16	PURGE PUMP TYPE OR BAILER: PP			
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons							
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons							
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 14:19	PURGING ENDED AT: 14:46	TOTAL VOLUME PURGED (gallons):	

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
14:23	0.5	0.5	0.1	10.7	6.96	25.4	2.59	0.96	0	clear	none
14:30	1.0	1.0	0.1	13.5	6.97	25.5	2.39	0.52	11	"	"
14:35	1.5	1.5	0.1	15.1	6.98	25.56	2.38	0.47	11	"	"
14:40	2.0	2.0	0.1	15.4	7.02	25.73	2.56	0.43	14	"	"

ORP
-195
-213
-219

mS/cm

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes (Arcadis)	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 14:41	SAMPLING ENDED AT: 14:45
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: µm	DUPLICATE: Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-209	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: *Purging as slow as possible to prevent drawdown*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallahassee, FL
WELL NO: MW-210	SAMPLE ID: MW-210 DATE: 1.23.08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 4"	WELL SCREEN INTERVAL DEPTH: 93 feet to 103 feet	STATIC DEPTH TO WATER (feet): 8.14	PURGE PUMP TYPE OR BAILER: PP							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X feet) + 0.25 gallons = gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1543	PURGING ENDED AT: 1604	TOTAL VOLUME PURGED (gallons): 2.1							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1552	0.9	0.9	0.1	10.78	7.42	24.76	0.95	0.58	0.97		
1555	0.3	1.2	↓	11.18	7.43	24.74	0.95	0.60	1.07		
1558	0.3	1.5	↓	11.46	7.43	24.65	0.95	0.79	1.17		
1601	0.3	1.8	↓	11.82	7.43	24.60	0.95	0.75	1.09		
1604	0.3	2.1	↓	11.96	7.43	24.60	0.95	0.72	1.03		
							µS/cm				
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

ORP
-232
-238
-241
-244
-245

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Ami Coats / Arcadis				SAMPLER(S) SIGNATURES: <i>Ami Coats</i>				SAMPLING INITIATED AT: 1606		SAMPLING ENDED AT: 1610	
PUMP OR TUBING DEPTH IN WELL (feet):				SAMPLE PUMP FLOW RATE (mL per minute): <100 (MDS 4/15/08)				TUBING MATERIAL CODE: PE			
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N				FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm				DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH					
MW-210	3	CG	40ml	HCl	—	—	8260		RF PP		
REMARKS:											
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)											
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)											

MDS
4/15/08

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC		SITE LOCATION: Tallerost, FL	
WELL NO: MW-211	SAMPLE ID: MW-211	DATE: 1.23.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 140 feet to 150 feet	STATIC DEPTH TO WATER (feet): 7.8	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + 0.0026 gallons/foot X feet) + 0.25 gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1509	PURGING ENDED AT: 1545	TOTAL VOLUME PURGED (gallons): 4.0
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1515	1.0	1.0	0.1	10.25	8.01	25.60	0.466	0.53	0	clear	none
1520	1.5	1.5	0.1	11.20	7.78	25.4	0.468	0.45	618	turbid	none
1525	2.0	2.0	0.1	12.2	7.73	25.4	0.467	0.39	341	turbid	none
1530	2.5	2.5	0.1	13.25	7.78	25.4	0.463	0.37	82	sl. turbid	none
1535	3.0	3.0	0.1	13.95	7.95	25.3	0.456	0.35	16	clear	none
1540	3.5	3.5	0.1	14.1	7.92	25.2	0.447	0.33	15	clear	none
1545	4.0	4.0	0.1	14.2	7.95	25.20	0.446	0.31	12	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Micah Forbes / Arcadis	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1546	SAMPLING ENDED AT: 15:50
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: 4100 µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-211	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallahassee, FL
WELL NO: MW-212	SAMPLE ID: MW-212 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 4"	WELL SCREEN INTERVAL DEPTH: 281 feet to 301 feet	STATIC DEPTH TO WATER (feet): 4.04	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
only fill out if applicable)

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= **0** gallons + (**0.0026** gallons/foot X **4** feet) + **0.25** gallons = **0.0104** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 1507	PURGING ENDED AT: 1518	PURGED VOLUME (gallons): 1.41
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1512	0.5	0.5	0.1	6.61	7.69	24.96	1.09	0.32	0.0	none	none
1515	0.3	0.8	↓	7.36	7.70	24.96	1.09	0.32	0.0	↓	↓
1518	0.3	1.1	↓	7.69	7.71	24.98	1.09	0.31	0.0	↓	↓

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amy Poats / Arcadis	SAMPLER(S) SIGNATURE: <i>Amy Poats</i>	SAMPLING INITIATED AT: 1519	SAMPLING ENDED AT: 1520
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 4100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-212	3	CG	40mL	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. **STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)**
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-213	SAMPLE ID: MW-213
DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 2.29	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (.0026 gallons/foot X 23 feet) + .25 gallons = gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 3:18	PURGING ENDED AT: 1:54	TOTAL VOLUME PURGED (gallons): 2.3

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
15:25	0.7	0.7	0.1	3.82	6.14	23.87	.755	2.06	.19	clear	none
15:28	0.3	1.0	0.1	4.09	6.18	24.0	.747	1.61	.30	clear	none
15:31	0.3	1.3	0.1	4.17	6.21	24.07	.740	1.49	2.06	clear	none
15:34	0.3	1.7	0.1	4.25	6.33	24.15	.740	1.38	0.29	clear	none
15:37	0.3	2.0	0.1	4.3	6.44	24.19	.736	1.33	0.57	clear	none
15:40	0.3	2.3	0.1	4.33	6.51	24.21	.739	1.31	0.43	clear	none

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Derrin Johnson	SAMPLER(S) SIGNATURES: Derrin Johnson	SAMPLING INITIATED AT: 15:41	SAMPLING ENDED AT: 15:46
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N FILTER SIZE: µm	DUPLICATE: Y N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-213	3	CG	90ml	HLL	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG



SITE NAME: Tallevast **MDS 9/15/08** SITE LOCATION: Sarasota, FL **MDS 9/15/08**
 WELL NO: MW-214 SAMPLE ID: MW-214 DATE: 1/31/08

PURGING DATA
 WELL SCREEN INTERVAL: 1/4 inch TUBING DIAMETER (inches): 1/4 inch
 STATIC DEPTH TO WATER (feet): 9.99 PURGE PUMP TYPE OR BAILER: PP
 WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 = (30 feet - 35 feet) X 0.0026 gallons/foot = 0 gallons
 EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 = 0 gallons + (0.0026 gallons/foot X 35 feet) + 0.25 gallons = 0.115 gallons
 INITIAL PUMP OR TUBING DEPTH IN WELL (feet): _____ FINAL PUMP OR TUBING DEPTH IN WELL (feet): _____
 PURGING INITIATED AT: 1511 PURGING ENDED AT: 1533 TOTAL VOLUME PURGED (gallons): 2.3 **MDS 9/15/08**

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	GRP
1516	0.5	0.5	0.1	10.27	6.9	25.0	91	4.6	1.23	clear	none	170
1519	0.3	0.8		10.28	6.8	25.1	91	3.5	1.46			151
1521	0.3	1.1		10.29	6.8	25.2	91	3.2	0.62			145
1524	0.3	1.4		10.28	6.9	25.1	90	3.1	0.34			140
1527	0.3	1.7		10.28	6.9	25.1	90	3.0	0.24			136
1530	0.3	2.0		10.28	6.9	25.0	90	2.8	0.28			129
1533	0.3	2.3			6.9	25.0	90	2.9	0.45			
1536												

Sampled @ 1535

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) AFFILIATION: [Signature] ARCADIS SAMPLER(S) SIGNATURES: Kate Kerwell
 PUMP OR TUBING DEPTH IN WELL (feet): _____ SAMPLE PUMP FLOW RATE (mL per minute): 400 **MDS 9/15/08** TUBING MATERIAL CODE: _____
 FIELD DECONTAMINATION: Y N **MDS 9/15/08** FIELD-FILTERED: Y N **MDS 9/15/08** FILTER SIZE: _____ µm
 Filtration Equipment Type: _____ DUPLICATE: Y N **MDS 9/15/08**
 SAMPLING INITIATED AT: 1535 SAMPLING ENDED AT: 1540

SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-214	3	CG	90ml	HCL	-	✓	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2);
 optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 4/15/08	SITE LOCATION: <u>Sarasota FL</u> MDS 4/15/08
WELL NO: <u>MW-215</u>	SAMPLE ID: <u>MW-215</u> DATE: <u>1/30/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>90</u> feet to <u>100</u> feet	STATIC DEPTH TO WATER (feet): <u>13.41</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = <u>0</u> gallons + (<u>00026</u> gallons/foot X <u>100</u> feet) + <u>.25</u> gallons = <u>.51</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>95</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>95</u>	PURGING INITIATED AT: <u>1419</u>	PURGING ENDED AT: <u>1439</u> MDS 4/15/08	TOTAL VOLUME PURGED (gallons): <u>2.0</u> MDS 4/15/08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1424	0.5	0.5	0.1	13.61	7.5	25.0	.78	4.0	46.6	cloudy	none	ORP 60
1427	0.8	0.3	0.1	13.64	7.5	24.8	.78	3.4	13.9			63
1430	1.1	0.3	0.1	13.64	7.3	24.8	.76	3.1	2.84	clear		36
1433	1.4	0.3	0.1	13.64	7.2	24.8	.77	3.0	1.10			24
1436	1.7	0.3	0.1	13.64	7.2	24.7	.77	3.0	.80			23
1439	2.0	0.3	0.1	13.64	7.2	24	.77	2.9	.47			

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasey Cornwell / ARCADIS</u>			SAMPLER(S) SIGNATURES:			SAMPLING INITIATED AT: <u>1445</u>		SAMPLING ENDED AT: <u>1450</u>	
PUMP OR TUBING DEPTH IN WELL (feet): <u>95</u>			SAMPLE PUMP FLOW RATE (mL per minute): <u><100</u> MDS 4/15/08			TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N			FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm			DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-215</u>	<u>3</u>	<u>LG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <i>Tellevesi</i>	SITE LOCATION: <i>Sarasota, FL</i>
WELL NO: <i>MW-216</i>	SAMPLE ID: <i>MW-216</i> DATE: <i>1/31/08</i>

PURGING DATA

WELL DIAMETER (inches): <i>2</i>	TUBING DIAMETER (inches): <i>1/2</i>	WELL SCREEN INTERVAL DEPTH: <i>140</i> feet to <i>150</i> feet	STATIC DEPTH TO WATER (feet): <i>14.72</i>	PURGE PUMP TYPE OR BAILER: <i>(APP) peristaltic</i>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (<i>0.022</i> gallons/foot X <i>160</i> feet) + <i>0.25</i> gallons = <i>(0.25)</i> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~145</i>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~145</i>	PURGING INITIATED AT: <i>1505</i>	PURGING ENDED AT: <i>1518</i>	TOTAL VOLUME PURGED (gallons): <i>1.30</i>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm) <i>(circle)</i>	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<i>1512</i>	<i>0.20</i>	<i>0.70</i>	<i>0.10</i>	<i>17.21</i>	<i>7.33</i>	<i>24.79</i>	<i>0.683</i>	<i>0.63</i>	<i>13.6</i>	<i>clear</i>	<i>yes</i>
<i>1515</i>	<i>0.30</i>	<i>1.00</i>	<i>0.10</i>	<i>18.42</i>	<i>7.36</i>	<i>24.77</i>	<i>0.680</i>	<i>0.58</i>	<i>16.4</i>	<i>clear</i>	<i>yes</i>
<i>1518</i>	<i>0.30</i>	<i>1.30</i>	<i>0.10</i>	<i>18.45</i>	<i>7.39</i>	<i>24.76</i>	<i>0.679</i>	<i>0.56</i>	<i>14.9</i>	<i>clear</i>	<i>yes</i>
<i>sw purged 1/31/08</i>											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Neil Swisher / ARCADIS</i>	SAMPLER(S) SIGNATURES: <i>Neil Swisher</i>	SAMPLING INITIATED AT: <i>1520</i>	SAMPLING ENDED AT: <i>1525</i>
PUMP OR TUBING DEPTH IN WELL (feet): <i>~145</i>	SAMPLE PUMP FLOW RATE (mL per minute): <i>400</i>	TUBING MATERIAL CODE: <i>1F</i>	
FIELD DECONTAMINATION: <i>Y</i> N	FIELD-FILTERED: <i>Y</i> N FILTER SIZE: _____ µm	DUPLICATE: <i>Y</i> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>MW-216</i>	<i>3</i>	<i>CG</i>	<i>400</i>	<i>HCL</i>	<i>-</i>	<i>6.2</i>	<i>8260 BETA 2000 SM RFPF</i>	

REMARKS: *Placed new lock & cap on well*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast MDS 4/15/08	SITE LOCATION: Sarasota FL MDS 4/15/08
WELL NO: MW-217	SAMPLE ID: MW-217 DATE: 11/31/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 280 feet to 300 feet	STATIC DEPTH TO WATER (feet): 13.03	PURGE PUMP TYPE OR BAILER: Peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY = (300 feet - 13.03 feet) X 0.75 gallons/foot = 0.78 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME = 0 gallons + (300 gallons/foot X 13.03 feet) + 0.25 gallons = 6.83 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 290	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 290	PURGING INITIATED AT: 1422	PURGING ENDED AT: 1619	TOTAL VOLUME PURGED (gallons): 5.24 MDS 4/15/08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1510	2.4	2.4	0.05	20.20	7.63	26.97	0.587	0.94	3.51	clear	—
1525	0.75	3.15	0.05	21.09	7.63	26.31	0.584	0.94	1.91	clear	—
1540	0.75	3.90	0.05	21.74	7.64	25.82	0.587	0.95	1.23	clear	—
1555	0.75	3.460	0.05	22.20	7.64	25.14	0.584	0.96	1.03	clear	—
1610	0.40	5.0	0.026	22.50	7.64	24.57	0.584	0.97	1.01	clear	—
1613	0.08	5.08	0.026	22.51	7.64	24.59	0.584	0.98	1.05	clear	—
1616	0.08	5.16	0.026	22.53	7.64	24.59	0.584	0.98	1.00	clear	—
1619	0.08	5.24	0.026	22.54	7.64	24.60	0.584	0.98	1.00	clear	—

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Nathan Smith		SAMPLER(S) SIGNATURES: Nath Smith		SAMPLING INITIATED AT: 1619	SAMPLING ENDED AT: 1625			
PUMP OR TUBING DEPTH IN WELL (feet): 290 MDS 4/15/08		SAMPLE PUMP FLOW RATE (mL per minute): 4100 MDS 4/15/08		TUBING MATERIAL CODE:				
FIELD DECONTAMINATION: Y		FIELD-FILTERED: Y		DUPLICATE: Y (N)				
SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION					
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
MW-217	3	LG	40ml	HCL	-	-	8260	RFPP
REMARKS: Initial purge water is silty gray-blue in color. water levels would not stabilize until around 1610.								
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)								
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)								

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

0-15
0.99

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast		SITE LOCATION: Tallevast FL	
WELL NO: MW-219	SAMPLE ID: MW-219	DATE: 1/25/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 22 feet to 27 feet	STATIC DEPTH TO WATER (feet): 3.60	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 30 feet) + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 24.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 24.5	PURGING INITIATED AT: 1233	PURGING ENDED AT: 1247	TOTAL VOLUME PURGED (gallons): 1.82

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1241	1.04	1.04	0.53	3.97	6.6	24.6	0.418	0.8	1.95	clear	none
1243	0.26	1.30	0.13	3.97	6.5	24.6	0.416	1.0	1.21	clear	none
1245	0.26	1.56	0.13	3.99	6.5	24.6	0.416	1.4	1.00	clear	none
1247	0.26	1.82	0.13	4.00	6.5	24.6	0.411	1.4	0.89	clear	none

ORP
-111
-109
-110
-110

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / Arcadis	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1248	SAMPLING ENDED AT: 1252
PUMP OR TUBING DEPTH IN WELL (feet): 24.5	SAMPLE PUMP FLOW RATE (mL per minute): 2492	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-219	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **missing 1 bolt for cover**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.11

SITE NAME: <u>Tallevast</u>		SITE LOCATION: <u>Tallevast, FL</u>	
WELL NO: <u>MW-220</u>	SAMPLE ID: <u>MW-220</u>	DATE: <u>1/25/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>39</u> feet to <u>44</u> feet	STATIC DEPTH TO WATER (feet): <u>595</u>	PURGE PUMP TYPE OR BAILER: <u>Geopump (PP)</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>47</u> feet) + <u>0.25</u> gallons = <u>0.37</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>41.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>41.5</u>	PURGING INITIATED AT: <u>1232</u>	PURGING ENDED AT: <u>1246</u>	TOTAL VOLUME PURGED (gallons): <u>1.94</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
8 1240	1.12	1.12	0.14	10.82	7.31	25.0	794	0.56	7.68	Clear	NONE	-228
1242	.28	1.40	0.14	12.71	7.31	25.2	808	0.60	2.12	Clear	NONE	-230
1244	.28	1.68	0.14	13.09	7.32	25.2	802	0.58	0.62	clear	NONE	-232
1246	.28	1.96	0.14	13.90	7.31	25.1	806	0.55	0.62	CLEAR	NONE	-231

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>BRIDGET STALL / ARCADIS</u>		SAMPLER(S) SIGNATURES: <u>Bridget Stall</u>		SAMPLING INITIATED AT: <u>1247</u>	SAMPLING ENDED AT: <u>1251</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>41.5</u>		SAMPLE PUMP FLOW RATE (mL per minute): <u>~530</u>		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N		FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm		DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-220	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: missing bolt

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.62

SITE NAME: TAIlevAST	SITE LOCATION: TAIlevAST, FL
WELL NO: mw-221	SAMPLE ID: mw-221 DATE: 1-25-08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 97 feet to 107 feet	STATIC DEPTH TO WATER (feet): 8.88	PURGE PUMP TYPE OR BAILER: PP Geo Pump
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable)
 = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = **0** gallons + (**0.0026** gallons/foot X **110** feet) + **.25** gallons = **0.51** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 102	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 102	PURGING INITIATED AT: 1326	PURGING ENDED AT: 1332	TOTAL VOLUME PURGED (gallons): 2.40
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm)	DISSOLVED OXYGEN (circle % or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1326											
1327	1.68	1.68	0.12	10.69	7.30	25.2	1.24	0.52	64.4	Cloudy	yes
1328	0.24	1.92	0.12	10.76	7.31	25.3	1.24	0.51	61.0	Cloudy	yes
1330	0.24	2.16	0.12	10.80	7.32	25.1	1.24	0.50	59.8	cloudy	yes
1332	0.24	2.40	0.12	10.82	7.32	25.1	1.24	0.49	57.7	Cloudy	yes

14

ORP
-283
-286
-288
-288

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1333	SAMPLING ENDED AT: 1343
PUMP OR TUBING DEPTH IN WELL (feet): 102	SAMPLE PUMP FLOW RATE (mL per minute): 455	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-221	3	CG	40ml	HCl	40ml	-	8260	RF PP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

1.89

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST, FL	
WELL NO: mw-222		SAMPLE ID: mw-222	
DATE: 1-08			

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 133 feet to 143 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 140 feet) + .25 gallons = 0.63 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 138	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 138	PURGING INITIATED AT: 1309	PURGING ENDED AT: 1334	TOTAL VOLUME PURGED (gallons): 2.64
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1327	1.98	1.98	0.11	17.65	7.5	25.7	1.21	0.4	5.28	clear	yes
1329	0.22	2.20	0.11	18.15	7.5	25.6	1.30	0.4	3.88	clear	yes
1331	0.22	2.42	0.11	18.32	7.5	25.7	1.29	0.5	2.65	clear	yes
1333	0.22	2.64	0.11	18.56	7.5	25.7	1.31	0.4	1.53	clear	yes

ORP
-29
-20
-220
-222

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 1337	SAMPLING ENDED AT: 1341
PUMP OR TUBING DEPTH IN WELL (feet): 138	SAMPLE PUMP FLOW RATE (mL per minute): ~416	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-222	3	CG	40ml	HCl	40ml	-	8260	RFPP

REMARKS: **Missing bolt**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tellawayt</u>	SITE LOCATION: <u>Jarasota, FL</u>
WELL NO: <u>MW-223</u>	SAMPLE ID: <u>MW-223</u> DATE: <u>1/25/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>4 1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>10</u> feet to <u>15</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>(pp) geopump</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
= gallons + (<u>0.0026</u> gallons/foot X <u>30</u> feet) + <u>0.50</u> gallons = <u>0.578</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~12.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~12.5</u>	PURGING INITIATED AT: <u>1040</u>	PURGING ENDED AT: <u>1052</u>	TOTAL VOLUME PURGED (gallons): <u>1.32</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (umhos/cm or uS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
3 1043	0.63	0.63	0.21	5.75	6.77	22.97	920	0.78	5.26	clear	yes
1046	0.21	0.84	0.07	5.77	6.78	22.76	937	0.64	3.45	clear	yes
1049	0.27	1.11	0.07	5.77	6.79	22.78	931	0.66	4.07	clear	yes
1052	0.21	1.32	0.07	5.77	6.79	22.81	932	0.64	3.01	clear	yes

ORP
-95
-104
-108
-111

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>N. Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>N. Smith</u>	SAMPLING INITIATED AT: <u>1052</u>	SAMPLING ENDED AT: <u>1100</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~12.5</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> <u>N</u>	FIELD-FILTERED: <u>Y</u> <u>N</u> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <u>N</u>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-223</u>	<u>3</u>	<u>CG</u>	<u>40mL</u>	<u>HCL</u>	<u>-</u>	<u>6.2</u>	<u>82003 (8260) M</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallahassee	SITE LOCATION: Sarasota, FL
WELL NO: MW-224	SAMPLE ID: MW-224 DATE: 1/25/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 25 feet to 30 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0024 gallons/foot X 34 feet) + 0.50 gallons = (0.60) 0.58 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~27.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~27.5	PURGING INITIATED AT: 0955	PURGING ENDED AT: 1012	TOTAL VOLUME PURGED (gallons): 2.38

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle (100) or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
5 1000	0.70	0.70	0.14	12.42	6.85	25.41	1630	0.54	3.05	clear	yes	OK -134
1003	0.42	1.12	0.14	13.61	6.91	25.37	1590	0.95	2.76	clear	yes	-123
1006	0.42	1.54	0.14	14.89	6.92	25.47	1650	1.65	1.45	clear	yes	-115
1009	0.42	1.96	0.14	15.75	6.92	25.42	1630	0.60	1.22	clear	yes	-115
1012	0.42	2.38	0.14	15.99	6.92	25.42	1630	0.69	1.00	clear	yes	-116

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Neil Suf Arcadis	SAMPLER(S) SIGNATURES: Neil Suf	SAMPLING INITIATED AT: 1014	SAMPLING ENDED AT: 1020
PUMP OR TUBING DEPTH IN WELL (feet): ~27.5	SAMPLE PUMP FLOW RATE (mL per minute): 1100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-224	3	CG	40m	HCL	-	6.2	RFPP-224-SM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Tellevert</i>		SITE LOCATION: <i>Sarasota, FL</i>	
WELL NO: <i>MW-225</i>	SAMPLE ID: <i>MW-225</i>	DATE: <i>1/25/08</i>	

PURGING DATA

WELL DIAMETER (inches): <i>2</i>	TUBING DIAMETER (inches): <i>1/4</i>	WELL SCREEN INTERVAL DEPTH: <i>40</i> feet to <i>45</i> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <i>PP</i> <i>MDS 9/15/08</i>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
= gallons + (<i>0.0026</i> gallons/foot X <i>50</i> feet) + <i>0.50</i> gallons = <i>0.63</i> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~43</i>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <i>~43</i>	PURGING INITIATED AT: <i>1052</i>	PURGING ENDED AT: <i>1109</i>	TOTAL VOLUME PURGED (gallons): <i>1.70</i>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<i>1100</i>	<i>0.80</i>	<i>0.80</i>	<i>0.10</i>	<i>6.82</i> <i>6.82</i>	<i>6.82</i>	<i>24.69</i>	<i>181</i>	<i>0.45</i>	<i>13.5</i>	<i>clear</i>	<i>yes</i>
<i>1103</i>	<i>0.30</i>	<i>1.10</i>	<i>0.10</i>	<i>12.41</i>	<i>6.84</i>	<i>24.65</i>	<i>180</i>	<i>0.45</i>	<i>7.4</i>	<i>clear</i>	<i>yes</i>
<i>1106</i>	<i>0.30</i>	<i>1.40</i>	<i>0.10</i>	<i>12.01</i>	<i>6.85</i>	<i>24.60</i>	<i>181</i>	<i>0.47</i>	<i>3.9</i>	<i>clear</i>	<i>yes</i>
<i>1109</i>	<i>0.70</i>	<i>1.70</i>	<i>0.10</i>	<i>13.47</i>	<i>6.85</i>	<i>24.53</i>	<i>181</i>	<i>0.46</i>	<i>2.4</i>	<i>clear</i>	<i>yes</i>
<i>Not done 1/25/08</i>											

-159
-155
-158
-15A

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>N. Smith / AREAS</i>	SAMPLER(S) SIGNATURES: <i>N. Smith</i>	SAMPLING INITIATED AT: <i>1110</i>	SAMPLING ENDED AT: <i>1120</i>
PUMP OR TUBING DEPTH IN WELL (feet): <i>~43</i>	SAMPLE PUMP FLOW RATE (mL per minute): <i>1000</i>	TUBING MATERIAL CODE: <i>PE</i>	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y FILTER SIZE: µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<i>MW-225</i>	<i>3</i>	<i>CG</i>	<i>40mL</i>	<i>HCL</i>	<i>-</i>	<i>~2</i>	<i>826B+826DM RFPP</i>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: <u>Telleart</u>	SITE LOCATION: <u>Sarasota FL</u>
WELL NO: <u>MW-226</u>	SAMPLE ID: <u>MW-226</u> DATE: <u>1/25/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>1/4</u>	WELL SCREEN INTERVAL DEPTH: <u>90</u> feet to <u>100</u> feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: <u>(pp) gasyne</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
= gallons + (<u>8.0026</u> gallons/foot X <u>98</u> feet) + <u>0.50</u> gallons = <u>0.75</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~95</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>~95</u>	PURGING INITIATED AT: <u>1130</u>	PURGING ENDED AT: <u>1147</u>	TOTAL VOLUME PURGED (gallons): <u>1.70</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or (S/cm))	DISSOLVED OXYGEN (circle DO or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
8 1138	0.80	0.80	0.10	6.31	7.34	23.15	101	0.49	10.34	clear	NO
1141	0.30	1.10	0.10	6.31	7.34	22.98	103	0.47	9.1	clear	NO
1144	0.30	1.40	0.10	6.31	7.33	22.10	103	0.44	8.21	clear	NO
1147	0.30	1.70	0.10		7.30	23.18	108	0.45	3.9	clear	NO

ORP
-181
-182
-184
-187

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>N. Smith / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>N. Smith</u>	SAMPLING INITIATED AT: <u>1148</u>	SAMPLING ENDED AT: <u>1155</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>~95</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE: <u>PE</u>	
FIELD DECONTAMINATION: <u>Y</u> N	FIELD-FILTERED: <u>Y</u> <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: <u>Y</u> <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-226</u>	<u>3</u>	<u>CG</u>	<u>40</u>	<u>HCL</u>	<u>-</u>	<u>< 2</u>	<u>8260B + 826 SM RFPF</u>	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEMIST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-227	SAMPLE ID: MW-227
DATE: 1/25/08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 160 feet) + 0.5 gallons = 0.92 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150	PURGING INITIATED AT: 0948	PURGING ENDED AT: 1024	TOTAL VOLUME PURGED (gallons): 2.16								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1004	0.96	0.96	0.06	12.33	6.88	21.03	1070	1.99	2.09	clear	yes	-170
1008	0.24	1.20	0.06	13.09	7.15	21.46	1060	1.88	1.38	clear	yes	-192
1012	0.24	1.44	0.06	13.45	7.20	21.56	1060	0.97	1.22	clear	yes	-200
1016	0.24	1.68	0.06	14.04	7.24	21.86	1050	0.67	2.12	clear	yes	-211
1020	0.24	1.92	0.06	14.66	7.26	21.97	1050	0.58	2.22	clear	yes	-218
1024	0.24	2.16	0.06	15.59	7.27	22.02	1050	0.53	0.94	clear	yes	-222

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/NS/ARADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 1026	SAMPLING ENDED AT: 1032
PUMP OR TUBING DEPTH IN WELL (feet): 150	SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-227	3	CG	40mL	HCl	40mL	—	8260 + 8260 SM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

**Form FD 9000-24
GROUNDWATER SAMPLING LOG**

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA, FL
WELL NO: MW-228	SAMPLE ID: MW-228
DATE: 1/23/08	

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 98 feet to 108 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = 0.5 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 113 feet) + 0.5 gallons = 0.80 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 103	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 103	PURGING INITIATED AT: 0925	PURGING ENDED AT: 0941	PURGE VOLUME PURGED (gallons): 1.92
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0932	0.84	0.84	0.12	18.70	7.27	24.58	920	0.42	17.91	clear	yes	-242
0935	0.36	1.20	0.12	20.94	7.34	24.55	990	0.42	11.10	clear	yes	-245
0938	0.36	1.56	0.12	21.50	7.29	24.67	999	0.47	6.78	clear	yes	-246
0941	0.36	1.92	0.12	22.25	7.34	24.65	999	0.41	4.44	clear	yes	-246

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NS/SW/ARCADIS	SAMPLER(S) SIGNATURES: <i>[Signature]</i>	SAMPLING INITIATED AT: 0944	SAMPLING ENDED AT: 0948
PUMP OR TUBING DEPTH IN WELL (feet): 103	SAMPLE PUMP FLOW RATE (mL per minute): 100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> N	
Filtration Equipment Type:			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-228	3	CG	40mL	HCl	40mL	—	8260 + 8260 SIM	RFPP
DUP-4	3	CG	40mL	HCl	40mL	—	8260 + 8260 SIM	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 4/15/08	SITE LOCATION: <u>Sarasota FL</u> MDS 9/15/08
WELL NO: <u>MW-229</u>	SAMPLE ID: <u>MW-229</u>
DATE: <u>1/29/08</u>	

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>0.25</u>	WELL SCREEN INTERVAL DEPTH: <u>17.5</u> feet to <u>22.5</u> feet	STATIC DEPTH TO WATER (feet): <u>6.43</u>	PURGE PUMP TYPE OR BAILER: <u>Peristaltic</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>6</u> gallons + (<u>0.0026</u> gallons/foot X <u>23</u> feet) + <u>0.25</u> gallons = <u>0.2</u> gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>21</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>21</u>	PURGING INITIATED AT: <u>10:20</u>	PURGING ENDED AT: <u>1042</u>	TOTAL VOLUME PURGED (gallons): <u>2.2</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1030		1.0	0.1	6.59	5.35	27.38	0.636	1.56	6.38	clear	-	20
1033	0.3	1.3	0.1	6.59	5.34	27.34	0.636	1.48	5.21	clear	-	30
1036	0.3	1.6	0.1	6.59	5.34	27.43	0.637	1.37	1.58	clear	-	27
1039	0.3	1.9	0.1	6.59	5.34	27.48	0.633	1.25	1.02	clear	-	22
1042	0.3	2.2	0.1	6.59	5.34	27.47	0.635	1.16	1.07	"	-	18
Sampled @ 1042												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith</u>	SAMPLER(S) SIGNATURES: <u>Nathan Smith</u>	SAMPLING INITIATED AT: <u>1042</u>	SAMPLING ENDED AT: <u>1048</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>21</u> MDS 9/15/08	SAMPLE PUMP FLOW RATE (mL per minute): <u>4100</u> MDS 9/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u>	FIELD-FILTERED: <u>Y</u> N	FILTER SIZE: <u> </u> µm	DUPLICATE: <u>Y</u> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-229	3	CG	90ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 9/15/08	SITE LOCATION: Suragoba FL MDS 9/15/08
WELL NO: MW-230	SAMPLE ID: MW-230 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 31 feet to 36 feet	STATIC DEPTH TO WATER (feet): 6.85	PURGE PUMP TYPE: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 36 feet) + 0.25 gallons = 0.34 gallons MDS 9/15/08				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 33	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 33	PURGING INITIATED AT: 1022	PURGING ENDED AT: 1039	TOTAL VOLUME PURGED (gallons): 1.4								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1027	.5	.5	0.1	9.36	6.9	26.6	.63	3.9	0.33	Clear	none	ORP 138
1030	.3	.8	↓	9.84	6.9	26.8	.62	2.7	1.26	↓	↓	72
1033	.3	1.1	↓	10.21	7.0	26.8	.62	2.7	1.83	↓	↓	48
1036	.3	1.4	↓	11.40	7.0	26.8	.63	2.6	1.94	↓	↓	38
1039	→ Started Sampling @ 1039 ←											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Rasey Cornwell / ARCADIS	SAMPLER(S) SIGNATURES: <i>Rasey Cornwell</i>	SAMPLING INITIATED AT: 1039	SAMPLING ENDED AT: 1045
PUMP OR TUBING DEPTH IN WELL (feet): 33	SAMPLE PUMP FLOW RATE (mL per minute): 2100 MDS 9/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-230	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/15/08	SITE LOCATION: Sarasota FL MDS 4/15/08
WELL NO: MW-231	SAMPLE ID: MW-231 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 96 feet to 107 feet	STATIC DEPTH TO WATER (feet): 20.65	PURGE PUMP TYPE OR BAILER: Peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (8 feet - 107 feet) X 50.25 gallons/foot = _____ gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0006 gallons/foot X 107 feet) + 0.25 gallons = 0.5 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 102	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 102	PURGING INITIATED AT: 1130	PURGING ENDED AT: 1036	TOTAL VOLUME PURGED (gallons): 1.6

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
1027	0.7	0.7	0.1	21.25	6.7	25.7	0.77	3.4	0.90	Clear	—	12
1030	0.3	1.0	0.1	21.25	6.7	25.7	0.77	2.9	1.50	"	—	2
1033	0.3	1.3	0.1	21.25	6.7	25.6	0.77	1.0-2.7	1.0	"	—	-6
1036	0.3	1.6	0.1	21.25	6.7	25.6	0.77	2.6	1.1	"	—	-10
MDS 4/15/08												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: ARLAJIS MDS 4/15/08	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1030	SAMPLING ENDED AT:
PUMP OR TUBING DEPTH IN WELL (feet): 102 MDS 4/15/08	SAMPLE PUMP FLOW RATE (mL per minute): 2100 MDS 4/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-231	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 40% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: SARASOTA
WELL NO: MW-232	SAMPLE ID: MW-232 DATE: 1/30/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 98 feet to 108 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 113 feet) + 0.5 gallons = 0.79 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 103	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 103	PURGING INITIATED AT: 1400	PURGING ENDED AT: 1455	TOTAL VOLUME PURGED (gallons): 3.32
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1427	1.62	1.64	0.06	14.11	7.31	25.60	700	5.27	1.48	clear	yes	-284
1431	0.24	1.88	0.06	14.11	7.31	25.62	701	5.11	1.08	clear	yes	-284
1435	0.24	2.12	0.06	14.03	7.32	25.65	701	3.88	0.86	clear	yes	-292
1439	0.24	2.36	0.06	14.15	7.30	25.63	703	2.28	17.5	clear	yes	-304
1443	0.24	2.60	0.06	14.11	7.42	25.60	699	2.96	5.19	clear	yes	-318
1447	0.24	2.84	0.06	14.15	7.30	25.38	699	0.89	4.42	clear	yes	-324
1451	0.24	3.08	0.06	14.18	7.27	25.35	699	0.59	2.79	clear	yes	-328
1455	0.24	3.32	0.06	14.19	7.26	25.31	698	0.57	2.95	clear	yes	-330

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SW/TH / ARCADIS	SAMPLER(S) SIGNATURES: [Signature]	SAMPLING INITIATED AT: 1457	SAMPLING ENDED AT: 1503
PUMP OR TUBING DEPTH IN WELL (feet): 103	SAMPLE PUMP FLOW RATE (mL per minute): ≤100	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y <input type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-232	3	CG	40mL	HCl	40mL	—	8260B+8260BSM	RFPP
DUP-24	3	CG	40mL	HCl	40mL	—	11	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

1.56

SITE NAME: Tallevast	SITE LOCATION: Sarasota FL
WELL NO: mw-233	SAMPLE ID: mw-233
DATE: 1-24-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 90 feet to 100 feet	STATIC DEPTH TO WATER (feet): 16.08	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable) = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable) = **0** gallons + (**.0026** gallons/foot X **103** feet) + **.25** gallons = **0.52** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 95	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 95	PURGING INITIATED AT: 1621	PURGING ENDED AT: 1706	TOTAL VOLUME PURGED (gallons): 1.80
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1700	1.56	1.56	0.04	28.69	7.52	24.9	.635	0.45	53.4	slightly cloudy	yes
1702	0.08	1.64	0.04	28.85	7.55	24.9	.637	0.46	53.4	slightly cloudy	yes
1704	0.08	1.72	0.04	28.99	7.56	24.9	.636	0.45	50.1	slightly cloudy	yes
1706	0.08	1.80	0.04	29.06	7.56	24.8	.640	0.45	48.9	slightly cloudy	yes

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: Jennifer King	SAMPLING INITIATED AT: 1708	SAMPLING ENDED AT: 1715
PUMP OR TUBING DEPTH IN WELL (feet): 95	SAMPLE PUMP FLOW RATE (mL per minute): 2150	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-233	3	CG	40mL	HeL	40mL	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/15/08	SITE LOCATION: Sarasota FL MDS 4/15/08
WELL NO: MW-234	SAMPLE ID: MW-234 DATE: 1/31/08

PURGING DATA			
WELL DIAMETER (inches):	TUBING DIAMETER (inches): 4	WELL SCREEN INTERVAL DEPTH: 15 feet to 20 feet	STATIC DEPTH TO WATER (feet): 2.58
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY			PURGE PUMP TYPE OR BAILER: PP
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME			
= 0 gallons + (0.0026 gallons/foot X 20 feet) + 1.25 gallons = 0.052 gallons			

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1023	0.5	0.5	0.1	3.36	6.54	25.09	382	1.50	11.8	Clear	none
1026	0.3	0.8		3.39	6.61	25.08	396	1.31	10.43		
1029	0.3	1.1		3.42	6.67	25.17	412	1.20	8.27		
1032	0.3	1.4		3.58	6.65	25.20	417	1.18	12.0		
1036	0.3	1.7		3.52	6.67	25.21	434	1.17	17.60	"muddy"	
1038	0.3	2.0		3.53	6.66	25.25	433	1.15	49.5	Cloudy	
1041	0.3	2.3		3.53	6.67	25.29	871	1.10	11.6	Clearish	
1044	0.3	2.6		3.53	6.67	25.35	0.900	1.09	4.22	Clear	
1047	0.3	2.9	✓	3.53	6.67	25.37	0.900	1.08	3.42	↓	↓
1050	0.3	3.2			6.69	25.52	0.900	1.07	3.50	↓	↓

WELL CAPACITY (gallons per foot) = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIAMETER (gallons per foot) = 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLED BY (PRINT) / AFFILIATION: 1 ARCADIS	SAMPLER(S) SIGNATURES: MDS 4/15/08	SAMPLING INITIATED AT: 1055	SAMPLING ENDED AT: 1100
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): L100 MDS 4/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: N	FIELD-FILTERED: Y MDS 4/15/08	DUPLICATE: Y N MDS 4/15/08	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-234	3	LG	40ml	HCl	-	-	8260	RFP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast	SITE LOCATION: Sarasota FL	
WELL NO: MW-235	SAMPLE ID: MW-235	DATE: 1/31/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 35 feet to 40 feet	STATIC DEPTH TO WATER (feet): 10.41	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
= gallons + (0.002 gallons/foot X 40 feet) + 0.25 gallons = 0.35 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 38	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 38	PURGING INITIATED AT: 1144	PURGING ENDED AT: 1146	TOTAL VOLUME PURGED (gallons): 1.50
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1130		0.75	0.05	12.71	7.15	25.71	0.544	0.98	0.29	clear	-
1133	0.15	0.90	0.05	12.71	7.15	25.64	0.556	0.97	0.29	clear	-
1137	0.15	1.05	0.05	12.71	7.15	25.70	0.840	0.95	0.29	clear	-
1140	0.15	1.20	0.05	12.71	7.15	25.84	0.900	0.93	0.18	clear	-
1143	0.15	1.35	0.05	12.71	7.16	25.85	0.940	0.92	0.18	clear	-
1146	0.15	1.50	0.05	12.71	7.16	25.86	0.950	0.92	0.16	clear	-

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Nathan Smith	SAMPLER(S) SIGNATURES: Nathan Smith	SAMPLING INITIATED AT: 1146	SAMPLING ENDED AT: 1150
PUMP OR TUBING DEPTH IN WELL (feet): 38	SAMPLE PUMP FLOW RATE (mL per minute): 2100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: N	FIELD-FILTERED: Y	DUPLICATE: Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-235	3	CG	40ml	HCl	-	-	8260	R.F.P.P.

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/15/08
SITE LOCATION: Talleast
WELL NO: MW-236 SAMPLE ID: MW-236 DATE: 1-31-08

PURGING DATA
WELL DIAMETER (inches): _____ TUBING DIAMETER (inches): 1/4 GREEN INTERVAL DEPTH: 90 feet STATIC DEPTH TO WATER (feet): 1556 PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
(only fill out if applicable) = (_____ feet - _____ feet) X _____ gallons/foot = _____ gallons
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable) = 0 gallons + (1.0076 gallons/foot X 103 feet) + 1.75 gallons = 0.52 gallons MDS 4/15/08

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
10:47	2.0	2.0	0.1	10.12	7.53	24.70	0.596	1.39	15.7	murky	none
10:50	0.3	2.3	0.1	10.12	7.54	24.65	0.610	1.29	19.3	↓	↓
10:53	0.3	2.6	0.1	10.12	7.62	24.76	0.612	1.15	46.9	↓	↓
10:56	0.3	2.9	0.1	10.12	7.65	24.78	0.622	1.14	18.7	↓	↓
10:59	0.3	3.2	0.1	10.10	7.59	24.83	0.592	1.08	15.7	↓	↓
11:02	0.3	3.5	0.1	10.10	7.48	24.72	0.601	1.06	19.9	↓	↓

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0028; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA
SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson SAMPLER(S) SIGNATURES: Darrin Johnson MDS 4/15/08
PUMP OR TUBING DEPTH IN WELL (feet): _____ SAMPLE PUMP FLOW RATE (mL per minute): <100 TUBING MATERIAL CODE: T
FIELD DECONTAMINATION: Y N FIELD-FILTERED: Y N FILTER SIZE: _____ µm DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-236</u>	<u>3</u>	<u>LG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS: believe per fell down well. Regular ball point.
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast MDS 9/15/08	SITE LOCATION: Sarasota FL MDS 9/15/08
WELL NO: MW-237	SAMPLE ID: MW-237 DATE: 1/31/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet): 14.15	PURGE PUMP TYPE OR BAILER: Peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = 0.6 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (0.0026 gallons/foot X 150 feet) + 0.25 gallons = 0.6 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150		FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150		PURGING INITIATED AT: 1030		PURGING ENDED AT: 1051		TOTAL VOLUME PURGED (gallons): 1.65				
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1040	1.0	1.0	0.1	17.85	7.2	25.0	.600	3.4	4.19	clear	—	39
1045	0.25	1.25	0.05	17.75	7.2	25.0	.690	3.3	3.58	clear	—	38
1048	0.25	1.40	0.03	17.75	7.2	25.0	.690	3.3	4.10	clear	—	38
1051	0.15	1.65	0.05	17.75	7.3	25.1	.690	3.2	2.62	clear	—	38
<p>Sampled @ 1051 DUP-30 taken @ 1051</p>												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN SAITH	SAMPLER(S) SIGNATURES: <i>Nathan P. Saith</i>	SAMPLING INITIATED AT: 1051	SAMPLING ENDED AT: 1100
PUMP OR TUBING DEPTH IN WELL (feet): 150 MDS 9/15/08	SAMPLE PUMP FLOW RATE (mL per minute): 400 MDS 9/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y MDS 9/15/08	FIELD-FILTERED: <input checked="" type="checkbox"/> Y MDS 9/15/08	DUPLICATE: <input checked="" type="checkbox"/> Y DUP-30	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH			
MW-237	3	CG	90ml	HCl	—	—	8260	RFPP	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump

EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 40% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC	SITE LOCATION: Tallevast
WELL NO: MW-238	SAMPLE ID: MW-238
DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: feet to 300 feet	STATIC DEPTH TO WATER (feet): 1286	PURGE PUMP TYPE OR BAILER: PP
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WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 only fill out if applicable)
 = (feet - feet) X gallons/foot = gallons

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 (only fill out if applicable)
 = **0** gallons + (**10026** gallons/foot X **303** feet) + **125** gallons = **1.03** gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 10:48	PURGING ENDED AT: 11:10	TOTAL VOLUME PURGED (gallons): 1.93
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
11:01	1.03	1.03	0.1	17.5	6.85	24.76	1299	1.11	1.58	amber	none
11:04	.3	1.33	0.1	19.3	6.84	24.8	1285	1.01	1.57	amber	none
11:07	.3	1.63	0.1	21.0	6.83	24.79	1280	1.86	1.40	↓	↓
11:10	.3	1.93	0.1	21.34	6.83	24.88	1279	1.82	1.75	↓	↓

169
151

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darrin Johnson	SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>	SAMPLING INITIATED AT: 11:11	SAMPLING ENDED AT: 11:16
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): 4100 ^{MDS} _{4/15/08}	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-238	3	CG	40ml	HCl	—	—	8260	RFPP

MDS
4/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings < 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: TAIlerAST SITE LOCATION: Sarasota FL MDC 2/15/08
WELL NO: mw-239 SAMPLE ID: mw-239 DATE: 1-23-08

PURGING DATA

WELL DIAMETER (inches): 2 TUBING DIAMETER (inches): 1/4 WELL SCREEN INTERVAL DEPTH: 99 feet to 108 feet STATIC DEPTH TO WATER (feet): 15.56 PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
(only fill out if applicable) = (99 feet - 15.56 feet) X 0.0026 gallons/foot = 0.59 gallons
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 111 feet) + .25 gallons = 0.59 gallons MDC 2/15/08
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 103 FINAL PUMP OR TUBING DEPTH IN WELL (feet): 103 PURGING INITIATED AT: 0948 PURGING ENDED AT: 1116 TOTAL VOLUME PURGED (gallons): 1.76

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1110	1.64	1.64	0.02	17.06	7.06	24.7	<u>inst</u> 0.618	0.47	2.95	clear	yes
1112	.04	1.68	0.02	17.01	7.06	24.6	.617	0.47	2.95	clear	yes
1114	.04	1.72	0.02	17.01	7.07	24.6	.625	.48	2.71	clear	yes
1116	.04	1.76	0.02	17.01	7.08	24.6	.624	.47	2.35	clear	yes

ORP
-332
-333
-333
-334

82
2

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS SAMPLER(S) SIGNATURES: Jennifer King SAMPLING INITIATED AT: 1117 SAMPLING ENDED AT: 1120
PUMP OR TUBING DEPTH IN WELL (feet): 103 SAMPLE PUMP FLOW RATE (mL per minute): 78 TUBING MATERIAL CODE: PE
FIELD DECONTAMINATION: Y N FIELD-FILTERED: Y N FILTER SIZE: _____ µm DUPLICATE: Y N
Filtration Equipment Type: _____

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>mw239</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCL</u>	<u>40ml</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: STA	SITE LOCATION: Sarasota, FL
WELL NO: MW-240	SAMPLE ID: MW-240 DATE: 1/23/08

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 156.5 feet to 116.5 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: (PP)geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 170 feet) + 0.50 gallons = 0.44 gallons (1.00)				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): ~161	FINAL PUMP OR TUBING DEPTH IN WELL (feet): ~161	PURGING INITIATED AT: 1350	PURGING ENDED AT: 1410	TOTAL VOLUME PURGED (gallons): 2.60
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle µg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	GRP
1358	1.04	1.04	0.13	20.72	7.32	26.13	622	6.04	1.61	clear	no	-182
1401	0.39	1.43	0.13	21.14	7.32	26.05	621	3.69	2.03	clear	no	-183
1404	0.39	1.82	0.13	21.45	7.31	26.09	619	8.24	2.42	clear	no	-182
1407	0.39	2.21	0.13	21.55	7.32	26.03	615	8.70	2.88	clear	no	-176
1410	0.39	2.60	0.13	21.58	7.31	25.99	614	8.48	2.96	clear	no	-175

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: N. SMITH / ARCADIS	SAMPLER(S) SIGNATURES: <i>N. Smith</i>	SAMPLING INITIATED AT: 1412	SAMPLING ENDED AT: 1422
PUMP OR TUBING DEPTH IN WELL (feet): ~161	SAMPLE PUMP FLOW RATE (mL per minute): SL00	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="checkbox"/>	
Filtration Equipment Type: _____			

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-240	3	CG	40ml	HCL	-	<2	82608-SIM	RFP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

SITE NAME: **TALLEVAET** SITE LOCATION: **SARASOTA, FL**
WELL NO: **MW-241** SAMPLE ID: **MW-241** DATE: **1/25/08**

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 281 feet to 301 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: PP								
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (feet - feet) X gallons/foot = gallons												
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 306 feet) + 0.5 gallons = 1.30 gallons												
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 290	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 290	PURGING INITIATED AT: 1125	PURGING ENDED AT: 1150	TOTAL VOLUME PURGED (gallons): 2.50								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1138	1.30	1.30	0.10	8.75	7.55	24.52	767	0.33	17.9	gray	yes	-243
1142	0.40	1.70	0.10	10.24	7.54	24.59	766	0.30	6.58	clear	yes	-246
1146	0.40	2.10	0.10	11.76	7.53	24.56	767	0.30	2.62	clear	yes	-249
1150	0.40	2.50	0.10	13.31	7.56	24.51	764	0.30	1.52	clear	yes	-253
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016												

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: SWINS/ARCADIS				SAMPLER(S) SIGNATURES: [Signature]			SAMPLING INITIATED AT: 1153		SAMPLING ENDED AT: 1157	
PUMP OR TUBING DEPTH IN WELL (feet): 290				SAMPLE PUMP FLOW RATE (mL per minute): ≤ 100			TUBING MATERIAL CODE: PE			
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N				FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm			DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N			
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH				
MW-241	3	CG	40mL	HCl	40mL	—	82608+82685M		RFPP	
REMARKS:										
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)										
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)										

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> MDS 4/15/08	SITE LOCATION: <u>Sarasota FL</u> MDS 4/15/08
WELL NO: <u>MW-221-242</u>	SAMPLE ID: <u>MW-221-242</u> DATE: <u>1/29/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>20</u> feet to <u>100</u> feet	STATIC DEPTH TO WATER (feet): <u>3.38</u>	PURGE PUMP TYPE: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
(only fill out if applicable)

= 0 gallons + (0.0026 gallons/foot X 20 feet) + 0.25 gallons = 0.052 gallons

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>20</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>20</u>	PURGING INITIATED AT: <u>1353</u>	PURGING ENDED AT: <u>1420</u>	TOTAL VOLUME PURGED (gallons): <u>2.3</u>
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<u>1358</u>	<u>0.5</u>	<u>0.5</u>	<u>0.1</u>	<u>4.18</u>	<u>6.69</u>	<u>25.01</u>	<u>614</u>	<u>1.33</u>	<u>6.16</u>	<u>Clear</u>	<u>None</u>
<u>1401</u>	<u>3</u>	<u>3.8</u>	<u>0.1</u>	<u>4.39</u>	<u>6.47</u>	<u>25.08</u>	<u>0.438</u>	<u>2.48</u>	<u>2.63</u>		
<u>1404</u>	<u>3</u>	<u>3.1</u>		<u>4.64</u>	<u>6.13</u>	<u>25.13</u>	<u>473</u>	<u>1.79</u>	<u>4.24</u>		
<u>1407</u>	<u>3</u>	<u>3.1</u>		<u>4.69</u>	<u>6.12</u>	<u>25.15</u>	<u>572</u>	<u>1.53</u>	<u>3.06</u>		
<u>1410</u>	<u>3</u>	<u>1.87</u>		<u>4.70</u>	<u>6.05</u>	<u>25.21</u>	<u>829</u>	<u>1.33</u>	<u>1.46</u>		
<u>1413</u>	<u>3</u>	<u>2.0</u>		<u>4.72</u>	<u>6.05</u>	<u>25.24</u>	<u>874</u>	<u>1.26</u>	<u>0.91</u>		
<u>1416</u>	<u>3</u>	<u>2.3</u>			<u>6.01</u>	<u>25.27</u>	<u>900</u>	<u>1.22</u>	<u>0.81</u>		
→ <u>Sampled @ 1420</u> →											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasey Cornwell / ARCADIS</u>	SAMPLER(S) SIGNATURES: <u>Kasey Cornwell</u>	SAMPLING INITIATED AT: <u>1420</u>	SAMPLING ENDED AT: <u>1425</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>22.5</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u> MDS 4/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u> <u>PE</u>	FIELD-FILTERED: <u>Y</u> <u>PE</u> FILTER SIZE: <u>0.1</u> µm	DUPLICATE: <u>Y</u> <u>N</u>	MDS 4/15/08

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-242</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	<u>RFPP</u>

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 9/15/08	SITE LOCATION: <u>Sarasota</u> MDS 9/15/08
WELL NO: <u>MW-243</u>	SAMPLE ID: <u>MW-243</u> DATE: <u>1/29/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2</u>	TUBING DIAMETER (inches): <u>0.25</u>	WELL SCREEN INTERVAL DEPTH: <u>33</u> feet to <u>38</u> feet	STATIC DEPTH TO WATER (feet): <u>11.83</u>
PURGE PUMP TYPE OR BAILER: <u>peristaltic</u>			
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable			
= (feet - feet) X gallons/foot = gallons			
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME. (only fill out if applicable)			
= <u>0</u> gallons + (<u>0.0026</u> gallons/foot X <u>38</u> feet) + <u>0.25</u> gallons = <u>0.35</u> gallons			
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>36</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>36</u>	PURGING INITIATED AT: <u>1353</u>	PURGING ENDED AT: <u>1422</u>
			TOTAL VOLUME PURGED (gallons): <u>0.9</u>

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	ORP
1359	—	0.6	0.1	12.42	7.3	25.0	0.760	3.0	448	light drusy	light drusy	29
1404	0.5	1.1	0.1	12.42	7.1	24.9	0.770	2.8	108.9	clearing up	slight odor	24
1407	0.3	1.4	0.1	12.42	7.1	24.9	0.770	2.7	42.4	clearing	" "	20
1410	0.3	1.7	0.1	12.42	7.0	24.9	0.780	2.6	33.4	clear	" "	14
1413	0.3	2.0	0.1	12.42	7.0	24.9	0.780	2.6	21.5	" "	" "	12
1416	0.3	2.3	0.1	12.42	7.0	24.8	0.780	2.5	14.8	" "	" "	7
1419	0.3	2.6	0.1	12.42	7.0	24.9	0.780	2.5	10.76	" "	" "	3
1422	0.3	2.9	0.1	12.42	7.0	24.9	0.780	2.4	6.43	" "	" "	
Sampled @ 1422												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Nathan Smith/ Arcadis</u>	SAMPLER(S) SIGNATURES: <u>Nathan S L</u>	SAMPLING INITIATED AT: <u>1422</u>	SAMPLING ENDED AT: <u>1426</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>36</u> MDS 9/15/08	SAMPLE PUMP FLOW RATE (mL per minute): <u>1100</u> MDS 9/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-243	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/15/08	SITE LOCATION: Sarasota FL MDS 4/15/08
WELL NO: MW-244	SAMPLE ID: MW-244 DATE: 1/20/08

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 106 feet to 116 feet	STATIC DEPTH TO WATER (feet): 12.78	PURGE PUMP TYPE OR BAILER: PP MDS 4/15/08
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 116. feet) + 0.25 gallons = 0.55 gallons				

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1437	0.7	0.7	0.1	13.87	6.8	24.7	0.72	3.30	22.0	light grey	—
1440	0.3	1.0	0.1	13.87	6.8	24.8	0.71	2.6	17.8	clear	—
1443	0.3	1.3	0.1	13.87	6.8	24.7	0.71	2.5	4.14	clear	—
1446	0.3	1.6	0.1	13.87	6.8	24.7	0.71	2.4	6.55	" "	—
1449	0.3	1.9	0.1	13.87	6.2	24.7	0.71	2.4	6.58	" "	—

ORR
47
47
43
40
37

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN SMITH	SAMPLER(S) SIGNATURES: Nathan P. Smith	SAMPLING INITIATED AT: 1449	SAMPLING ENDED AT: 1458
PUMP OR TUBING DEPTH IN WELL (feet): 114 MDS 4/15/08	SAMPLE PUMP FLOW RATE (mL per minute): 4100 MDS 4/15/08	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: 20 µm	DUPLICATE: <input type="checkbox"/> Y <input checked="" type="checkbox"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-244	3	CG	40mL	HCl	—	—	8260	RFPP

MDS 4/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/15/08	SITE LOCATION: Sarasota MDS 4/15/08
WELL NO: MW-245	SAMPLE ID: MW-245 DATE: 1/29/08

PURGING DATA			
WELL DIAMETER (inches):	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 17 feet to 22 feet	STATIC DEPTH TO WATER (feet): 6.15
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)		PURGE PUMP TYPE OR BAILER: PP MDS 4/15/08	
= (20 feet -		feet) X gallons/foot = gallons	
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)			
= 0 gallons + (0.0026 gallons/foot X 20 feet) + 0.25 gallons = 0.3 gallons			

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 9	PURGING INITIATED AT: 9:15	PURGING ENDED AT: 9:33	TOTAL VOLUME PURGED (gallons): 2.0							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0924	0.25	0.75	0.08	6.60	6.88	23.76	0.95	1.12	0.98	Clear	—
0927	0.08	1.0	0.08	6.60	6.88	23.92	0.94	1.04	1.01	Clear	—
0930	0.25	1.25	0.08	6.60	6.88	23.92	0.94	1.04	0.68	Clear	—
0933	0.25	2.0	0.08	6.60	6.88	24.04	0.94	1.02	0.67	Clear	—
Sampled @ 0933											

-267
-272
-271

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA	
SAMPLED BY (PRINT) / AFFILIATION: Nathan Smith	SAMPLER(S) SIGNATURES: <i>Nathan Smith</i>
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLING INITIATED AT: 0933 SAMPLING ENDED AT: 0940
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: 0.45µm MDS 4/15/08
TUBING MATERIAL CODE: _____	
DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-245	3	CG	40ml	HCl	—	—	8260	RFPP

MDS 4/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump

EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MOS 4/15/08	SITE LOCATION: <u>Sarasota</u> MOS 4/15/08
WELL NO: <u>MW-246</u>	SAMPLE ID: <u>MW-246</u> DATE: <u>1/29/08</u>

PURGING DATA

WELL DIAMETER (inches): <u>2"</u>	TUBING DIAMETER (inches): <u>1/4"</u>	WELL SCREEN INTERVAL DEPTH: <u>40</u> feet	STATIC DEPTH TO WATER (feet): <u>15.44</u>	PURGE PUMP TYPE OR BAILER: <u>PP</u>
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= <u>0</u> gallons + (0.0026 gallons/foot X <u>45'</u> feet) + <u>0.25</u> gallons = <u>0.37</u> gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): <u>42.5</u>	FINAL PUMP OR TUBING DEPTH IN WELL (feet): <u>42.5</u>	PURGING INITIATED AT: <u>0833</u>	PURGING ENDED AT: <u>0900</u>	TOTAL VOLUME PURGED (gallons): <u>2.0</u>								
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	OPP 97 47 17 7 0 0
0838	0.5	0.5	0.1	16.01	6.5	25.2	0.11	3.0	2.17	clear	none	
0841	0.3	0.8	0.1	17.31	6.6	25.4	0.11	2-8	1.11			
0844	0.3	1.1	0.1	20.45	6.6	25.5	0.11	4.3	7.89			
0849	0.3	1.4	0.1	21.58	6.6	25.4	0.11	4.3	0.22			
0852	0.3	1.7	0.1	22.46	6.7	25.5	0.10	4.1	0.50			
0855	0.3	2.0	0.1	25.15	6.7	25.5	0.10	3.9	0.77	↓	↓	
0900	→ Sampled @ 0900 ←											

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasey Cornwell (KLC) / Armas</u>	SAMPLER(S) SIGNATURES: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>0900</u>	SAMPLING ENDED AT: <u>0905</u>
PUMP OR TUBING DEPTH IN WELL (feet): <u>42.5</u>	SAMPLE PUMP FLOW RATE (mL per minute): <u>1100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N	FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>MW-246</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>

REMARKS: Initial purge clear no noticeable odor

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast MDS 4/15/08	SITE LOCATION: Sarasota MDS 4/15/08
WELL NO: MW-247	SAMPLE ID: MW-247 DATE: 1/29/08

PURGING DATA

WELL DIAMETER (inches): 2.5	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 98 feet to 100 feet	STATIC DEPTH TO WATER (feet): 13.18	PURGE PUMP TYPE OR BAILER: Peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (108 feet - 13.18 feet) X 0.75 gallons/foot = 74.82 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 103 feet) + 0.25 gallons = 0.5 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 108		FINAL PUMP OR TUBING DEPTH IN WELL (feet): 103		PURGING INITIATED AT: 8:25		PURGING ENDED AT: 8:47		TOTAL VOLUME PURGED (gallons): 28				
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0838	-	1.3	350.0	14.54	7.24	24.98	0.97	1.06	3.08	clear	-	-329
0841	0.5	1.8	0.1	14.54	7.27	24.82	0.96	1.06	1.28	clear	-	-337
0844	0.5	2.3	0.1	14.54	7.27	25.13	0.96	1.05	1.02	clear	-	-338
0847	0.5	2.8	0.1	14.54	7.27	24.99	0.96	1.04	1.52	clear	-	-340

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN SMITH		SAMPLER(S) SIGNATURES: <i>Nathan P. Smith</i>		SAMPLING INITIATED AT: 08:50	SAMPLING ENDED AT: 08:59
PUMP OR TUBING DEPTH IN WELL (feet): 103 MDS 4/15/08		SAMPLE PUMP FLOW RATE (mL per minute): 1100 MDS 4/15/08		TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N		FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N FILTER SIZE: 1 µm		DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N → DUP-18	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-247	3	CG	40ml	HCl	-	-	8260	RFPP
DUP-18	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-248	SAMPLE ID: MW-248
DATE: 1.28.08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 103 feet to 113 feet	STATIC DEPTH TO WATER (feet): 13.9	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)				
= (103 feet - 13.9 feet) X 0.026 gallons/foot = 2.5 gallons				

EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 113 feet) + 0.25 gallons = 0.54 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 108	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 108	PURGING INITIATED AT: 11:55	PURGING ENDED AT: 12:20	TOTAL VOLUME PURGED (gallons): 2.5
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
12:00	0.5	0.5	0.1	13.9	7.21	23.37	0.584	0.89	9	clear	none
12:05	0.5	1.0	0.1	13.9	7.12	23.4	0.596	0.36	3	"	"
12:10	0.5	1.5	0.1	13.9	7.08	23.4	0.582	0.34	2.63	"	"
12:15	0.5	2.0	0.1	14.8	6.98	23.3	0.578	0.28	0	"	"
12:20	0.5	2.5	0.1	14.8	6.85	23.4	0.574	0.28	0	"	"

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Nican Forbes / ARCADIS	SAMPLER(S) SIGNATURES: <i>Nican Forbes</i>	SAMPLING INITIATED AT: 12:21	SAMPLING ENDED AT: 12:25
PUMP OR TUBING DEPTH IN WELL (feet): 108	SAMPLE PUMP FLOW RATE (mL per minute): 400	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: Y	FIELD-FILTERED: Y	DUPLICATE: Y	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-248	3	CG	40ml	HCl	—	—	8260	RFPP
Dup-21	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **collect Dup-21 here**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC		SITE LOCATION: Tallevast	
WELL NO: MW-249	SAMPLE ID: MW-249	DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 4	WELL SCREEN INTERVAL DEPTH: 88 feet to 98 feet	STATIC DEPTH TO WATER (feet): 11.10	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (.0026 gallons/foot X 101 feet) + .25 gallons = gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 98	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 93	PURGING INITIATED AT: 13:55	PURGING ENDED AT: 14:22	TOTAL VOLUME PURGED (gallons): 2.6

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
14:03	.8	.8	0.1	11.36	6.98	25.18	.464	2.59	.24	clear	none
14:07	.3	1.1	0.1	11.36	6.95	24.74	.470	1.67	.94	clear	none
14:10	.3	1.4	0.1	11.38	6.94	25.18	.466	1.46	1.83	clear	none
14:13	.3	1.7	0.1	11.38	6.94	24.82	.471	1.41	1.76	clear	none
14:16	.3	2.0	0.1	11.38	6.95	25.11	.468	1.33	1.56	clear	none
14:19	.3	2.3	0.1	11.38	6.97	25.12	.468	1.29	1.78	clear	none
14:22	.3	2.6	0.1	11.38	6.98	25.12	.469	1.28	1.83	clear	none

MDS 4/15/08
19
14
12
10
9
8

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Darin Johnson	SAMPLER(S) SIGNATURES: <i>Darin Johnson</i>	SAMPLING INITIATED AT: 14:23	SAMPLING ENDED AT: 14:28
PUMP OR TUBING DEPTH IN WELL (feet): 93	SAMPLE PUMP FLOW RATE (mL per minute): 4100	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>	FIELD-FILTERED: Y <input type="checkbox"/> N <input checked="" type="checkbox"/> FILTER SIZE: 0.45 µm	DUPLICATE: Y <input type="checkbox"/> N <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-249	3	CG	90ml	HCl	-	-	8260	RFPP

MDS 4/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

2

GROUNDWATER SAMPLING LOG

SITE NAME: Former ABC Facility	SITE LOCATION: Tallahassee, FL
WELL NO: MW-250	SAMPLE ID: MW-250 DATE: 1-31-08

PURGING DATA

WELL DIAMETER (inches): 2"	TUBING DIAMETER (inches): 1/4"	WELL SCREEN INTERVAL DEPTH: 90 feet to 100 feet	STATIC DEPTH TO WATER (feet): 1301	PURGE PUMP TYPE OR BAILER: PP
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 100 feet) + 0.25 gallons = 0.51 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: 1342
				PURGING ENDED AT: 1404
				TOTAL VOLUME PURGED (gallons): 2.1

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1352	1.0	1.0	0.1	13.00	7.38	24.86	0.045	1.69	0.06	clear	none noted
1355	0.3	1.3	0.1	13.00	7.39	24.85	0.049	1.47	0.29		
1358	0.3	1.6	0.1	13.01	7.39	24.87	0.053	1.22	0.37		
							0.053 S				
1401	0.3	1.9	0.1	13.00	7.41	24.81	0.054	1.21	0.29		
1404	0.3	2.1	0.1	13.00	7.45	24.80	0.060	1.20	0.25		

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: AMY COOK / Arcadis	SAMPLER(S) SIGNATURES:	SAMPLING INITIATED AT: 1405	SAMPLING ENDED AT: 1410
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <100	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N	FILTER SIZE: 0.45µm	DUPLICATE: <input checked="" type="checkbox"/> Y <input type="checkbox"/> N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-250	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **pumped some before reaching a state. Took dtw this morning**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

- NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevoast	SITE LOCATION: Sarasota FL
WELL NO: MW-251	SAMPLE ID: MW-251
DATE: 1/20/08	

PURGING DATA

WELL DIAMETER (inches): 2.5	TUBING DIAMETER (inches): 0.25	WELL SCREEN INTERVAL DEPTH: 380 feet to 400 feet	STATIC DEPTH TO WATER (feet): 13.20	PURGE PUMP TYPE OR BAILER: Peristaltic
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (4 feet - 13.20 feet) X 0.0026 gallons/foot = 0.026 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (400 gallons/foot X 390 feet) + 0.25 gallons = 1.3 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 390	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 390	PURGING INITIATED AT: 09:30	PURGING ENDED AT: 09:57	PURGED (gallons): 2.7

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)	
0945	0.5	1.5	0.1	13.22	7.33	26.59	1.26	1.17	3.69	clear	slight odor	-297
0948	1.8	0.3	0.1	13.22	7.33	26.57	1.25	1.15	3.21	clear	" "	-296
0951	2.1	0.3	0.1	13.22	7.34	26.56	1.26	1.14	2.64	clear	" "	-296
0954	2.4	0.3	0.1	13.22	7.34	26.58	1.26	1.12	1.55	clear	" "	-295
0957	2.7	0.3	0.1	13.22	7.34	26.62	1.25	1.11	1.51	clear	" "	-296
Sampled @ 0957												

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: NATHAN SMITH		SAMPLER(S) SIGNATURES: <i>Nathan P. Smith</i>		SAMPLING INITIATED AT: 0957	SAMPLING ENDED AT: 1005			
PUMP OR TUBING DEPTH IN WELL (feet): 390		SAMPLE PUMP FLOW RATE (mL per minute): 4100		TUBING MATERIAL CODE:				
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> Y		FIELD-FILTERED: <input checked="" type="checkbox"/> Y		DUPLICATE: <input checked="" type="checkbox"/> Y				
SAMPLE CONTAINER SPECIFICATION		SAMPLE PRESERVATION		INTENDED ANALYSIS AND/OR METHOD				
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
MW-251	3	CG	40ml	HCL	—	—	8260	RFPP
REMARKS:								
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)								
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)								

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

DEP-SOP-001/01
Form FD 9000-24
GROUNDWATER SAMPLING LOG

P. 98

SITE NAME: Tallevast		SITE LOCATION: Tallevast, FL	
WELL NO: MW-252	SAMPLE ID: MW-252	DATE: 1-31-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 145 feet to 155 feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER: Geopump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.024 gallons/foot X 158 feet) + .25 gallons = .66 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 150	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 150	PURGING INITIATED AT: 1530	PURGING ENDED AT: 1605	TOTAL VOLUME PURGED (gallons): 2.45

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1559	2.03	2.03	0.07	21.16	7.50	24.7	.505	.66	23.9	clear	yes
1601	.14	2.17	0.07	21.19	7.49	24.7	.503	.64	23.7	clear	yes
1603	.14	2.31	0.07	21.20	7.49	24.6	.503	.63	22.3	clear	yes
1605	.14	2.45	0.07	21.20	7.47	24.6	.504	.61	21.4	clear	yes

ORP
-248
-249
-250
-251

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIDGET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Bridget Stahl</i>	SAMPLING INITIATED AT: 1607	SAMPLING ENDED AT: 1617
PUMP OR TUBING DEPTH IN WELL (feet): 150	SAMPLE PUMP FLOW RATE (mL per minute): ~280	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <input checked="" type="radio"/> Y <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y <input type="radio"/> N FILTER SIZE: _____ µm	DUPLICATE: <input type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-252	3	CG	40ml	HCl	—	—	8260	RFPP

REMARKS: **placeel 158' teflon tubing in mw**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

153

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST	
WELL NO: mw-253	SAMPLE ID: mw-253	DATE: F31 08 21-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 100 feet to 110 feet	STATIC DEPTH TO WATER (feet): 15.85	PURGE PUMP TYPE OR BAILER: PP Gas Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0026 gallons/foot X 113 feet) + .25 gallons = .54 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 105	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 105	PURGING INITIATED AT: 0903	PURGING ENDED AT: 0923	TOTAL VOLUME PURGED (gallons): 2.40
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TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or m/cm)	DISSOLVED OXYGEN (circled mg/l or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0917	1.68	1.68	.12	17.00	7.0	24.0	.967	0.9	4.97	clear	yo
0919	0.24	1.92	.12	17.00	7.0	24.0	.973	0.9	4.09	clear	yo
0921	0.24	2.16	.12	17.0	7.0	24.0	.977	0.9	4.64	clear	yo
0923	0.24	2.40	.12	17.0	7.0	24.1	.977	0.8	2.54	clear	yo

ORP
-201
-206
-200
-214

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King/Alcorbis	SAMPLER(S) SIGNATURES: Jennifer King	SAMPLING INITIATED AT: 0924	SAMPLING ENDED AT: 0930
PUMP OR TUBING DEPTH IN WELL (feet): 105	SAMPLE PUMP FLOW RATE (mL per minute): 1.440	TUBING MATERIAL CODE: T	
FIELD DECONTAMINATION: <input type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y	FILTER SIZE: _____ µm	DUPLICATE: <input checked="" type="radio"/> Y

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-253	3	CG	40ml	Hel	-	-	P260	RPPP

REMARKS: **ADDED 113' teflon tubing; missing two bolts**

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RPPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

99

SITE NAME: TALLEVAST		SITE LOCATION: TALLEVAST, FL	
WELL NO: MW-254	SAMPLE ID: MW-254	DATE: 12/08/2010	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 24 feet to 29 feet	STATIC DEPTH TO WATER (feet): 498	PURGE PUMP TYPE OR BAILER: PP Gas Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= 0 gallons + (0.0024 gallons/foot X 32 feet) + 0.25 gallons = 0.33 gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 26.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 26.5	PURGING INITIATED AT: 823	PURGING ENDED AT: 0835	TOTAL VOLUME PURGED (gallons): 2.16

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mS/cm)	DISSOLVED OXYGEN (circle % or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
6 0829	1.08	1.08	0.18	6.30	5.3	29.0	812	0.8	4.00	Clear	NO
0831	0.36	1.44	0.18	6.34	5.2	28.7	871	0.7	7.19	Clear	NO
0833	0.36	1.80	0.18	6.35	5.2	28.8	898	0.6	6.21	Clear	NO
0835	0.36	2.16	0.18	6.35	5.2	28.8	903	0.6	7.95	Clear	NO

ORP
17
7
-2
-5

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer Key / Arcadis	SAMPLER(S) SIGNATURES: Jennifer Key	SAMPLING INITIATED AT: 0836	SAMPLING ENDED AT: 0840
PUMP OR TUBING DEPTH IN WELL (feet): 26.5	SAMPLE PUMP FLOW RATE (mL per minute): 700	TUBING MATERIAL CODE: 7	
FIELD DECONTAMINATION: Y N	FIELD-FILTERED: Y N	FILTER SIZE: _____ µm	DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
MW-254	3	CG	4ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Talleast</u> ^{MDS} <u>4/15/08</u>	SITE LOCATION: <u>Sarasota FL</u> ^{MDS} <u>4/15/08</u>
WELL NO: <u>PW-7</u>	SAMPLE ID: <u>PW-7</u> DATE: <u>2/1/08</u>

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable)											
= (feet - feet) X gallons/foot = gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)											
= gallons + (gallons/foot X feet) + gallons = gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>849</u>	PURGING ENDED AT: <u>854</u>	TOTAL VOLUME PURGED (gallons):							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88											
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>ARLADIS</u>		SAMPLER(S) SIGNATURES: ^{MDS} <u>4/15/08</u>		SAMPLING INITIATED AT: <u>854</u>	SAMPLING ENDED AT: <u>857</u>			
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u>100</u>	TUBING MATERIAL CODE:						
FIELD DECONTAMINATION: <input checked="" type="checkbox"/> N	FIELD-FILTERED: <input checked="" type="checkbox"/> Y	FILTER SIZE: <u>0.45</u> µm	DUPLICATE: <input checked="" type="checkbox"/> Y					
SAMPLE CONTAINER SPECIFICATION		SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE		
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED			TOTAL VOL ADDED IN FIELD (mL)	FINAL pH
<u>PW-7</u>	<u>3</u>	<u>LG</u>	<u>40ml</u>	<u>HCL</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>RFPP</u>
REMARKS:								
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)								
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)								

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast **MDS 4/15/08** SITE LOCATION: Sarasota FL **MDS 4/15/08**
 WELL NO: PW-38 SAMPLE ID: PW-38 DATE: 2/1/08

PURGING DATA

WELL DIAMETER (inches): 1.75 TUBING DIAMETER (inches): 1/4 WELL SCREEN INTERVAL DEPTH: NA feet to NA feet STATIC DEPTH TO WATER (feet): 13.85 OR BAILER: PP
 WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY
 = (140 feet - 13.85 feet) X 0.02 gallons/foot = 2.8 gallons
 EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME
 = 0 gallons + (0.0006 gallons/foot X 140 feet) + .25 gallons = .61 gallons

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
1040	2.8	2.8	0.1	13.85	7.02	23.6	0.636	1.8	25.9	Cloudy none	-251
1042	0.2	3.0			7.02	23.6	0.638	1.4	17.7		-252
1044	0.2	3.2			7.01	23.5	0.637	1.4	15.5		-254
1046	0.2	3.4			7.02	23.6	0.640	1.4	24.6		-258
1048	0.2	3.6			7.02	23.5	0.635	1.3	15.7	Clear	-259
1050	0.2	3.8			7.02	23.6	0.635	1.3	17.2		-262
1052	0.2	4.0			7.02	23.6	0.632	1.3	12.4		
1055	→ Sampled			@ 1055							

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: KLC/NPS/ARCADIS SAMPLER(S) SIGNATURES: [Signature] SAMPLING INITIATED AT: 1055 SAMPLING ENDED AT:
 PUMP OR TUBING DEPTH IN WELL (feet): SAMPLE PUMP FLOW RATE (mL per minute): < 100 TUBING MATERIAL CODE:
 FIELD DECONTAMINATION: N FIELD-FILTERED: Y FILTER SIZE: 5 µm DUPLICATE: Y
 Filtration Equipment Type: N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
PW-28	3	CG	90ml	HCl	-	-	8260	

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Private Well # 47

DEP-SOP-001/01
Form FD 9000-24

GROUNDWATER SAMPLING LOG

SITE NAME: Talleast MDS 4/15/08	SITE LOCATION: Sarasota FL 4/15/08
WELL NO: HW-47	SAMPLE ID: HW-47 DATE: 4/29/1/30/08

* Southern most on figure.

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable											
= (feet - feet) X gallons/foot = gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)											
= gallons + (gallons/foot X feet) + gallons = gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 0855	PURGING ENDED AT: 0900	TOTAL VOLUME PURGED (gallons):							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
Table content is crossed out with a large X.											
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88											
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: WATSON/ARCADIS			SAMPLER(S) SIGNATURES:			SAMPLING INITIATED AT: 0900		SAMPLING ENDED AT: 0903			
PUMP OR TUBING DEPTH IN WELL (feet):			SAMPLE PUMP FLOW RATE (mL per minute) 100 MDS 4/15/08			TUBING MATERIAL CODE:					
FIELD DECONTAMINATION: N MDS 4/15/08			FIELD-FILTERED: Y MDS 4/15/08			FILTRATION EQUIPMENT TYPE: µm		DUPLICATE: Y N MDS 4/15/08			
SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	8260				
PW-47	3	CG	90ml	HCl	-	-					
REMARKS:											
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)											
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)											

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* Spigot/well/sprinkler located on the driving range side of the netting, near driving range monitoring wells.

Private Well ~~48~~ 47 KT 4/29/08

DEP-SOP-001/01
Form FD 9000-24

GROUNDWATER SAMPLING LOG

MDS
4/15/08

SITE NAME: Talleast MDS 4/15/08
 WELL NO: 47 KT PRIVATE WELL SAMPLE ID: 48
 SITE LOCATION: Sarasota FL DATE: 1/30/08

* may not need to sample

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable											
= (feet - feet) X gallons/foot = gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)											
= gallons + (gallons/foot X feet) + gallons = gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>0850</u>	PURGING ENDED AT: <u>0855</u>	TOTAL VOLUME PURGED (gallons):							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
Table content is crossed out with a large X											
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88											
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>Kasei Comwell / ARCADIS</u>			SAMPLER(S) SIGNATURES: <u>[Signature]</u>			SAMPLING INITIATED AT: <u>0855</u>		SAMPLING ENDED AT: <u>0858</u>	
PUMP OR TUBING DEPTH IN WELL (feet):			SAMPLE PUMP FLOW RATE (mL per minute): <u>400</u> MDS 4/15/08			TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <u>N</u> MDS 4/15/08			FIELD-FILTERED: <u>Y</u> MDS 4/15/08			DUPLICATE: <u>Y</u>		<u>N</u> MDS 4/15/08	
SAMPLE CONTAINER SPECIFICATION					SAMPLE PRESERVATION				
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE	
<u>PW-48</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>-</u>	<u>-</u>	<u>8260</u>	<u>0 - sampled at spigot</u>	
REMARKS:									

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
 SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

* Spigot / Just east of clubhouse
 North of PW-47. near sidewalk / parking lot.
 Revision Date: February 1, 2004

GROUNDWATER SAMPLING LOG

SITE NAME: Tallevast	MDS 4/15/08	SITE LOCATION: Sarasota FL	MDS 4/15/08
WELL NO: PW-57	SAMPLE ID: PW-57	DATE: 2/1/08	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: — feet to — feet	STATIC DEPTH TO WATER (feet): —	PURGE PUMP TYPE OR BAILER: —							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)											
= (— feet - — feet) X — gallons/foot = — gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)											
= — gallons + (— gallons/foot X — feet) + — gallons = — gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: 815	PURGING ENDED AT: 820	TOTAL VOLUME PURGED (gallons):							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Kasey Cornwall / Arcadis			SAMPLER(S) SIGNATURES: <i>[Signature]</i>			SAMPLING INITIATED AT: 820	SAMPLING ENDED AT: 823
PUMP OR TUBING DEPTH IN WELL (feet):			SAMPLE PUMP FLOW RATE (mL per minute): 400			TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: N			FIELD-FILTERED: Y			DUPLICATE: Y	
SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	SAMPLING EQUIPMENT CODE
PW-57	3	CG	40ml	HCL	—	—	8260
							MDS 4/15/08
REMARKS:							

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u>	MDS <u>9/15/08</u>	SITE LOCATION: <u>Sarasota FL</u>	MDS <u>9/15/08</u>
WELL NO: <u>PW-59</u>	SAMPLE ID: <u>PW-59</u>	DATE: <u>2/1/08</u>	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)											
= (feet - feet) X gallons/foot = gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)											
= gallons + (gallons/foot X feet) + gallons = gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT: <u>900</u>	PURGING ENDED AT: <u>905</u>	TOTAL VOLUME PURGED (gallons):							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>KLC/UPS / ARCADIS</u>			SAMPLER(S) SIGNATURES: <u>[Signature]</u>			SAMPLING INITIATED AT: <u>905</u>	SAMPLING ENDED AT: <u>908</u>
PUMP OR TUBING DEPTH IN WELL (feet):			SAMPLE PUMP FLOW RATE (mL per minute): <u>4100</u>			TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>Y</u>			FIELD-FILTERED: <u>Y</u>			DUPLICATE: <u>Y</u>	
SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	SAMPLING EQUIPMENT CODE	
<u>PW-59</u>	<u>3</u>	<u>LG</u>	<u>HCL</u>	<u>-</u>	<u>-</u>	<u>8260</u>	
REMARKS:							
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)							
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)							

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Former ABC</i>		SITE LOCATION: <i>Tallevast</i>	
WELL NO: <i>Private Well 64</i>	SAMPLE ID: PW-64 <i>PW-64</i>	DATE: <i>2-4-08</i>	

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable = (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons				
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT:	PURGING ENDED AT:	TOTAL VOLUME PURGED (gallons):

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Darrin Johnson</i>	SAMPLER(S) SIGNATURE(S): <i>[Signature]</i>	SAMPLING INITIATED AT: <i>9:55</i>	SAMPLING ENDED AT: <i>9:55</i>				
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <i>4100</i>	TUBING MATERIAL CODE:					
FIELD DECONTAMINATION: <i>N</i>	FIELD-FILTERED: <i>Y</i>	FILTRATION EQUIPMENT TYPE: <i>N</i>	DUPLICATE: <i>Y</i>				
SAMPLE CONTAINER SPECIFICATION: <i>N</i>	SAMPLE PRESERVATION:		INTENDED ANALYSIS AND/OR METHOD: <i>8260</i>				
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH	SAMPLING EQUIPMENT CODE
<i>PW-64</i>	<i>3</i>	<i>CG</i>	<i>40ml</i>	<i>HCl</i>	<i>—</i>	<i>—</i>	<i>[Signature]</i>

REMARKS: *This sample was taken from sprinkler system at 1107 Tallevast Rd*

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
 EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: <i>Arcadis</i>		SITE LOCATION: <i>Talleast</i>	
WELL NO: <i>Schmit private well</i>		SAMPLE ID: <i>84</i>	DATE: <i>1-24-08</i>

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY (only fill out if applicable) = (feet - feet) X gallons/foot = gallons											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = gallons + (gallons/foot X feet) + gallons = gallons											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT:	PURGING ENDED AT:	TOTAL VOLUME PURGED (gallons):					
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
<i>none taken private well!</i>											
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <i>Darrin Johnson</i>			SAMPLER(S) SIGNATURES: <i>Darrin Johnson</i>				SAMPLING INITIATED AT: <i>15:40</i>		SAMPLING ENDED AT: <i>15:40</i>	
PUMP OR TUBING DEPTH IN WELL (feet):			SAMPLE PUMP FLOW RATE (mL per minute): <i>4100</i> (MDS 8-715/08)				TUBING MATERIAL CODE:			
FIELD DECONTAMINATION: <i>Y</i>			FIELD-FILTERED: <i>Y</i> (MDS 8-715/08)				DUPLICATE: <i>Y</i> N			
SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD		SAMPLING EQUIPMENT CODE	
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH				
<i>PIW-84</i>	<i>3</i>	<i>LG</i>	<i>40ml</i>	<i>HCl</i>	<i>—</i>	<i>—</i>	<i>8260</i>			
<i>DUP-7</i>	<i>3</i>	<i>LG</i>	<i>40ml</i>	<i>HCl</i>	<i>—</i>	<i>—</i>	<i>8260</i>			
<i>MDS 8-715/08</i>										
REMARKS: <i>DUP-7-012408</i>										
MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)										
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)										

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: FUNNY ABC SITE LOCATION: Talbotsh PL
 WELL NO: PW-85 MDS 9/15/08 SAMPLE ID: Private well 85 DATE: 1-31-08

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:							
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable)											
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)											
INITIAL PUMP OR TUBING DEPTH IN WELL (feet):	FINAL PUMP OR TUBING DEPTH IN WELL (feet):	PURGING INITIATED AT:	PURGING ENDED AT:	TOTAL VOLUME PURGED (gallons):							
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88 TUBING INSIDE DIA. CAPACITY (Gal/FL): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016											

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Amy Coats SAMPLER(S) SIGNATURES: [Signature] SAMPLING INITIATED AT: 15:00 SAMPLING ENDED AT: 15:05

PUMP OR TUBING DEPTH IN WELL (feet): _____ SAMPLE PUMP FLOW RATE (mL per minute): 400 MDS 9/15/08 TUBING MATERIAL CODE: spigot

FIELD DECONTAMINATION: Y N FIELD-FILTERED: Y N FILTER SIZE: _____ µm DUPLICATE: Y N

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>PW-85</u>	<u>3</u>	<u>CG</u>	<u>40ml</u>	<u>HCl</u>	<u>—</u>	<u>—</u>	<u>8260</u>	

REMARKS: spigot pulls from 80 gallon tank. Purged ~4.5 gallons then sampled. 4.5 gall is more than enough to purge line. Cannot purge tank

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFPF = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

Private well Desenberg

GROUNDWATER SAMPLING LOG

SITE NAME: <u>Tallevast</u> MDS 9/15/08	SITE LOCATION: <u>Sarasota FL</u> MDS 9/15/08
WELL NO: <u>PW-102</u>	SAMPLE ID: <u>PW-102</u> DATE: <u>2/1/08</u>

PURGING DATA

WELL DIAMETER (inches):	TUBING DIAMETER (inches):	WELL SCREEN INTERVAL DEPTH: feet to feet	STATIC DEPTH TO WATER (feet):	PURGE PUMP TYPE OR BAILER:
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable				
= (feet - feet) X gallons/foot = gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable)				
= gallons + (gallons/foot X feet) + gallons = gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet):		FINAL PUMP OR TUBING DEPTH IN WELL (feet):		PURGING INITIATED AT: <u>1110</u>	PURGING ENDED AT: <u>1115</u>	TOTAL VOLUME PURGED (gallons):					
TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or µS/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: <u>KASEL / ARCADIS</u>	SAMPLER(S) SIGNATURE: <u>[Signature]</u>	SAMPLING INITIATED AT: <u>11215</u>	SAMPLING ENDED AT: <u>1118</u>
PUMP OR TUBING DEPTH IN WELL (feet):	SAMPLE PUMP FLOW RATE (mL per minute): <u>2100</u>	TUBING MATERIAL CODE:	
FIELD DECONTAMINATION: <u>N</u>	FIELD-FILTERED: <u>N</u> FILTER SIZE: <u>5µm</u>	DUPLICATE: <u>Y</u>	MDS 9/15/08

SAMPLE CONTAINER SPECIFICATION			SAMPLE PRESERVATION				INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
<u>PW-102</u>	<u>3</u>	<u>CG</u>	<u>90ml</u>	<u>HCL</u>	<u>-</u>	<u>-</u>	<u>8260</u>	MDS 9/15/08

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
EQUIPMENT CODES: RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: TALLEVAST
WELL NO: DW-84B	SAMPLE ID: MW-84B
DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 1.5	TUBING DIAMETER (inches): 1.5	WELL SCREEN INTERVAL DEPTH: 7 feet to 12 feet	STATIC DEPTH TO WATER (feet): 6.56	PURGE PUMP TYPE OR BAILER: PP Geo Pump
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY = (12 feet - 6.56 feet) X .14 gallons/foot = .87 gallons				
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME = 0 gallons + (0.0026 gallons/foot X 30 feet) + .25 gallons = 0.33 gallons				

INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 9.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 9.5	PURGING INITIATED AT: 0932	PURGING ENDED AT: 0946	TOTAL VOLUME PURGED (gallons): 1.68
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8
4
2
2

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or mg/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0940	0.10	0.10	.12	6.75	5.6	22.7	.113	.5	12.7	Clear	ND
0942	0.24	1.2	.12	6.74	5.6	22.7	.115	.5	12.0	Clear	N
0944	0.24	1.44	.12	6.74	5.6	22.9	.115	.5	9.27	Clear	N
0946	0.24	1.68	.12	6.76	5.6	22.9	.115	.6	8.05	Clear	N

ORP
44
34
25
20

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
 TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: Jennifer King / ARCADIS	SAMPLER(S) SIGNATURES: <i>Jennifer King</i>	SAMPLING INITIATED AT: 0947	SAMPLING ENDED AT: 0952
PUMP OR TUBING DEPTH IN WELL (feet): 9.5	SAMPLE PUMP FLOW RATE (mL per minute): ~450	TUBING MATERIAL CODE: PE <small>MS 4/11/02</small>	
FIELD DECONTAMINATION: 0 N	FIELD-FILTERED: Y <input checked="" type="checkbox"/> FILTER SIZE: _____ µm	DUPLICATE: Y <input checked="" type="checkbox"/>	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-84B	3	CG	40ml	HCl	-	-	8260	RFP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)

SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailer; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump; RFP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.
 2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)
 pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: ± 5% Dissolved Oxygen: all readings ≤ 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

GROUNDWATER SAMPLING LOG

SITE NAME: TALLEVAST	SITE LOCATION: TALLEVAST, FL
WELL NO: AW-84A	SAMPLE ID: mw-84A
DATE: 1-29-08	

PURGING DATA

WELL DIAMETER (inches): 2	TUBING DIAMETER (inches): 1/4	WELL SCREEN INTERVAL DEPTH: 7 feet to 12 feet	STATIC DEPTH TO WATER (feet): 6.49
PURGE PUMP TYPE OR BAILER: PP Geo Pump			
WELL VOLUME PURGE: 1 WELL VOLUME = (TOTAL WELL DEPTH - STATIC DEPTH TO WATER) X WELL CAPACITY only fill out if applicable) = (12 feet - 6.49 feet) X 0.16 gallons/foot = 0.88 gallons			
EQUIPMENT VOLUME PURGE: 1 EQUIPMENT VOL. = PUMP VOLUME + (TUBING CAPACITY X TUBING LENGTH) + FLOW CELL VOLUME (only fill out if applicable) = 0 gallons + (0.0026 gallons/foot X 530 feet) + .25 gallons = 1.33 gallons			
INITIAL PUMP OR TUBING DEPTH IN WELL (feet): 9.5	FINAL PUMP OR TUBING DEPTH IN WELL (feet): 9.5	PURGING INITIATED AT: 0940	PURGING ENDED AT: 0954
TOTAL VOLUME PURGED (gallons): 1.68			

TIME	VOLUME PURGED (gallons)	CUMUL. VOLUME PURGED (gallons)	PURGE RATE (gpm)	DEPTH TO WATER (feet)	pH (standard units)	TEMP. (°C)	COND. (µmhos/cm or M/S/cm)	DISSOLVED OXYGEN (circle mg/L or % saturation)	TURBIDITY (NTUs)	COLOR (describe)	ODOR (describe)
0948	0.96	0.96	0.12	6.52	6.04	22.3	1.37	0.52	12.6	clear	yes -180
0950	0.24	1.20	0.12	6.52	6.02	22.5	1.38	0.50	11.0	clear	yes -180
0952	0.24	1.44	0.12	6.52	6.01	22.5	1.38	0.51	11.6	clear	yes -180
0954	0.24	1.68	0.12	6.52	6.00	22.5	1.38	0.49	12.1	clear	yes -181

WELL CAPACITY (Gallons Per Foot): 0.75" = 0.02; 1" = 0.04; 1.25" = 0.06; 2" = 0.16; 3" = 0.37; 4" = 0.65; 5" = 1.02; 6" = 1.47; 12" = 5.88
TUBING INSIDE DIA. CAPACITY (Gal./Ft.): 1/8" = 0.0006; 3/16" = 0.0014; 1/4" = 0.0026; 5/16" = 0.004; 3/8" = 0.006; 1/2" = 0.010; 5/8" = 0.016

SAMPLING DATA

SAMPLED BY (PRINT) / AFFILIATION: BRIQUET STAHL / ARCADIS	SAMPLER(S) SIGNATURES: <i>Briquet Stahl</i>	SAMPLING INITIATED AT: 0955	SAMPLING ENDED AT: 0957
PUMP OR TUBING DEPTH IN WELL (feet): 9.5	SAMPLE PUMP FLOW RATE (mL per minute): 460	TUBING MATERIAL CODE: PE	
FIELD DECONTAMINATION: <input checked="" type="radio"/> N	FIELD-FILTERED: <input checked="" type="radio"/> Y Filtration Equipment Type: <input checked="" type="radio"/>	FILTER SIZE: _____ µm	
SAMPLE CONTAINER SPECIFICATION		DUPLICATE: <input checked="" type="radio"/> Y <input checked="" type="radio"/> N	

SAMPLE CONTAINER SPECIFICATION				SAMPLE PRESERVATION			INTENDED ANALYSIS AND/OR METHOD	SAMPLING EQUIPMENT CODE
SAMPLE ID CODE	# CONTAINERS	MATERIAL CODE	VOLUME	PRESERVATIVE USED	TOTAL VOL ADDED IN FIELD (mL)	FINAL pH		
mw-84A	3	CG	40ml	HCl	-	-	8260	RFPP

REMARKS:

MATERIAL CODES: AG = Amber Glass; CG = Clear Glass; PE = Polyethylene; PP = Polypropylene; S = Silicone; T = Teflon; O = Other (Specify)
SAMPLING/PURGING EQUIPMENT CODES: APP = After Peristaltic Pump; B = Bailor; BP = Bladder Pump; ESP = Electric Submersible Pump; PP = Peristaltic Pump
RFPP = Reverse Flow Peristaltic Pump; SM = Straw Method (Tubing Gravity Drain); VT = Vacuum Trap; O = Other (Specify)

NOTES: 1. The above do not constitute all of the information required by Chapter 62-160, F.A.C.

2. STABILIZATION CRITERIA FOR RANGE OF VARIATION OF LAST THREE CONSECUTIVE READINGS (SEE FS 2212, SECTION 3)

pH: ± 0.2 units Temperature: ± 0.2 °C Specific Conductance: + 5% Dissolved Oxygen: all readings < 20% saturation (see Table FS 2200-2); optionally, ± 0.2 mg/L or ± 10% (whichever is greater) Turbidity: all readings ≤ 20 NTU; optionally ± 5 NTU or ± 10% (whichever is greater)

ARCADIS

Appendix B

Chain-of-Custody Forms

586 660-20738

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

660-20738

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

TestAmerica Tampa
6712 Benjamin Road, Suite 100
Tampa, FL 33634

Alternate Laboratory Name/Location

Website: www.testamericainc.com
Phone: (813) 885-7427
Fax: (813) 885-7049

Serial Number 00525

PROJECT REFERENCE Vallewest		PROJECT NO. BOSSOSS	PROJECT LOCATION (STATE) FL	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 1 OF 4
SAMPLER'S SIGNATURE <i>Steve Skimble</i>		P.O. NUMBER	CONTRACT NO.	COMPOSITE (S) OR GRAB (G) INDICATE	STANDARD REPORT DELIVERY	DATE DUE
CLIENT (SITE) PM Guy Yeminski		CLIENT PHONE 813-915-1216	CLIENT FAX	AQUEOUS (WATER)	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE
CLIENT NAME ARCADIS		CLIENT EMAIL		SOLID OR SEMISOLID	NUMBER OF COOLERS SUBMITTED PER SHIPMENT:	
CLIENT ADDRESS 350 Buschwood Park Drive		Sikelofsky		AIR	REMARKS	
COMPANY CONTRACTING THIS WORK (if applicable) 33618				NONAQUEOUS LIQUID (OIL, SOLVENT...)		
DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED		REMARKS
1/31/07	948	MW-228		3		
		MW-204		3		
		MW-205		3		
		MW-240		3		
		MW-169		3		
		MW-182		3		
		MW-167		3		
		MW-168		3		
		MW-170		3		
		MW-154		3		
		MW-151		3		
		MW-121		3		
RELINQUISHED BY: (SIGNATURE)		DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
<i>Steve Skimble</i>		1/31/07	815			
RECEIVED BY: (SIGNATURE)		DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
<i>Steve Skimble</i>		1/31/08	815			
EMPTY CONTAINERS						
RECEIVED FOR LABORATORY BY: (SIGNATURE)		DATE	TIME	CUSTODY INTACT	CUSTODY SEAL NO.	TAMPA LOG NO.
<i>Angy Akrun</i>		1/31/08	945	YES <input type="radio"/> NO <input type="radio"/>		20738
LABORATORY USE ONLY						
COOLER TEMP UPON RECEIPT						
LABORATORY REMARKS						

8260+8260 SIM

PRESEPTATIVE

Serial Number 00524

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

0600-20738

TestAmerica Tampa
6712 Benjamin Road, Suite 100
Tampa, FL 33634

Website: www.testamericainc.com
Phone: (813) 885-7427
Fax: (813) 885-7049

Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE	OF
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE		2	4
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)			
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID			
CLIENT ADDRESS			AIR			
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)			
DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS	
1/23/08	15:10	MW-172		3		
		1552		3		
		1413		3		
		1002		3		
		1120		3		
		1141		3		
		1415		3		
		1417		3		
		1484		3		
		1485		3		
		1534		3		
		1840		3		
RELINQUISHED BY (SIGNATURE)	DATE	TIME	RELINQUISHED BY (SIGNATURE)	DATE	TIME	RELINQUISHED BY (SIGNATURE)
<i>[Signature]</i>	1/19/08	07:30	<i>[Signature]</i>	1/24/07	815	
RECEIVED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)
<i>[Signature]</i>	1/24/08	815	<i>[Signature]</i>	1/24/08	815	

WTS
0928-0928
PRESERVATIVE

RECEIVED FOR LABORATORY BY: *[Signature]* DATE: 1/24/08 TIME: 745 CUSTODY INTACT: YES NO LABORATORY USE ONLY: TAMPA LOG NO: 200738 COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

6600-20738

TestAmerica Tampa
6712 Benjamin Road, Suite 100
Tampa, FL 33634
Alternate Laboratory Name/Location

Website: www.testamericainc.com
Phone: (813) 885-7427
Fax: (813) 885-7049

Phone:
Fax:

PROJECT REFERENCE PROJECT NO. PROJECT LOCATION (STATE) CONTRACT NO.

SAMPLER'S SIGNATURE P.O. NUMBER CLIENT PHONE CLIENT FAX

CLIENT (SITE) PM CLIENT NAME CLIENT E-MAIL

CLIENT ADDRESS

COMPANY CONTRACTING THIS WORK (if applicable)

MATRIX TYPE

REQUIRED ANALYSIS

PAGE 3 OF 4

STANDARD REPORT DELIVERY DATE DUE

EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE

NUMBER OF COOLERS SUBMITTED PER SHIPMENT: _____

COMPOSITE (C) OR GRAB (G) INDICATE	DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT	REMARKS
AQUEOUS (WATER)			
SOLID OR SEMISOLID			
AIR			
NONAQUEOUS LIQUID (OIL, SOLVENT...)			

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	MATRIX TYPE	REQUIRED ANALYSIS	PAGE	STANDARD REPORT DELIVERY DATE DUE	EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT	REMARKS
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1/23/08	1100	mw-188	g			3				
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	1205	mw-192				3				
--	------	--------	--	--	--	---	--	--	--	--

	1150	mw-191				3				
--	------	--------	--	--	--	---	--	--	--	--

	1085	mw-189				3				
--	------	--------	--	--	--	---	--	--	--	--

	1520	mw-212				3				
--	------	--------	--	--	--	---	--	--	--	--

	1440	mw-208				3				
--	------	--------	--	--	--	---	--	--	--	--

	1550	mw-211				3				
--	------	--------	--	--	--	---	--	--	--	--

	1010	mw-210				3				
--	------	--------	--	--	--	---	--	--	--	--

	1445	mw-209				3				
--	------	--------	--	--	--	---	--	--	--	--

		DUP2				3				
--	--	------	--	--	--	---	--	--	--	--

		DUP3				3				
--	--	------	--	--	--	---	--	--	--	--

		DUP4				3				
--	--	------	--	--	--	---	--	--	--	--

RELINQUISHED BY: (SIGNATURE) DATE TIME RECEIVED BY: (SIGNATURE) DATE TIME

RECEIVED BY: (SIGNATURE) DATE TIME

EMPTY CONTAINERS

EMPTY CONTAINERS

RECEIVED BY: (SIGNATURE) DATE TIME

RELINQUISHED BY: (SIGNATURE) DATE TIME

RECEIVED BY: (SIGNATURE) DATE TIME

LABORATORY USE ONLY

LABORATORY REMARKS

COOLER TEMP UPON RECEIPT

LABORATORY REMARKS

Serial Number 00526

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

TestAmerica
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660-20738

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Serial Number 00527

Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 4 OF 4
SAMPLER'S SIGNATURE	CONTRACT NO.	CLIENT (SITE) PM	COMPOSITE (C) OR GRAB (G) INDICATE	STANDARD REPORT DELIVERY	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID	NONAQUEOUS LIQUID (OIL, SOLVENT...)	
CLIENT ADDRESS			AIR		
COMPANY CONTRACTING THIS WORK (if applicable)					
DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/23/08	10:30	DUP-5		3	STK
		DUP-2		2	STK
		DUP-3		2	STK
		DUP-4		2	STK
		DUP-5		2	STK
		TB-2 (SIL)		2	STK
		TB-3 (MTE)		2	STK
		TB-4 (SW/AS)		2	STK
		TB-5 (VLE/DUPS)		2	STK
		DUPA		3	STK
1/23/08 10:44		MW-152			Added per client N4 12-8-08
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS			RECEIVED BY: (SIGNATURE)		
RECEIVED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS			RECEIVED BY: (SIGNATURE)		
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	YES <input type="radio"/> NO <input type="radio"/>	
LABORATORY USE ONLY	TAMPA LOG NO.	COOLER TEMP. UPON RECEIPT	LABORATORY REMARKS		
	AD938	815			

SDG 660-20784

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Fax:

Serial Number **00528**

PROJECT REFERENCE: Tallowast PROJECT NO. _____ PROJECT LOCATION (STATE): _____ CONTRACT NO. _____

CLIENT'S SIGNATURE: [Signature] P.O. NUMBER _____ CLIENT PHONE _____ CLIENT FAX _____

CLIENT (SITE) PM: Gov Yaminiski CLIENT E-MAIL _____

CLIENT NAME: Attacalis

CLIENT ADDRESS: 350 Besenwood Park Drive, Suite 100, Tampa, FL 33618

COMPANY CONTRACTING THIS WORK (if applicable): _____

SAMPLE DATE	SAMPLE TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT...)	MATRIX TYPE	REQUIRED ANALYSIS					PAGE 1 OF 5			
									STANDARD REPORT DELIVERY	DATE DUE	EXPEDITED REPORT DELIVERY (SUCH AS HARGE)	DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT				
1/24/08	1630	MW-94															
1/24/08	1547	MW-95															
		TR-1															
	930	RW-2															
	1035	MW-12															
	1130	MW-5															
	1420	MW-114															
	1525	MW-83															
	1540	Private Mail 84															
	1012	MW-32															
	1054	MW-75															
		DO07															

RELINQUISHED BY (SIGNATURE): _____ DATE: _____ TIME: _____

EMPTY CONTAINERS: _____ DATE: _____ TIME: _____

RECEIVED BY (SIGNATURE): _____ DATE: _____ TIME: _____

EMPTY CONTAINERS: _____ DATE: _____ TIME: _____

RELINQUISHED BY (SIGNATURE): _____ DATE: _____ TIME: _____

RECEIVED BY (SIGNATURE): _____ DATE: _____ TIME: _____

RECEIVED FOR LABORATORY BY (SIGNATURE): Carol Mc Nutty DATE: 1/23/08 TIME: 0915

CUSTODY INTACT: YES NO

LABORATORY USE ONLY: TAMPA LOG NO. 20784 COOLER TEMP. UPON RECEIPT: _____

LABORATORY REMARKS: _____

500

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Serial Number

00529

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 2 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE		STANDARD REPORT DELIVERY <input type="checkbox"/>
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)		DATE DUE _____
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID		EXPEDITED REPORT DELIVERY (SUCH AS CHARGE) <input type="checkbox"/>
CLIENT ADDRESS			AIR		DATE DUE _____
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		NUMBER OF COOLERS SUBMITTED PER SHIPMENT: _____
SAMPLE DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/24/08	1117	MW-7D	S	1	
1/31/08	1515	MW-15A		1	
1/31/08	1432	MW-157		1	
1/31/08	1515	MW-110D		1	
1/31/08	1435	SK MW-156		1	
1/31/08	1105	MW-11		1	
1/31/08	1628	EB-7		1	
1/31/08	1628	TR-7		1	
1/31/08	1628	TR-6		1	
1/31/08	1628	EB-6		1	
1/31/08	1450	MW-107		1	
1/31/08	1045	MW-153		1	
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME

RECEIVED FOR LABORATORY BY: (SIGNATURE) *Muel McSwatky* DATE: 1/31/08 TIME: 0915

RECEIVED BY: (SIGNATURE) *Amil Roberts* DATE: 1/31/08 TIME: 754

LABORATORY USE ONLY

CUSTODY SEAL NO. _____ TAMPA LOG NO. _____ COOLER TEMP UPON RECEIPT _____

LABORATORY REMARKS _____

Serial Number 00530

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 3 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE	STANDARD REPORT DELIVERY	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID	NUMBER OF COOLERS SUBMITTED PER SHIPMENT	
CLIENT ADDRESS			AIR		
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		
SAMPLE		SAMPLE IDENTIFICATION		REMARKS	
DATE	TIME			NUMBER OF CONTAINERS SUBMITTED	
1/24/08	1130	MW-185	<input checked="" type="checkbox"/>	3	
	1534	MW-161	<input checked="" type="checkbox"/>	5	
	1535	MW-158	<input checked="" type="checkbox"/>		
	1650	ER-9	<input checked="" type="checkbox"/>		
	930	MW-124	<input checked="" type="checkbox"/>		
	1010	MW-126	<input checked="" type="checkbox"/>		
	1018	MW-106	<input checked="" type="checkbox"/>		
	1045	MW-112	<input checked="" type="checkbox"/>		
	1214	MW-195	<input checked="" type="checkbox"/>		
	1245	MW-194	<input checked="" type="checkbox"/>		
	1248	MW-193	<input checked="" type="checkbox"/>		
		DUOL	<input checked="" type="checkbox"/>		
REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS			AMM BATS	1/25/08	754
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS				1/25/08	754

RECEIVED FOR LABORATORY BY (SIGNATURE)	DATE	TIME	CUSTODY SEAL NO.	TAMPA LOG NO.	COOLER TEMP. UPON RECEIPT	LABORATORY REMARKS
And the truth	1/25/08	0915				

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ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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Alternate Laboratory Name/Location

Phone:
Fax:

Serial Number

00531

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE	OF		
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE		STANDARD REPORT DELIVERY			
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)		DATE DUE			
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID		EXPEDITED REPORT DELIVERY (SURCHARGE)			
CLIENT ADDRESS			AIR		DATE DUE			
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		NUMBER OF COOLERS SUBMITTED PER SHIPMENT			
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS	DATE	TIME	
12/10/08	15:10	MW-184	<input checked="" type="checkbox"/>	1				
12/10/08	15:19	MW-183	<input checked="" type="checkbox"/>	1				
12/10/08	16:00	MW-185	<input checked="" type="checkbox"/>	1				
12/10/08	16:10	MW-187	<input checked="" type="checkbox"/>	1				
12/10/08	16:40	MW-186	<input checked="" type="checkbox"/>	1	Changed per email from sampler			
		DUP-9	<input checked="" type="checkbox"/>	1				
		TR-9	<input checked="" type="checkbox"/>	1				
		TR-8	<input checked="" type="checkbox"/>	1				
		EB-8	<input checked="" type="checkbox"/>	1				
		DUP-8	<input checked="" type="checkbox"/>	1				
		MW-185	<input checked="" type="checkbox"/>	1				
		MW-150	<input checked="" type="checkbox"/>	1				
REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
RECEIVED BY: (SIGNATURE)	DATE	TIME	LABORATORY USE ONLY	DATE	TIME	LABORATORY USE ONLY	DATE	TIME
RECEIVED FOR LABORATORY BY (SIGNATURE)	DATE	TIME	CUSTODY SEAL NO.	TAMPA LOG NO.	COOLER TEMP UPON RECEIPT	LABORATORY REMARKS		
Carol McInally	12/29/08	09:15			754			

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Serial Number

00532

Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 5 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE	STANDARD REPORT DELIVERY	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID		
CLIENT ADDRESS			AIR		
GENERALLY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		
SAMPLE DATE	SAMPLE TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/25/08	1012	MW-14S		3	
	1009	MW-14D		3	
	1132	MW-77		3	
	1211	MW-20		3	
	1408	MW-95		3	
	1400	MW-9D		3	
	1445	MW-76		3	
	1530	MW-25		3	
	1452	MW-85		3	
	1715	MW-233		3	
TOTAL OF 58 Samples					
REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTOMER CONTACT YES/NO	LABORATORY/USE ONLY	COOLER TEMP. UPON RECEIPT
Carol Mc Nulty	1/25/08	0915	YES	TAMPA LOG NO.	

SDG-660-20815

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660-20815

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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Fax: (813) 885-7049

Serial Number

00533

Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE
Taleo West

SAMPLER'S SIGNATURE
Tina Skunk

CLIENT (SITE) PM
Guy Kaminski

CLIENT NAME
Arcadis.

CLIENT ADDRESS
3350 Brownwood Park Drive Suite 100 Tampa, FL 33618

COMPANY CONTRACTING THIS WORK (if applicable)

PROJECT NO.
60038055

PO. NUMBER

CLIENT PHONE
813-915-1720

CLIENT E-MAIL
guy.kaminski@arcadis-us.com

PROJECT LOCATION (STATE)
FL

CONTRACT NO.

CLIENT FAX

MATRIX TYPE

REQUIRED ANALYSIS

PAGE 1 OF 9

STANDARD REPORT DELIVERY

DATE DUE

EXPEDITED REPORT DELIVERY (SURCHARGE)

DATE DUE

NUMBER OF COOLERS SUBMITTED PER SHIPMENT:

REMARKS

COMPOSITE (S) OR GRAB (G) INDICATE

AQUEOUS (WATER)

SOLID OR SEMISOLID

AIR

NONAQUEOUS LIQUID (OIL, SOLVENT...)

NUMBER OF CONTAINERS SUBMITTED

REMARKS

DATE

TIME

SAMPLE IDENTIFICATION

DATE

TIME

REINQUISHED BY: (SIGNATURE)

DATE

TIME

RECEIVED BY: (SIGNATURE)

DATE

TIME

RECEIVED FOR LABORATORY BY (SIGNATURE)	DATE	TIME	CUSTODY INTACT (YES/NO)	LABORATORY SEAL NO.	TAMPA LOG NO.	COOLER TEMP. UPON RECEIPT	LABORATORY REMARKS
<i>[Signature]</i>	1/28/08	936	YES		660	4040	

SAMPLE	DATE	TIME	SAMPLE IDENTIFICATION	REINQUISHED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	REMARKS
105108	903		MW-117	<i>[Signature]</i>	1/28/08	936				
	840		MW-118							
	806		MW-119							
	750		MW-120							
	1020		MW-224							
	1032		MW-227							
	1100		MW-223							
	1120		MW-225							
	1155		MW-226							
	1157		MW-241							
	1300		MW-100							
	1303		MW-99							

Serial Number

00534

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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6600-20815

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Alternate Laboratory Name/location

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PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 2 OF 4
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE		STANDARD REPORT DELIVERY <input type="checkbox"/>
CLIENT (SITE) PM	PHONE	CLIENT FAX	AQUEOUS (WATER)		DATE DUE
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID		EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="checkbox"/>
CLIENT ADDRESS			AIR		DATE DUE
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		NUMBER OF COOLERS SUBMITTED PER SHIPMENT
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/28/08	13:23	EB-10	<input checked="" type="checkbox"/>	1	
		1343	<input checked="" type="checkbox"/>	1	
		1121	<input checked="" type="checkbox"/>	1	
		1041	<input checked="" type="checkbox"/>	1	
		1252	<input checked="" type="checkbox"/>	1	
		1251	<input checked="" type="checkbox"/>	1	
		1341	<input checked="" type="checkbox"/>	1	
		1345	<input checked="" type="checkbox"/>	1	
		1455	<input checked="" type="checkbox"/>	1	
		1410	<input checked="" type="checkbox"/>	1	
		DUP-11	<input checked="" type="checkbox"/>	1	
		DUP-10	<input checked="" type="checkbox"/>	1	
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
<i>Diana Salinas</i>	1/28/08	9:36	<i>Mary Stiles</i>	1/28/08	9:36
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	LABORATORY USE ONLY	LABORATORY REMARKS
<i>Mary Stiles</i>	1-28-08		YES <input type="checkbox"/> NO <input type="checkbox"/>	TAMPA LOG NO. 6600	COOLER TEMP. UPON RECEIPT 40.4 °C

Serial Number

00535

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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660-2085

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THE LEADER IN ENVIRONMENTAL TESTING

Alternate Laboratory Name/Location

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PROJECT REFERENCE: PROJECT NO.: PROJECT LOCATION (STATE): MATRIX TYPE: REQUIRED ANALYSIS: PAGE 3 OF 4

SAMPLER'S SIGNATURE: P.O. NUMBER: CONTRACT NO. STANDARD REPORT DELIVERY DATE DUE: 0

CLIENT (SITE) PM: CLIENT NAME: CLIENT E-MAIL: CLIENT FAX: EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE: 0

CLIENT ADDRESS: COMPANY CONTRACTING THIS WORK (if applicable): NUMBER OF COOLERS SUBMITTED PER SHIPMENT: 0

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/25/08	1115	MW-103						3	
	1030	MW-102						4	
	855	MW-190							
	1508	EB-13							
	1317	MW-88							
	1339	MW-3							
	1500	MW-84							
	1345	MW-4							
		DP-13							
		TR-13							
	840	MW-174							
	1125	MW-179							

REINQUISHED BY: (SIGNATURE) DATE: TIME: RELINQUISHED BY: (SIGNATURE) DATE: TIME:

EMPTY CONTAINERS: RECEIVED BY: (SIGNATURE) DATE: TIME:

EMPTY CONTAINERS: RECEIVED BY: (SIGNATURE) DATE: TIME:

RECEIVED FOR LABORATORY BY: DATE: TIME: CUSTODY INTACT: YES NO

LABORATORY USE ONLY: TAMPA LOG NO. COOLER TEMP. UPON RECEIPT: 40.4

LABORATORY REMARKS:

TAL8240-580 (1207)

Serial Number

00536

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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6600-20815

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Alternate Laboratory Name/Location

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PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 4 OF 4
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SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	STANDARD REPORT DELIVERY
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CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	DATE DUE
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CLIENT NAME	CLIENT E-MAIL	EXPEDITED REPORT DELIVERY (SURCHARGE)
-------------	---------------	---------------------------------------

CLIENT ADDRESS	DATE DUE
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COMPANY CONTRACTING THIS WORK (if applicable)	NUMBER OF COOLERS SUBMITTED PER SHIPMENT
---	--

SAMPLE DATE	SAMPLE TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
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1/25/08	1420	MW-116	9	1	
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	925	MW-122		1	
--	-----	--------	--	---	--

	1115	MW-178		1	
--	------	--------	--	---	--

	1410	MW-180		1	
--	------	--------	--	---	--

	1510	MW181		1	
--	------	-------	--	---	--

		Tamp Blank			
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REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME
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EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
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RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
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RECEIVED FOR LABORATORY USE ONLY	DATE	TIME	LABORATORY USE ONLY	DATE	TIME
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RECEIVED BY: (SIGNATURE)	DATE	TIME	COOLER TEMP UPON RECEIPT	LABORATORY REMARKS
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SDC 660-20822

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660-20822

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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Alternate Laboratory Name/Location

Phone:
Fax:

Serial Number

00537

PROJECT REFERENCE: Tallevest PROJECT NO.: B0088055 PROJECT LOCATION (STATE): FL CONTRACT NO.: 33618

SAMPLER'S SIGNATURE: [Signature] PO. NUMBER: W150998-0028

CLIENT (SITE) PM: Guy Leminski CLIENT PHONE: 813-915-4720 CLIENT FAX: 813-915-4720

CLIENT NAME: Arcaadis CLIENT E-MAIL: guy.leminski@arcaadis-us.com

CLIENT ADDRESS: 3350 Buschwood Park Drive, Suite 100, Tampa, FL 33618

COMPANY CONTRACTING THIS WORK (if applicable): 33618

REQUIREMENT ANALYSIS: W150998-0028

MATRIX TYPE: W150998-0028

COMPOSITE (C) OR GRAB (G) INDICATE: G

AQUEOUS (WATER): W150998-0028

SOLID OR SEMISOLID: W150998-0028

AIR: W150998-0028

NONAQUEOUS LIQUID (OIL, SOLVENT...): W150998-0028

REQUIRED ANALYSIS: PRESERVATIVE

DATE DUE: 01/29/08

STANDARD REPORT DELIVERY: 0

EXPEDITED REPORT DELIVERY (SURCHARGE): 0

DATE DUE: 01/29/08

PAGE 1 OF 4

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT...)	REMARKS
1/28/08	1305	MW-31						
		MW-59						
		MW-68						
		MW-60						
		MW-109						
		MW-21						
		MW-86						
		MW-110						
		MW-107						
		MW-105						
		MW-135						
		MW-93						

RELINQUISHED BY: (SIGNATURE) [Signature] DATE: 1/29/08 TIME: 8:48

RECEIVED BY: (SIGNATURE) [Signature] DATE: 1/29/08 TIME: 8:17

RECEIVED FOR LABORATORY BY: (SIGNATURE) [Signature] DATE: 1/29/08 TIME: 1:00

CUSTODY INTACT: YES NO

CUSTODY SEAL NO. 129108

TAMPA LOG NO. 017

COOLER TEMP UPON RECEIPT: 4

LABORATORY REMARKS: LABORATORY USE ONLY

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Fax:

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Serial Number

00803

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 2 OF 4
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE		STANDARD REPORT DELIVERY <input type="radio"/>
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)		DATE DUE _____
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID		EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="radio"/>
CLIENT ADDRESS			AIR		DATE DUE _____
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		NUMBER OF COOLERS SUBMITTED PER SHIPMENT: _____
DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/20/08	1318	MW-50	<input checked="" type="checkbox"/>	3	
	1700	MW-111		3	
	1618	MW-113		3	
	1640	MW-51		3	
	1503	MW-26		3	
	1319	MW-44		3	
	1345	MW-70		3	
	1050	MW-188		3	
	1035	MW-18D		3	
	935	MW-91		3	
	1110	MW-29		3	
	937	MW-133		3	
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					

PRESERVATIVE

020-2005DM

LABORATORY USE ONLY
RECEIVED FOR LABORATORY BY: (SIGNATURE) DATE TIME
CUSTODY SEAL NO. TAMPA LOG NO. COOLER TEMP. UPON RECEIPT
LABORATORY REMARKS

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD
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Alternate Laboratory Name/Location

Phone:
 Fax:

Serial Number **03804**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 3 OF 4
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	STANDARD REPORT DELIVERY <input type="radio"/>	EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="radio"/>	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	DATE DUE		
CLIENT NAME	CLIENT E-MAIL		DATE DUE		
CLIENT ADDRESS			DATE DUE		
COMPANY CONTRACTING THIS WORK (if applicable)					

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT,...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/28/08	1140	MW-52						3	
		MW-104						3	
		MW-28						3	
		MW-53						3	
		MW-54						3	
		MW-27						3	
		MW-104						3	
		MW-105						3	
		MW-248						3	
		Skate-2 Dup-14						3	
		Dup15						3	
		Dup16						3	

REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
LABORATORY USE ONLY					
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT <input type="radio"/>	LABORATORY SEAL NO.	TAMPA LOG NO.
			YES <input type="radio"/>		
			NO <input type="radio"/>		
COOLER TEMP UPON RECEIPT			LABORATORY REMARKS		

TAL 8240-690 (1207)

Serial Number

00805

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE PROJECT NO. PROJECT LOCATION (STATE) CONTRACT NO.

SAMPLER'S SIGNATURE P.O. NUMBER CLIENT PHONE CLIENT FAX

CLIENT (SITE) PM CLIENT NAME CLIENT E-MAIL

CLIENT ADDRESS

COMPANY CONTRACTING THIS WORK (if applicable)

DATE TIME SAMPLE IDENTIFICATION

COMPOSITE (C) OR GRAB (G) INDICATE

AQUEOUS (WATER)

SOLID OR SEMISOLID

AIR

NONAQUEOUS LIQUID (OIL, SOLVENT...)

MATRIX TYPE

REQUIRED ANALYSIS

PAGE 4 OF 4

STANDARD REPORT DELIVERY

DATE DUE

EXPEDITED REPORT DELIVERY (SURCHARGE)

DATE DUE

NUMBER OF COOLERS SUBMITTED PER SHIPMENT

REMARKS

DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)
1/28/08	14	[Signature]	1/28/08	8:47	[Signature]
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			
	14	[Signature]			
	15	[Signature]			
	16	[Signature]			
	17	[Signature]			

49 Samples total

RELINQUISHED BY (SIGNATURE)

DATE

TIME

RECEIVED BY (SIGNATURE)

DATE

TIME

RECEIVED BY (SIGNATURE)

DATE

TIME

RECEIVED BY (SIGNATURE)

DATE

TIME

RECEIVED BY (SIGNATURE)

DATE

TIME

RECEIVED BY (SIGNATURE)

DATE

TIME

RECEIVED BY (SIGNATURE)

DATE

TIME

70602-099 SDC

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Alternate Laboratory Name/Location
 Phone:
 Fax:

Serial Number **00811**

PROJECT REFERENCE: Tallevest PROJECT NO: 80038055 PROJECT LOCATION (STATE): FL CONTRACT NO. F1 MATRIX TYPE: WTS 0098-098 REQUIRED ANALYSIS: PRESERVATIVE PAGE 1 OF 5

SAMPLES SIGNATURE: [Signature] P.O. NUMBER: _____ STANDARD REPORT DELIVERY: DATE DUE: _____

CLIENT (SITE) PM: GUY KAMINSKI CLIENT PHONE: 813-915-4726 CLIENT FAX: _____ EXPEDITED REPORT DELIVERY (SURCHARGE): DATE DUE: _____

CLIENT NAME: Arcaadis CLIENT E-MAIL: guy.kaminski@arcaadis-us.com NUMBER OF COOLERS SUBMITTED PER SHIPMENT: _____

CLIENT ADDRESS: 3350 Bosch Road, Raleigh Drive, Suite 100, Tampa, FL 33606

COMPANY CONTRACTING THIS WORK (if applicable): _____

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR SUB (S) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/30/08	1340	MW-71							
	1349	MW-149							
	1431	MW-129							
	1436	MW-58							
	922	MW-130							
	930	MW-57							
	10210	MW-80							
	1027	MW-72							
	1118	MW-6							
	1620	MW-30							
	1206	MW-48							
	1503	MW-232							

RELINQUISHED BY: (SIGNATURE) DATE TIME RECEIVED BY: (SIGNATURE) DATE TIME

EMPTY CONTAINERS: _____ DATE TIME

RECEIVED BY: (SIGNATURE) DATE TIME

EMPTY CONTAINERS: _____ DATE TIME

RECEIVED FOR LABORATORY BY: [Signature] DATE: 1-31-08 TIME: 0820 CUSTODY INTACT: YES NO

LABORATORY USE ONLY: TAMPA LOG NO: 40-20907 COOLER TEMP. UPON RECEIPT: 49° LABORATORY REMARKS: _____

LABORATORY SEAL NO. _____

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Serial Number

00814

Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 2 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	STANDARD REPORT DELIVERY	DATE DUE	
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE	
CLIENT NAME	CLIENT ADDRESS		DATE DUE		
COMPANY CONTRACTING THIS WORK (if applicable)					
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/30/08	—	DUP-22 EGG- EW-110	X AQUEOUS (WATER)	3	
		1530		3	
		1538		3	
		1908		3	
		1213		3	
		1447		3	
		1450		3	
		1251		3	
		1445		3	
		1440		3	
		1440		3	
		DUP-25		3	
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	LABORATORY SEAL NO.	TAMPA LOG NO.
	1/30/08	1000	YES <input type="radio"/> NO <input type="radio"/>		660-20907
COOLER TEMP. UPON RECEIPT	LABORATORY USE ONLY		COOLER TEMP. UPON RECEIPT	LABORATORY REMARKS	
40			40		

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Fax: (813) 885-7049

Alternate Laboratory Name/Location

Phone:
Fax:

Serial Number

00815

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 3 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTACT NO.	REQUIRED ANALYSIS	STANDARD REPORT DELIVERY <input type="checkbox"/>	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	REQUIRED ANALYSIS	EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="checkbox"/>	DATE DUE
CLIENT NAME	CLIENT E-MAIL		REQUIRED ANALYSIS	DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT:
CLIENT ADDRESS			REQUIRED ANALYSIS		
COMPANY CONTACTING THIS WORK (if applicable)			REQUIRED ANALYSIS		
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/30/08	1430	MW-103	AQUEOUS (WATER)		
		MW-102	SOLID OR SEMISOLID		
		MW-97	AIR		
		MW-87	NONAQUEOUS LIQUID (OIL, SOLVENT....)		
		MW-75			
		MW-34			
		MW-35			
		MW-251			
		MW-109			
		MW-101			
		MW-196			
		MW-197			
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME

RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT YES <input type="checkbox"/> NO <input type="checkbox"/>	LABORATORY SEAL NO.	TAMPA LOG NO.	COOLER TEMP. UPON RECEIPT	LABORATORY REMARKS
M. Williams	1-31-08	1000			600-20907	49°	

Serial Number 00812

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE: PROJECT NO.: PROJECT LOCATION (STATE): MATRIX TYPE: REQUIRED ANALYSIS: PAGE 5 OF 5

SAMPLER'S SIGNATURE: P.O. NUMBER: CONTRACT NO.: STANDARD REPORT DELIVERY:

CLIENT (SITE) PM: CLIENT PHONE: CLIENT FAX: DATE DUE:

CLIENT NAME: CLIENT E-MAIL: EXPEDITED REPORT DELIVERY (SURCHARGE):

CLIENT ADDRESS: DATE DUE: NUMBER OF COOLERS SUBMITTED PER SHIPMENT:

COMPANY CONTRACTING THIS WORK (if applicable):

SAMPLE DATE: SAMPLE TIME: SAMPLE IDENTIFICATION: NUMBER OF CONTAINERS SUBMITTED: REMARKS:

1/30/08 1705 MW-47
11060 MW-81
IB-24
Dup-24

RELINQUISHED BY: (SIGNATURE) DATE: TIME: RELINQUISHED BY: (SIGNATURE) DATE: TIME:

EMPTY CONTAINERS: RECEIVED BY: (SIGNATURE) DATE: TIME: RECEIVED BY: (SIGNATURE) DATE: TIME:

EMPTY CONTAINERS: RECEIVED BY: (SIGNATURE) DATE: TIME: RECEIVED BY: (SIGNATURE) DATE: TIME:

RECEIVED FOR LABORATORY BY: (SIGNATURE) DATE: TIME: CUSTODY INTACT: YES NO

LABORATORY SEAL NO.: TAMPA LOG NO.: COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

LABORATORY USE ONLY: DATE: TIME: RECEIVED BY: (SIGNATURE) DATE: TIME:

LABORATORY SEAL NO.: TAMPA LOG NO.: COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

LABORATORY SEAL NO.: TAMPA LOG NO.: COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

LABORATORY SEAL NO.: TAMPA LOG NO.: COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

LABORATORY SEAL NO.: TAMPA LOG NO.: COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

LABORATORY SEAL NO.: TAMPA LOG NO.: COOLER TEMP UPON RECEIPT: LABORATORY REMARKS:

SDG 660-20941

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Serial Number 00823

Alternate Laboratory Name/Location

Phone:
Fax:

PAGE 1 OF 2

PROJECT REFERENCE: Valleerst PROJECT NO: 60388055 PROJECT LOCATION (STATE): FL CONTRACT NO. F1 MATRIX TYPE: PM REQUIRED ANALYSIS: PM

SAMPLES SIGNATURE: [Signature] P.O. NUMBER: 813-915-4700 CLIENT FAX: 813-915-4700 CLIENT PHONE: 813-915-4700 CLIENT E-MAIL: gy.kaminski@arccadis-us.com

CLIENT NAME: Arccadis CLIENT ADDRESS: 250 Bosch Woodward Park Drive Suite 100 Tampa COMPANY CONTRACTING THIS WORK (if applicable): F1 33618

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR TAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT, ...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS
8/1/08	810	MW-202						3	
	1132	MW-134						3	
	945	DW-1						3	
	1050	MW-42						3	
	1103	MW-43						3	
	915	IWT-1						3	
	1025	IWT-2						3	
	1000	EW-UACG-1						3	
	940	MW-38						3	
	1010	MW-39						3	
	930	MW-253						3	
	840	MW-254						3	

RECEIVED BY: (SIGNATURE) [Signature]
DATE: 8/1/08 TIME: 1518

RECEIVED BY: (SIGNATURE) [Signature]
DATE: 8/1/08 TIME: 1518

RECEIVED BY: (SIGNATURE) _____
DATE: _____ TIME: _____

RECEIVED FOR LABORATORY BY: (SIGNATURE) _____ DATE: _____ TIME: _____
CUSTODY: INTACT YES NO
LABORATORY USE ONLY: CUSTODY SEAL NO. _____ TAMPA LOG NO. _____ COOLER TEMP. UPON RECEIPT: _____ LABORATORY REMARKS: _____

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Fax:

Serial Number **00822**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE OF
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	AQUEOUS (WATER)		2 OF 2
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	SOLID OR SEMISOLID		
CLIENT NAME	SHEET/EMAIL		AIR		
CLIENT ADDRESS			NONAQUEOUS LIQUID (OIL, SOLVENT...)		
COMPANY CONTRACTING THIS WORK (if applicable)					
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
2/1/08	906	MW-81D	X	3	
	921	MW-123		3	
	910	RW-1		3	
		DOD-31		2	
		EB-31		2	
		TR-31		2	
	854	Private Well 7		3	
	820	Private Well 57		3	
	908	Private Well 59		3	
	1055	Private Well 38		3	
	115	Private Well 102		3	
		TR-32		2	
REINQUISHED BY: (SIGNATURE)	DATE <td>TIME <td>REINQUISHED BY: (SIGNATURE)</td> <td>DATE <td>TIME </td></td></td>	TIME <td>REINQUISHED BY: (SIGNATURE)</td> <td>DATE <td>TIME </td></td>	REINQUISHED BY: (SIGNATURE)	DATE <td>TIME </td>	TIME
EMPTY CONTAINERS			<i>[Signature]</i>	2/1/07	1578
RECEIVED BY: (SIGNATURE)	DATE <td>TIME <td>RECEIVED BY: (SIGNATURE)</td> <td>DATE <td>TIME </td></td></td>	TIME <td>RECEIVED BY: (SIGNATURE)</td> <td>DATE <td>TIME </td></td>	RECEIVED BY: (SIGNATURE)	DATE <td>TIME </td>	TIME
EMPTY CONTAINERS			<i>[Signature]</i>	2/1/08	1578
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE <td>TIME <td>CUSTODY SEAL NO.</td> <td>TAMPA LOG NO. <td>COOLER TEMP. UPON RECEIPT</td> </td></td>	TIME <td>CUSTODY SEAL NO.</td> <td>TAMPA LOG NO. <td>COOLER TEMP. UPON RECEIPT</td> </td>	CUSTODY SEAL NO.	TAMPA LOG NO. <td>COOLER TEMP. UPON RECEIPT</td>	COOLER TEMP. UPON RECEIPT
LABORATORY USE ONLY					
LABORATORY REMARKS					

PRESERVATIVE

Serial Number 43659

SEVERN
TRENT

STL
660-21028

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

STL Tampa
6712 Benjamin Road, Suite 100
Tampa, FL 33634

Website: www.stl-inc.com
Phone: (813) 885-7427
Fax: (813) 885-7049

Alternate Laboratory Name/Location

Phone:
Fax:

PROJECT REFERENCE: 1 letterest PROJECT NO.: 38055.003 PROJECT LOCATION (STATE): FL MATRIX TYPE: REQUIRED ANALYSIS: PAGE 1 OF 1

SAMPLER'S SIGNATURE: [Signature] P.O. NUMBER: 38055 CONTRACT NO.:

CLIENT (SITE) P# [Signature] CLIENT PHONE: 813-933-0691 CLIENT FAX: 813-932-9574

CLIENT NAME: Arcadis CLIENT E-MAIL: Sara.Klimmek@arcadis-us.com

CLIENT ADDRESS: 3350 Buschwood Park Dr, Tampa, FL 33618

COMPANY CONTRACTING THIS WORK (if applicable): LMC

COMPOSITE (C) OR GRAB (G) INDICATE: AQUEOUS (WATER) SOLID OR SEMISOLID AIR NONAQUEOUS LIQUID (OIL, SOLVENT, ...)

DATE: 2-4-08 TIME: 9:55 SAMPLE IDENTIFICATION: PW-64

DATE: 2-4-08 TIME: 11:18 SAMPLE IDENTIFICATION: PW-65

Table with columns: DATE, TIME, RELINQUISHED BY (SIGNATURE), RECEIVED BY (SIGNATURE), DATE, TIME, RELINQUISHED BY (SIGNATURE), RECEIVED BY (SIGNATURE). Includes handwritten signatures and dates.

RECEIVED FOR LABORATORY BY: [Signature] DATE: 2-7-08 TIME: 1:00 CUSTODY SEAL NO.: 660- LABORATORY USE ONLY: STL TAMPA LOG NO.: 660- LABORATORY REMARKS: 4 cooler temp

SDG 660 21028

SDG 660-20857

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Serial Number **00806**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	CONTRACT NO.	MATRIX TYPE	REQUIRED ANALYSIS	PAGE	STANDARD REPORT DELIVERY	DATE DUE	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT	REMARKS
Vallejo East						1	<input type="radio"/>		<input type="radio"/>			
SAMPLER'S SIGNATURE <i>Steve Plimack</i>	P.O. NUMBER	CLIENT PHONE 813-915-1720		CLIENT FAX		<input type="radio"/> STANDARD REPORT DELIVERY <input type="radio"/> EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE _____						
CLIENT (SITE) PM Gurkamski	CLIENT PHONE	CLIENT E-MAIL gurkamski@arcadis.com		CLIENT FAX		<input type="radio"/> STANDARD REPORT DELIVERY <input type="radio"/> EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE _____						
CLIENT NAME Arcadis	CLIENT PHONE	CLIENT E-MAIL		CLIENT FAX		<input type="radio"/> STANDARD REPORT DELIVERY <input type="radio"/> EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE _____						
CLIENT ADDRESS 3350 Bosque Boulevard Company Contracting This Work (if applicable)	CLIENT PHONE	CLIENT E-MAIL		CLIENT FAX		<input type="radio"/> STANDARD REPORT DELIVERY <input type="radio"/> EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE _____						
DATE <th>SAMPLE TIME</th> <th>SAMPLE IDENTIFICATION</th> <th>COMPOSITE (C) OR GRAB (G) INDICATE</th> <th>AQUEOUS (WATER)</th> <th>SOLID OR SEMISOLID</th> <th>AIR</th> <th>NONAQUEOUS LIQUID (OIL, SOLVENT,...)</th> <th>NUMBER OF CONTAINERS SUBMITTED</th> <th>REMARKS</th>	SAMPLE TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT,...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS			
1/29/08	9110	MW-24						3				
	1149	MW-1410						3				
	1557	MW-56						3				
	1519	MW-88						3				
	952	MW-84B						3				
	957	MW-84A						3				
	1450	MW-82						3				
	1622	MW-74						3				
	-	DUP-19						3				
	-	DUP-20						3				
	1605	MW-103						3				
	1130	MW-145						3				
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME	
<i>Steve Plimack</i>	1/30/08	840	<i>Steve Plimack</i>	1/30/08	546							
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	

RECEIVED FOR LABORATORY BY: *Martha Williams* DATE: 1/30/08 TIME: 1145

LABORATORY USE ONLY

COOLERS TEMP. (IF ON RECEIPT)

LABORATORY REMARKS

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Serial Number **UU8U1**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 2 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	AQUEOUS (WATER)		STANDARD REPORT DELIVERY <input type="radio"/>
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	SOLID OR SEMISOLID		DATE DUE _____
CLIENT NAME	CLIENT EMAIL		AIR		EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="radio"/>
CLIENT ADDRESS			NONAQUEOUS LIQUID (OIL, SOLVENT...)		DATE DUE _____
COMPANY CONTRACTING THIS WORK (if applicable)			COMPOSITE (C) OR GRAB (G) INDICATE		NUMBER OF COOLERS SUBMITTED PER SHIPMENT
			<input checked="" type="checkbox"/> AQUEOUS (WATER)		
			<input type="checkbox"/> SOLID OR SEMISOLID		
			<input type="checkbox"/> AIR		
			<input type="checkbox"/> NONAQUEOUS LIQUID (OIL, SOLVENT...)		
SAMPLE	SAMPLE IDENTIFICATION			NUMBER OF CONTAINERS SUBMITTED	REMARKS
DATE	TIME				
1/30/08	1315	MW-143		W	
	1010	MW-140		W	
	850	MW-136		W	
	1415	MW-207		W	
	1535	MW-160		W	
	1405	MW-206		W	
	1315	MW-142		W	
	1350	MW-141		W	
	1200	MW-144		W	
	1115	MW-910		W	
	1005	MW-139		W	
	915	MW-137		W	
RELINQUISHED BY: (SIGNATURE) EMPTY CONTAINERS	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
				1/30/08	846
RECEIVED BY: (SIGNATURE) EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
				1/30/08	846
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	TAMP LOG NO.	COOLER TEMP UPON RECEIPT
			<input type="radio"/> YES <input checked="" type="radio"/> NO	98857	

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 Fax:

Serial Number **00808**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 3 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	<input checked="" type="checkbox"/> COMPOSITE (C) OR GRAB (G) INDICATE <input type="checkbox"/> AQUEOUS (WATER) <input type="checkbox"/> SOLID OR SEMISOLID <input type="checkbox"/> AIR <input type="checkbox"/> NONAQUEOUS LIQUID (OIL, SOLVENT, ...)	STANDARD REPORT DELIVERY <input type="radio"/>	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX		EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="radio"/>	DATE DUE
CLIENT NAME	CLIENT E-MAIL				
CLIENT ADDRESS					
COMPAWK CONTRACTING THIS WORK (if applicable)					
DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/29/08	845	MW-138		3	
	930	MW-176		5	
	1130	MW-175			
	1114	MW-125			
	1305	MW-135			
	1344	MW-131			
	1353	MW-45			
	1005	MW-115			
	928	MW-177			
	1422	MW-243			
	1449	MW-244			
		Duo-18			

RELINQUISHED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME
<i>[Signature]</i>	1/30/08	846	<i>[Signature]</i>	1/30/08	846
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME

RECEIVED FOR LABORATORY BY (SIGNATURE)	DATE	TIME	CUSTODY INTACT	LABORATORY USE ONLY	LABORATORY SIGNATURE
<i>[Signature]</i>			YES <input type="radio"/> NO <input type="radio"/>	AMPER LOG NO. <i>20857</i>	
				COOL/RT TEMP. UPON RECEIPT	

Serial Number 00809

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PROJECT REFERENCE: PROJECT NO. PROJECT LOCATION (STATE) CONTRACT NO. MATRIX TYPE REQUIRED ANALYSIS PAGE 4 OF 5

SAMPLER'S SIGNATURE: P.O. NUMBER CLIENT PHONE CLIENT FAX STANDARD REPORT DELIVERY DATE DUE EXPEDITED REPORT DELIVERY (SURCHARGE) DATE DUE

CLIENT (SITE) PM CLIENT PHONE CLIENT FAX CLIENT NAME CLIENT ADDRESS COMPANY CONTRACTING THIS WORK (if applicable)

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/31/08	1030	MW-231						3	
	1042	MW-229						3	
	933	MW-245						3	
	850	MW-247						3	
	900	MW-246						3	
	1420	MW-242						3	
	1039	MW-230						3	
	1527	MW-108						3	
	1025	EB-19						2	
	1250	EB-21						2	
	1452	EB-20						2	
	1000	EB-18						2	

RELINQUISHED BY (SIGNATURE) DATE TIME RECEIVED BY (SIGNATURE) DATE TIME

RELINQUISHED BY (SIGNATURE) DATE TIME RECEIVED BY (SIGNATURE) DATE TIME

RELINQUISHED BY (SIGNATURE) DATE TIME RECEIVED BY (SIGNATURE) DATE TIME

RELINQUISHED BY (SIGNATURE) DATE TIME RECEIVED BY (SIGNATURE) DATE TIME

RECEIVED FOR LABORATORY BY (SIGNATURE) DATE TIME CUSTODY SEAL NO. TAMPA LIC. NO. COOLER TEMP UPON RECEIPT LABORATORY REMARKS

820-820-820
PRESERVATIVE

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Serial Number **00810**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 5	5 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	<input checked="" type="checkbox"/> AQUEOUS (WATER) <input type="checkbox"/> SOLID OR SEMISOLID <input type="checkbox"/> AIR <input type="checkbox"/> NONAQUEOUS LIQUID (OIL, SOLVENT,...)	MS 098-098 8260-8260 HELI PRESERVATIVE	STANDARD REPORT DELIVERY <input type="checkbox"/>	DATE DUE _____ EXPEDITED REPORT DELIVERY (SUPCHARGE) <input type="checkbox"/> DATE DUE _____
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	NUMBER OF CONTAINERS SUBMITTED			
CLIENT NAME	CLIENT E-MAIL		REMARKS			
CLIENT ADDRESS	COMPANY CONTRACTING THIS WORK (if applicable)					
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE			
1/29/08	845	MWJ-101	W			
		DUP-19	W			
		MWJ-55	W			
		MWJ-150	W			
		MWJ-147	W			
		MWJ-148	W			
		TB-19	W			
		TB-18	W			
		TB-20	W			
		TB-21	W			
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)
EMPTY CONTAINERS				1/30/08	846	
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)
EMPTY CONTAINERS				1/30/08	846	
RECEIVED FOR LABORATORY BY (SIGNATURE)	DATE	TIME	CUSTODY INTACT YES <input type="radio"/> NO <input type="radio"/>	LABORATORY SEAL NO.	TAMPA LOG NO.	COOLER TEMP. (°C) RECEIPT
					8857	
LABORATORY USE ONLY						
LABORATORY BENMARKS						

SDX 660 - 20926

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Serial Number

00816

Alternate Laboratory Name/Location

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Fax:

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE	STANDARD REPORT DELIVERY	DATE DUE	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT	
Talbot St	BDD38055	FL	Soil	ITS 00928-0928	1	5	OF				
SAMPLER'S SIGNATURE <i>Sue Skimal</i>	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE								
CLIENT (SITE) PM <i>Guy Kaminski</i>	CLIENT PHONE <i>813-915-1700</i>	CLIENT FAX	AQUEOUS (WATER)								
CLIENT NAME <i>Accadis</i>	CLIENT E-MAIL <i>guy.kaminski@accadis-us.com</i>		SOLID OR SEMISOLID								
CLIENT ADDRESS <i>3350 Buschway, Real Drive, Suite 100, Tampa, FL 33608</i>			AIR								
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)								
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	AQUEOUS (WATER)	SOLID OR SEMISOLID	AIR	NONAQUEOUS LIQUID (OIL, SOLVENT...)	NUMBER OF CONTAINERS SUBMITTED	REMARKS		
1/31/08	1428	MW-249						3			
	1410	MW-250						3			
	1540	MW-214						3			
	1450	MW-215						3			
	1525	MW-216						3			
	1025	MW-217						3			
	1100	MW-231						3			
	1150	MW-235						3			
	1108	MW-236						3			
	1100	MW-237						3			
	1110	MW-238						3			
	1546	MW-213						3			
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
<i>Sue Skimal</i>	2/1/08	846	<i>Guy Kaminski</i>	2/1/08	840						
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME

RECEIVED FOR LABORATORY BY: *MARLE* DATE: 2/1/08 TIME: 1000

LABORATORY USE ONLY

CUSTODY SEAL NO. 20926

COOLER TEMP UPON RECEIPT: 4.4

LABORATORY REMARKS:

Serial Number

00817

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 2 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	AQUEOUS (WATER)	STANDARD REPORT DELIVERY	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	SOLID OR SEMISOLID	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE
CLIENT NAME	CLIENT EMAIL		AIR		
CLIENT ADDRESS			NONAQUEOUS LIQUID (OIL, SOLVENT...)		
COMPANY CONTRACTING THIS WORK (if applicable)					
DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/21/08	10:25	MW-102	<input checked="" type="checkbox"/>	3	
	9:20	MW-92		3	
	1:10	MW-96		3	
	9:55	MW-103		3	
	9:20	MW-131		3	
	1:10	P2-LSAS-1		3	
	1:15	P2-LSAS-2		3	
	1:20	P2-LSAS-3		3	
	1:20	P2-LSAS-4		3	
	1:25	P2-LSAS-5		3	
	1:25	P2-LSAS-6		3	
	1:25	P2-LSAS-7		3	
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
<i>[Signature]</i>	2/1/08	8:40	<i>[Signature]</i>	2/1/08	8:40
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	YES <input type="radio"/> NO <input type="radio"/>	LABORATORY SEAL NO.
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	LABORATORY USE ONLY	TAMPA LOG INO.	COOLER TEMP UPON RECEIPT
					LABORATORY REMARKS

TAL 8240-660 (1207)

Serial Number

00818

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 3 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER 1	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE	STANDARD REPORT DELIVERY	DATE DUE
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID		
CLIENT ADDRESS			AIR		
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)		
DATE	TIME	SAMPLE IDENTIFICATION		NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/31/08		DUP-30	<input checked="" type="checkbox"/>	3	
		EB-30		3	
		FB-30		3	
		DUP-27		3	
		EB-27		3	
		FB-27		3	
		MW-40		3	
		MW-41		3	
		MW-33		3	
		EW-102		3	
		EW-107		3	
		EW-108		3	
RELINQUISHED BY: (SIGNATURE)	DATE	TIME	RELINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	LABORATORY SEAL NO.	TAMPA LOG NO.
			YES <input type="radio"/> NO <input type="radio"/>		
LABORATORY USE ONLY					
COOLER TEMP UPON RECEIPT			LABORATORY REMARKS		

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Serial Number **00819**

PROJECT REFERENCE	PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE	STANDARD REPORT DELIVERY	DATE DUE	EXPEDITED REPORT DELIVERY (SURCHARGE)	DATE DUE	NUMBER OF COOLERS SUBMITTED PER SHIPMENT	REMARKS
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.	COMPOSITE (C) OR GRAB (G) INDICATE		4	<input type="radio"/>		<input type="radio"/>			
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX	AQUEOUS (WATER)			<input type="radio"/>		<input type="radio"/>			
CLIENT NAME	CLIENT E-MAIL		SOLID OR SEMISOLID			<input type="radio"/>		<input type="radio"/>			
CLIENT ADDRESS			AIR			<input type="radio"/>		<input type="radio"/>			
COMPANY CONTRACTING THIS WORK (if applicable)			NONAQUEOUS LIQUID (OIL, SOLVENT...)			<input type="radio"/>		<input type="radio"/>			
SAMPLE DATE	SAMPLE TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS	RECEIVED BY (SIGNATURE)	DATE	TIME	RELINQUISHED BY (SIGNATURE)	DATE	TIME
1/31/08	1030	MW-10	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	840	[Signature]		
	1120	MW-36	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	842	[Signature]		
	1038	MW-37	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	840	[Signature]		
	1221	MW-127	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	840	[Signature]		
	1220	MW-128	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	840	[Signature]		
		DUP-28	<input checked="" type="checkbox"/>	2		[Signature]	2/1/08	840	[Signature]		
		ERB-28	<input checked="" type="checkbox"/>	2		[Signature]	2/1/08	840	[Signature]		
		FB-28	<input checked="" type="checkbox"/>	2		[Signature]	2/1/08	840	[Signature]		
		DUP-29	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	840	[Signature]		
		ERB-29	<input checked="" type="checkbox"/>	2		[Signature]	2/1/08	840	[Signature]		
		FB-29	<input checked="" type="checkbox"/>	2		[Signature]	2/1/08	840	[Signature]		
		Private well	<input checked="" type="checkbox"/>	3		[Signature]	2/1/08	840	[Signature]		
RELINQUISHED BY (SIGNATURE)	DATE	TIME	RELINQUISHED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME	RELINQUISHED BY (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS											
RECEIVED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME	RECEIVED BY (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS											

Serial Number

00820

ANALYSIS REQUEST AND CHAIN OF CUSTODY RECORD

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Alternate Laboratory Name/Location

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Fax:

PROJECT REFERENCE		PROJECT NO.	PROJECT LOCATION (STATE)	MATRIX TYPE	REQUIRED ANALYSIS	PAGE 5 OF 5
SAMPLER'S SIGNATURE	P.O. NUMBER	CONTRACT NO.		AQUEOUS (WATER)		STANDARD REPORT DELIVERY <input type="radio"/>
CLIENT (SITE) PM	CLIENT PHONE	CLIENT FAX		SOLID OR SEMISOLID		DATE DUE <input type="radio"/>
CLIENT NAME	CLIENT E-MAIL			AIR		EXPEDITED REPORT DELIVERY (SURCHARGE) <input type="radio"/>
CLIENT ADDRESS				NONAQUEOUS LIQUID (OIL, SOLVENT....)		DATE DUE
COMPANY CONTRACTING THIS WORK (if applicable)						NUMBER OF COOLERS SUBMITTED PER SHIPMENT

DATE	TIME	SAMPLE IDENTIFICATION	COMPOSITE (C) OR GRAB (G) INDICATE	NUMBER OF CONTAINERS SUBMITTED	REMARKS
1/31/08	16:13	MW-19	→	3	
		1617			

REINQUISHED BY: (SIGNATURE)	DATE	TIME	REINQUISHED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					
RECEIVED BY: (SIGNATURE)	DATE	TIME	RECEIVED BY: (SIGNATURE)	DATE	TIME
EMPTY CONTAINERS					

RECEIVED FOR LABORATORY BY: (SIGNATURE)	DATE	TIME	CUSTODY INTACT	LABORATORY/USE ONLY	COOLER TEMP. UPON RECEIPT	LABORATORY REMARKS
			YES <input type="radio"/> NO <input type="radio"/>	CUSTODY SEAL NO.	TAMPA LOG NO.	

PRESERVATIVE

TAL 8240-880 (1207)

ARCADIS

Appendix C

Validated Laboratory Analytical
Reports and Data Assessment
Reports

Lockheed Martin Corporation

**Appendix C - Data
Validation/Quality Control
Summary and Data Review**

**2008 Groundwater Monitoring
Report**

Former American Beryllium Company Site

July 7, 2008

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Attachments

- 1 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-20738
- 2 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-20784
- 3 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-20815
- 4 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-20822
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- 7 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-20926
- 8 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-20941
- 9 Data Review for Lockheed Martin - Tallevast, Tallevast, FL - SDG #660-21028

1. Data Validation/ Quality Control Summary

This report summarizes the analytical data and quality assurance/quality control (QA/QC) evaluation associated with the January/February 2008 annual sampling event at the former American Beryllium Company (ABC) facility. A summary of the groundwater data with any applicable data qualification and data verification reports are provided in the Sample Delivery Group (SDG) Data Reviews included in Attachments 1-9 of this appendix.

The samples were analyzed by TestAmerica Laboratories, of Tampa, Florida for volatile organic compounds (VOCs) and 1,4-dioxane. Analyses were performed according to United States Environmental Protection Agency (USEPA) SW-846 Method 8260B Total Ion Monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution (ID). Data review and verification were performed in accordance with *Organic USEPA National Functional Guidelines* (October 1999 and January 2005) and *USEPA Region II SOP HW-24, revision 2 (October 2006) Validating Volatile Organic Compounds by SW-846 Method 8260B* was also used to supplement the data review process.

Below is an overall summary of the sample analyses conducted during monitoring activities.

Field Sample Collection

Parameter	Samples	Field Duplicate	Equipment Blanks	Trip Blanks	Total
VOCs	295	27	27	31	380
Total	295	27	27	31	380

Laboratory QC samples prepared and analyzed with the field samples included method blanks, site-specific matrix spike (MS) pairs, and laboratory control samples (LCSs). Field QC samples included field blanks, trip blanks and field duplicates.

The sections below present a summary of the results of the data quality review completed by ARCADIS. The following subsections describe the various types of QA/QC parameter deviations identified during the data validation process. The complete sets of detailed data validation reports are included in the SDG Data Reviews included in Attachments 1-9 of this appendix.

1.1 Sample Preservation and Holding Time

Samples were preserved correctly with hydrochloric acid at the time of collection. All sample analyses were completed within holding time limits with the exception of the 24 analyses presented in the following table.

Sample Locations	Holding Time	Criteria
MW-87 (SIM ID) DL	Analysis Completed	23 Days
PZ-LSAS-2 (SIM ID) DL PZ-LSAS-4 (SIM ID) DL Dup-30 (SIM ID)		18 Days
MW-37 DL MW-37 (SIM ID) DL MW-127 DL MW-127 (SIM ID) DL		19 Days
PZ-LSAS-5 DL PZ-LSAS-5 (SIM ID) PZ-LSAS-6 (SIM ID) PZ-LSAS-7 DL PZ-LSAS-7 (SIM ID) Dup-27 (SIM ID) EB-27 (SIM ID) MW-40 (SIM ID) MW-41 (SIM ID) MW-33 DL MW-33 (SIM ID) EW-102 (SIM ID) DL EW-107 (SIM ID)		21 Days
MW-213 EW-108 (SIM ID)		22 Days
PZ-LSAS-6 DL		25 Days

Sample results reported from the diluted associated sample locations analyzed by analytical method USEPA SW-846 8260 were qualified, as specified in the table below.

Criteria	Qualification	
	Detected Analytes	Non-detect Analytes
Analysis completed less than two times holding time	J	UJ

J = Estimated value.

UJ = The analyte was analyzed for, but not detected. The reporting limit is an estimated value.

1.2 Gas Chromatograph Instrument Performance

Mass spectrometer tuning performance was acceptable. System performance and column resolution were acceptable.

The initial 1,4-dioxane (SIM ID) analysis for six samples (MW-213, PZ-LSAS-4, Dup-30, Dup-27, EB-27 and EW-107) were analyzed outside of the 12-hour tune clock. These samples were reanalyzed outside of holding time (as noted in Section 1.1) and reported. The reported results were qualified as estimated using “J” and “UJ” flags as described in Section 1.1.

The initial 8260B TIM analysis for sample location MW-213 was analyzed outside of the 12-hour tune clock. This sample was reanalyzed outside of holding time (as noted in Section 1.1) and reported. The reported results were qualified as estimated using a “UJ” flag as described in Section 1.1.

1.3 Blank Contamination

Blank samples were utilized to evaluate the potential introduction of contaminants into the field samples. Method blanks, equipment blanks and trip blanks were prepared and analyzed in conjunction with the field samples to monitor potential contamination impacts during collection, shipment and analysis. Method blanks were included in each analytical batch to ensure that potential interferences associated with the analytical system, reagents, and/or laboratory glassware were identified. Trip blanks

were prepared by the laboratory and shipped with sample bottles to the site, then returned to the laboratory in each cooler to evaluate potential impacts to samples during transport and storage. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

Target compounds were detected in the several QA blanks. Sample results associated with blank contamination greater than the BAL and/or non-detect did not result in any qualification of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
MW-192 MW-211	Toluene	Detected sample results <RL and <BAL	“U” at the PQL
MW-112 MW-195 MW-20 MW-233	Toluene	Detected sample results <RL and <BAL	“U” at the PQL
MW-120 MW-224 MW-226 MW-220 MW-181	Toluene	Detected sample results <RL and <BAL	“U” at the PQL
MW-99 MW-22 DUP-10 MW-103 MW-102 MW-4	o-Xylene	Detected sample results <RL and <BAL	“U” at the PQL
MW-238 MW-128 Dup-28	Acetone	Detected sample results <RL and <BAL	“U” at the PQL
EW-108	Acetone	Detected sample results >RL and <BAL	“U” at detected sample concentration
MW-131 EW-102	Trichloroethene	Detected sample results >RL and <BAL	“U” at detected sample concentration

Sample Locations	Analytes	Sample Result	Qualification
MW-10			

RL = reporting limit
PQL = practical quantitation limit

1.4 Analytical Instrument Calibrations

USEPA SW-846 Method 8260B and 8260C specify percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

Initial Calibration Criteria

All target compounds analyzed by USEPA SW-846 Method 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by USEPA SW-846 Method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by USEPA SW-846 Method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

Continuing Calibration Criteria

All target compounds analyzed by USEPA SW-846 Method 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by USEPA SW-846 Method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by USEPA SW-846 Method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

Calibration Assessment

Many Target Compound List (TCL) compounds associated with the continuing calibrations exhibited percent deviations (%D) that were greater than the 20% criteria resulting in the qualification of the associated sample results as estimated using a “J” or “UJ” qualifier, as appropriate. A detailed description of the individual qualification of

the data can be found in the SDG Data Reviews included in Attachments 1-9 of this appendix. Below is a summary of the continuing calibrations that exhibited %D greater than the 90% criteria resulting in the qualification of associated sample results as rejected (“R” qualifier). All samples were non-detect prior to rejection.

Sample Locations	Initial/Continuing	Compound	Criteria
MW-124 MW-126 MW-106 MW-112 MW-195 MW-194 MW-193 Dup6 MW-184 MW-183 MW-185 MW-187 MW-186 Dup9 TB-9 TB-8 EB-8 Dup-8 MW-15S MW-15D	CCV %D	Chloroethane	92.7%

1.5 System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

Sample locations associated with surrogates exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Surrogate	Recovery
MW-26	Toluene-d8	AC
	4-Bromofluorobenzene	> UL
	Dibromofluoromethane	AC
DUP21	Toluene-d8	> UL
	4-Bromofluorobenzene	> UL
	Dibromofluoromethane	> UL
EB-22	Toluene-d8	> UL
	4-Bromofluorobenzene	AC
	Dibromofluoromethane	AC
Dup-24	Toluene-d8	AC
	4-Bromofluorobenzene	> UL
	Dibromofluoromethane	AC
MW-143 MW-140 MW-142 MW-141	Toluene-d8	> UL
	4-Bromofluorobenzene	AC
	Dibromofluoromethane	AC
MW-33	Toluene-d8	> UL
	4-Bromofluorobenzene	AC
	Dibromofluoromethane	AC

Upper control limit (UL)
Acceptable (AC)

The criteria used to evaluate the surrogate recoveries are presented in the following table. In the case of a surrogate deviation, the sample results associated with the deviant fraction are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	No Action
	Detect	J
< LL but > 10%	Non-detect	J
	Detect	J
< 10%	Non-detect	R
	Detect	J
	Detect	

Lower control limit (LL)

1.6 Internal Standard Performance

Internal standard performance criteria insure that the Gas Chromatography/Mass Spectrometry (GC/MS) sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard. All internal standard areas and retention times were within established limits. Details of internal standard performance are provided in the SDG Data Reviews included in Attachments 1-9 of this appendix.

1.7 Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

LCSs/LCSDs were prepared and analyzed by the laboratory to evaluate method performance and analytical accuracy utilizing chemical standards prepared from a source different than the calibration standards in a clean matrix similar to the field samples included in an analytical batch. The LCSs/LCSDs establish control of the method within the laboratory environment from sample preparation through instrument performance without potential bias from field matrices. Percent recoveries were utilized to evaluate and qualify associated field samples based on laboratory established control limits. All LCSs/LCSDs recoveries were within laboratory control limits and acceptable relative percent difference (RPD) between the LCS/LCSD recoveries. Details of LCSs/LCSDs are provided in the SDG Data Reviews included in Attachments 1-9 of this appendix.

1.8 Matrix Spike Samples

Laboratory MS and matrix spike duplicate (MSD) analyses were performed to determine precision and accuracy of the analytical method on site-specific matrix and to demonstrate acceptable compound recovery by the laboratory at the time of sample analysis. Accuracy was evaluated based on the percent recoveries of the spiked compounds. Precision was based on the RPD between the MS and MSD recoveries. The MS/MSD results alone were not used to evaluate the precision and accuracy of data due to a variety of sampling and analytical conditions including heterogeneity of site samples, variability in constituent concentrations, various matrix effects, analytical batching and sample collection conditions. MS/MSD data were used in conjunction with other available QC information to formulate professional judgments and to apply qualification relative to the validity and usability of the analytical results. Laboratory

established control limits, as defined in the method protocols, were utilized to evaluate MS/MSD performance. Data qualification was based on professional judgment and overall compliance with control limits.

Two site-specific sample locations (MW-182 located in SDG 660-20738 and MW-159 located in SDG 660-20784) were selected and evaluated for MS/MSD performance. The MS/MSD performed on site sample locations exhibited acceptable recoveries and RPD between the MS/MSD recoveries. Detailed discussion of the matrix spike evaluation is provided in the data validation reports in the SDG Data Reviews included in Attachments 1-9 of this appendix.

1.9 Field Duplicates

Site-specific precision was also monitored through the collection of field duplicates at the rate of approximately one per twenty field samples. Duplicate samples are defined as samples collected simultaneously from the location under identical conditions. Duplicate aqueous samples were collected by splitting aliquots of sample between sample bottles. Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Field Duplicate Samples			
Sample ID	Parent Sample	Sample Collection Date	Sample Delivery Group
Dup2	MW-132	1/23/2008	660-20738
Dup3	MW-189	1/23/2008	660-20738
Dup4	MW-228	1/23/2008	660-20738
Dup6	MW-153	1/24/2008	660-20784
Dup7	Private Well 84	1/24/2008	660-20784
Dup8	MW-77	1/24/2008	660-20784
Dup9	MW-195	1/24/2008	660-20784
Dup10	MW-99	1/25/2008	660-20815

Field Duplicate Samples			
Sample ID	Parent Sample	Sample Collection Date	Sample Delivery Group
Dup11	MW-23	1/25/2008	660-20815
Dup13	MW-179	1/25/2008	660-20815
Dup14	MW-21	1/28/2008	660-20822
Dup15	MW-135	1/28/2008	660-20822
Dup16	MW-133	1/28/2008	660-20822
Dup17	MW-53	1/28/2008	660-20822
Dup21	MW-248	1/28/2008	660-20822
Dup22	MW-166	1/30/2008	660-20907
Dup23	MW-58	1/30/2008	660-20907
Dup24	MW-232	1/30/2008	660-20907
Dup25	MW-61	1/30/2008	660-20907
Dup31	MW-43	2/1/2008	660-20941
Dup18	MW-247	1/29/2008	660-20857
Dup19	MW-55	1/29/2008	660-20857
Dup20	MW-175	1/29/2008	660-20857
Dup27	MW-46	1/31/2008	660-20926
Dup28	MW-128	1/31/2008	660-20926
Dup29	EW-107	1/31/2008	660-20926
Dup30	MW-237	1/31/2008	660-20926

All field duplicate RPDs for values above the quantitative reporting limits were within the 50% control limit for this sampling event with the exception of the compounds carbon disulfide and 1,4-dioxane for associated sample locations MW-248 and DUP21. The reported results for these associated sample locations were qualified as estimated using a “J” qualifier.

1.10 System Performance

Manual integration was performed on TCL compound 1,4-dioxane and internal standard 1,4-dioxane-d8 associated sample locations found in SDGs/validation reports (660-20738/8094R and 660-20784/8107R). All manual integrations performed by the laboratory were consistent from baseline to baseline of the primary quantitation ion; therefore detected sample results were not qualified. Additionally, non-detect sample

results were reviewed and there was no evidence of a detectable peak of the primary quantitation ion within the retention time window of 1,4-dioxane; therefore all non-detect results were confirmed and are usable as reported.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

1.11 Data Quality Assessment Conclusion

All constituents of concern (COC) data associated with the January/February 2008 sampling event are usable for the intended purpose. Overall, the quality control data, as defined in the USEPA SW-846 Method 8260B/8260C and laboratory performance criteria were within the guidelines specified in the method with the exception of those deviations specifically mentioned in this review. Specifically, several non-detected concentrations of chloroethene were qualified as rejected using an "R" qualifier as described in Section 1.4.

2. References

United States Environmental Protection Agency (USEPA), 2005. Organic USEPA National Functional Guidelines. January 2005.

United States Environmental Protection Agency (USEPA), 1999. Organic USEPA National Functional Guidelines. October 1999.

United States Environmental Protection Agency (USEPA), Region II, 2006. USEPA Region II SOP HW-24, revision 2 (October 2006) Validating Volatile Organic Compounds by SW-846 Method 8260B.

ARCADIS

Attachment 1

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20738

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8094R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20738 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-228	660-20738-1	Water	1/23/2008	X				
MW-154	660-20738-10	Water	1/23/2008	X				
MW-151	660-20738-11	Water	1/23/2008	X				
MW-121	660-20738-12	Water	1/23/2008	X				
MW-172	660-20738-13	Water	1/23/2008	X				
MW-173	660-20738-14	Water	1/23/2008	X				
MW-171	660-20738-15	Water	1/23/2008	X				
MW-90	660-20738-16	Water	1/23/2008	X				
MW-239	660-20738-17	Water	1/23/2008	X				
MW-49	660-20738-18	Water	1/23/2008	X				
MW-66	660-20738-19	Water	1/23/2008	X				
MW-204	660-20738-2	Water	1/23/2008	X				
MW-89	660-20738-20	Water	1/23/2008	X				
MW-17S	660-20738-21	Water	1/23/2008	X				
MW-17D	660-20738-22	Water	1/23/2008	X				
MW-79	660-20738-23	Water	1/23/2008	X				
MW-132	660-20738-24	Water	1/23/2008	X				
MW-188	660-20738-25	Water	1/23/2008	X				
MW-192	660-20738-26	Water	1/23/2008	X				
MW-191	660-20738-27	Water	1/23/2008	X				
MW-189	660-20738-28	Water	1/23/2008	X				
MW-212	660-20738-29	Water	1/23/2008	X				
MW-205	660-20738-3	Water	1/23/2008	X				
MW-208	660-20738-30	Water	1/23/2008	X				
MW-211	660-20738-31	Water	1/23/2008	X				
MW-210	660-20738-32	Water	1/23/2008	X				
MW-209	660-20738-33	Water	1/23/2008	X				
Dup2	660-20738-34	Water	1/23/2008	X				
Dup3	660-20738-35	Water	1/23/2008	X				
Dup4	660-20738-36	Water	1/23/2008	X				
Equipment-2	660-20738-37	Water	1/23/2008	X				
Equipment-3	660-20738-38	Water	1/23/2008	X				
Equipment-5	660-20738-39	Water	1/23/2008	X				
MW-240	660-20738-4	Water	1/23/2008	X				
Trip Blank 2	660-20738-40	Water	1/23/2008	X				
Trip Blank 3	660-20738-41	Water	1/23/2008	X				
Trip Blank 4	660-20738-42	Water	1/23/2008	X				
Trip Blank 5	660-20738-43	Water	1/23/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-152	660-20738-44	Water	1/23/2008	X				
MW-169	660-20738-5	Water	1/23/2008	X				
MW-182	660-20738-6	Water	1/23/2008	X				
MW-167	660-20738-7	Water	1/23/2008	X				
MW-168	660-20738-8	Water	1/23/2008	X				
MW-170	660-20738-9	Water	1/23/2008	X				

Notes:

1. Matrix spike/matrix spike duplicate (MS/MSD) analyses performed on sample location MW-182.
2. Sample location Dup2 is the field duplicate of parent sample location MW-132.
3. Sample location Dup3 is the field duplicate of parent sample location MW-189.
4. Sample location Dup4 is the field duplicate of parent sample location MW-228.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

Target compounds were detected in the associated QA blanks. Sample results associated with blank contamination that were greater than the BAL and/or non-detect did not result in any qualification of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
MW-192 MW-211	Toluene	Detected sample results <RL and <BAL	"U" at the PQL

RL = reporting limit

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of

acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
MW-170 MW-79 MW-132 MW-212 MW-208 MW-211 MW-210 MW-209 Dup2 Dup3 Dup4 Equipment-2 Equipment-3 Equipment-5 Trip Blank 2 Trip Blank 3 Trip Blank 4 Trip Blank 5 MW-152	CCV %D	Methyl tert-butyl ether	-20.3%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing	RRF <0.05	Non-detect	R
		Detect	J

Initial/Continuing	Criteria	Sample Result	Qualification
Calibration	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

The MS/MSD exhibited acceptable recoveries and RPD between the MS/MSD recoveries.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-132/Dup2	1,1-Dichloroethene	2.6	0.9 l	AC
	Trichloroethene	3.7	1.7	AC
	1,4-Dioxane (SIM ID)	2.8	2.1	AC
MW-189/Dup3	All Compounds	ND	ND	AC
MW-228/Dup4	All Compounds	ND	ND	AC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

All identified compounds met the specified criteria.

11. System Performance and Overall Assessment

Manual integration was performed on TCL compound 1,4-Dioxane and internal standard 1,4-Dioxane-d8 for associated sample locations found within this SDG. All manual integrations performed by the laboratory were consistent from baseline to baseline of the primary quantitation ion; therefore detected sample results were not qualified.

Additionally, non-detect sample results were reviewed and there is no detected peak of the primary quantitation ion within the retention time window of 1,4-Dioxane; therefore all non-detect results were confirmed and are usable as reported.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-228

Sdg Number: 660-20738

Lab Sample ID: 660-20738-1

Date Sampled: 01/23/2008 0948

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2527.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2052		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2052		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-228

Sdg Number: 660-20738

Lab Sample ID: 660-20738-1

Date Sampled: 01/23/2008 0948

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2527.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2052		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2052		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-204

Lab Sample ID: 660-20738-2

Date Sampled: 01/23/2008 1040

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-56731 Instrument ID: BVMJ GC/MS
Preparation: 5030B Lab File ID: 1JA2528.D
Dilution: 1.0 Initial Weight/Volume: 5 mL
Date Analyzed: 01/25/2008 2114 Final Weight/Volume: 5 mL
Date Prepared: 01/25/2008 2114

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-204

Lab Sample ID: 660-20738-2

Date Sampled: 01/23/2008 1040

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56731

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2528.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/25/2008 2114

Final Weight/Volume: 5 mL

Date Prepared: 01/25/2008 2114

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	99		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-205

Sdg Number: 660-20738

Lab Sample ID: 660-20738-3

Date Sampled: 01/23/2008 1044

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2529.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2136		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2136		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-205

Sdg Number: 660-20738

Lab Sample ID: 660-20738-3

Date Sampled: 01/23/2008 1044

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2529.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2136		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2136		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-240

Lab Sample ID: 660-20738-4

Date Sampled: 01/23/2008 1422

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-56731 Instrument ID: BVMJ GC/MS
Preparation: 5030B Lab File ID: 1JA2530.D
Dilution: 1.0 Initial Weight/Volume: 5 mL
Date Analyzed: 01/25/2008 2159 Final Weight/Volume: 5 mL
Date Prepared: 01/25/2008 2159

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-240

Sdg Number: 660-20738

Lab Sample ID: 660-20738-4

Date Sampled: 01/23/2008 1422

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2530.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2159		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2159		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	103	74 - 126
Dibromofluoromethane	111	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-169

Sdg Number: 660-20738

Lab Sample ID: 660-20738-5

Date Sampled: 01/23/2008 1515

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2531.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2221		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2221		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-169

Lab Sample ID: 660-20738-5

Date Sampled: 01/23/2008 1515

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2531.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2221		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2221		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	28		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	117	77 - 122
4-Bromofluorobenzene	108	74 - 126
Dibromofluoromethane	117	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-182

Lab Sample ID: 660-20738-6

Date Sampled: 01/23/2008 1515

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2532.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2243		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2243		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-182

Sdg Number: 660-20738

Lab Sample ID: 660-20738-6

Date Sampled: 01/23/2008 1515

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2532.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2243		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2243		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-167

Lab Sample ID: 660-20738-7

Date Sampled: 01/23/2008 1600

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56731

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2533.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/25/2008 2305

Final Weight/Volume: 5 mL

Date Prepared: 01/25/2008 2305

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-167

Sdg Number: 660-20738

Lab Sample ID: 660-20738-7

Date Sampled: 01/23/2008 1600

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56731

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2533.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/25/2008 2305

Final Weight/Volume: 5 mL

Date Prepared: 01/25/2008 2305

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	117	77 - 122
4-Bromofluorobenzene	107	74 - 126
Dibromofluoromethane	121	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-168

Sdg Number: 660-20738

Lab Sample ID: 660-20738-8

Date Sampled: 01/23/2008 1647

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2534.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2328		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2328		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-168

Lab Sample ID: 660-20738-8

Date Sampled: 01/23/2008 1647

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56731	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2534.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/25/2008 2328		Final Weight/Volume: 5 mL
Date Prepared:	01/25/2008 2328		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	101	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-170

Sdg Number: 660-20738

Lab Sample ID: 660-20738-9

Date Sampled: 01/23/2008 1705

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2612.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1528		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1528		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-170

Lab Sample ID: 660-20738-9

Date Sampled: 01/23/2008 1705

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2612.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1528		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1528		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	103	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-154

Sdg Number: 660-20738

Lab Sample ID: 660-20738-10

Date Sampled: 01/23/2008 1000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2547.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0416		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0416		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: **MW-154**

Sdg Number: 660-20738

Lab Sample ID: 660-20738-10

Date Sampled: 01/23/2008 1000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2547.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0416		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0416		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	103	74 - 126
Dibromofluoromethane	114	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-151

Lab Sample ID: 660-20738-11

Date Sampled: 01/23/2008 1050

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2548.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0439		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0439		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-151

Lab Sample ID: 660-20738-11

Date Sampled: 01/23/2008 1050

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2548.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0439		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0439		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	89	77 - 122
4-Bromofluorobenzene	81	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-121

Lab Sample ID: 660-20738-12

Date Sampled: 01/23/2008 1615

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2549.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0501		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0501		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-121

Sdg Number: 660-20738

Lab Sample ID: 660-20738-12

Date Sampled: 01/23/2008 1615

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56734

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2549.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0501

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0501

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	12	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	98	77 - 122
4-Bromofluorobenzene	89	74 - 126
Dibromofluoromethane	98	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-172

Sdg Number: 660-20738

Lab Sample ID: 660-20738-13

Date Sampled: 01/23/2008 1510

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2550.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0523		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0523		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-172

Sdg Number: 660-20738

Lab Sample ID: 660-20738-13

Date Sampled: 01/23/2008 1510

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2550.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0523		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0523		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	113		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-173

Sdg Number: 660-20738

Lab Sample ID: 660-20738-14

Date Sampled: 01/23/2008 1552

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2551.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0545		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0545		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-173

Sdg Number: 660-20738

Lab Sample ID: 660-20738-14

Date Sampled: 01/23/2008 1552

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2551.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0545		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0545		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	113		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-171

Sdg Number: 660-20738

Lab Sample ID: 660-20738-15

Date Sampled: 01/23/2008 1433

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2552.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0607		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0607		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-171

Sdg Number: 660-20738

Lab Sample ID: 660-20738-15

Date Sampled: 01/23/2008 1433

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2552.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0607		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0607		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	111		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: **MW-90**

Sdg Number: 660-20738

Lab Sample ID: 660-20738-16

Date Sampled: 01/23/2008 1002

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2553.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0630		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0630		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-90

Sdg Number: 660-20738

Lab Sample ID: 660-20738-16

Date Sampled: 01/23/2008 1002

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2553.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0630		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0630		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-239

Lab Sample ID: 660-20738-17

Date Sampled: 01/23/2008 1120

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2554.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0652		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0652		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	2.1		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: MW-239

Lab Sample ID: 660-20738-17
Client Matrix: Water

Date Sampled: 01/23/2008 1120
Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 01/26/2008 0652
Date Prepared: 01/26/2008 0652

Analysis Batch: 660-56734

Instrument ID: BVMJ GC/MS
Lab File ID: 1JA2554.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2.9		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate				
Toluene-d8 (Surr)	110			77 - 122
4-Bromofluorobenzene	97			74 - 126
Dibromofluoromethane	109			70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-49

Lab Sample ID: 660-20738-18

Date Sampled: 01/23/2008 1141

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2555.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0714		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0714		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-49

Lab Sample ID: 660-20738-18

Date Sampled: 01/23/2008 1141

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2555.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0714		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0714		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	111		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-66

Sdg Number: 660-20738

Lab Sample ID: 660-20738-19

Date Sampled: 01/23/2008 1415

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56734

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2556.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0736

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0736

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-66

Sdg Number: 660-20738

Lab Sample ID: 660-20738-19

Date Sampled: 01/23/2008 1415

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2556.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0736		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0736		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	120		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	116		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-89

Lab Sample ID: 660-20738-20

Date Sampled: 01/23/2008 1417

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-56734 Instrument ID: BVMJ GC/MS
Preparation: 5030B Lab File ID: 1JA2557.D
Dilution: 1.0 Initial Weight/Volume: 5 mL
Date Analyzed: 01/26/2008 0758 Final Weight/Volume: 5 mL
Date Prepared: 01/26/2008 0758

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-89

Sdg Number: 660-20738

Lab Sample ID: 660-20738-20

Date Sampled: 01/23/2008 1417

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2557.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0758		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0758		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec			Acceptance Limits
Toluene-d8 (Surr)	108			77 - 122
4-Bromofluorobenzene	97			74 - 126
Dibromofluoromethane	107			70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-17S

Sdg Number: 660-20738

Lab Sample ID: 660-20738-21

Date Sampled: 01/23/2008 1454

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2558.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0821		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0821		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-17S

Sdg Number: 660-20738

Lab Sample ID: 660-20738-21

Date Sampled: 01/23/2008 1454

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56734

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2558.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0821

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0821

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-17D

Lab Sample ID: 660-20738-22

Date Sampled: 01/23/2008 1455

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2559.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0843		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0843		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-17D

Sdg Number: 660-20738

Lab Sample ID: 660-20738-22

Date Sampled: 01/23/2008 1455

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2559.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0843		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0843		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec	Acceptance Limits		
Toluene-d8 (Surr)	114	77 - 122		
4-Bromofluorobenzene	99	74 - 126		
Dibromofluoromethane	113	70 - 130		

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-79

Sdg Number: 660-20738

Lab Sample ID: 660-20738-23

Date Sampled: 01/23/2008 1534

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2613.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1550		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 1550		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	4.0		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	14		0.45	1.0
cis-1,2-Dichloroethene	140		0.65	1.0
trans-1,2-Dichloroethene	3.8		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-79

Sdg Number: 660-20738

Lab Sample ID: 660-20738-23

Date Sampled: 01/23/2008 1534

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 01/26/2008 1550
Date Prepared: 01/26/2008 1550

Analysis Batch: 660-56869

Instrument ID: BVMJ GC/MS
Lab File ID: 1JA2613.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	100		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	94		77 - 122	
4-Bromofluorobenzene	87		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-132

Sdg Number: 660-20738

Lab Sample ID: 660-20738-24

Date Sampled: 01/23/2008 1540

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2614.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1613		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 1613		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	2.6	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-132

Sdg Number: 660-20738

Lab Sample ID: 660-20738-24

Date Sampled: 01/23/2008 1540

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2614.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1613		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 1613		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	3.7		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	101	74 - 126
Dibromofluoromethane	115	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-188

Sdg Number: 660-20738

Lab Sample ID: 660-20738-25

Date Sampled: 01/23/2008 1100

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2562.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0949		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0949		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-188

Sdg Number: 660-20738

Lab Sample ID: 660-20738-25

Date Sampled: 01/23/2008 1100

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2562.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 0949		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 0949		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-192

Sdg Number: 660-20738

Lab Sample ID: 660-20738-26

Date Sampled: 01/23/2008 1205

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56874	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JA2929.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/30/2008 0052		Final Weight/Volume: 5 mL
Date Prepared:	01/30/2008 0052		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-192

Lab Sample ID: 660-20738-26

Date Sampled: 01/23/2008 1205

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B	Analysis Batch: 660-56874	Instrument ID: BVMJ GC/MS	
Preparation: 5030B		Lab File ID: 2JA2929.D	
Dilution: 1.0		Initial Weight/Volume: 5 mL	
Date Analyzed: 01/30/2008 0052		Final Weight/Volume: 5 mL	
Date Prepared: 01/30/2008 0052			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.81 1.0	U + U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	12	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	43		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-191

Sdg Number: 660-20738

Lab Sample ID: 660-20738-27

Date Sampled: 01/23/2008 1150

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56734

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2564.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 1034

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 1034

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-191

Sdg Number: 660-20738

Lab Sample ID: 660-20738-27

Date Sampled: 01/23/2008 1150

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2564.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1034		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 1034		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.89	I	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-189

Sdg Number: 660-20738

Lab Sample ID: 660-20738-28

Date Sampled: 01/23/2008 1055

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2565.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1056		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 1056		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-189

Lab Sample ID: 660-20738-28

Date Sampled: 01/23/2008 1055

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56734	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2565.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1056		Final Weight/Volume: 5 mL
Date Prepared:	01/26/2008 1056		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	120	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-212

Lab Sample ID: 660-20738-29

Date Sampled: 01/23/2008 1520

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56869

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2615.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 1635

Final Weight/Volume: mL

Date Prepared: 01/26/2008 1635

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-212

Lab Sample ID: 660-20738-29

Date Sampled: 01/23/2008 1520

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2615.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1635		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1635		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	2.0		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-208

Lab Sample ID: 660-20738-30

Date Sampled: 01/23/2008 1440

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56874	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2930.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0114		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0114			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	11		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-208

Sdg Number: 660-20738

Lab Sample ID: 660-20738-30

Date Sampled: 01/23/2008 1440

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56874	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JA2930.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/30/2008 0114		Final Weight/Volume: 5 mL
Date Prepared:	01/30/2008 0114		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	109		74 - 126	
Dibromofluoromethane	115		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-211

Sdg Number: 660-20738

Lab Sample ID: 660-20738-31

Date Sampled: 01/23/2008 1550

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2617.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1719		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1719		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.64	I	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.64	I	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	1.1	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-211

Sdg Number: 660-20738

Lab Sample ID: 660-20738-31

Date Sampled: 01/23/2008 1550

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2617.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1719		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1719		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.93 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.53	I	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	16		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	113		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	114		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-210

Lab Sample ID: 660-20738-32

Date Sampled: 01/23/2008 1610

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-56869 Instrument ID: BVMJ GC/MS
Preparation: 5030B Lab File ID: 1JA2618.D
Dilution: 1.0 Initial Weight/Volume: 5 mL
Date Analyzed: 01/26/2008 1742 Final Weight/Volume: mL
Date Prepared: 01/26/2008 1742

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-210

Lab Sample ID: 660-20738-32

Date Sampled: 01/23/2008 1610

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2618.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1742		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1742		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-209

Lab Sample ID: 660-20738-33

Date Sampled: 01/23/2008 1445

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2619.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1804		Final Weight/Volume:	mL
Date Prepared:	01/26/2008 1804			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-209

Sdg Number: 660-20738

Lab Sample ID: 660-20738-33

Date Sampled: 01/23/2008 1445

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2619.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1804		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1804		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Dup2

Sdg Number: 660-20738

Lab Sample ID: 660-20738-34

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2620.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1826		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1826		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.90	I	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Dup2

Sdg Number: 660-20738

Lab Sample ID: 660-20738-34

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2620.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1826		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1826		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1.7		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	116		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Dup3

Sdg Number: 660-20738

Lab Sample ID: 660-20738-35

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2621.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1849		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1849		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: Dup3

Lab Sample ID: 660-20738-35

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56869

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2621.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 1849

Final Weight/Volume: mL

Date Prepared: 01/26/2008 1849

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	115		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Dup4

Sdg Number: 660-20738

Lab Sample ID: 660-20738-36

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2622.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1911		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1911		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Dup4

Sdg Number: 660-20738

Lab Sample ID: 660-20738-36

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation: 5030B		Lab File ID: 1JA2622.D
Dilution: 1.0		Initial Weight/Volume: 5 mL
Date Analyzed: 01/26/2008 1911		Final Weight/Volume: mL
Date Prepared: 01/26/2008 1911		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec	Acceptance Limits		
Toluene-d8 (Surr)	116	77 - 122		
4-Bromofluorobenzene	103	74 - 126		
Dibromofluoromethane	120	70 - 130		

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: Equipment-2

Lab Sample ID: 660-20738-37

Date Sampled: 01/23/2008 1600

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56869

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2623.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 1933

Final Weight/Volume: mL

Date Prepared: 01/26/2008 1933

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: **Equipment-2**

Sdg Number: 660-20738

Lab Sample ID: 660-20738-37
Client Matrix: Water

Date Sampled: 01/23/2008 1600
Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation: 5030B		Lab File ID: 1JA2623.D
Dilution: 1.0		Initial Weight/Volume: 5 mL
Date Analyzed: 01/26/2008 1933		Final Weight/Volume: mL
Date Prepared: 01/26/2008 1933		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.87	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Equipment-3

Sdg Number: 660-20738

Lab Sample ID: 660-20738-38

Date Sampled: 01/23/2008 1630

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2624.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1955		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1955		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: Equipment-3

Lab Sample ID: 660-20738-38
Client Matrix: Water

Date Sampled: 01/23/2008 1630
Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2624.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 1955		Final Weight/Volume: mL
Date Prepared:	01/26/2008 1955		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.99	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	121	77 - 122
4-Bromofluorobenzene	106	74 - 126
Dibromofluoromethane	119	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Equipment-5

Sdg Number: 660-20738

Lab Sample ID: 660-20738-39

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2625.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2018		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2018		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: Equipment-5

Lab Sample ID: 660-20738-39
Client Matrix: Water

Date Sampled: 01/23/2008 0000
Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2625.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2018		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2018		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.98	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	84		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: Trip Blank 2

Lab Sample ID: 660-20738-40

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56869

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2626.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 2040

Final Weight/Volume: mL

Date Prepared: 01/26/2008 2040

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: Trip Blank 2

Lab Sample ID: 660-20738-40
Client Matrix: Water

Date Sampled: 01/23/2008 0000
Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2626.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2040		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2040		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	118		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	117		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 3

Sdg Number: 660-20738

Lab Sample ID: 660-20738-41

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2627.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2102		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2102		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: Trip Blank 3

Lab Sample ID: 660-20738-41

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2627.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2102		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2102		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	110	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 4

Sdg Number: 660-20738

Lab Sample ID: 660-20738-42

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2628.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2125		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2125		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 4

Sdg Number: 660-20738

Lab Sample ID: 660-20738-42

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2628.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2125		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2125		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	122	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	117	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 5

Sdg Number: 660-20738

Lab Sample ID: 660-20738-43

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2629.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2147		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2147		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 5

Sdg Number: 660-20738

Lab Sample ID: 660-20738-43

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2629.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2147		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2147		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: MW-152

Lab Sample ID: 660-20738-44
Client Matrix: Water

Date Sampled: 01/23/2008 1044
Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2630.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2209		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2209		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-152

Lab Sample ID: 660-20738-44

Date Sampled: 01/23/2008 1044

Client Matrix: Water

Date Received: 01/24/2008 0945

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56869	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2630.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/26/2008 2209		Final Weight/Volume: mL
Date Prepared:	01/26/2008 2209		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-228

Sdg Number: 660-20738

Lab Sample ID: 660-20738-1

Date Sampled: 01/23/2008 0948

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2527.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2052		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2052			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-204

Sdg Number: 660-20738

Lab Sample ID: 660-20738-2

Date Sampled: 01/23/2008 1040

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2528.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2114		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2114			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-205

Sdg Number: 660-20738

Lab Sample ID: 660-20738-3

Date Sampled: 01/23/2008 1044

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2529.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2136		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2136			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-240

Lab Sample ID: 660-20738-4

Date Sampled: 01/23/2008 1422

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2530.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2159		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2159			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-169

Sdg Number: 660-20738

Lab Sample ID: 660-20738-5

Date Sampled: 01/23/2008 1515

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2531.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2221		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2221			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-182

Lab Sample ID: 660-20738-6

Date Sampled: 01/23/2008 1515

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2532.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2243		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2243			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-167

Sdg Number: 660-20738

Lab Sample ID: 660-20738-7

Date Sampled: 01/23/2008 1600

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2533.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2305		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2305			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-168

Sdg Number: 660-20738

Lab Sample ID: 660-20738-8

Date Sampled: 01/23/2008 1647

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2534.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/25/2008 2328		Final Weight/Volume:	5 mL
Date Prepared:	01/25/2008 2328			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-170

Sdg Number: 660-20738

Lab Sample ID: 660-20738-9

Date Sampled: 01/23/2008 1705

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 01/26/2008 1528
Date Prepared: 01/26/2008 1528

Analysis Batch: 660-56870

Instrument ID: BVMJ GC/MS

Lab File ID: 1JA2612.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-154

Sdg Number: 660-20738

Lab Sample ID: 660-20738-10

Date Sampled: 01/23/2008 1000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2547.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0416		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0416			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-151

Lab Sample ID: 660-20738-11

Date Sampled: 01/23/2008 1050

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2548.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0439

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0439

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.8		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-121

Sdg Number: 660-20738

Lab Sample ID: 660-20738-12

Date Sampled: 01/23/2008 1615

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2549.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0501		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0501			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-172

Lab Sample ID: 660-20738-13

Date Sampled: 01/23/2008 1510

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2550.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0523

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0523

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-173

Sdg Number: 660-20738

Lab Sample ID: 660-20738-14

Date Sampled: 01/23/2008 1552

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2551.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0545		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0545			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-171

Sdg Number: 660-20738

Lab Sample ID: 660-20738-15

Date Sampled: 01/23/2008 1433

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2552.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0607		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0607			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-90

Lab Sample ID: 660-20738-16

Date Sampled: 01/23/2008 1002

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2553.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0630

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0630

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-239

Lab Sample ID: 660-20738-17

Date Sampled: 01/23/2008 1120

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2554.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0652

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0652

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	6.7		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-49

Lab Sample ID: 660-20738-18

Date Sampled: 01/23/2008 1141

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2555.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0714		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0714			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-66

Sdg Number: 660-20738

Lab Sample ID: 660-20738-19

Date Sampled: 01/23/2008 1415

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2556.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0736		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0736			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-89

Lab Sample ID: 660-20738-20

Date Sampled: 01/23/2008 1417

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2557.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0758

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0758

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.9		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-17S

Sdg Number: 660-20738

Lab Sample ID: 660-20738-21

Date Sampled: 01/23/2008 1454

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2558.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/26/2008 0821

Final Weight/Volume: 5 mL

Date Prepared: 01/26/2008 0821

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: MW-17D

Lab Sample ID: 660-20738-22
Client Matrix: Water

Date Sampled: 01/23/2008 1455
Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2559.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0843		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0843			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-79

Sdg Number: 660-20738

Lab Sample ID: 660-20738-23

Date Sampled: 01/23/2008 1534

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2613.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1550		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1550			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	52		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-132

Lab Sample ID: 660-20738-24

Date Sampled: 01/23/2008 1540

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2561.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0927		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0927			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.8		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: MW-188

Lab Sample ID: 660-20738-25
Client Matrix: Water

Date Sampled: 01/23/2008 1100
Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2562.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 0949		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 0949			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-192

Lab Sample ID: 660-20738-26

Date Sampled: 01/23/2008 1205

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2563.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1011		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1011			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-191

Sdg Number: 660-20738

Lab Sample ID: 660-20738-27

Date Sampled: 01/23/2008 1150

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2564.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1034		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1034			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: MW-189

Lab Sample ID: 660-20738-28
Client Matrix: Water

Date Sampled: 01/23/2008 1055
Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2565.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1056		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1056			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-212

Lab Sample ID: 660-20738-29

Date Sampled: 01/23/2008 1520

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2615.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1635		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1635			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-208

Sdg Number: 660-20738

Lab Sample ID: 660-20738-30

Date Sampled: 01/23/2008 1440

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2616.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1657		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1657			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-211

Sdg Number: 660-20738

Lab Sample ID: 660-20738-31

Date Sampled: 01/23/2008 1550

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2617.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1719		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1719			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-210

Lab Sample ID: 660-20738-32

Date Sampled: 01/23/2008 1610

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3055.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1230

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1230

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: MW-209

Sdg Number: 660-20738

Lab Sample ID: 660-20738-33

Date Sampled: 01/23/2008 1445

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57011	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3056.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1252		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1252			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Dup2

Sdg Number: 660-20738

Lab Sample ID: 660-20738-34

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2620.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1826		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1826			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.1		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: Dup3

Lab Sample ID: 660-20738-35

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2621.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1849		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1849			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: Dup4

Lab Sample ID: 660-20738-36

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57011	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3057.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1314		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1314			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Equipment-2

Sdg Number: 660-20738

Lab Sample ID: 660-20738-37

Date Sampled: 01/23/2008 1600

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2623.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 1933		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 1933			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Equipment-3

Sdg Number: 660-20738

Lab Sample ID: 660-20738-38

Date Sampled: 01/23/2008 1630

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57011	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3058.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1336		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1336			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Equipment-5

Sdg Number: 660-20738

Lab Sample ID: 660-20738-39

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2625.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 2018		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 2018			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 2

Sdg Number: 660-20738

Lab Sample ID: 660-20738-40

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2626.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 2040		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 2040			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 3

Sdg Number: 660-20738

Lab Sample ID: 660-20738-41

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2627.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 2102		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 2102			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Client Sample ID: Trip Blank 4

Sdg Number: 660-20738

Lab Sample ID: 660-20738-42

Date Sampled: 01/23/2008 0000

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2628.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 2125		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 2125			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1
Sdg Number: 660-20738

Client Sample ID: Trip Blank 5

Lab Sample ID: 660-20738-43
Client Matrix: Water

Date Sampled: 01/23/2008 0000
Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2629.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 2147		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 2147			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20738-1

Sdg Number: 660-20738

Client Sample ID: MW-152

Lab Sample ID: 660-20738-44

Date Sampled: 01/23/2008 1044

Client Matrix: Water

Date Received: 01/24/2008 0945

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56870	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2630.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/26/2008 2209		Final Weight/Volume:	5 mL
Date Prepared:	01/26/2008 2209			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20738	1/23/2008	SW-846	MW-228	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-154	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-151	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-121	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-172	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-173	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-171	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-90	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-239	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-49	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-66	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-204	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-89	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-17S	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-17D	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-79	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-132	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-188	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-192	Water	No	--	--	--	--	VOC – Associated Blank
660-20738	1/23/2008	SW-846	MW-191	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-189	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-212	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-205	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-208	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-211	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20738	1/23/2008	SW-846	MW-210	Water	No	--	--	--	--	VOC – CCAL %D

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20738	1/23/2008	SW-846	MW-209	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Dup2	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Dup3	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Dup4	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Equipment-2	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Equipment-3	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Equipment-5	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-240	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	Trip Blank 2	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Trip Blank 3	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Trip Blank 4	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	Trip Blank 5	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-152	Water	No	--	--	--	--	VOC – CCAL %D
660-20738	1/23/2008	SW-846	MW-169	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-182	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-167	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-168	Water	Yes	--	--	--	--	
660-20738	1/23/2008	SW-846	MW-170	Water	No	--	--	--	--	VOC – CCAL %D

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 2

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20784

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8107R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20784 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-94	660-20784-1	Water	1/24/2008	X				
MW-32	660-20784-10	Water	1/24/2008	X				
MW-7S	660-20784-11	Water	1/24/2008	X				
Dup7	660-20784-12	Water	1/24/2008	X				
MW-7D	660-20784-13	Water	1/24/2008	X				
MW-159	660-20784-14	Water	1/24/2008	X				
MW-157	660-20784-15	Water	1/24/2008	X				
MW-160	660-20784-16	Water	1/24/2008	X				
MW-156	660-20784-17	Water	1/24/2008	X				
MW-11	660-20784-18	Water	1/24/2008	X				
EB-7	660-20784-19	Water	1/24/2008	X				
MW-95	660-20784-2	Water	1/24/2008	X				
TB-7	660-20784-20	Water	1/24/2008	X				
TB-6	660-20784-21	Water	1/24/2008	X				
EB-6	660-20784-22	Water	1/24/2008	X				
MW-107	660-20784-23	Water	1/24/2008	X				
MW-153	660-20784-24	Water	1/24/2008	X				
MW-155	660-20784-25	Water	1/24/2008	X				
MW-161	660-20784-26	Water	1/24/2008	X				
MW-158	660-20784-27	Water	1/24/2008	X				
EB-9	660-20784-28	Water	1/24/2008	X				
MW-124	660-20784-29	Water	1/24/2008	X				
TB-1	660-20784-3	Water	1/24/2008	X				
MW-126	660-20784-30	Water	1/24/2008	X				
MW-106	660-20784-31	Water	1/24/2008	X				
MW-112	660-20784-32	Water	1/24/2008	X				
MW-195	660-20784-33	Water	1/24/2008	X				
MW-194	660-20784-34	Water	1/24/2008	X				
MW-193	660-20784-35	Water	1/24/2008	X				
Dup6	660-20784-36	Water	1/24/2008	X				
MW-184	660-20784-37	Water	1/24/2008	X				
MW-183	660-20784-38	Water	1/24/2008	X				
MW-185	660-20784-39	Water	1/24/2008	X				
RW-2	660-20784-4	Water	1/24/2008	X				
MW-187	660-20784-40	Water	1/24/2008	X				
MW-186	660-20784-41	Water	1/24/2008	X				
Dup9	660-20784-42	Water	1/24/2008	X				
TB-9	660-20784-43	Water	1/24/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
TB-8	660-20784-44	Water	1/24/2008	X				
EB-8	660-20784-45	Water	1/24/2008	X				
Dup-8	660-20784-46	Water	1/24/2008	X				
MW-15S	660-20784-47	Water	1/24/2008	X				
MW-15D	660-20784-48	Water	1/24/2008	X				
MW-14S	660-20784-49	Water	1/24/2008	X				
MW-12	660-20784-5	Water	1/24/2008	X				
MW-14D	660-20784-50	Water	1/24/2008	X				
MW-77	660-20784-51	Water	1/24/2008	X				
MW-20	660-20784-52	Water	1/24/2008	X				
MW-9S	660-20784-53	Water	1/24/2008	X				
MW-9D	660-20784-54	Water	1/24/2008	X				
MW-76	660-20784-55	Water	1/24/2008	X				
MW-25	660-20784-56	Water	1/24/2008	X				
MW-85	660-20784-57	Water	1/24/2008	X				
MW-233	660-20784-58	Water	1/24/2008	X				
MW-5	660-20784-6	Water	1/24/2008	X				
MW-114	660-20784-7	Water	1/24/2008	X				
MW-83	660-20784-8	Water	1/24/2008	X				
Private Well 84	660-20784-9	Water	1/24/2008	X				

Notes:

1. Matrix spike/matrix spike duplicate (MS/MSD) analyses performed on sample location MW-159.
2. Sample location Dup6 is the field duplicate of parent sample location MW-153.
3. Sample location Dup7 is the field duplicate of parent sample location Private Well 84.
4. Sample location Dup8 is the field duplicate of parent sample location MW-77.
5. Sample location Dup9 is the field duplicate of parent sample location MW-195.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

Target compounds were detected in the associated QA blanks. Sample results associated with blank contamination that were greater than the BAL and/or non-detect did not result in any qualification of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
MW-112 MW-195 MW-20 MW-233	Toluene	Detected sample results <RL and <BAL	"U" at the PQL

RL = reporting limit

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
MW-124 MW-126 MW-106 MW-112 MW-195 MW-194 MW-193 Dup6 MW-184 MW-183 MW-185 MW-187 MW-186 Dup9 TB-9 TB-8 EB-8 Dup-8 MW-15S MW-15D	CCV %D	Chloroethane	92.7%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

1. RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

The MS/MSD exhibited acceptable recoveries and RPD between the MS/MSD recoveries.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-77/Dup8	1,1-Dichloroethane	39	45	14.2 %
	1,1-Dichloroethene	18	20	10.5 %
	cis-1,2-Dichloroethene	3.5	3.8	AC
	Tetrachloroethene	1.3	1.4	AC
	Trichloroethene	24	26	8.0 %
	1,4-Dioxane (SIM ID)	150	220	37.8 %
MW-153/Dup6	All Compounds	ND	ND	AC
Private Well 84/Dup7	All Compounds	ND	ND	AC
MW-195/Dup9	All Compounds	ND	ND	AC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
MW-12	Trichloroethene	180 E	190 D	190 D

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
MW-9D	Acetone	760 E	580 D	580 D

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

11. System Performance and Overall Assessment

Manual integration was performed on TCL compound 1,4-Dioxane and internal standard 1,4-Dioxane-d8 for associated sample locations found within this SDG. All manual integrations performed by the laboratory were consistent from baseline to baseline of the primary quantitation ion; therefore detected sample results were not qualified.

Additionally, non-detect sample results were reviewed and there is no detected peak of the primary quantitation ion within the retention time window of 1,4-Dioxane; therefore all non-detect results were confirmed and are usable as reported.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-94

Sdg Number: 660-20784

Lab Sample ID: 660-20784-1

Date Sampled: 01/24/2008 1630

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 01/27/2008 0747
 Date Prepared: 01/27/2008 0747

Analysis Batch: 660-56879

Instrument ID: BVMJ GC/MS
 Lab File ID: 1JA2656.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

TestAmerica Tampa

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-94

Sdg Number: 660-20784

Lab Sample ID: 660-20784-1

Date Sampled: 01/24/2008 1630

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56879

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2656.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 0747

Final Weight/Volume: 5 mL

Date Prepared: 01/27/2008 0747

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	83		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-95

Sdg Number: 660-20784

Lab Sample ID: 660-20784-2

Client Matrix: Water

Date Sampled: 01/24/2008 1547

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56879

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2657.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 0809

Date Prepared: 01/27/2008 0809

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-95

Sdg Number: 660-20784

Lab Sample ID: 660-20784-2

Date Sampled: 01/24/2008 1547

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56879

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2657.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 0809

Final Weight/Volume: 5 mL

Date Prepared: 01/27/2008 0809

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	89		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-1

Sdg Number: 660-20784

Lab Sample ID: 660-20784-3

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-56879 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JA2658.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 01/27/2008 0832 Final Weight/Volume: 5 mL
 Date Prepared: 01/27/2008 0832

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	1.0
n-Butylbenzene	0.67	U	0.67	5.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-1

Sdg Number: 660-20784

Lab Sample ID: 660-20784-3

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA2658.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0832			Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0832				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	116	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	118	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: RW-2

Sdg Number: 660-20784

Lab Sample ID: 660-20784-4

Date Sampled: 01/24/2008 0930

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2659.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0854		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0854			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	4.1	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.56	I	0.45	1.0
cis-1,2-Dichloroethene	15	U	0.65	1.0
trans-1,2-Dichloroethene	0.46	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: RW-2

Sdg Number: 660-20784

Lab Sample ID: 660-20784-4

Date Sampled: 01/24/2008 0930

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56879

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2659.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 0854

Final Weight/Volume: 5 mL

Date Prepared: 01/27/2008 0854

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	21		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	9.7		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	160		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	90	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-12

Sdg Number: 660-20784

Lab Sample ID: 660-20784-5

Date Sampled: 01/24/2008 1035

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2660.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0916		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0916			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	1.0
n-Butylbenzene	0.67	U	0.67	5.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	1.0
Chloroform	0.90	U	0.90	5.0
Chloromethane	1.0	U	1.0	1.0
2-Chlorotoluene	0.65	U	1.0	4.0
4-Chlorotoluene	0.52	U	0.65	1.0
Chlorodibromomethane	0.34	U	0.52	1.0
1,2-Dibromo-3-Chloropropane	0.34	U	0.34	1.0
Ethylene Dibromide	2.5	U	2.5	5.0
Dibromomethane	0.50	U	0.50	1.0
1,2-Dichlorobenzene	0.41	U	0.41	1.0
1,3-Dichlorobenzene	0.44	U	0.44	1.0
1,4-Dichlorobenzene	0.64	U	0.64	1.0
Dichlorodifluoromethane	0.52	U	0.52	1.0
1,1-Dichloroethane	2.5	U	2.5	5.0
1,2-Dichloroethane	10		0.52	1.0
1,1-Dichloroethene	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	6.9		0.45	1.0
trans-1,2-Dichloroethene	5.1		0.65	1.0
1,2-Dichloropropane	0.44	U	0.44	1.0
1,3-Dichloropropane	0.52	U	0.52	1.0
2,2-Dichloropropane	0.39	U	0.39	1.0
1,1-Dichloropropene	0.36	U	0.36	1.0
cis-1,3-Dichloropropene	0.31	U	0.31	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.14	U	0.14	1.0
Hexachlorobutadiene	0.44	U	0.44	1.0
Isopropylbenzene	0.40	U	0.40	1.0
4-Isopropyltoluene	0.19	U	0.19	1.0
Methylene Chloride	0.69	U	0.69	1.0
Naphthalene	4.0	U	4.0	5.0
N-Propylbenzene	2.5	U	2.5	5.0
Styrene	0.59	U	0.59	1.0
1,1,1,2-Tetrachloroethane	0.98	U	0.98	2.0
	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-12

Sdg Number: 660-20784

Lab Sample ID: 660-20784-5

Date Sampled: 01/24/2008 1035

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2660.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0916		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0916			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	58		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.7		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	97		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	108		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-12

Sdg Number: 660-20784

Lab Sample ID: 660-20784-5

Date Sampled: 01/24/2008 1035

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3059.D

Dilution: 2.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1358

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1358

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	190	D	1.0	2.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-5

Sdg Number: 660-20784

Lab Sample ID: 660-20784-6

Date Sampled: 01/24/2008 1130

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA2661.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0938			Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0938				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-5

Sdg Number: 660-20784

Lab Sample ID: 660-20784-6

Date Sampled: 01/24/2008 1130

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2661.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0938		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0938			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	117		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	116		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-114

Sdg Number: 660-20784

Lab Sample ID: 660-20784-7

Date Sampled: 01/24/2008 1420

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA2662.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 1000			Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 1000				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-114

Sdg Number: 660-20784

Lab Sample ID: 660-20784-7

Date Sampled: 01/24/2008 1420

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56879	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA2662.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/27/2008 1000		Final Weight/Volume: 5 mL
Date Prepared:	01/27/2008 1000		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	90		74 - 126	
Dibromofluoromethane	111		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-83

Sdg Number: 660-20784

Lab Sample ID: 660-20784-8

Date Sampled: 01/24/2008 1525

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA2663.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 1022			Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 1022				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-83

Sdg Number: 660-20784

Lab Sample ID: 660-20784-8

Date Sampled: 01/24/2008 1525

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA2663.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 1022			Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 1022				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	110	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Private Well 84

Sdg Number: 660-20784

Lab Sample ID: 660-20784-9

Date Sampled: 01/24/2008 1540

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-56879	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2664.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 1045		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 1045			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Private Well 84

Sdg Number: 660-20784

Lab Sample ID: 660-20784-9

Date Sampled: 01/24/2008 1540

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-56879

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2664.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 1045

Final Weight/Volume: 5 mL

Date Prepared: 01/27/2008 1045

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	109	77 - 122
4-Bromofluorobenzene	89	74 - 126
Dibromofluoromethane	112	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-32

Sdg Number: 660-20784

Lab Sample ID: 660-20784-10

Date Sampled: 01/24/2008 1012

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3060.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1420			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1420				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	9.9		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	8.6		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-32

Sdg Number: 660-20784

Lab Sample ID: 660-20784-10

Date Sampled: 01/24/2008 1012

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3060.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1420		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1420			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	3.5		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	8.3		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	17	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	104	74 - 126
Dibromofluoromethane	111	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-7S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-11

Date Sampled: 01/24/2008 1054

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2948.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0756		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0756			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-7S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-11

Date Sampled: 01/24/2008 1054

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2948.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0756			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0756				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	107	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup7

Sdg Number: 660-20784

Lab Sample ID: 660-20784-12

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2949.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 0818

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 0818

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup7

Sdg Number: 660-20784

Lab Sample ID: 660-20784-12

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2949.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0818			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0818				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	4.1	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-7D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-13

Date Sampled: 01/24/2008 1117

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2950.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0840			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0840				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-7D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-13

Date Sampled: 01/24/2008 1117

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JA2950.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/30/2008 0840		Final Weight/Volume: 5 mL
Date Prepared:	01/30/2008 0840		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	18	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	103	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-159

Sdg Number: 660-20784

Lab Sample ID: 660-20784-14

Date Sampled: 01/24/2008 1515

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2960.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1223			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1223				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-159

Sdg Number: 660-20784

Lab Sample ID: 660-20784-14

Date Sampled: 01/24/2008 1515

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2960.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1223

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1223

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	92	77 - 122
4-Bromofluorobenzene	84	74 - 126
Dibromofluoromethane	94	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-157

Sdg Number: 660-20784

Lab Sample ID: 660-20784-15

Date Sampled: 01/24/2008 1422

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2961.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1245		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1245			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-157

Sdg Number: 660-20784

Lab Sample ID: 660-20784-15

Date Sampled: 01/24/2008 1422

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2961.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1245		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1245			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	91		74 - 126	
Dibromofluoromethane	108		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-160

Sdg Number: 660-20784

Lab Sample ID: 660-20784-16

Date Sampled: 01/24/2008 1515

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2953.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0947			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0947				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-160

Lab Sample ID: 660-20784-16

Client Matrix: Water

Date Sampled: 01/24/2008 1515

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2953.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 0947

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 0947

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	114		70 - 130	

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-156

Lab Sample ID: 660-20784-17

Date Sampled: 01/24/2008 1425

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2954.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1009			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1009				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-156

Lab Sample ID: 660-20784-17

Client Matrix: Water

Date Sampled: 01/24/2008 1425

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2954.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1009

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1009

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	114	70 - 130

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-11

Lab Sample ID: 660-20784-18

Date Sampled: 01/24/2008 1105

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2955.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1031			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1031				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	9.1		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.5		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-11

Lab Sample ID: 660-20784-18

Date Sampled: 01/24/2008 1105

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2955.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1031			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1031				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.4		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	17		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	24		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	115	70 - 130



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: EB-7

Lab Sample ID: 660-20784-19

Date Sampled: 01/24/2008 1628

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2956.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1054

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1054

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: EB-7

Lab Sample ID: 660-20784-19

Client Matrix: Water

Date Sampled: 01/24/2008 1628

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2956.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1054			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1054				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.2		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	104		70 - 130	



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: TB-7

Lab Sample ID: 660-20784-20

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2957.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1116

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1116

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: TB-7

Lab Sample ID: 660-20784-20

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2957.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1116			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1116				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	112	70 - 130



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: TB-6

Lab Sample ID: 660-20784-21

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2958.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1138

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1138

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: TB-6

Lab Sample ID: 660-20784-21

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2958.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1138

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1138

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	104		70 - 130	



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: EB-6

Lab Sample ID: 660-20784-22

Date Sampled: 01/24/2008 1645

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57058

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2959.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1200

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1200

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: EB-6

Lab Sample ID: 660-20784-22

Client Matrix: Water

Date Sampled: 01/24/2008 1645

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2959.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1200			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1200				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.2		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	113		70 - 130	



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-107

Lab Sample ID: 660-20784-23

Date Sampled: 01/24/2008 1450

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2962.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1307			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1307				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-107

Lab Sample ID: 660-20784-23

Client Matrix: Water

Date Sampled: 01/24/2008 1450

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2962.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1307			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1307				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	94	77 - 122
4-Bromofluorobenzene	87	74 - 126
Dibromofluoromethane	92	70 - 130



THE LEADER IN ENVIRONMENTAL TESTING

Client: ARCADIS

Analytical Data

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-153

Lab Sample ID: 660-20784-24

Date Sampled: 01/24/2008 1045

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2963.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1329			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1329				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-153

Sdg Number: 660-20784

Lab Sample ID: 660-20784-24

Date Sampled: 01/24/2008 1045

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2963.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1329		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1329			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	92		77 - 122	
4-Bromofluorobenzene	90		74 - 126	
Dibromofluoromethane	95		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-155

Sdg Number: 660-20784

Lab Sample ID: 660-20784-25

Date Sampled: 01/24/2008 1130

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2964.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1352			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1352				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-155

Sdg Number: 660-20784

Lab Sample ID: 660-20784-25

Date Sampled: 01/24/2008 1130

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2964.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1352			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1352				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	104	77 - 122
4-Bromofluorobenzene	101	74 - 126
Dibromofluoromethane	106	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-161

Sdg Number: 660-20784

Lab Sample ID: 660-20784-26

Date Sampled: 01/24/2008 1554

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2965.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1414			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1414				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-161

Sdg Number: 660-20784

Lab Sample ID: 660-20784-26

Date Sampled: 01/24/2008 1554

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2965.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1414		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1414			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	109	77 - 122
4-Bromofluorobenzene	104	74 - 126
Dibromofluoromethane	113	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-158

Sdg Number: 660-20784

Lab Sample ID: 660-20784-27

Date Sampled: 01/24/2008 1535

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JA2966.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/30/2008 1436		Final Weight/Volume: 5 mL
Date Prepared:	01/30/2008 1436		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	4.7		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	4.6		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-158

Sdg Number: 660-20784

Lab Sample ID: 660-20784-27

Date Sampled: 01/24/2008 1535

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2966.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1436		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1436			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-28

Date Sampled: 01/24/2008 1650

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2967.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1459		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1459			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-28

Date Sampled: 01/24/2008 1650

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57058	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JA2967.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1459			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1459				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1
Sdg Number: 660-20784

Client Sample ID: MW-124

Lab Sample ID: 660-20784-29
Client Matrix: WaterDate Sampled: 01/24/2008 0930
Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3011.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2003		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2003			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-124

Sdg Number: 660-20784

Lab Sample ID: 660-20784-29

Date Sampled: 01/24/2008 0930

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA3011.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/30/2008 2003		Final Weight/Volume: 5 mL
Date Prepared:	01/30/2008 2003		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	103		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-126

Sdg Number: 660-20784

Lab Sample ID: 660-20784-30

Date Sampled: 01/24/2008 1010

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3012.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2026			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2026				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	R 2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-126

Sdg Number: 660-20784

Lab Sample ID: 660-20784-30

Date Sampled: 01/24/2008 1010

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3012.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2026			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2026				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	16	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	5.3		0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-106

Sdg Number: 660-20784

Lab Sample ID: 660-20784-31

Date Sampled: 01/24/2008 1018

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3013.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2048			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2048				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-106

Sdg Number: 660-20784

Lab Sample ID: 660-20784-31

Date Sampled: 01/24/2008 1018

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3013.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2048			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2048				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-112

Sdg Number: 660-20784

Lab Sample ID: 660-20784-32

Date Sampled: 01/24/2008 1045

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3014.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2111			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2111				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-112

Sdg Number: 660-20784

Lab Sample ID: 660-20784-32

Date Sampled: 01/24/2008 1045

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3014.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2111		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2111			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.53 / 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	119		77 - 122	
4-Bromofluorobenzene	114		74 - 126	
Dibromofluoromethane	125		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-195

Sdg Number: 660-20784

Lab Sample ID: 660-20784-33

Date Sampled: 01/24/2008 1214

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3015.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2133

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2133

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-195

Sdg Number: 660-20784

Lab Sample ID: 660-20784-33

Date Sampled: 01/24/2008 1214

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3015.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2133		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2133			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.61 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-194

Sdg Number: 660-20784

Lab Sample ID: 660-20784-34

Date Sampled: 01/24/2008 1245

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3016.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2155			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2155				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	R 2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-194

Sdg Number: 660-20784

Lab Sample ID: 660-20784-34

Date Sampled: 01/24/2008 1245

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3016.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2155		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2155			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-193

Sdg Number: 660-20784

Lab Sample ID: 660-20784-35

Date Sampled: 01/24/2008 1248

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3017.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2218		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2218			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	R 2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-193

Sdg Number: 660-20784

Lab Sample ID: 660-20784-35

Date Sampled: 01/24/2008 1248

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3017.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2218			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2218				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Sum)	113	77 - 122
4-Bromofluorobenzene	102	74 - 126
Dibromofluoromethane	119	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: Dup6

Lab Sample ID: 660-20784-36

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3018.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2240

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2240

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup6

Sdg Number: 660-20784

Lab Sample ID: 660-20784-36

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57062 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JA3018.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 01/30/2008 2240 Final Weight/Volume: 5 mL
 Date Prepared: 01/30/2008 2240

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	113		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-184

Sdg Number: 660-20784

Lab Sample ID: 660-20784-37

Date Sampled: 01/24/2008 1510

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3019.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2302

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2302

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-184

Sdg Number: 660-20784

Lab Sample ID: 660-20784-37

Date Sampled: 01/24/2008 1510

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3019.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2302		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2302			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	106		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-183

Lab Sample ID: 660-20784-38

Date Sampled: 01/24/2008 1519

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3020.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2325

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2325

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-183

Lab Sample ID: 660-20784-38

Date Sampled: 01/24/2008 1519

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3020.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2325		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2325			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-185

Sdg Number: 660-20784

Lab Sample ID: 660-20784-39

Date Sampled: 01/24/2008 1600

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3021.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2347		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2347			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	<i>R</i> 2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-185

Sdg Number: 660-20784

Lab Sample ID: 660-20784-39

Date Sampled: 01/24/2008 1600

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3021.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2347			Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2347				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	107	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-187

Sdg Number: 660-20784

Lab Sample ID: 660-20784-40

Date Sampled: 01/24/2008 1610

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3022.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0009			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0009				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-187

Sdg Number: 660-20784

Lab Sample ID: 660-20784-40

Date Sampled: 01/24/2008 1610

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3022.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0009		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0009			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	21		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	3.4	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	115		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-186

Sdg Number: 660-20784

Lab Sample ID: 660-20784-41

Date Sampled: 01/24/2008 1640

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3023.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0032

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0032

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	<i>R</i> 2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-186

Sdg Number: 660-20784

Lab Sample ID: 660-20784-41

Date Sampled: 01/24/2008 1640

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3023.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0032		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0032			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	113		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-42

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57062 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JA3024.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 01/31/2008 0054 Final Weight/Volume: 5 mL
 Date Prepared: 01/31/2008 0054

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-42

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3024.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0054		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0054			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	119		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-43

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3025.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0116			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0116				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-43

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3025.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0116		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0116			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	102		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-44

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3026.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0139			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0139				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: TB-8

Lab Sample ID: 660-20784-44

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3026.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0139

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0139

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-45

Date Sampled: 01/24/2008 1630

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3027.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0201			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0201				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: EB-8

Lab Sample ID: 660-20784-45

Date Sampled: 01/24/2008 1630

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3027.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0201		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0201			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	111	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-46

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3028.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0223		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0223			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	R 2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	45		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	20		0.45	1.0
cis-1,2-Dichloroethene	3.8		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-46

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57062	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3028.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0223			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0223				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.4		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	26		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	118		77 - 122	
4-Bromofluorobenzene	108		74 - 126	
Dibromofluoromethane	120		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-15S

Lab Sample ID: 660-20784-47

Date Sampled: 01/24/2008 0918

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3029.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0246

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0246

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-15S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-47

Date Sampled: 01/24/2008 0918

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3029.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0246

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0246

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	104	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	113	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-15D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-48

Date Sampled: 01/24/2008 0910

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3030.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0308

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0308

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	4.9	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	3.0	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-15D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-48

Date Sampled: 01/24/2008 0910

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57062

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3030.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0308

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0308

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	4.0		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	103	74 - 126
Dibromofluoromethane	114	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-14S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-49

Date Sampled: 01/24/2008 1012

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA3045.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/31/2008 0842		Final Weight/Volume: 5 mL
Date Prepared:	01/31/2008 0842		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-14S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-49

Date Sampled: 01/24/2008 1012

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3045.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0842		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0842			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-14D

Lab Sample ID: 660-20784-50

Date Sampled: 01/24/2008 1009

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3046.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0904

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0904

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-14D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-50

Date Sampled: 01/24/2008 1009

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3046.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0904		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0904			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	106	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-77

Sdg Number: 660-20784

Lab Sample ID: 660-20784-51

Date Sampled: 01/24/2008 1122

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3047.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0927			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0927				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	39		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	18		0.45	1.0
cis-1,2-Dichloroethene	3.5		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-77

Sdg Number: 660-20784

Lab Sample ID: 660-20784-51

Date Sampled: 01/24/2008 1122

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3047.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0927		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0927			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.3		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	24		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	113		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-20

Sdg Number: 660-20784

Lab Sample ID: 660-20784-52

Date Sampled: 01/24/2008 1211

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3048.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0949		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0949			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.96	I	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	1.1		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.94	I	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-20

Lab Sample ID: 660-20784-52

Date Sampled: 01/24/2008 1211

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JA3048.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	01/31/2008 0949		Final Weight/Volume: 5 mL
Date Prepared:	01/31/2008 0949		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.73 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1.3		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	114		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-9S

Lab Sample ID: 660-20784-53

Date Sampled: 01/24/2008 1408

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3049.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1017			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1017				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	6.4		2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-9S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-53

Date Sampled: 01/24/2008 1408

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3049.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1017		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1017			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-9D

Lab Sample ID: 660-20784-54

Date Sampled: 01/24/2008 1400

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3050.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1040

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1040

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-9D

Lab Sample ID: 660-20784-54

Date Sampled: 01/24/2008 1400

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3050.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1040		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1040			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	119		77 - 122	
4-Bromofluorobenzene	110		74 - 126	
Dibromofluoromethane	128		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-9D

Lab Sample ID: 660-20784-54

Date Sampled: 01/24/2008 1400

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57130

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0133.D

Dilution: 2.0


Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2133

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2133

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Acetone	580 		20	40

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-76

Sdg Number: 660-20784

Lab Sample ID: 660-20784-55

Date Sampled: 01/24/2008 1445

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3051.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1102		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1102			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	3.5	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	3.1	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-76

Lab Sample ID: 660-20784-55

Date Sampled: 01/24/2008 1445

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3051.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1102

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1102

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	3.1		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	51		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	23		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	98	77 - 122
4-Bromofluorobenzene	87	74 - 126
Dibromofluoromethane	103	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-25

Lab Sample ID: 660-20784-56

Date Sampled: 01/24/2008 1550

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3052.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1124

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1124

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	14	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	7.5	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-25

Sdg Number: 660-20784

Lab Sample ID: 660-20784-56

Date Sampled: 01/24/2008 1550

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3052.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1124		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1124			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2.3		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	109	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	117	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-85

Sdg Number: 660-20784

Lab Sample ID: 660-20784-57

Date Sampled: 01/24/2008 1452

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3053.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1146

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1146

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	98		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	92		0.45	1.0
cis-1,2-Dichloroethene	34		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: MW-85

Lab Sample ID: 660-20784-57

Date Sampled: 01/24/2008 1452

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JA3053.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1146			Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1146				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	7.6		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	95		77 - 122	
4-Bromofluorobenzene	89		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-233

Sdg Number: 660-20784

Lab Sample ID: 660-20784-58

Date Sampled: 01/24/2008 1715

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57065	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3054.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1208		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1208			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-233

Sdg Number: 660-20784

Lab Sample ID: 660-20784-58

Date Sampled: 01/24/2008 1715

Client Matrix: Water

Date Received: 01/25/2008 0915

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57065

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3054.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1208

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1208

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.58 <i>1.0</i>	<i>+ U</i>	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	106	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	112	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-94

Sdg Number: 660-20784

Lab Sample ID: 660-20784-1

Date Sampled: 01/24/2008 1630

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56873	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2656.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0747		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0747			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	19		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-95

Sdg Number: 660-20784

Lab Sample ID: 660-20784-2

Date Sampled: 01/24/2008 1547

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56873	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2657.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0809		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0809			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	35		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-1

Sdg Number: 660-20784

Lab Sample ID: 660-20784-3

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56873	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2658.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0832		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0832			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: RW-2

Sdg Number: 660-20784

Lab Sample ID: 660-20784-4

Date Sampled: 01/24/2008 0930

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 01/27/2008 0854
Date Prepared: 01/27/2008 0854

Analysis Batch: 660-56873

Instrument ID: BVMJ GC/MS

Lab File ID: 1JA2659.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-12

Sdg Number: 660-20784

Lab Sample ID: 660-20784-5

Date Sampled: 01/24/2008 1035

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56873	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2660.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 0916		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 0916			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	23		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-5

Sdg Number: 660-20784

Lab Sample ID: 660-20784-6

Date Sampled: 01/24/2008 1130

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56873

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2661.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 0938

Final Weight/Volume: 5 mL

Date Prepared: 01/27/2008 0938

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-114

Sdg Number: 660-20784

Lab Sample ID: 660-20784-7

Date Sampled: 01/24/2008 1420

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-56873

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA2662.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/27/2008 1000

Final Weight/Volume: 5 mL

Date Prepared: 01/27/2008 1000

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.1		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-83

Sdg Number: 660-20784

Lab Sample ID: 660-20784-8

Date Sampled: 01/24/2008 1525

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56873	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2663.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 1022		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 1022			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Private Well 84

Sdg Number: 660-20784

Lab Sample ID: 660-20784-9

Date Sampled: 01/24/2008 1540

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-56873	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA2664.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/27/2008 1045		Final Weight/Volume:	5 mL
Date Prepared:	01/27/2008 1045			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-32

Sdg Number: 660-20784

Lab Sample ID: 660-20784-10

Client Matrix: Water

Date Sampled: 01/24/2008 1012

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3060.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1420

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1420

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-7S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-11

Date Sampled: 01/24/2008 1054

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2948.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 0756

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 0756

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup7

Sdg Number: 660-20784

Lab Sample ID: 660-20784-12

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2949.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 0818

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 0818

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-7D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-13

Client Matrix: Water

Date Sampled: 01/24/2008 1117

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2950.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 0840

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 0840

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-159

Sdg Number: 660-20784

Lab Sample ID: 660-20784-14

Date Sampled: 01/24/2008 1515

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2960.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1223

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1223

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-157

Sdg Number: 660-20784

Lab Sample ID: 660-20784-15

Date Sampled: 01/24/2008 1422

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2961.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1245

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1245

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-160

Sdg Number: 660-20784

Lab Sample ID: 660-20784-16

Date Sampled: 01/24/2008 1515

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2953.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 0947		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 0947			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-156

Sdg Number: 660-20784

Lab Sample ID: 660-20784-17

Date Sampled: 01/24/2008 1425

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2954.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1009

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1009

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-11

Sdg Number: 660-20784

Lab Sample ID: 660-20784-18

Date Sampled: 01/24/2008 1105

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2955.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1031		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1031			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-7

Sdg Number: 660-20784

Lab Sample ID: 660-20784-19

Date Sampled: 01/24/2008 1628

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2956.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1054		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1054			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-7

Sdg Number: 660-20784

Lab Sample ID: 660-20784-20

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2957.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1116		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1116			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Sdg Number: 660-20784

Client Sample ID: TB-6

Lab Sample ID: 660-20784-21

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57039

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JA2958.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 1138

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 1138

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-6

Sdg Number: 660-20784

Lab Sample ID: 660-20784-22

Date Sampled: 01/24/2008 1645

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2959.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1200		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1200			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-107

Sdg Number: 660-20784

Lab Sample ID: 660-20784-23

Date Sampled: 01/24/2008 1450

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2962.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1307		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1307			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-153

Sdg Number: 660-20784

Lab Sample ID: 660-20784-24

Date Sampled: 01/24/2008 1045

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2963.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1329		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1329			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-155

Sdg Number: 660-20784

Lab Sample ID: 660-20784-25

Date Sampled: 01/24/2008 1130

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2964.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1352		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1352			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-161

Sdg Number: 660-20784

Lab Sample ID: 660-20784-26

Date Sampled: 01/24/2008 1554

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2965.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1414		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1414			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-158

Sdg Number: 660-20784

Lab Sample ID: 660-20784-27

Date Sampled: 01/24/2008 1535

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2966.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1436		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1436			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	84		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-28

Date Sampled: 01/24/2008 1650

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57039	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JA2967.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 1459		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 1459			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-124

Sdg Number: 660-20784

Lab Sample ID: 660-20784-29

Date Sampled: 01/24/2008 0930

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3061.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1442

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1442

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.3		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-126

Sdg Number: 660-20784

Lab Sample ID: 660-20784-30

Date Sampled: 01/24/2008 1010

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3012.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2026		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2026			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-106

Sdg Number: 660-20784

Lab Sample ID: 660-20784-31

Date Sampled: 01/24/2008 1018

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3013.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2048		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2048			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-112

Sdg Number: 660-20784

Lab Sample ID: 660-20784-32

Date Sampled: 01/24/2008 1045

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3014.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2111		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2111			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-195

Sdg Number: 660-20784

Lab Sample ID: 660-20784-33

Date Sampled: 01/24/2008 1214

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3015.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2133		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2133			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-194

Sdg Number: 660-20784

Lab Sample ID: 660-20784-34

Date Sampled: 01/24/2008 1245

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3016.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2155

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-193

Sdg Number: 660-20784

Lab Sample ID: 660-20784-35

Date Sampled: 01/24/2008 1248

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3017.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2218		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2218			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup6

Sdg Number: 660-20784

Lab Sample ID: 660-20784-36

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3018.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/30/2008 2240		Final Weight/Volume:	5 mL
Date Prepared:	01/30/2008 2240			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-184

Sdg Number: 660-20784

Lab Sample ID: 660-20784-37

Date Sampled: 01/24/2008 1510

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3019.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2302

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2302

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-183

Sdg Number: 660-20784

Lab Sample ID: 660-20784-38

Date Sampled: 01/24/2008 1519

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3020.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2325

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2325

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-185

Sdg Number: 660-20784

Lab Sample ID: 660-20784-39

Date Sampled: 01/24/2008 1600

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3021.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/30/2008 2347

Final Weight/Volume: 5 mL

Date Prepared: 01/30/2008 2347

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-187

Sdg Number: 660-20784

Lab Sample ID: 660-20784-40

Date Sampled: 01/24/2008 1610

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3022.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0009		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0009			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-186

Sdg Number: 660-20784

Lab Sample ID: 660-20784-41

Date Sampled: 01/24/2008 1640

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3023.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0032

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0032

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-42

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3024.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0054

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0054

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-9

Sdg Number: 660-20784

Lab Sample ID: 660-20784-43

Client Matrix: Water

Date Sampled: 01/24/2008 0000

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3025.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0116

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0116

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: TB-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-44

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3026.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0139		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0139			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: EB-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-45

Date Sampled: 01/24/2008 1630

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3027.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0201		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0201			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: Dup-8

Sdg Number: 660-20784

Lab Sample ID: 660-20784-46

Date Sampled: 01/24/2008 0000

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57263	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0632.D
Dilution:	4.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/06/2008 1958		Final Weight/Volume:	5 mL
Date Prepared:	02/06/2008 1958			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	220		4.0	4.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-15S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-47

Date Sampled: 01/24/2008 0918

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 01/31/2008 0246
Date Prepared: 01/31/2008 0246

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS
Lab File ID: 1JA3029.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-15D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-48

Date Sampled: 01/24/2008 0910

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3030.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0308

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0308

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.5		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-14S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-49

Date Sampled: 01/24/2008 1012

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3045.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0842

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0842

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-14D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-50

Date Sampled: 01/24/2008 1009

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3046.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 0904

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 0904

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-77

Sdg Number: 660-20784

Lab Sample ID: 660-20784-51

Date Sampled: 01/24/2008 1122

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57011	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3063.D
Dilution:	2.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1526		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1526			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	150		2.0	2.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-20

Sdg Number: 660-20784

Lab Sample ID: 660-20784-52

Date Sampled: 01/24/2008 1211

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57011	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3048.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 0949		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 0949			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	6.1		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-9S

Sdg Number: 660-20784

Lab Sample ID: 660-20784-53

Date Sampled: 01/24/2008 1408

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57011	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JA3049.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	01/31/2008 1017		Final Weight/Volume:	5 mL
Date Prepared:	01/31/2008 1017			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-9D

Sdg Number: 660-20784

Lab Sample ID: 660-20784-54

Date Sampled: 01/24/2008 1400

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3050.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1040

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1040

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-76

Sdg Number: 660-20784

Lab Sample ID: 660-20784-55

Client Matrix: Water

Date Sampled: 01/24/2008 1445

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3051.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1102

Final Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1102

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	9.9		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-25

Sdg Number: 660-20784

Lab Sample ID: 660-20784-56

Date Sampled: 01/24/2008 1550

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3052.D

Dilution: 1.0

Date Analyzed: 01/31/2008 1124

Initial Weight/Volume: 5 mL

Date Prepared: 01/31/2008 1124

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.7		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-85

Sdg Number: 660-20784

Lab Sample ID: 660-20784-57

Date Sampled: 01/24/2008 1452

Client Matrix: Water

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57263	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0633.D
Dilution:	10		Initial Weight/Volume:	5 mL
Date Analyzed:	02/06/2008 2021		Final Weight/Volume:	5 mL
Date Prepared:	02/06/2008 2021			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	590		10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20784-1

Client Sample ID: MW-233

Sdg Number: 660-20784

Lab Sample ID: 660-20784-58

Client Matrix: Water

Date Sampled: 01/24/2008 1715

Date Received: 01/25/2008 0915

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57011

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JA3054.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 01/31/2008 1208

Final Weight/Volume: mL

Date Prepared: 01/31/2008 1208

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20784	1/24/2008	SW-846	MW-94	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-32	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-7S	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	Dup7	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-7D	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-159	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-157	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-160	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-156	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-11	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	EB-7	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-95	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	TB-7	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	TB-6	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	EB-6	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-107	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-153	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-155	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-161	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-158	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	EB-9	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-124	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	TB-1	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-126	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-106	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-112	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20784	1/24/2008	SW-846	MW-195	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20784	1/24/2008	SW-846	MW-194	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-193	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	Dup6	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-184	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-183	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-185	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	RW-2	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-187	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-186	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	Dup9	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	TB-9	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	TB-8	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	EB-8	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	Dup-8	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-15S	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-15D	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-14S	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-12	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-14D	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-77	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-20	Water	No	--	--	--	--	VOC – Associated Blank
660-20784	1/24/2008	SW-846	MW-9S	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-9D	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-76	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-25	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-85	Water	Yes	--	--	--	--	
660-20784	1/24/2008	SW-846	MW-233	Water	No	--	--	--	--	VOC – Associated Blank

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20784	1/24/2008	SW-846	MW-5	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-114	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	MW-83	Water	No	--	--	--	--	VOC – CCAL %D
660-20784	1/24/2008	SW-846	Private Well 84	Water	No	--	--	--	--	VOC – CCAL %D

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 3

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20815

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8114R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20815 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-117	660-20815-1	Water	1/25/2008	X				
MW-241	660-20815-10	Water	1/25/2008	X				
MW-100	660-20815-11	Water	1/25/2008	X				
MW-99	660-20815-12	Water	1/25/2008	X				
EB-10	660-20815-13	Water	1/25/2008	X				
MW-221	660-20815-14	Water	1/25/2008	X				
MW-22	660-20815-15	Water	1/25/2008	X				
MW-23	660-20815-16	Water	1/25/2008	X				
MW-219	660-20815-17	Water	1/25/2008	X				
MW-220	660-20815-18	Water	1/25/2008	X				
MW-222	660-20815-19	Water	1/25/2008	X				
MW-118	660-20815-2	Water	1/25/2008	X				
EB-11	660-20815-20	Water	1/25/2008	X				
MW-73	660-20815-21	Water	1/25/2008	X				
MW-98	660-20815-22	Water	1/25/2008	X				
DUP-11	660-20815-23	Water	1/25/2008	X				
DUP-10	660-20815-24	Water	1/25/2008	X				
MW-103	660-20815-25	Water	1/25/2008	X				
MW-102	660-20815-26	Water	1/25/2008	X				
MW-190	660-20815-27	Water	1/25/2008	X				
EB-13	660-20815-28	Water	1/25/2008	X				
MW-8S	660-20815-29	Water	1/25/2008	X				
MW-119	660-20815-3	Water	1/25/2008	X				
MW-3	660-20815-30	Water	1/25/2008	X				
MW-84	660-20815-31	Water	1/25/2008	X				
MW-4	660-20815-32	Water	1/25/2008	X				
DUP-13	660-20815-33	Water	1/25/2008	X				
TB-13	660-20815-34	Water	1/25/2008	X				
MW-174	660-20815-35	Water	1/25/2008	X				
MW-179	660-20815-36	Water	1/25/2008	X				
MW-116	660-20815-37	Water	1/25/2008	X				
MW-122	660-20815-38	Water	1/25/2008	X				
MW-178	660-20815-39	Water	1/25/2008	X				
MW-120	660-20815-4	Water	1/25/2008	X				
MW-180	660-20815-40	Water	1/25/2008	X				
MW-181	660-20815-41	Water	1/25/2008	X				
TB-10	660-20815-43	Water	1/25/2008	X				
TB-11	660-20815-44	Water	1/25/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
TB-14	660-20815-45	Water	1/25/2008	X				
MW-224	660-20815-5	Water	1/25/2008	X				
MW-227	660-20815-6	Water	1/25/2008	X				
MW-223	660-20815-7	Water	1/25/2008	X				
MW-225	660-20815-8	Water	1/25/2008	X				
MW-226	660-20815-9	Water	1/25/2008	X				

Notes:

1. Sample location DUP-10 is the field duplicate of parent sample location MW-99.
2. Sample location DUP-11 is the field duplicate of parent sample location MW-23.
3. Sample location DUP-13 is the field duplicate of parent sample location MW-179.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

Target compounds were detected in the associated QA blanks. Sample results associated with blank contamination that were greater than the BAL and/or non-detect did not result in any qualification of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
MW-120 MW-224 MW-226 MW-220 MW-181	Toluene	Detected sample results <RL and <BAL	"U" at the PQL
MW-99 MW-22 DUP-10 MW-103 MW-102 MW-4	o-Xylene	Detected sample results <RL and <BAL	"U" at the PQL

RL = reporting limit

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
MW-224 MW-227 MW-223 MW-225 MW-226 MW-241 MW-100 MW-99 EB-10 MW-221 MW-22 MW-23 MW-219 MW-220	CCV %D	1,2,3-Trichlorobenzene	-28.6%
		1,2,4-Trichlorobenzene	-28.6%
		Bromomethane	-26.0%
		Hexachlorobutadiene	-21.7%
		Naphthalene	-20.8%
MW-178 MW-180 MW-181 TB-10 TB-11 TB-14	CCV %D	trans-1,2-Dichloroethene	-21.5%
		1,2,4-Trichlorobenzene	-24.7%
		2,2-Dichloropropane	-22.9%
		4-Methyl-2-pentanone	21.0%
		Benzene	-21.3%
		Bromomethane	-22.4%

Sample Locations	Initial/Continuing	Compound	Criteria
		Chlorobromomethane	-20.2%
		Chloroethane	54.4%
		Hexachlorobutadiene	-22.7%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-99/DUP-10	m-Xylene & p-Xylene	ND(1.0)	0.71 l	AC
MW-23/DUP-11	Acetone	15 l	ND(20)	AC
	1,4-Dioxane (SIM ID)	4.0	4.9	AC
MW-179/DUP-13	All Compounds	ND	ND	AC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
MW-98	1,1-Dichloroethane	160 E	170 D	170 D
	1,1-Dichloroethene	350 E	480 D	480 D

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-117

Sdg Number: 660-20815

Lab Sample ID: 660-20815-1

Date Sampled: 01/25/2008 0903

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57130

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0134.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2155

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-117

Sdg Number: 660-20815

Lab Sample ID: 660-20815-1

Date Sampled: 01/25/2008 0903

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57130

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0134.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2155

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.2		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-118

Sdg Number: 660-20815

Lab Sample ID: 660-20815-2

Date Sampled: 01/25/2008 0840

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57130 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB0135.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/01/2008 2217 Final Weight/Volume: 5 mL
 Date Prepared: 02/01/2008 2217

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-118

Sdg Number: 660-20815

Lab Sample ID: 660-20815-2

Date Sampled: 01/25/2008 0840

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57130	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0135.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/01/2008 2217		Final Weight/Volume:	5 mL
Date Prepared:	02/01/2008 2217			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-119

Sdg Number: 660-20815

Lab Sample ID: 660-20815-3

Date Sampled: 01/25/2008 0806

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57130

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0136.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2240

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2240

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-119

Sdg Number: 660-20815

Lab Sample ID: 660-20815-3

Date Sampled: 01/25/2008 0806

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57130

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0136.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2240

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2240

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	32		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	89		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-120

Sdg Number: 660-20815

Lab Sample ID: 660-20815-4

Date Sampled: 01/25/2008 0750

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57307

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0634.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/06/2008 2043

Final Weight/Volume: 5 mL

Date Prepared: 02/06/2008 2043

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-120

Sdg Number: 660-20815

Lab Sample ID: 660-20815-4

Date Sampled: 01/25/2008 0750

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57307

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0634.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/06/2008 2043

Final Weight/Volume: 5 mL

Date Prepared: 02/06/2008 2043

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.77 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	103		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-224

Sdg Number: 660-20815

Lab Sample ID: 660-20815-5

Date Sampled: 01/25/2008 1020

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0155.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0541		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0541			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-224

Sdg Number: 660-20815

Lab Sample ID: 660-20815-5

Date Sampled: 01/25/2008 1020

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0155.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/02/2008 0541		Final Weight/Volume: 5 mL
Date Prepared:	02/02/2008 0541		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.53 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	100		77 - 122	
4-Bromofluorobenzene	87		74 - 126	
Dibromofluoromethane	102		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-227

Sdg Number: 660-20815

Lab Sample ID: 660-20815-6

Date Sampled: 01/25/2008 1032

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0156.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0603		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0603			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-227

Sdg Number: 660-20815

Lab Sample ID: 660-20815-6

Date Sampled: 01/25/2008 1032

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0156.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0603		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0603			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	92		77 - 122	
4-Bromofluorobenzene	81		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-223

Sdg Number: 660-20815

Lab Sample ID: 660-20815-7

Date Sampled: 01/25/2008 1100

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0157.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0625

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0625

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-223

Sdg Number: 660-20815

Lab Sample ID: 660-20815-7

Date Sampled: 01/25/2008 1100

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0157.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0625		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0625			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U J	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U J	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-225

Sdg Number: 660-20815

Lab Sample ID: 660-20815-8

Date Sampled: 01/25/2008 1120

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57307	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0635.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/06/2008 2105		Final Weight/Volume:	5 mL
Date Prepared:	02/06/2008 2105			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-225

Sdg Number: 660-20815

Lab Sample ID: 660-20815-8

Date Sampled: 01/25/2008 1120

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57307

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0635.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/06/2008 2105

Final Weight/Volume: 5 mL

Date Prepared: 02/06/2008 2105

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-226

Lab Sample ID: 660-20815-9

Date Sampled: 01/25/2008 1155

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0159.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0709		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0709			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-226

Sdg Number: 660-20815

Lab Sample ID: 660-20815-9

Date Sampled: 01/25/2008 1155

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0159.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0709		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0709			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.55 1.0	+ U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U J	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U J	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	104	77 - 122
4-Bromofluorobenzene	89	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-241

Sdg Number: 660-20815

Lab Sample ID: 660-20815-10

Date Sampled: 01/25/2008 1157

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0160.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0731		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0731			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-241

Sdg Number: 660-20815

Lab Sample ID: 660-20815-10

Date Sampled: 01/25/2008 1157

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0160.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0731

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0731

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	91	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-100

Sdg Number: 660-20815

Lab Sample ID: 660-20815-11

Date Sampled: 01/25/2008 1300

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0161.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0754

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0754

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	7.5	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	2.2	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-100

Sdg Number: 660-20815

Lab Sample ID: 660-20815-11

Date Sampled: 01/25/2008 1300

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0161.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0754

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0754

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	87		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-99

Sdg Number: 660-20815

Lab Sample ID: 660-20815-12

Date Sampled: 01/25/2008 1303

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0162.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0816

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0816

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-99

Sdg Number: 660-20815

Lab Sample ID: 660-20815-12

Date Sampled: 01/25/2008 1303

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0162.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0816

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0816

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U J	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U J	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.54 1.0	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	89	74 - 126
Dibromofluoromethane	113	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: EB-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-13

Date Sampled: 01/25/2008 1323

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0163.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0838

Final Weight/Volume: mL

Date Prepared: 02/02/2008 0838

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: EB-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-13

Date Sampled: 01/25/2008 1323

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0163.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0838		Final Weight/Volume:	mL
Date Prepared:	02/02/2008 0838			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	100	77 - 122
4-Bromofluorobenzene	85	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-221

Sdg Number: 660-20815

Lab Sample ID: 660-20815-14

Date Sampled: 01/25/2008 1343

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0164.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0900		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0900			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-221

Sdg Number: 660-20815

Lab Sample ID: 660-20815-14

Date Sampled: 01/25/2008 1343

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0164.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0900			Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0900				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	89	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-22

Sdg Number: 660-20815

Lab Sample ID: 660-20815-15

Date Sampled: 01/25/2008 1121

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0165.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0922

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0922

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-22

Sdg Number: 660-20815

Lab Sample ID: 660-20815-15
Client Matrix: WaterDate Sampled: 01/25/2008 1121
Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0165.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/02/2008 0922		Final Weight/Volume: 5 mL
Date Prepared:	02/02/2008 0922		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.52 1.0	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	100	77 - 122
4-Bromofluorobenzene	86	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-23

Sdg Number: 660-20815

Lab Sample ID: 660-20815-16

Date Sampled: 01/25/2008 1041

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0166.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0944

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0944

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-23

Lab Sample ID: 660-20815-16

Date Sampled: 01/25/2008 1041

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0166.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/02/2008 0944		Final Weight/Volume: 5 mL
Date Prepared:	02/02/2008 0944		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	15	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	92		77 - 122	
4-Bromofluorobenzene	82		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-219

Sdg Number: 660-20815

Lab Sample ID: 660-20815-17

Date Sampled: 01/25/2008 1252

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0167.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 1006

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 1006

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-219

Sdg Number: 660-20815

Lab Sample ID: 660-20815-17

Date Sampled: 01/25/2008 1252

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0167.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 1006

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 1006

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	33		0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	113		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-220

Sdg Number: 660-20815

Lab Sample ID: 660-20815-18

Date Sampled: 01/25/2008 1251

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57134	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0168.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 1028		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 1028			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-220

Lab Sample ID: 660-20815-18

Date Sampled: 01/25/2008 1251

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57134

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0168.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 1028

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 1028

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.53 ^{1.0}	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	85	74 - 126
Dibromofluoromethane	106	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-222

Lab Sample ID: 660-20815-19

Date Sampled: 01/25/2008 1341

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0235.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0058			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0058				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-222

Sdg Number: 660-20815

Lab Sample ID: 660-20815-19

Date Sampled: 01/25/2008 1341

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0235.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0058

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0058

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	1.1	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: EB-11

Lab Sample ID: 660-20815-20

Date Sampled: 01/25/2008 1345

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0236.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0120		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0120			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: EB-11

Lab Sample ID: 660-20815-20

Date Sampled: 01/25/2008 1345

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0236.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0120		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0120			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.99	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	93		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	90		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-73

Sdg Number: 660-20815

Lab Sample ID: 660-20815-21

Date Sampled: 01/25/2008 1455

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0237.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0143

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0143

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	53	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	91	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-73

Sdg Number: 660-20815

Lab Sample ID: 660-20815-21

Date Sampled: 01/25/2008 1455

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0237.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0143		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0143			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	7.9		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	11		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	89	77 - 122
4-Bromofluorobenzene	86	74 - 126
Dibromofluoromethane	94	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-98

Sdg Number: 660-20815

Lab Sample ID: 660-20815-22

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0238.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0205			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0205				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,2-Dichloroethane	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	3.6		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	13		0.50	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-98

Sdg Number: 660-20815

Lab Sample ID: 660-20815-22

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0238.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0205

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0205

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	47		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.6		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	88		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	91		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-98

Sdg Number: 660-20815

Lab Sample ID: 660-20815-22

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0654.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0408

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0408

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethane	170		10	20
1,1-Dichloroethene	480	SP	9.0	20

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: DUP-11

Lab Sample ID: 660-20815-23

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0239.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0227		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0227			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-11

Sdg Number: 660-20815

Lab Sample ID: 660-20815-23

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0239.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0227

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0227

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	90	77 - 122
4-Bromofluorobenzene	88	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-24

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0240.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0250		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0250			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-24

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0240.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0250

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0250

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.59 1.0	+U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.71	I	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	97	77 - 122
4-Bromofluorobenzene	97	74 - 126
Dibromofluoromethane	97	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-103

Sdg Number: 660-20815

Lab Sample ID: 660-20815-25

Date Sampled: 01/25/2008 1115

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0241.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0312		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0312			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-103

Sdg Number: 660-20815

Lab Sample ID: 660-20815-25

Date Sampled: 01/25/2008 1115

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0241.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0312		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0312			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.53 1.0	+ U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	89	77 - 122
4-Bromofluorobenzene	91	74 - 126
Dibromofluoromethane	87	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW102

Sdg Number: 660-20815

Lab Sample ID: 660-20815-26

Date Sampled: 01/25/2008 1030

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0242.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0334		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0334			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW102

Sdg Number: 660-20815

Lab Sample ID: 660-20815-26

Date Sampled: 01/25/2008 1030

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0242.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0334		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0334			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.56 1.0	+ U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	115		74 - 126	
Dibromofluoromethane	111		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-190

Sdg Number: 660-20815

Lab Sample ID: 660-20815-27

Date Sampled: 01/25/2008 0855

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0243.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0357

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0357

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-190

Sdg Number: 660-20815

Lab Sample ID: 660-20815-27

Date Sampled: 01/25/2008 0855

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0243.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0357		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0357			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.66	I	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	95	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	97	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: EB-13

Lab Sample ID: 660-20815-28

Date Sampled: 01/25/2008 1508

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0244.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0419		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0419			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: EB-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-28

Date Sampled: 01/25/2008 1508

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0244.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0419		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0419			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.55	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.51	I	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	77		77 - 122	
4-Bromofluorobenzene	81		74 - 126	
Dibromofluoromethane	76		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-8S

Sdg Number: 660-20815

Lab Sample ID: 660-20815-29

Date Sampled: 01/25/2008 1317

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0245.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0442

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0442

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-8S

Sdg Number: 660-20815

Lab Sample ID: 660-20815-29

Date Sampled: 01/25/2008 1317

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0245.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0442

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0442

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	92		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	92		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-3

Sdg Number: 660-20815

Lab Sample ID: 660-20815-30

Date Sampled: 01/25/2008 1354

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0246.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0504		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0504			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-3

Sdg Number: 660-20815

Lab Sample ID: 660-20815-30

Date Sampled: 01/25/2008 1354

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0246.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0504		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0504			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	93		77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	91		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-84

Lab Sample ID: 660-20815-31

Date Sampled: 01/25/2008 1500

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0247.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0526		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0526			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-84

Lab Sample ID: 660-20815-31

Date Sampled: 01/25/2008 1500

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0247.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0526

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0526

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-4

Sdg Number: 660-20815

Lab Sample ID: 660-20815-32

Date Sampled: 01/25/2008 1345

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0248.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0549

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0549

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	6.1	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.7	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.50	I	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-4

Sdg Number: 660-20815

Lab Sample ID: 660-20815-32

Date Sampled: 01/25/2008 1345

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0248.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0549

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0549

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	5.9		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.60 1.0	TV	0.50	1.0
Acetone	31		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.68	I	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	96	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	91	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-33

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0249.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0611		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0611			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-33

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0249.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0611

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0611

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	107	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: TB-13

Lab Sample ID: 660-20815-34

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0250.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0633		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0633			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-34

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0250.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0633		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0633			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	98		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	93		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-174

Lab Sample ID: 660-20815-35

Date Sampled: 01/25/2008 0840

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0251.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0656

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0656

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-174

Sdg Number: 660-20815

Lab Sample ID: 660-20815-35

Date Sampled: 01/25/2008 0840

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0251.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0656		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0656			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	98		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-179

Lab Sample ID: 660-20815-36

Date Sampled: 01/25/2008 1125

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0252.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/03/2008 0718		Final Weight/Volume: 5 mL
Date Prepared:	02/03/2008 0718		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-179

Lab Sample ID: 660-20815-36

Date Sampled: 01/25/2008 1125

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0252.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0718		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0718			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	89	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	85	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-116

Lab Sample ID: 660-20815-37

Date Sampled: 01/25/2008 1420

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57197	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0253.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0740		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0740			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-116

Lab Sample ID: 660-20815-37

Date Sampled: 01/25/2008 1420

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0253.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0740

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0740

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	88	77 - 122
4-Bromofluorobenzene	90	74 - 126
Dibromofluoromethane	83	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-122

Lab Sample ID: 660-20815-38

Date Sampled: 01/25/2008 0925

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0254.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0803

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0803

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-122

Sdg Number: 660-20815

Lab Sample ID: 660-20815-38

Date Sampled: 01/25/2008 0925

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57197

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0254.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0803

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0803

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	89		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	84		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-178

Sdg Number: 660-20815

Lab Sample ID: 660-20815-39

Date Sampled: 01/25/2008 1115

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0314.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1321		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1321			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U J	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U J	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U J	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-178

Sdg Number: 660-20815

Lab Sample ID: 660-20815-39

Date Sampled: 01/25/2008 1115

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0314.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/03/2008 1321		Final Weight/Volume: 5 mL
Date Prepared:	02/03/2008 1321		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.1		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	98	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	90	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-180

Lab Sample ID: 660-20815-40

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0315.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1343		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1343			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U ↓	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U ↓	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U ↓	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U ↓	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U ↓	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U ↓	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-180

Sdg Number: 660-20815

Lab Sample ID: 660-20815-40

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0315.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1343		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1343			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	92	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	86	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-181

Sdg Number: 660-20815

Lab Sample ID: 660-20815-41

Date Sampled: 01/25/2008 1510

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0316.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1406		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1406			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.79	+ J	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U J	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U J	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-181

Lab Sample ID: 660-20815-41

Date Sampled: 01/25/2008 1510

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0316.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1406

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1406

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.88 1.0	+U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U J	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	3.5		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	103	74 - 126
Dibromofluoromethane	90	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-43

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0317.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1428			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1428				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U J	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U J	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U J	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-43

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0317.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1428		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1428			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	116		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-11

Sdg Number: 660-20815

Lab Sample ID: 660-20815-44

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0318.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1450

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1450

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U J	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U J	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U J	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: TB-11

Lab Sample ID: 660-20815-44

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0318.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1450		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1450			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	94		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	86		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: TB-14

Lab Sample ID: 660-20815-45

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0319.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1513

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1513

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U J	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U J	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U J	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: TB-14

Lab Sample ID: 660-20815-45

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0319.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1513

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1513

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	96		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-117

Sdg Number: 660-20815

Lab Sample ID: 660-20815-1

Date Sampled: 01/25/2008 0903

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57131

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0134.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2155

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-118

Lab Sample ID: 660-20815-2

Date Sampled: 01/25/2008 0840

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/01/2008 2217
Date Prepared: 02/01/2008 2217

Analysis Batch: 660-57131

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0135.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.3		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-119

Lab Sample ID: 660-20815-3

Date Sampled: 01/25/2008 0806

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57131

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0136.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/01/2008 2240

Final Weight/Volume: 5 mL

Date Prepared: 02/01/2008 2240

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-120

Sdg Number: 660-20815

Lab Sample ID: 660-20815-4

Date Sampled: 01/25/2008 0750

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57263

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0634.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/06/2008 2043

Final Weight/Volume: 5 mL

Date Prepared: 02/06/2008 2043

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-224

Sdg Number: 660-20815

Lab Sample ID: 660-20815-5

Date Sampled: 01/25/2008 1020

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0155.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0541

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0541

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-227

Lab Sample ID: 660-20815-6

Date Sampled: 01/25/2008 1032

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57198	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0156.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0603		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0603			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-223

Lab Sample ID: 660-20815-7

Date Sampled: 01/25/2008 1100

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0157.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0625

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0625

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-225

Lab Sample ID: 660-20815-8

Date Sampled: 01/25/2008 1120

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57263

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0635.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/06/2008 2105

Final Weight/Volume: 5 mL

Date Prepared: 02/06/2008 2105

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-226

Sdg Number: 660-20815

Lab Sample ID: 660-20815-9

Date Sampled: 01/25/2008 1155

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0159.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0709

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0709

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-241

Lab Sample ID: 660-20815-10

Date Sampled: 01/25/2008 1157

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57198	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0160.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0731		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0731			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-100

Lab Sample ID: 660-20815-11

Date Sampled: 01/25/2008 1300

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0161.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0754

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0754

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	28		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-99

Sdg Number: 660-20815

Lab Sample ID: 660-20815-12

Date Sampled: 01/25/2008 1303

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0162.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0816

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0816

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: EB-10

Lab Sample ID: 660-20815-13

Date Sampled: 01/25/2008 1323

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/02/2008 0838
Date Prepared: 02/02/2008 0838

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0163.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-221

Sdg Number: 660-20815

Lab Sample ID: 660-20815-14

Date Sampled: 01/25/2008 1343

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0164.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0900

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0900

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-22

Sdg Number: 660-20815

Lab Sample ID: 660-20815-15

Date Sampled: 01/25/2008 1121

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0165.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 0922

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 0922

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-23

Sdg Number: 660-20815

Lab Sample ID: 660-20815-16

Date Sampled: 01/25/2008 1041

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57198	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0166.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/02/2008 0944		Final Weight/Volume:	5 mL
Date Prepared:	02/02/2008 0944			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	4.0		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-219

Lab Sample ID: 660-20815-17

Date Sampled: 01/25/2008 1252

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0167.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 1006

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 1006

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-220

Sdg Number: 660-20815

Lab Sample ID: 660-20815-18

Date Sampled: 01/25/2008 1251

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57198

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0168.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/02/2008 1028

Final Weight/Volume: 5 mL

Date Prepared: 02/02/2008 1028

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-222

Lab Sample ID: 660-20815-19

Date Sampled: 01/25/2008 1341

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0235.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0058

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0058

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: EB-11

Lab Sample ID: 660-20815-20

Date Sampled: 01/25/2008 1345

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0236.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0120

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0120

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-73

Sdg Number: 660-20815

Lab Sample ID: 660-20815-21

Date Sampled: 01/25/2008 1455

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/03/2008 0143
 Date Prepared: 02/03/2008 0143

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS
 Lab File ID: 1JB0237.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	84		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-98

Sdg Number: 660-20815

Lab Sample ID: 660-20815-22

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 20
Date Analyzed: 02/07/2008 0408
Date Prepared: 02/07/2008 0408

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0654.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	710		20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-11

Sdg Number: 660-20815

Lab Sample ID: 660-20815-23

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0239.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0227

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0227

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	4.9		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: DUP-10

Lab Sample ID: 660-20815-24

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0240.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0250

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0250

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-103

Sdg Number: 660-20815

Lab Sample ID: 660-20815-25

Date Sampled: 01/25/2008 1115

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/03/2008 0312
 Date Prepared: 02/03/2008 0312

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0241.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW102

Sdg Number: 660-20815

Lab Sample ID: 660-20815-26

Date Sampled: 01/25/2008 1030

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0242.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0334

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0334

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-190

Sdg Number: 660-20815

Lab Sample ID: 660-20815-27

Date Sampled: 01/25/2008 0855

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0243.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0357

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0357

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: EB-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-28

Date Sampled: 01/25/2008 1508

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0244.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0419

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0419

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-8S

Sdg Number: 660-20815

Lab Sample ID: 660-20815-29

Date Sampled: 01/25/2008 1317

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0245.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0442

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0442

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Sdg Number: 660-20815

Client Sample ID: MW-3

Lab Sample ID: 660-20815-30

Date Sampled: 01/25/2008 1354

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57112	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0246.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0504		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0504			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-84

Sdg Number: 660-20815

Lab Sample ID: 660-20815-31

Date Sampled: 01/25/2008 1500

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0247.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0526

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0526

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-4

Sdg Number: 660-20815

Lab Sample ID: 660-20815-32

Date Sampled: 01/25/2008 1345

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57112	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0248.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 0549		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 0549			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: DUP-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-33

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0249.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0611

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0611

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-13

Sdg Number: 660-20815

Lab Sample ID: 660-20815-34

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0250.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0633

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0633

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-174

Sdg Number: 660-20815

Lab Sample ID: 660-20815-35

Date Sampled: 01/25/2008 0840

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0251.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0656

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0656

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-179

Sdg Number: 660-20815

Lab Sample ID: 660-20815-36

Date Sampled: 01/25/2008 1125

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0252.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0718

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0718

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-116

Sdg Number: 660-20815

Lab Sample ID: 660-20815-37

Date Sampled: 01/25/2008 1420

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/03/2008 0740
 Date Prepared: 02/03/2008 0740

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0253.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-122

Sdg Number: 660-20815

Lab Sample ID: 660-20815-38

Date Sampled: 01/25/2008 0925

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57112

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0254.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 0803

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 0803

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-178

Sdg Number: 660-20815

Lab Sample ID: 660-20815-39

Date Sampled: 01/25/2008 1115

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0314.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1321

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1321

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-180

Sdg Number: 660-20815

Lab Sample ID: 660-20815-40

Date Sampled: 01/25/2008 1410

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0315.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1343

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1343

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: MW-181

Sdg Number: 660-20815

Lab Sample ID: 660-20815-41

Date Sampled: 01/25/2008 1510

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0316.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1406		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1406			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-10

Sdg Number: 660-20815

Lab Sample ID: 660-20815-43

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0317.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1428

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1428

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-11

Sdg Number: 660-20815

Lab Sample ID: 660-20815-44

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0318.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1450

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1450

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20815-1

Client Sample ID: TB-14

Sdg Number: 660-20815

Lab Sample ID: 660-20815-45

Date Sampled: 01/25/2008 0000

Client Matrix: Water

Date Received: 01/28/2008 1254

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/03/2008 1513
Date Prepared: 02/03/2008 1513

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS
Lab File ID: 1JB0319.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20815	1/25/2008	SW-846	MW-117	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-241	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-100	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-99	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20815	1/25/2008	SW-846	EB-10	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-221	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-22	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20815	1/25/2008	SW-846	MW-23	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-219	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-220	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20815	1/25/2008	SW-846	MW-222	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-118	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	EB-11	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-73	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-98	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	DUP-11	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	DUP-10	Water	No	--	--	--	--	VOC – Associated Blank
660-20815	1/25/2008	SW-846	MW-103	Water	No	--	--	--	--	VOC – Associated Blank
660-20815	1/25/2008	SW-846	MW-102	Water	No	--	--	--	--	VOC – Associated Blank
660-20815	1/25/2008	SW-846	MW-190	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	EB-13	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-8S	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-119	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-3	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-84	Water	Yes	--	--	--	--	

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20815	1/25/2008	SW-846	MW-4	Water	No	--	--	--	--	VOC – Associated Blank
660-20815	1/25/2008	SW-846	DUP-13	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	TB-13	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-174	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-179	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-116	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-122	Water	Yes	--	--	--	--	
660-20815	1/25/2008	SW-846	MW-178	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-120	Water	No	--	--	--	--	VOC – Associated Blank
660-20815	1/25/2008	SW-846	MW-180	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-181	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20815	1/25/2008	SW-846	TB-10	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	TB-11	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	TB-14	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-224	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20815	1/25/2008	SW-846	MW-227	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-223	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-225	Water	No	--	--	--	--	VOC – CCAL %D
660-20815	1/25/2008	SW-846	MW-226	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 4

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20822

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8118R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20822 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-31	660-20822-1	Water	1/28/2008	X				
MW-65	660-20822-10	Water	1/28/2008	X				
MW-135	660-20822-11	Water	1/28/2008	X				
MW-93	660-20822-12	Water	1/28/2008	X				
MW-50	660-20822-13	Water	1/28/2008	X				
MW-111	660-20822-14	Water	1/28/2008	X				
MW-113	660-20822-15	Water	1/28/2008	X				
MW-51	660-20822-16	Water	1/28/2008	X				
MW-26	660-20822-17	Water	1/28/2008	X				
MW-44	660-20822-18	Water	1/28/2008	X				
MW-78	660-20822-19	Water	1/28/2008	X				
MW-59	660-20822-2	Water	1/28/2008	X				
MW-18S	660-20822-20	Water	1/28/2008	X				
MW-18D	660-20822-21	Water	1/28/2008	X				
MW-91	660-20822-22	Water	1/28/2008	X				
MW-29	660-20822-23	Water	1/28/2008	X				
MW-133	660-20822-24	Water	1/28/2008	X				
MW-52	660-20822-25	Water	1/28/2008	X				
MW-64	660-20822-26	Water	1/28/2008	X				
MW-28	660-20822-27	Water	1/28/2008	X				
MW-53	660-20822-28	Water	1/28/2008	X				
MW-54	660-20822-29	Water	1/28/2008	X				
MW-68	660-20822-3	Water	1/28/2008	X				
MW-27	660-20822-30	Water	1/28/2008	X				
MW-104	660-20822-31	Water	1/28/2008	X				
MW-105	660-20822-32	Water	1/28/2008	X				
MW-248	660-20822-33	Water	1/28/2008	X				
DUP14	660-20822-34	Water	1/28/2008	X				
DUP15	660-20822-35	Water	1/28/2008	X				
DUP16	660-20822-36	Water	1/28/2008	X				
DUP17	660-20822-37	Water	1/28/2008	X				
TB-14	660-20822-38	Water	1/28/2008	X				
TB-15	660-20822-39	Water	1/28/2008	X				
MW-60	660-20822-4	Water	1/28/2008	X				
TB-16	660-20822-40	Water	1/28/2008	X				
TB-17	660-20822-41	Water	1/28/2008	X				
EB-14	660-20822-42	Water	1/28/2008	X				
EB-15	660-20822-43	Water	1/28/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
EB-16	660-20822-44	Water	1/28/2008	X				
MW-201	660-20822-45	Water	1/28/2008	X				
MW-199	660-20822-46	Water	1/28/2008	X				
MW-200	660-20822-47	Water	1/28/2008	X				
MW-198	660-20822-48	Water	1/28/2008	X				
DUP21	660-20822-49	Water	1/28/2008	X				
MW-109	660-20822-5	Water	1/28/2008	X				
EB-21	660-20822-50	Water	1/28/2008	X				
MW-21	660-20822-6	Water	1/28/2008	X				
MW-86	660-20822-7	Water	1/28/2008	X				
MW-110	660-20822-8	Water	1/28/2008	X				
MW-67	660-20822-9	Water	1/28/2008	X				

Notes:

1. Sample location DUP14 is the field duplicate of parent sample location MW-21.
2. Sample location DUP15 is the field duplicate of parent sample location MW-135.
3. Sample location DUP16 is the field duplicate of parent sample location MW-133.
4. Sample location DUP17 is the field duplicate of parent sample location MW-53.
5. Sample location DUP21 is the field duplicate of parent sample location MW-248.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No compounds were detected in the associated blanks.

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria	
MW-59	CCV %D	trans-1,2-Dichloroethene	-21.5%	
MW-68		1,2,4-Trichlorobenzene	-24.7%	
MW-60		2,2-Dichloropropane	-22.9%	
MW-109		4-Methyl-2-pentanone	21.0%	
MW-21		Benzene	-21.3%	
MW-86		Bromomethane	-22.4%	
MW-110		Chlorobromomethane	-20.2%	
MW-67		Chloroethane	54.4%	
MW-65		Hexachlorobutadiene	-22.7%	
MW-135				
MW-93				
MW-50		CCV %D	trans-1,2-Dichloroethene	-21.0%
MW-111	2-Hexanone		31.6%	
MW-113	4-Methyl-2-pentanone		33.7%	
MW-51	Acetone		34.6%	
MW-26	Benzene		-23.2%	
MW-44	Chlorobromomethane		-27.4%	
MW-78	MTBE		-24.0%	
MW-18S	Trichloroethene		-25.8%	
MW-18D	Tetrachloroethene		-27.2%	
MW-91				
MW-29				
MW-133				
MW-52				
MW-64				
MW-28				
MW-53				
MW-54				
MW-27				
MW-104				
MW-105				

Sample Locations	Initial/Continuing	Compound	Criteria
MW-248 DUP14 DUP15 DUP16 DUP17 TB-15 TB-16 TB-17 EB-14 EB-15 EB-16 MW-201 MW-199 MW-200 MW-198 DUP21 EB-21	CCV %D	Bromomethane	77.5%
Chloroethane		26.5%	
TB-14	CCV %D	Bromomethane	25.3%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

Sample locations associated with surrogates exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Surrogate	Recovery
MW-26	Toluene-d8	AC
	4-Bromofluorobenzene	> UL
	Dibromofluoromethane	AC
DUP21	Toluene-d8	> UL
	4-Bromofluorobenzene	> UL
	Dibromofluoromethane	> UL

Upper control limit (UL)

Lower control limit (LL)

Diluted (D)

Acceptable (AC)

The criteria used to evaluate the surrogate recoveries are presented in the following table. In the case of a surrogate deviation, the sample results associated with the deviant fraction are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	No Action
	Detect	J
< LL but > 10%	Non-detect	J
	Detect	J
< 10%	Non-detect	R
	Detect	J
Surrogates diluted below the calibration curve due to the high concentration of a target compounds	Non-detect	No Action
	Detect	

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-21/DUP14	Carbon disulfide	1.2	ND(0.85)	AC
	1,4-Dioxane (SIM ID)	ND(1.0)	2.6	AC
MW-135/DUP15	1,1-Dichloroethane	0.72 l	0.98 l	AC
	1,1-Dichloroethene	1.2	2.1	AC
	cis-1,2-Dichloroethene	63	93	38.4%
	trans-1,2-Dichloroethene	0.5 l	0.93 l	AC
	Trichloroethene	80	120	40.0%
	1,4-Dioxane (SIM ID)	29	33	12.9%
MW-133/DUP16	Acetone	180	130	32.2%
MW-53/DUP17	All Compounds	ND	ND	AC
MW-248/DUP21	1,1-Dichloroethane	1.1	ND(0.52)	AC
	1,1-Dichloroethene	0.68 l	ND(0.45)	AC
	Acetone	14 l	ND(9.9)	AC
	Carbon disulfide	3.6	ND(0.85)	NC
	1,4-Dioxane (SIM ID)	12	ND(1.0)	NC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The compounds carbon disulfide and 1,4-dioxane associated with samples MW-248 and DUP21 exhibited a field duplicate RPD greater than the control limit. The associated sample results from sample locations for the listed compound were qualified as estimated.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
MW-78	1,1-Dichloroethane	240 E	340 D	340 D
	1,1-Dichloroethene	180 E	320 D	320 D

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-31

Sdg Number: 660-20822

Lab Sample ID: 660-20822-1

Date Sampled: 01/28/2008 1305

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57668	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0821.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/08/2008 2248		Final Weight/Volume: 5 mL
Date Prepared:	02/08/2008 2248		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-31

Sdg Number: 660-20822

Lab Sample ID: 660-20822-1

Date Sampled: 01/28/2008 1305

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57668	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0821.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2248		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2248			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.70	I	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	87		77 - 122	
4-Bromofluorobenzene	91		74 - 126	
Dibromofluoromethane	90		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-59

Sdg Number: 660-20822

Lab Sample ID: 660-20822-2

Date Sampled: 01/28/2008 1240

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/03/2008 1557
Date Prepared: 02/03/2008 1557

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0321.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U J	0.50	1.0
Bromobenzene	0.58	U J	0.58	1.0
Chlorobromomethane	0.58	U J	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U J	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U J	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-59

Lab Sample ID: 660-20822-2

Date Sampled: 01/28/2008 1240

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0321.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1557		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1557			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	2.0		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	95		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	84		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-68

Sdg Number: 660-20822

Lab Sample ID: 660-20822-3

Date Sampled: 01/28/2008 1350

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57098 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB0322.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/03/2008 1620 Final Weight/Volume: 5 mL
 Date Prepared: 02/03/2008 1620

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.8	U	0.45	1.0
cis-1,2-Dichloroethene	1.6	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-68

Lab Sample ID: 660-20822-3

Date Sampled: 01/28/2008 1350

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0322.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1620

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1620

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	5.2	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	91		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	80		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-60

Sdg Number: 660-20822

Lab Sample ID: 660-20822-4

Date Sampled: 01/28/2008 1627

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0323.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1642			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1642				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-60

Sdg Number: 660-20822

Lab Sample ID: 660-20822-4

Date Sampled: 01/28/2008 1627

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0323.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1642		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1642			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-109

Sdg Number: 660-20822

Lab Sample ID: 660-20822-5

Date Sampled: 01/28/2008 1528

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0324.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1704		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1704			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.77	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-109

Sdg Number: 660-20822

Lab Sample ID: 660-20822-5

Date Sampled: 01/28/2008 1528

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0324.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1704

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1704

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	96		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	82		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-6

Date Sampled: 01/28/2008 1121

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0325.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1727		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1727			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-6

Date Sampled: 01/28/2008 1121

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0325.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1727

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1727

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.2		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	98		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	89		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-86

Sdg Number: 660-20822

Lab Sample ID: 660-20822-7

Date Sampled: 01/28/2008 1028

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0326.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1749

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1749

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	1.6	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.8	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-86

Sdg Number: 660-20822

Lab Sample ID: 660-20822-7

Date Sampled: 01/28/2008 1028

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0326.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1749			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1749				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	88		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-110

Sdg Number: 660-20822

Lab Sample ID: 660-20822-8

Date Sampled: 01/28/2008 1112

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0327.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1811		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1811			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	15	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	15	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-110

Sdg Number: 660-20822

Lab Sample ID: 660-20822-8

Date Sampled: 01/28/2008 1112

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0327.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1811

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1811

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	87		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-67

Sdg Number: 660-20822

Lab Sample ID: 660-20822-9

Date Sampled: 01/28/2008 1657

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0328.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1834

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1834

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	25	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	41	U	0.45	1.0
cis-1,2-Dichloroethene	2.3	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-67

Sdg Number: 660-20822

Lab Sample ID: 660-20822-9

Date Sampled: 01/28/2008 1657

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0328.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1834		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1834			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	14		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	107		74 - 126	
Dibromofluoromethane	87		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-65

Sdg Number: 660-20822

Lab Sample ID: 660-20822-10

Date Sampled: 01/28/2008 1012

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0329.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1857		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1857			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-65

Lab Sample ID: 660-20822-10

Date Sampled: 01/28/2008 1012

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0329.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1857			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1857				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	101	77 - 122
4-Bromofluorobenzene	107	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-135

Sdg Number: 660-20822

Lab Sample ID: 660-20822-11

Date Sampled: 01/28/2008 1105

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0330.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1921		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1921			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.72	I	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.2		0.45	1.0
cis-1,2-Dichloroethene	63		0.65	1.0
trans-1,2-Dichloroethene	0.50	+	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-135

Sdg Number: 660-20822

Lab Sample ID: 660-20822-11

Date Sampled: 01/28/2008 1105

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57098	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0330.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1921			Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1921				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	80		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	101	77 - 122
4-Bromofluorobenzene	106	74 - 126
Dibromofluoromethane	88	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-93

Lab Sample ID: 660-20822-12

Date Sampled: 01/28/2008 1055

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57098	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0331.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/03/2008 1946		Final Weight/Volume: 5 mL
Date Prepared:	02/03/2008 1946		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-93

Lab Sample ID: 660-20822-12

Date Sampled: 01/28/2008 1055

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57098

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0331.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1946

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1946

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	93		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	82		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-50

Sdg Number: 660-20822

Lab Sample ID: 660-20822-13

Date Sampled: 01/28/2008 1318

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0428.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2200		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2200			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-50

Sdg Number: 660-20822

Lab Sample ID: 660-20822-13

Date Sampled: 01/28/2008 1318

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0428.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2200		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2200			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	94		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	79		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-111

Sdg Number: 660-20822

Lab Sample ID: 660-20822-14

Date Sampled: 01/28/2008 1700

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57469 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB0429.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/04/2008 2222 Final Weight/Volume: 5 mL
 Date Prepared: 02/04/2008 2222

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-111

Sdg Number: 660-20822

Lab Sample ID: 660-20822-14

Date Sampled: 01/28/2008 1700

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0429.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/04/2008 2222

Final Weight/Volume: 5 mL

Date Prepared: 02/04/2008 2222

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	94		77 - 122	
4-Bromofluorobenzene	107		74 - 126	
Dibromofluoromethane	75		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-113

Sdg Number: 660-20822

Lab Sample ID: 660-20822-15

Date Sampled: 01/28/2008 1618

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0430.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2244		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2244			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-113

Lab Sample ID: 660-20822-15

Date Sampled: 01/28/2008 1618

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0430.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2244			Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2244				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	121	74 - 126
Dibromofluoromethane	88	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-51

Sdg Number: 660-20822

Lab Sample ID: 660-20822-16

Date Sampled: 01/28/2008 1640

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0431.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2307			Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2307				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-51

Sdg Number: 660-20822

Lab Sample ID: 660-20822-16

Date Sampled: 01/28/2008 1640

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0431.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2307		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2307			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	117		74 - 126	
Dibromofluoromethane	86		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-26

Sdg Number: 660-20822

Lab Sample ID: 660-20822-17

Date Sampled: 01/28/2008 1503

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0432.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2329		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2329			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-26

Sdg Number: 660-20822

Lab Sample ID: 660-20822-17

Date Sampled: 01/28/2008 1503

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0432.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/04/2008 2329

Final Weight/Volume: 5 mL

Date Prepared: 02/04/2008 2329

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	127	J1	74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-44

Sdg Number: 660-20822

Lab Sample ID: 660-20822-18

Date Sampled: 01/28/2008 1319

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57469 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB0433.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/04/2008 2351 Final Weight/Volume: 5 mL
 Date Prepared: 02/04/2008 2351

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	5.1	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	2.9	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-44

Sdg Number: 660-20822

Lab Sample ID: 660-20822-18

Date Sampled: 01/28/2008 1319

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0433.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2351			Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2351				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	123		74 - 126	
Dibromofluoromethane	86		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-78

Sdg Number: 660-20822

Lab Sample ID: 660-20822-19

Date Sampled: 01/28/2008 1345

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0434.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0013

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0013

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,2-Dichloroethane	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	110		0.65	1.0
trans-1,2-Dichloroethene	3.7		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	6.4	U	0.50	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-78

Sdg Number: 660-20822

Lab Sample ID: 660-20822-19

Date Sampled: 01/28/2008 1345

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0434.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0013			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0013				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	48	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.2		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	120		74 - 126	
Dibromofluoromethane	87		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-78

Sdg Number: 660-20822

Lab Sample ID: 660-20822-19

Date Sampled: 01/28/2008 1345

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0667.D

Dilution: 5.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0858

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0858

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethane	340	D	2.6	5.0
1,1-Dichloroethene	320	D	2.2	5.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-18S

Sdg Number: 660-20822

Lab Sample ID: 660-20822-20

Date Sampled: 01/28/2008 1050

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0435.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0036			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0036				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-18S

Lab Sample ID: 660-20822-20

Date Sampled: 01/28/2008 1050

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0435.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0036		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0036			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	116		74 - 126	
Dibromofluoromethane	88		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-18D

Sdg Number: 660-20822

Lab Sample ID: 660-20822-21

Date Sampled: 01/28/2008 1035

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0436.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/05/2008 0058		Final Weight/Volume: 5 mL
Date Prepared:	02/05/2008 0058		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.74	I	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	2.1		0.65	1.0
trans-1,2-Dichloroethene	3.1		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-18D

Sdg Number: 660-20822

Lab Sample ID: 660-20822-21

Date Sampled: 01/28/2008 1035

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0436.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0058			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0058				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.52	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	117		74 - 126	
Dibromofluoromethane	86		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-91

Sdg Number: 660-20822

Lab Sample ID: 660-20822-22

Date Sampled: 01/28/2008 0935

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0437.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0120		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0120			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	34		2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	3.8		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	3.1		0.45	1.0
cis-1,2-Dichloroethene	7.0		0.65	1.0
trans-1,2-Dichloroethene	0.92	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-91

Sdg Number: 660-20822

Lab Sample ID: 660-20822-22

Date Sampled: 01/28/2008 0935

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0437.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/05/2008 0120		Final Weight/Volume: 5 mL
Date Prepared:	02/05/2008 0120		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	16	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	79		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-29

Sdg Number: 660-20822

Lab Sample ID: 660-20822-23

Date Sampled: 01/28/2008 1116

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0438.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0142		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0142			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	10	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	7.5	U	0.45	1.0
cis-1,2-Dichloroethene	12	U	0.65	1.0
trans-1,2-Dichloroethene	3.2	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-29

Sdg Number: 660-20822

Lab Sample ID: 660-20822-23

Date Sampled: 01/28/2008 1116

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0438.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0142

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0142

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	91	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	124	74 - 126
Dibromofluoromethane	87	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-133

Sdg Number: 660-20822

Lab Sample ID: 660-20822-24

Date Sampled: 01/28/2008 0937

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0439.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0205

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0205

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-133

Sdg Number: 660-20822

Lab Sample ID: 660-20822-24

Date Sampled: 01/28/2008 0937

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0439.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0205			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0205				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	180	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	101	77 - 122
4-Bromofluorobenzene	115	74 - 126
Dibromofluoromethane	82	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-52

Sdg Number: 660-20822

Lab Sample ID: 660-20822-25

Date Sampled: 01/28/2008 1140

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0440.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0227

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0227

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-52

Sdg Number: 660-20822

Lab Sample ID: 660-20822-25

Date Sampled: 01/28/2008 1140

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0440.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0227			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0227				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	100		77 - 122	
4-Bromofluorobenzene	120		74 - 126	
Dibromofluoromethane	83		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-64

Lab Sample ID: 660-20822-26

Date Sampled: 01/28/2008 1205

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0441.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0249

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0249

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	1.7		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.79	I	0.45	1.0
cis-1,2-Dichloroethene	2.0		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-64

Lab Sample ID: 660-20822-26

Date Sampled: 01/28/2008 1205

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0441.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0249		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0249			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	10	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	114		74 - 126	
Dibromofluoromethane	81		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-28

Sdg Number: 660-20822

Lab Sample ID: 660-20822-27

Date Sampled: 01/28/2008 1255

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0442.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0311

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0311

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	3.1	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.4	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-28

Sdg Number: 660-20822

Lab Sample ID: 660-20822-27

Date Sampled: 01/28/2008 1255

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0442.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0311		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0311			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	111		74 - 126	
Dibromofluoromethane	77		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-53

Sdg Number: 660-20822

Lab Sample ID: 660-20822-28

Date Sampled: 01/28/2008 1215

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0443.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0333

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0333

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-53

Lab Sample ID: 660-20822-28

Date Sampled: 01/28/2008 1215

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0443.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/05/2008 0333		Final Weight/Volume: 5 mL
Date Prepared:	02/05/2008 0333		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	125		74 - 126	
Dibromofluoromethane	89		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-54

Sdg Number: 660-20822

Lab Sample ID: 660-20822-29

Date Sampled: 01/28/2008 1115

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57469

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0444.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0356

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0356

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-54

Sdg Number: 660-20822

Lab Sample ID: 660-20822-29

Date Sampled: 01/28/2008 1115

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0444.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0356		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0356			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	98		77 - 122	
4-Bromofluorobenzene	113		74 - 126	
Dibromofluoromethane	80		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-27

Sdg Number: 660-20822

Lab Sample ID: 660-20822-30

Date Sampled: 01/28/2008 1040

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0445.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/05/2008 0418		Final Weight/Volume: 5 mL
Date Prepared:	02/05/2008 0418		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	69		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	120		0.45	1.0
cis-1,2-Dichloroethene	11		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-27

Sdg Number: 660-20822

Lab Sample ID: 660-20822-30

Date Sampled: 01/28/2008 1040

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0445.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0418			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0418				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	79	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	23	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	125	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-104

Sdg Number: 660-20822

Lab Sample ID: 660-20822-31

Date Sampled: 01/28/2008 1115

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0446.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0440			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0440				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	23		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	25		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-104

Sdg Number: 660-20822

Lab Sample ID: 660-20822-31

Date Sampled: 01/28/2008 1115

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0446.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0440			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0440				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.85	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	110		74 - 126	
Dibromofluoromethane	83		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-105

Sdg Number: 660-20822

Lab Sample ID: 660-20822-32

Date Sampled: 01/28/2008 1010

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0447.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0502		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0502			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-105

Lab Sample ID: 660-20822-32

Date Sampled: 01/28/2008 1010

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57469	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0447.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0502			Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0502				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	100	77 - 122
4-Bromofluorobenzene	118	74 - 126
Dibromofluoromethane	84	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-248

Sdg Number: 660-20822

Lab Sample ID: 660-20822-33

Date Sampled: 01/28/2008 1225

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0649.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0217

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0217

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	1.1	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.68	I	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-248

Sdg Number: 660-20822

Lab Sample ID: 660-20822-33

Date Sampled: 01/28/2008 1225

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0649.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0217

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0217

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	14	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	3.6	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-34

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0650.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0240			Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0240				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-34

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0650.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0240		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0240			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-35

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0651.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0302		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0302			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.98	I	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	2.1		0.45	1.0
cis-1,2-Dichloroethene	93		0.65	1.0
trans-1,2-Dichloroethene	0.93	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-35

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0651.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0302		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0302			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	120		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-36

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0652.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0324

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0324

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-36

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0652.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0324		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0324			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	130		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP17

Sdg Number: 660-20822

Lab Sample ID: 660-20822-37

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0653.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0346

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0346

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP17

Sdg Number: 660-20822

Lab Sample ID: 660-20822-37

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0653.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0346		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0346			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	115	77 - 122
4-Bromofluorobenzene	103	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-38

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57397

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0913.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 1455

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 1455

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: TB-14

Lab Sample ID: 660-20822-38

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57397	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0913.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 1455		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 1455			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	86		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	84		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: TB-15

Lab Sample ID: 660-20822-39

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0655.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0431

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0431

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-39

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0655.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0431		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0431			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-40

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0656.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0453		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0453			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-40

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0656.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0453

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0453

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	111	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	106	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-17

Sdg Number: 660-20822

Lab Sample ID: 660-20822-41

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0657.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0515		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0515			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-17

Sdg Number: 660-20822

Lab Sample ID: 660-20822-41

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0657.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0515		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0515			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	90		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-42

Date Sampled: 01/28/2008 1130

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0658.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0538

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0538

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-42

Date Sampled: 01/28/2008 1130

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0658.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0538			Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0538				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-43

Date Sampled: 01/28/2008 1715

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0659.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0600			Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0600				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-43

Date Sampled: 01/28/2008 1715

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB0659.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0600			Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0600				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-44

Date Sampled: 01/28/2008 1625

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0660.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0622		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0622			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: EB-16

Lab Sample ID: 660-20822-44

Date Sampled: 01/28/2008 1625

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0660.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0622		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0622			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-201

Lab Sample ID: 660-20822-45

Date Sampled: 01/28/2008 1550

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB0661.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/07/2008 0644		Final Weight/Volume: 5 mL
Date Prepared:	02/07/2008 0644		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-201

Sdg Number: 660-20822

Lab Sample ID: 660-20822-45

Date Sampled: 01/28/2008 1550

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0661.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0644

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0644

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-199

Sdg Number: 660-20822

Lab Sample ID: 660-20822-46

Date Sampled: 01/28/2008 1550

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0662.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0706		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0706			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-199

Sdg Number: 660-20822

Lab Sample ID: 660-20822-46

Date Sampled: 01/28/2008 1550

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0662.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0706		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0706			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	106		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-200

Sdg Number: 660-20822

Lab Sample ID: 660-20822-47

Date Sampled: 01/28/2008 1440

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0663.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0729		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0729			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-200

Sdg Number: 660-20822

Lab Sample ID: 660-20822-47

Date Sampled: 01/28/2008 1440

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0663.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0729

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0729

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	119	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	119	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Sdg Number: 660-20822

Client Sample ID: MW-198

Lab Sample ID: 660-20822-48

Date Sampled: 01/28/2008 1445

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0664.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0751

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0751

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-198

Sdg Number: 660-20822

Lab Sample ID: 660-20822-48

Date Sampled: 01/28/2008 1445

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0664.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0751

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0751

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-49

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0665.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0813		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0813			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-49

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0665.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0813		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0813			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	151	J1	77 - 122	
4-Bromofluorobenzene	126	J1	74 - 126	
Dibromofluoromethane	147	J1	70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-50

Date Sampled: 01/28/2008 1650

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57315

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0666.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0835

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0835

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-50

Date Sampled: 01/28/2008 1650

Client Matrix: Water

Date Received: 01/29/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57315	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0666.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0835		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0835			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	87		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-31

Sdg Number: 660-20822

Lab Sample ID: 660-20822-1

Date Sampled: 01/28/2008 1305

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0320.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1535

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1535

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-59

Sdg Number: 660-20822

Lab Sample ID: 660-20822-2

Date Sampled: 01/28/2008 1240

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0321.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1557		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1557			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-68

Sdg Number: 660-20822

Lab Sample ID: 660-20822-3

Date Sampled: 01/28/2008 1350

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57684	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0822.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2311		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2311			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	16		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-60

Sdg Number: 660-20822

Lab Sample ID: 660-20822-4

Date Sampled: 01/28/2008 1627

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0323.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1642

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1642

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.5		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-109

Sdg Number: 660-20822

Lab Sample ID: 660-20822-5

Date Sampled: 01/28/2008 1528

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/03/2008 1704
Date Prepared: 02/03/2008 1704

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0324.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	15		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-6

Date Sampled: 01/28/2008 1121

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/03/2008 1727
 Date Prepared: 02/03/2008 1727

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0325.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-86

Sdg Number: 660-20822

Lab Sample ID: 660-20822-7

Date Sampled: 01/28/2008 1028

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0326.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1749		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1749			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	13		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-110

Sdg Number: 660-20822

Lab Sample ID: 660-20822-8

Date Sampled: 01/28/2008 1112

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0327.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1811		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1811			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	62		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-67

Sdg Number: 660-20822

Lab Sample ID: 660-20822-9

Date Sampled: 01/28/2008 1657

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0328.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1834		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1834			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	54		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-65

Sdg Number: 660-20822

Lab Sample ID: 660-20822-10

Date Sampled: 01/28/2008 1012

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57097

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0329.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/03/2008 1857

Final Weight/Volume: 5 mL

Date Prepared: 02/03/2008 1857

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-135

Sdg Number: 660-20822

Lab Sample ID: 660-20822-11

Date Sampled: 01/28/2008 1105

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0330.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1921		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1921			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	29		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-93

Sdg Number: 660-20822

Lab Sample ID: 660-20822-12

Date Sampled: 01/28/2008 1055

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57097	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0331.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/03/2008 1946		Final Weight/Volume:	5 mL
Date Prepared:	02/03/2008 1946			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-50

Sdg Number: 660-20822

Lab Sample ID: 660-20822-13

Date Sampled: 01/28/2008 1318

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0428.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/04/2008 2200

Final Weight/Volume: 5 mL

Date Prepared: 02/04/2008 2200

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-111

Sdg Number: 660-20822

Lab Sample ID: 660-20822-14

Date Sampled: 01/28/2008 1700

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57215	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0429.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2222		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2222			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-113

Sdg Number: 660-20822

Lab Sample ID: 660-20822-15

Date Sampled: 01/28/2008 1618

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/04/2008 2244
 Date Prepared: 02/04/2008 2244

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0430.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-51

Sdg Number: 660-20822

Lab Sample ID: 660-20822-16

Date Sampled: 01/28/2008 1640

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0431.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/04/2008 2307

Final Weight/Volume: 5 mL

Date Prepared: 02/04/2008 2307

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-26

Sdg Number: 660-20822

Lab Sample ID: 660-20822-17

Date Sampled: 01/28/2008 1503

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57215	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0432.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/04/2008 2329		Final Weight/Volume:	5 mL
Date Prepared:	02/04/2008 2329			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-44

Sdg Number: 660-20822

Lab Sample ID: 660-20822-18

Date Sampled: 01/28/2008 1319

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0433.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/04/2008 2351

Final Weight/Volume: 5 mL

Date Prepared: 02/04/2008 2351

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	4.8		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-78

Sdg Number: 660-20822

Lab Sample ID: 660-20822-19

Date Sampled: 01/28/2008 1345

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0434.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0013

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0013

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	99		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-18S

Sdg Number: 660-20822

Lab Sample ID: 660-20822-20

Date Sampled: 01/28/2008 1050

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57215	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0435.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0036		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0036			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-18D

Sdg Number: 660-20822

Lab Sample ID: 660-20822-21

Date Sampled: 01/28/2008 1035

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0436.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0058

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0058

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-91

Sdg Number: 660-20822

Lab Sample ID: 660-20822-22

Date Sampled: 01/28/2008 0935

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57215	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0437.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0120		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0120			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	85		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-29

Sdg Number: 660-20822

Lab Sample ID: 660-20822-23

Date Sampled: 01/28/2008 1116

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/05/2008 0142
Date Prepared: 02/05/2008 0142

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0438.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	91		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-133

Sdg Number: 660-20822

Lab Sample ID: 660-20822-24

Date Sampled: 01/28/2008 0937

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0439.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0205

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0205

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-52

Sdg Number: 660-20822

Lab Sample ID: 660-20822-25

Date Sampled: 01/28/2008 1140

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57215	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0440.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0227		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0227			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-64

Sdg Number: 660-20822

Lab Sample ID: 660-20822-26

Date Sampled: 01/28/2008 1205

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/05/2008 0249
Date Prepared: 02/05/2008 0249

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0441.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	19		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-28

Sdg Number: 660-20822

Lab Sample ID: 660-20822-27

Date Sampled: 01/28/2008 1255

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0442.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0311

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0311

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	28		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-53

Sdg Number: 660-20822

Lab Sample ID: 660-20822-28

Date Sampled: 01/28/2008 1215

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0443.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0333

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0333

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-54

Sdg Number: 660-20822

Lab Sample ID: 660-20822-29

Date Sampled: 01/28/2008 1115

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0444.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0356

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0356

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-27

Sdg Number: 660-20822

Lab Sample ID: 660-20822-30

Date Sampled: 01/28/2008 1040

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/05/2008 0418
 Date Prepared: 02/05/2008 0418

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0445.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	69		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-104

Sdg Number: 660-20822

Lab Sample ID: 660-20822-31

Date Sampled: 01/28/2008 1115

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57215	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0446.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/05/2008 0440		Final Weight/Volume:	5 mL
Date Prepared:	02/05/2008 0440			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	91		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-105

Sdg Number: 660-20822

Lab Sample ID: 660-20822-32

Date Sampled: 01/28/2008 1010

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57215

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0447.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/05/2008 0502

Final Weight/Volume: 5 mL

Date Prepared: 02/05/2008 0502

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-248

Sdg Number: 660-20822

Lab Sample ID: 660-20822-33

Date Sampled: 01/28/2008 1225

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0649.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0217

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0217

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	12	J	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-34

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/07/2008 0240
 Date Prepared: 02/07/2008 0240

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0650.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-35

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0651.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0302

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0302

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	33		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-36

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0652.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0324

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0324

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP17

Sdg Number: 660-20822

Lab Sample ID: 660-20822-37

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0653.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0346

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0346

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-38

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57645	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	3JB0710.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 1841		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 1841			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-39

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0655.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0431

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0431

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-40

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0656.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0453

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0453

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: TB-17

Sdg Number: 660-20822

Lab Sample ID: 660-20822-41

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0657.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0515

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0515

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-14

Sdg Number: 660-20822

Lab Sample ID: 660-20822-42

Date Sampled: 01/28/2008 1130

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0658.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0538

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0538

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-15

Sdg Number: 660-20822

Lab Sample ID: 660-20822-43

Date Sampled: 01/28/2008 1715

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/07/2008 0600
Date Prepared: 02/07/2008 0600

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0659.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-16

Sdg Number: 660-20822

Lab Sample ID: 660-20822-44

Date Sampled: 01/28/2008 1625

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0660.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0622

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0622

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-201

Sdg Number: 660-20822

Lab Sample ID: 660-20822-45

Date Sampled: 01/28/2008 1550

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57317	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB0661.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/07/2008 0644		Final Weight/Volume:	5 mL
Date Prepared:	02/07/2008 0644			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-199

Sdg Number: 660-20822

Lab Sample ID: 660-20822-46

Date Sampled: 01/28/2008 1550

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0662.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0706

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0706

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-200

Sdg Number: 660-20822

Lab Sample ID: 660-20822-47

Date Sampled: 01/28/2008 1440

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/07/2008 0729
 Date Prepared: 02/07/2008 0729

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB0663.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: MW-198

Sdg Number: 660-20822

Lab Sample ID: 660-20822-48

Date Sampled: 01/28/2008 1445

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0664.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0751

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0751

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: DUP21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-49

Date Sampled: 01/28/2008 0000

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0665.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0813

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0813

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20822-1

Client Sample ID: EB-21

Sdg Number: 660-20822

Lab Sample ID: 660-20822-50

Date Sampled: 01/28/2008 1650

Client Matrix: Water

Date Received: 01/29/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57317

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB0666.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/07/2008 0835

Final Weight/Volume: 5 mL

Date Prepared: 02/07/2008 0835

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20822	1/28/2008	SW-846	MW-31	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-65	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-135	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-93	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-50	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-111	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-113	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-51	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-26	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-44	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-78	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-59	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-18S	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-18D	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-91	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-29	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-133	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-52	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-64	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-28	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-53	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-54	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-68	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-27	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-104	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-105	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-248	Water	No	--	--	--	--	VOC – Field Duplicate RPD

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20822	1/28/2008	SW-846	DUP14	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	DUP15	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	DUP16	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	DUP17	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	TB-14	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	TB-15	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-60	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	TB-16	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	TB-17	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	EB-14	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	EB-15	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	EB-16	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-201	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-199	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-200	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-198	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	DUP21	Water	No	--	--	--	--	VOC – Field Duplicate RPD
660-20822	1/28/2008	SW-846	MW-109	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	EB-21	Water	Yes	--	--	--	--	
660-20822	1/28/2008	SW-846	MW-21	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-86	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-110	Water	No	--	--	--	--	VOC – CCAL %D
660-20822	1/28/2008	SW-846	MW-67	Water	No	--	--	--	--	VOC – CCAL %D

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 5

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20857

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8147R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20857 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-24	660-20857-1	Water	1/29/2008	X				
DUP-20	660-20857-10	Water	1/29/2008	X				
MW-16S	660-20857-11	Water	1/29/2008	X				
MW-145	660-20857-12	Water	1/29/2008	X				
MW-143	660-20857-13	Water	1/29/2008	X				
MW-140	660-20857-14	Water	1/29/2008	X				
MW-136	660-20857-15	Water	1/29/2008	X				
MW-207	660-20857-16	Water	1/29/2008	X				
MW-16D	660-20857-17	Water	1/29/2008	X				
MW-206	660-20857-18	Water	1/29/2008	X				
MW-142	660-20857-19	Water	1/29/2008	X				
MW-146	660-20857-2	Water	1/29/2008	X				
MW-141	660-20857-20	Water	1/29/2008	X				
MW-144	660-20857-21	Water	1/29/2008	X				
MW-96	660-20857-22	Water	1/29/2008	X				
MW-139	660-20857-23	Water	1/29/2008	X				
MW-137	660-20857-24	Water	1/29/2008	X				
MW-138	660-20857-25	Water	1/29/2008	X				
MW-176	660-20857-26	Water	1/29/2008	X				
MW-175	660-20857-27	Water	1/29/2008	X				
MW-125	660-20857-28	Water	1/29/2008	X				
MW-13S	660-20857-29	Water	1/29/2008	X				
MW-56	660-20857-3	Water	1/29/2008	X				
MW-13D	660-20857-30	Water	1/29/2008	X				
MW-45	660-20857-31	Water	1/29/2008	X				
MW-115	660-20857-32	Water	1/29/2008	X				
MW-177	660-20857-33	Water	1/29/2008	X				
MW-243	660-20857-34	Water	1/29/2008	X				
MW-244	660-20857-35	Water	1/29/2008	X				
DUP-18	660-20857-36	Water	1/29/2008	X				
MW-231	660-20857-37	Water	1/29/2008	X				
MW-229	660-20857-38	Water	1/29/2008	X				
MW-245	660-20857-39	Water	1/29/2008	X				
MW-88	660-20857-4	Water	1/29/2008	X				
MW-247	660-20857-40	Water	1/29/2008	X				
MW-246	660-20857-41	Water	1/29/2008	X				
MW-242	660-20857-42	Water	1/29/2008	X				
MW-230	660-20857-43	Water	1/29/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-108	660-20857-44	Water	1/29/2008	X				
EB-19	660-20857-45	Water	1/29/2008	X				
EB-21	660-20857-46	Water	1/29/2008	X				
EB-20	660-20857-47	Water	1/29/2008	X				
EB-18	660-20857-48	Water	1/29/2008	X				
MW-101	660-20857-49	Water	1/29/2008	X				
MW-84B	660-20857-5	Water	1/29/2008	X				
MW-55	660-20857-50	Water	1/29/2008	X				
MW-150	660-20857-51	Water	1/29/2008	X				
MW-147	660-20857-52	Water	1/29/2008	X				
MW-148	660-20857-53	Water	1/29/2008	X				
TB-19	660-20857-54	Water	1/29/2008	X				
TB-18	660-20857-55	Water	1/29/2008	X				
TB-20	660-20857-56	Water	1/29/2008	X				
TB-21	660-20857-57	Water	1/29/2008	X				
MW-84A	660-20857-6	Water	1/29/2008	X				
MW-82	660-20857-7	Water	1/29/2008	X				
MW-74	660-20857-8	Water	1/29/2008	X				
DUP-19	660-20857-9	Water	1/29/2008	X				

Notes:

1. Sample location DUP-18 is the field duplicate of parent sample location MW-247.
2. Sample location DUP-19 is the field duplicate of parent sample location MW-55.
3. Sample location DUP-20 is the field duplicate of parent sample location MW-175.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No compounds were detected in the associated blanks.

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria		
MW-24 MW-146 MW-56 MW-88 MW-84B MW-84A EB-19 EB-21 EB-18 TB-19 TB-18 TB-20 TB-21	CCV %D	2,2-Dichloropropane	-43.1%		
		4-Methyl-2-pentanone	27.2%		
		Benzene	-22.9%		
		Bromomethane	-71.8%		
		Carbon disulfide	-24.2%		
		Chloromethane	35.1%		
		Methylene chloride	-24.7%		
		trans-1,2-Dichloroethene	-24.5%		
		DUP-20 MW-230 MW-108 MW-101 MW-55 MW-150 MW-147 MW-148	CCV %D	1,2,3-Trichlorobenzene	-28.0%
				2,2-Dichloropropane	50.5%
Bromomethane	-66.3%				
Methylene chloride	-21.8%				
Naphthalene	-30.5%				
Tetrachloroethene	-26.7%				
MW-13D MW-45 MW-115 MW-177 MW-243 MW-244 DUP-18 MW-231 MW-229 MW-245 MW-247 MW-246 MW-242 EB-20	CCV %D	1,2,3-Trichlorobenzene	-20.2%		
		2,2-Dichloropropane	54.4%		
		Naphthalene	-26.2%		

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

Sample locations associated with surrogates exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Surrogate	Recovery
MW-143	Toluene-d8	> UL
MW-140	4-Bromofluorobenzene	AC
MW-142	Dibromofluoromethane	AC
MW-141		

Upper control limit (UL)

Lower control limit (LL)

Diluted (D)

Acceptable (AC)

The criteria used to evaluate the surrogate recoveries are presented in the following table. In the case of a surrogate deviation, the sample results associated with the deviant fraction are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	No Action
	Detect	J
< LL but > 10%	Non-detect	J
	Detect	J

Control Limit	Sample Result	Qualification
< 10%	Non-detect	R
	Detect	J
Surrogates diluted below the calibration curve due to the high concentration of a target compounds	Non-detect	No Action
	Detect	

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-247/DUP-18	All Compounds	ND	ND	AC

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-55/DUP-19	1,1-Dichloroethane	1.9	1.6	AC
	1,4-Dioxane (SIM ID)	14	15	6.8%
MW-175/DUP-20	All Compounds	ND	ND	AC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

All identified compounds met the specified criteria.

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-24

Sdg Number: 660-20857

Lab Sample ID: 660-20857-1

Date Sampled: 01/29/2008 0916

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0825.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1942		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1942			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U J	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U J	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U J	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U J	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U J	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-24

Sdg Number: 660-20857

Lab Sample ID: 660-20857-1

Date Sampled: 01/29/2008 0916

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0825.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1942		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1942			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	89		74 - 126	
Dibromofluoromethane	77		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-146

Sdg Number: 660-20857

Lab Sample ID: 660-20857-2

Date Sampled: 01/29/2008 1149

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB0826.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/08/2008 2005		Final Weight/Volume: 5 mL
Date Prepared:	02/08/2008 2005		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-146

Sdg Number: 660-20857

Lab Sample ID: 660-20857-2

Date Sampled: 01/29/2008 1149

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0826.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2005		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2005			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	88	74 - 126
Dibromofluoromethane	71	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-56

Sdg Number: 660-20857

Lab Sample ID: 660-20857-3

Date Sampled: 01/29/2008 1557

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0827.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2029		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2029			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-56

Sdg Number: 660-20857

Lab Sample ID: 660-20857-3

Date Sampled: 01/29/2008 1557

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0827.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2029			Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2029				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	25		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	74		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-88

Sdg Number: 660-20857

Lab Sample ID: 660-20857-4

Date Sampled: 01/29/2008 1519

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0828.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2052			Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2052				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-88

Sdg Number: 660-20857

Lab Sample ID: 660-20857-4

Date Sampled: 01/29/2008 1519

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0828.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2052		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2052			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	180		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	100	77 - 122
4-Bromofluorobenzene	88	74 - 126
Dibromofluoromethane	76	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-84B

Sdg Number: 660-20857

Lab Sample ID: 660-20857-5

Date Sampled: 01/29/2008 0952

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0829.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2117		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2117			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-84B

Lab Sample ID: 660-20857-5

Date Sampled: 01/29/2008 0952

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57398

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0829.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 2117

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 2117

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	94	77 - 122
4-Bromofluorobenzene	84	74 - 126
Dibromofluoromethane	72	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-84A

Sdg Number: 660-20857

Lab Sample ID: 660-20857-6

Date Sampled: 01/29/2008 0957

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0830.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2141		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2141			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-84A

Sdg Number: 660-20857

Lab Sample ID: 660-20857-6

Date Sampled: 01/29/2008 0957

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB0830.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/08/2008 2141		Final Weight/Volume: 5 mL
Date Prepared:	02/08/2008 2141		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	95		77 - 122	
4-Bromofluorobenzene	84		74 - 126	
Dibromofluoromethane	71		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-82

Sdg Number: 660-20857

Lab Sample ID: 660-20857-7

Date Sampled: 01/29/2008 1450

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0925.D
Dilution:	10			Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2013			Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2013				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	5.0	U	5.0	10
Bromobenzene	5.8	U	5.8	10
Chlorobromomethane	5.8	U	5.8	10
Dichlorobromomethane	3.5	U	3.5	10
Bromoform	5.8	U	5.8	10
Bromomethane	24	U	24	50
n-Butylbenzene	6.7	U	6.7	10
sec-Butylbenzene	6.3	U	6.3	10
tert-Butylbenzene	8.4	U	8.4	10
Carbon tetrachloride	4.2	U	4.2	10
Chlorobenzene	6.3	U	6.3	10
Chloroethane	24	U	24	50
Chloroform	9.0	U	9.0	10
Chloromethane	10	U	10	40
2-Chlorotoluene	6.4	U	6.4	10
4-Chlorotoluene	5.2	U	5.2	10
Chlorodibromomethane	3.4	U	3.4	10
1,2-Dibromo-3-Chloropropane	24	U	24	50
Ethylene Dibromide	5.0	U	5.0	10
Dibromomethane	4.1	U	4.1	10
1,2-Dichlorobenzene	4.4	U	4.4	10
1,3-Dichlorobenzene	6.4	U	6.4	10
1,4-Dichlorobenzene	5.2	U	5.2	10
Dichlorodifluoromethane	24	U	24	50
1,1-Dichloroethane	73		5.2	10
1,2-Dichloroethane	5.7	U	5.7	10
1,1-Dichloroethene	17		4.4	10
cis-1,2-Dichloroethene	6.4	U	6.4	10
trans-1,2-Dichloroethene	4.4	U	4.4	10
1,2-Dichloropropane	5.2	U	5.2	10
1,3-Dichloropropane	3.9	U	3.9	10
2,2-Dichloropropane	3.6	U	3.6	10
1,1-Dichloropropene	3.1	U	3.1	10
cis-1,3-Dichloropropene	1.4	U	1.4	10
trans-1,3-Dichloropropene	1.4	U	1.4	10
Ethylbenzene	4.4	U	4.4	10
Hexachlorobutadiene	4.0	U	4.0	10
Isopropylbenzene	1.9	U	1.9	10
4-Isopropyltoluene	6.9	U	6.9	10
Methylene Chloride	40	U	40	50
Naphthalene	24	U	24	50
N-Propylbenzene	5.9	U	5.9	10
Styrene	9.8	U	9.8	20
1,1,1,2-Tetrachloroethane	6.3	U	6.3	10

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-82

Sdg Number: 660-20857

Lab Sample ID: 660-20857-7

Date Sampled: 01/29/2008 1450

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0925.D
Dilution:	10			Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2013			Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2013				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	1.5	U	1.5	10
Tetrachloroethene	5.0	U	5.0	10
Toluene	5.1	U	5.1	10
1,2,3-Trichlorobenzene	7.7	U	7.7	10
1,2,4-Trichlorobenzene	5.8	U	5.8	10
1,1,1-Trichloroethane	4.6	U	4.6	10
1,1,2-Trichloroethane	4.7	U	4.7	10
Trichloroethene	5.0	U	5.0	10
Trichlorofluoromethane	24	U	24	50
1,2,3-Trichloropropane	1.8	U	1.8	10
1,2,4-Trimethylbenzene	8.6	U	8.6	10
1,3,5-Trimethylbenzene	5.4	U	5.4	10
Vinyl chloride	5.0	U	5.0	10
o-Xylene	5.0	U	5.0	10
Acetone	99	U	99	200
2-Butanone (MEK)	84	U	84	100
4-Methyl-2-pentanone (MIBK)	38	U	38	100
Carbon disulfide	8.4	U	8.4	10
2-Hexanone	44	U	44	100
Methyl tert-butyl ether	4.4	U	4.4	10
m-Xylene & p-Xylene	6.0	U	6.0	20
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	98		77 - 122	
4-Bromofluorobenzene	88		74 - 126	
Dibromofluoromethane	90		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-74

Sdg Number: 660-20857

Lab Sample ID: 660-20857-8

Date Sampled: 01/29/2008 1622

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0926.D
Dilution:	5.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2037			Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2037				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	2.4	U	2.4	5.0
Bromobenzene	2.9	U	2.9	5.0
Chlorobromomethane	2.9	U	2.9	5.0
Dichlorobromomethane	1.8	U	1.8	5.0
Bromoform	2.9	U	2.9	5.0
Bromomethane	13	U	13	24
n-Butylbenzene	3.4	U	3.4	5.0
sec-Butylbenzene	3.2	U	3.2	5.0
tert-Butylbenzene	4.2	U	4.2	5.0
Carbon tetrachloride	2.1	U	2.1	5.0
Chlorobenzene	3.2	U	3.2	5.0
Chloroethane	13	U	13	24
Chloroform	4.4	U	4.4	5.0
Chloromethane	5.0	U	5.0	20
2-Chlorotoluene	3.3	U	3.3	5.0
4-Chlorotoluene	2.6	U	2.6	5.0
Chlorodibromomethane	1.7	U	1.7	5.0
1,2-Dibromo-3-Chloropropane	13	U	13	24
Ethylene Dibromide	2.4	U	2.4	5.0
Dibromomethane	2.1	U	2.1	5.0
1,2-Dichlorobenzene	2.2	U	2.2	5.0
1,3-Dichlorobenzene	3.2	U	3.2	5.0
1,4-Dichlorobenzene	2.6	U	2.6	5.0
Dichlorodifluoromethane	13	U	13	24
1,1-Dichloroethane	82		2.6	5.0
1,2-Dichloroethane	2.9	U	2.9	5.0
1,1-Dichloroethene	37		2.3	5.0
cis-1,2-Dichloroethene	3.3	U	3.3	5.0
trans-1,2-Dichloroethene	2.2	U	2.2	5.0
1,2-Dichloropropane	2.6	U	2.6	5.0
1,3-Dichloropropane	2.0	U	2.0	5.0
2,2-Dichloropropane	1.8	U	1.8	5.0
1,1-Dichloropropene	1.6	U	1.6	5.0
cis-1,3-Dichloropropene	0.70	U	0.70	5.0
trans-1,3-Dichloropropene	0.70	U	0.70	5.0
Ethylbenzene	2.2	U	2.2	5.0
Hexachlorobutadiene	2.0	U	2.0	5.0
Isopropylbenzene	0.95	U	0.95	5.0
4-Isopropyltoluene	3.5	U	3.5	5.0
Methylene Chloride	20	U	20	24
Naphthalene	13	U	13	24
N-Propylbenzene	3.0	U	3.0	5.0
Styrene	4.9	U	4.9	10
1,1,1,2-Tetrachloroethane	3.2	U	3.2	5.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-74

Sdg Number: 660-20857

Lab Sample ID: 660-20857-8

Date Sampled: 01/29/2008 1622

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0926.D
Dilution:	5.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2037		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2037			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.75	U	0.75	5.0
Tetrachloroethene	2.4	U	2.4	5.0
Toluene	2.5	U	2.5	5.0
1,2,3-Trichlorobenzene	3.9	U	3.9	5.0
1,2,4-Trichlorobenzene	2.9	U	2.9	5.0
1,1,1-Trichloroethane	2.3	U	2.3	5.0
1,1,2-Trichloroethane	2.4	U	2.4	5.0
Trichloroethene	2.4	U	2.4	5.0
Trichlorofluoromethane	13	U	13	24
1,2,3-Trichloropropane	0.90	U	0.90	5.0
1,2,4-Trimethylbenzene	4.3	U	4.3	5.0
1,3,5-Trimethylbenzene	2.7	U	2.7	5.0
Vinyl chloride	2.4	U	2.4	5.0
o-Xylene	2.4	U	2.4	5.0
Acetone	50	U	50	100
2-Butanone (MEK)	42	U	42	50
4-Methyl-2-pentanone (MIBK)	19	U	19	50
Carbon disulfide	4.3	U	4.3	5.0
2-Hexanone	22	U	22	50
Methyl tert-butyl ether	2.2	U	2.2	5.0
m-Xylene & p-Xylene	3.0	U	3.0	10

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	97	77 - 122
4-Bromofluorobenzene	90	74 - 126
Dibromofluoromethane	85	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: DUP-19

Sdg Number: 660-20857

Lab Sample ID: 660-20857-9

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0927.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2100		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2100			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	1.6	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: DUP-19

Sdg Number: 660-20857

Lab Sample ID: 660-20857-9

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0927.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2100		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2100			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	117		77 - 122	
4-Bromofluorobenzene	109		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: DUP-20

Lab Sample ID: 660-20857-10

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1142.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0117		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0117			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: DUP-20

Sdg Number: 660-20857

Lab Sample ID: 660-20857-10

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57495

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1142.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0117

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0117

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	97	74 - 126
Dibromofluoromethane	91	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-16S

Sdg Number: 660-20857

Lab Sample ID: 660-20857-11

Date Sampled: 01/29/2008 1605

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0929.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2147		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2147			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-16S

Sdg Number: 660-20857

Lab Sample ID: 660-20857-11

Date Sampled: 01/29/2008 1605

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB0929.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/09/2008 2147		Final Weight/Volume: 5 mL
Date Prepared:	02/09/2008 2147		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	109		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-145

Lab Sample ID: 660-20857-12

Date Sampled: 01/29/2008 1130

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB0930.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/09/2008 2210		Final Weight/Volume: 5 mL
Date Prepared:	02/09/2008 2210		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-145

Sdg Number: 660-20857

Lab Sample ID: 660-20857-12

Date Sampled: 01/29/2008 1130

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0930.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2210			Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2210				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	119		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-143

Sdg Number: 660-20857

Lab Sample ID: 660-20857-13

Date Sampled: 01/29/2008 1315

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57501 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB0931.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/09/2008 2234 Final Weight/Volume: 5 mL
 Date Prepared: 02/09/2008 2234

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-143

Sdg Number: 660-20857

Lab Sample ID: 660-20857-13

Date Sampled: 01/29/2008 1315

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0931.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2234		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2234			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	127	J1	77 - 122	
4-Bromofluorobenzene	110		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-140

Sdg Number: 660-20857

Lab Sample ID: 660-20857-14

Date Sampled: 01/29/2008 1010

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0932.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2257		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2257			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-140

Sdg Number: 660-20857

Lab Sample ID: 660-20857-14

Date Sampled: 01/29/2008 1010

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0932.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2257		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2257			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	125	J1	77 - 122	
4-Bromofluorobenzene	110		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-136

Lab Sample ID: 660-20857-15

Date Sampled: 01/29/2008 0850

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0933.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2321			Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2321				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-136

Sdg Number: 660-20857

Lab Sample ID: 660-20857-15

Date Sampled: 01/29/2008 0850

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0933.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2321		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2321			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	108		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-207

Sdg Number: 660-20857

Lab Sample ID: 660-20857-16

Date Sampled: 01/29/2008 1415

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0934.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2344		Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2344			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-207

Sdg Number: 660-20857

Lab Sample ID: 660-20857-16

Date Sampled: 01/29/2008 1415

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0934.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/09/2008 2344			Final Weight/Volume:	5 mL
Date Prepared:	02/09/2008 2344				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	14	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	4.8		0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	120		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-16D

Sdg Number: 660-20857

Lab Sample ID: 660-20857-17

Date Sampled: 01/29/2008 1535

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0935.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 0007		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 0007			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-16D

Lab Sample ID: 660-20857-17

Date Sampled: 01/29/2008 1535

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57501	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0935.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 0007		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 0007			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	122		77 - 122	
4-Bromofluorobenzene	109		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-206

Sdg Number: 660-20857

Lab Sample ID: 660-20857-18

Date Sampled: 01/29/2008 1405

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1019.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 1945		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 1945			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-206

Sdg Number: 660-20857

Lab Sample ID: 660-20857-18

Date Sampled: 01/29/2008 1405

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 2HB1019.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/10/2008 1945		Final Weight/Volume: 5 mL
Date Prepared:	02/10/2008 1945		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.7		0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	117	77 - 122
4-Bromofluorobenzene	105	74 - 126
Dibromofluoromethane	106	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-142

Sdg Number: 660-20857

Lab Sample ID: 660-20857-19

Date Sampled: 01/29/2008 1315

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1020.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2009		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2009			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-142

Sdg Number: 660-20857

Lab Sample ID: 660-20857-19

Date Sampled: 01/29/2008 1315

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57531

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1020.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/10/2008 2009

Final Weight/Volume: 5 mL

Date Prepared: 02/10/2008 2009

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	130	J1	77 - 122
4-Bromofluorobenzene	112		74 - 126
Dibromofluoromethane	118		70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-141

Sdg Number: 660-20857

Lab Sample ID: 660-20857-20

Date Sampled: 01/29/2008 1250

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 2HB1021.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/10/2008 2032		Final Weight/Volume: 5 mL
Date Prepared:	02/10/2008 2032		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-141

Lab Sample ID: 660-20857-20

Date Sampled: 01/29/2008 1250

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1021.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2032		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2032			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	130	J1	77 - 122
4-Bromofluorobenzene	113		74 - 126
Dibromofluoromethane	121		70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-144

Sdg Number: 660-20857

Lab Sample ID: 660-20857-21

Date Sampled: 01/29/2008 1220

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57531 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 2HB1022.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/10/2008 2055 Final Weight/Volume: 5 mL
 Date Prepared: 02/10/2008 2055

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-144

Lab Sample ID: 660-20857-21

Date Sampled: 01/29/2008 1220

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1022.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2055		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2055			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-96

Sdg Number: 660-20857

Lab Sample ID: 660-20857-22

Date Sampled: 01/29/2008 1115

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1023.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2119		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2119			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-96

Sdg Number: 660-20857

Lab Sample ID: 660-20857-22

Date Sampled: 01/29/2008 1115

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1023.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2119		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2119			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-139

Sdg Number: 660-20857

Lab Sample ID: 660-20857-23

Date Sampled: 01/29/2008 1005

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1024.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2142		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2142			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-139

Lab Sample ID: 660-20857-23

Date Sampled: 01/29/2008 1005

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1024.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2142		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2142			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	94	74 - 126
Dibromofluoromethane	91	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-137

Lab Sample ID: 660-20857-24

Date Sampled: 01/29/2008 0915

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1025.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2206		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2206			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-137

Lab Sample ID: 660-20857-24

Date Sampled: 01/29/2008 0915

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 2HB1025.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/10/2008 2206		Final Weight/Volume: 5 mL
Date Prepared:	02/10/2008 2206		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	111	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	102	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-138

Sdg Number: 660-20857

Lab Sample ID: 660-20857-25

Date Sampled: 01/29/2008 0845

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1026.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2229		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2229			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.1	I	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-138

Sdg Number: 660-20857

Lab Sample ID: 660-20857-25

Date Sampled: 01/29/2008 0845

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1026.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2229		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2229			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	96	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-176

Lab Sample ID: 660-20857-26

Date Sampled: 01/29/2008 0920

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1027.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2252		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2252			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-176

Lab Sample ID: 660-20857-26

Date Sampled: 01/29/2008 0920

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 2HB1027.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/10/2008 2252		Final Weight/Volume: 5 mL
Date Prepared:	02/10/2008 2252		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	106		77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	95		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-175

Lab Sample ID: 660-20857-27

Date Sampled: 01/29/2008 1130

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1028.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2316		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2316			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-175

Lab Sample ID: 660-20857-27

Date Sampled: 01/29/2008 1130

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1028.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2316		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2316			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	102		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-125

Lab Sample ID: 660-20857-28

Date Sampled: 01/29/2008 1114

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	2HB1029.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2339			Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2339				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-125

Lab Sample ID: 660-20857-28

Date Sampled: 01/29/2008 1114

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57531	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1029.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2339		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2339			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-13S

Lab Sample ID: 660-20857-29

Date Sampled: 01/29/2008 1305

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/11/2008 0002
 Date Prepared: 02/11/2008 0002

Analysis Batch: 660-57531

Instrument ID: BVMH GC/MS
 Lab File ID: 2HB1030.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	9.2		0.90	1.0
Chloromethane	1.1	I	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1
Sdg Number: 660-20857

Client Sample ID: MW-13S

Lab Sample ID: 660-20857-29
Client Matrix: WaterDate Sampled: 01/29/2008 1305
Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/11/2008 0002
Date Prepared: 02/11/2008 0002

Analysis Batch: 660-57531

Instrument ID: BVMH GC/MS
Lab File ID: 2HB1030.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	117	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	100	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-13D

Lab Sample ID: 660-20857-30

Date Sampled: 01/29/2008 1244

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1113.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1342		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1342			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-13D

Lab Sample ID: 660-20857-30

Date Sampled: 01/29/2008 1244

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1113.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1342		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1342		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	100	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-45

Lab Sample ID: 660-20857-31

Date Sampled: 01/29/2008 1353

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1114.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1406		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1406			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-45

Lab Sample ID: 660-20857-31

Date Sampled: 01/29/2008 1353

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1114.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1406		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1406			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	111	77 - 122
4-Bromofluorobenzene	94	74 - 126
Dibromofluoromethane	95	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-115

Sdg Number: 660-20857

Lab Sample ID: 660-20857-32

Date Sampled: 01/29/2008 1005

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1115.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1429		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1429		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-115

Lab Sample ID: 660-20857-32

Date Sampled: 01/29/2008 1005

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57504

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1115.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1429

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1429

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	91	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-177

Lab Sample ID: 660-20857-33

Date Sampled: 01/29/2008 0928

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1116.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1453		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1453			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-177

Sdg Number: 660-20857

Lab Sample ID: 660-20857-33

Date Sampled: 01/29/2008 0928

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1116.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1453		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1453		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-243

Lab Sample ID: 660-20857-34

Date Sampled: 01/29/2008 1422

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1117.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1516		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1516			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-243

Lab Sample ID: 660-20857-34

Date Sampled: 01/29/2008 1422

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1117.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1516		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1516			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	95	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-244

Lab Sample ID: 660-20857-35

Date Sampled: 01/29/2008 1449

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1118.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1539		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1539			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-244

Lab Sample ID: 660-20857-35

Date Sampled: 01/29/2008 1449

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/11/2008 1539
 Date Prepared: 02/11/2008 1539

Analysis Batch: 660-57504

Instrument ID: BVMH GC/MS
 Lab File ID: 1HB1118.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	96	74 - 126
Dibromofluoromethane	97	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: DUP-18

Lab Sample ID: 660-20857-36

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1119.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1603			Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1603				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: DUP-18

Lab Sample ID: 660-20857-36

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1119.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1603		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1603			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	96	74 - 126
Dibromofluoromethane	96	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-231

Lab Sample ID: 660-20857-37

Date Sampled: 01/29/2008 1036

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1120.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1626		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1626			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-231

Lab Sample ID: 660-20857-37

Date Sampled: 01/29/2008 1036

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1120.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1626		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1626			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	109	77 - 122
4-Bromofluorobenzene	94	74 - 126
Dibromofluoromethane	90	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-229

Lab Sample ID: 660-20857-38

Date Sampled: 01/29/2008 1042

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1121.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1650		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1650		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-229

Lab Sample ID: 660-20857-38

Date Sampled: 01/29/2008 1042

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1121.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1650		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1650			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-245

Sdg Number: 660-20857

Lab Sample ID: 660-20857-39

Date Sampled: 01/29/2008 0933

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1122.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1713			Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1713				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-245

Sdg Number: 660-20857

Lab Sample ID: 660-20857-39

Date Sampled: 01/29/2008 0933

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1122.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1713			Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1713				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	103	77 - 122
4-Bromofluorobenzene	90	74 - 126
Dibromofluoromethane	89	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-247

Sdg Number: 660-20857

Lab Sample ID: 660-20857-40

Date Sampled: 01/29/2008 0850

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57504

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1123.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1737

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1737

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-247

Lab Sample ID: 660-20857-40

Date Sampled: 01/29/2008 0850

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1123.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1737		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1737		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	94	74 - 126
Dibromofluoromethane	94	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-246

Lab Sample ID: 660-20857-41

Date Sampled: 01/29/2008 0900

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1124.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1800		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1800			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-246

Lab Sample ID: 660-20857-41

Date Sampled: 01/29/2008 0900

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1124.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1800		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1800		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	93		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-242

Sdg Number: 660-20857

Lab Sample ID: 660-20857-42

Date Sampled: 01/29/2008 1420

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1125.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1824		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1824			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-242

Sdg Number: 660-20857

Lab Sample ID: 660-20857-42

Date Sampled: 01/29/2008 1420

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1125.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1824		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1824			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	91		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-230

Lab Sample ID: 660-20857-43

Date Sampled: 01/29/2008 1039

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1143.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0140		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0140			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-230

Sdg Number: 660-20857

Lab Sample ID: 660-20857-43

Date Sampled: 01/29/2008 1039

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57495 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 2HB1143.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/12/2008 0140 Final Weight/Volume: 5 mL
 Date Prepared: 02/12/2008 0140

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	92		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-108

Sdg Number: 660-20857

Lab Sample ID: 660-20857-44

Date Sampled: 01/29/2008 1527

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1144.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0203		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0203			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	1.3	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-108

Lab Sample ID: 660-20857-44

Date Sampled: 01/29/2008 1527

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57495

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1144.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0203

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0203

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U ↓	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U ↓	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	92	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: EB-19

Sdg Number: 660-20857

Lab Sample ID: 660-20857-45

Date Sampled: 01/29/2008 1625

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0814.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1522		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1522			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: EB-19

Sdg Number: 660-20857

Lab Sample ID: 660-20857-45

Date Sampled: 01/29/2008 1625

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0814.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1522		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1522			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	89		74 - 126	
Dibromofluoromethane	83		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-21

Lab Sample ID: 660-20857-46

Date Sampled: 01/29/2008 1550

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0815.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1545		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1545			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: EB-21

Sdg Number: 660-20857

Lab Sample ID: 660-20857-46

Date Sampled: 01/29/2008 1550

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0815.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1545		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1545			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	81		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-20

Lab Sample ID: 660-20857-47

Date Sampled: 01/29/2008 1452

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1126.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/11/2008 1847		Final Weight/Volume: 5 mL
Date Prepared:	02/11/2008 1847		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-20

Lab Sample ID: 660-20857-47

Date Sampled: 01/29/2008 1452

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57504	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1126.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1847		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1847			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	104	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	89	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: EB-18

Sdg Number: 660-20857

Lab Sample ID: 660-20857-48

Date Sampled: 01/29/2008 1600

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0817.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1632		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1632			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-18

Lab Sample ID: 660-20857-48

Date Sampled: 01/29/2008 1600

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0817.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1632		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1632			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	99		77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	77		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-101

Sdg Number: 660-20857

Lab Sample ID: 660-20857-49

Date Sampled: 01/29/2008 0845

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57495

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1145.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0227

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0227

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-101

Lab Sample ID: 660-20857-49

Date Sampled: 01/29/2008 0845

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1145.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0227		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0227			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	90		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-55

Lab Sample ID: 660-20857-50

Date Sampled: 01/29/2008 0855

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	2HB1146.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0250			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0250				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	1.9		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-55

Lab Sample ID: 660-20857-50

Date Sampled: 01/29/2008 0855

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1146.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0250		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0250			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U J	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U J	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	115	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	94	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-150

Sdg Number: 660-20857

Lab Sample ID: 660-20857-51

Date Sampled: 01/29/2008 1240

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1147.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0314		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0314			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-150

Sdg Number: 660-20857

Lab Sample ID: 660-20857-51

Date Sampled: 01/29/2008 1240

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1147.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0314		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0314			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-147

Lab Sample ID: 660-20857-52

Date Sampled: 01/29/2008 1059

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	2HB1148.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0337			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0337				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-147

Lab Sample ID: 660-20857-52

Date Sampled: 01/29/2008 1059

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1148.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0337		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0337			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	116	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	90	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-148

Sdg Number: 660-20857

Lab Sample ID: 660-20857-53

Date Sampled: 01/29/2008 1108

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1149.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0400		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0400			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-148

Sdg Number: 660-20857

Lab Sample ID: 660-20857-53

Date Sampled: 01/29/2008 1108

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57495	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1149.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0400		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0400			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	117		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	95		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-19

Lab Sample ID: 660-20857-54

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57398 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB0818.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/08/2008 1655 Final Weight/Volume: 5 mL
 Date Prepared: 02/08/2008 1655

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-19

Lab Sample ID: 660-20857-54

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0818.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1655			Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1655				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	92	74 - 126
Dibromofluoromethane	75	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: TB-18

Sdg Number: 660-20857

Lab Sample ID: 660-20857-55

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0819.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1718		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1718			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-18

Lab Sample ID: 660-20857-55

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB0819.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1718			Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1718				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	98		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	77		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: TB-20

Sdg Number: 660-20857

Lab Sample ID: 660-20857-56

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57398 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB0820.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/08/2008 1742 Final Weight/Volume: 5 mL
 Date Prepared: 02/08/2008 1742

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-20

Lab Sample ID: 660-20857-56

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB0820.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/08/2008 1742		Final Weight/Volume: 5 mL
Date Prepared:	02/08/2008 1742		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	90	74 - 126
Dibromofluoromethane	75	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-21

Lab Sample ID: 660-20857-57

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57398 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB0824.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/08/2008 1915 Final Weight/Volume: 5 mL
 Date Prepared: 02/08/2008 1915

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.4	U	2.4	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.4	U	2.4	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.64	U	0.64	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.4	U	2.4	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.4	U	2.4	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.44	U	0.44	1.0
cis-1,2-Dichloroethene	0.64	U	0.64	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.4	U	2.4	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: TB-21

Sdg Number: 660-20857

Lab Sample ID: 660-20857-57

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57398	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0824.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 1915		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 1915			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.4	U	2.4	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.84	U	0.84	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	95	77 - 122
4-Bromofluorobenzene	90	74 - 126
Dibromofluoromethane	74	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-24

Lab Sample ID: 660-20857-1

Date Sampled: 01/29/2008 0916

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0825.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1942

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1942

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-146

Lab Sample ID: 660-20857-2

Date Sampled: 01/29/2008 1149

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57510	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB0826.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/08/2008 2005		Final Weight/Volume:	5 mL
Date Prepared:	02/08/2008 2005			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: **MW-56**

Lab Sample ID: 660-20857-3

Date Sampled: 01/29/2008 1557

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0827.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 2029

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 2029

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-88

Lab Sample ID: 660-20857-4

Date Sampled: 01/29/2008 1519

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0828.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 2052

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 2052

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-84B

Lab Sample ID: 660-20857-5

Date Sampled: 01/29/2008 0952

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0829.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 2117

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 2117

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-84A

Lab Sample ID: 660-20857-6

Date Sampled: 01/29/2008 0957

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0830.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 2141

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 2141

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-82

Lab Sample ID: 660-20857-7

Date Sampled: 01/29/2008 1450

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1139.D

Dilution: 10

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0006

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0006

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	580		10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-74

Lab Sample ID: 660-20857-8

Date Sampled: 01/29/2008 1622

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1140.D

Dilution: 5.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0030

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0030

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	130		5.0	5.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: DUP-19

Lab Sample ID: 660-20857-9

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/12/2008 0053
Date Prepared: 02/12/2008 0053

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS
Lab File ID: 2HB1141.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	15		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: DUP-20

Lab Sample ID: 660-20857-10

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1142.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0117

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0117

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-16S

Lab Sample ID: 660-20857-11

Date Sampled: 01/29/2008 1605

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57509

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0929.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 2147

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 2147

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-145

Lab Sample ID: 660-20857-12

Date Sampled: 01/29/2008 1130

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57509

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0930.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 2210

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 2210

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-143

Lab Sample ID: 660-20857-13

Date Sampled: 01/29/2008 1315

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57509

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0931.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 2234

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 2234

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: **MW-140**

Lab Sample ID: 660-20857-14

Date Sampled: 01/29/2008 1010

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57509

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0932.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 2257

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 2257

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-136

Lab Sample ID: 660-20857-15

Date Sampled: 01/29/2008 0850

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57509

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0933.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 2321

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 2321

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-207

Lab Sample ID: 660-20857-16

Date Sampled: 01/29/2008 1415

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57509

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0934.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/09/2008 2344

Final Weight/Volume: 5 mL

Date Prepared: 02/09/2008 2344

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-16D

Lab Sample ID: 660-20857-17

Date Sampled: 01/29/2008 1535

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57645

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 3JB0727.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 0100

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 0100

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	6.0		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-206

Lab Sample ID: 660-20857-18

Date Sampled: 01/29/2008 1405

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57645

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 3JB0728.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 0122

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 0122

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-142

Lab Sample ID: 660-20857-19

Date Sampled: 01/29/2008 1315

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57645

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 3JB0729.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 0145

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 0145

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-141

Lab Sample ID: 660-20857-20

Date Sampled: 01/29/2008 1250

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1166.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1700

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1700

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-144

Lab Sample ID: 660-20857-21

Date Sampled: 01/29/2008 1220

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1167.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1722

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1722

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-96

Lab Sample ID: 660-20857-22

Date Sampled: 01/29/2008 1115

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1168.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1745

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1745

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-139

Lab Sample ID: 660-20857-23

Date Sampled: 01/29/2008 1005

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1209.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2105

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2105

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-137

Lab Sample ID: 660-20857-24

Date Sampled: 01/29/2008 0915

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57649	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	2HB1025.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/10/2008 2206		Final Weight/Volume:	5 mL
Date Prepared:	02/10/2008 2206			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-138

Lab Sample ID: 660-20857-25

Date Sampled: 01/29/2008 0845

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57649

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1026.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/10/2008 2229

Final Weight/Volume: 5 mL

Date Prepared: 02/10/2008 2229

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-176

Lab Sample ID: 660-20857-26

Date Sampled: 01/29/2008 0920

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57649

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1027.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/10/2008 2252

Final Weight/Volume: 5 mL

Date Prepared: 02/10/2008 2252

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-175

Lab Sample ID: 660-20857-27

Date Sampled: 01/29/2008 1130

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57649

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1028.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/10/2008 2316

Final Weight/Volume: 5 mL

Date Prepared: 02/10/2008 2316

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-125

Lab Sample ID: 660-20857-28

Date Sampled: 01/29/2008 1114

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57649

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1029.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/10/2008 2339

Final Weight/Volume: 5 mL

Date Prepared: 02/10/2008 2339

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-13S

Lab Sample ID: 660-20857-29

Date Sampled: 01/29/2008 1305

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57649

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1030.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 0002

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 0002

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	4.5		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1
Sdg Number: 660-20857**Client Sample ID: MW-13D**Lab Sample ID: 660-20857-30
Client Matrix: WaterDate Sampled: 01/29/2008 1244
Date Received: 01/30/2008 1145**8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution**

Method:	8260C SIM/ID	Analysis Batch: 660-57506	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1113.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1342		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1342			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-45

Lab Sample ID: 660-20857-31

Date Sampled: 01/29/2008 1353

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1114.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1406

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1406

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-115

Lab Sample ID: 660-20857-32

Date Sampled: 01/29/2008 1005

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1115.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1429

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1429

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-177

Lab Sample ID: 660-20857-33

Date Sampled: 01/29/2008 0928

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1116.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1453

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1453

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-243

Lab Sample ID: 660-20857-34

Date Sampled: 01/29/2008 1422

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1117.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1516

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1516

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-244

Lab Sample ID: 660-20857-35

Date Sampled: 01/29/2008 1449

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1118.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1539

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1539

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: DUP-18

Lab Sample ID: 660-20857-36

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1119.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1603

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1603

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-231

Lab Sample ID: 660-20857-37

Date Sampled: 01/29/2008 1036

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57506	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1120.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/11/2008 1626		Final Weight/Volume:	5 mL
Date Prepared:	02/11/2008 1626			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-229

Lab Sample ID: 660-20857-38

Date Sampled: 01/29/2008 1042

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1121.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1650

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1650

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-245

Lab Sample ID: 660-20857-39

Date Sampled: 01/29/2008 0933

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1122.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1713

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1713

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-247

Lab Sample ID: 660-20857-40

Date Sampled: 01/29/2008 0850

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1123.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1737

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1737

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-246

Lab Sample ID: 660-20857-41

Date Sampled: 01/29/2008 0900

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1124.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1800

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1800

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: **MW-242**

Lab Sample ID: 660-20857-42

Date Sampled: 01/29/2008 1420

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57506

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1125.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/11/2008 1824

Final Weight/Volume: 5 mL

Date Prepared: 02/11/2008 1824

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-230

Lab Sample ID: 660-20857-43

Date Sampled: 01/29/2008 1039

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1143.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0140

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0140

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: **MW-108**

Lab Sample ID: 660-20857-44

Date Sampled: 01/29/2008 1527

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1144.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0203

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0203

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	12		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-19

Lab Sample ID: 660-20857-45

Date Sampled: 01/29/2008 1625

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0814.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1522

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1522

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-21

Lab Sample ID: 660-20857-46

Date Sampled: 01/29/2008 1550

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0815.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1545

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1545

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-20

Lab Sample ID: 660-20857-47

Date Sampled: 01/29/2008 1452

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0816.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1608

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1608

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: EB-18

Lab Sample ID: 660-20857-48

Date Sampled: 01/29/2008 1600

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0817.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1632

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1632

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-101

Lab Sample ID: 660-20857-49

Date Sampled: 01/29/2008 0845

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1145.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0227

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0227

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.5		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Client Sample ID: MW-55

Sdg Number: 660-20857

Lab Sample ID: 660-20857-50

Date Sampled: 01/29/2008 0855

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1146.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0250

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0250

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	14		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-150

Lab Sample ID: 660-20857-51

Date Sampled: 01/29/2008 1240

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1147.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0314

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0314

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-147

Lab Sample ID: 660-20857-52

Date Sampled: 01/29/2008 1059

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1148.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0337

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0337

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: MW-148

Lab Sample ID: 660-20857-53

Date Sampled: 01/29/2008 1108

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57499

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 2HB1149.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0400

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0400

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-19

Lab Sample ID: 660-20857-54

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0818.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1655

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1655

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-18

Lab Sample ID: 660-20857-55

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0819.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1718

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1718

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-20

Lab Sample ID: 660-20857-56

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0820.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1742

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1742

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20857-1

Sdg Number: 660-20857

Client Sample ID: TB-21

Lab Sample ID: 660-20857-57

Date Sampled: 01/29/2008 0000

Client Matrix: Water

Date Received: 01/30/2008 1145

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57510

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB0824.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/08/2008 1915

Final Weight/Volume: 5 mL

Date Prepared: 02/08/2008 1915

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20857	1/29/2008	SW-846	MW-24	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	DUP-20	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-16S	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-145	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-143	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-140	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-136	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-207	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-16D	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-206	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-142	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-146	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-141	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-144	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-96	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-139	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-137	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-138	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-176	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-175	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-125	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-13S	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-56	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-13D	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-45	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-115	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-177	Water	No	--	--	--	--	VOC – CCAL %D

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20857	1/29/2008	SW-846	MW-243	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-244	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	DUP-18	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-231	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-229	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-245	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-88	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-247	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-246	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-242	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-230	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-108	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	EB-19	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	EB-21	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	EB-20	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	EB-18	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-101	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-84B	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-55	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-150	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-147	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-148	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	TB-19	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	TB-18	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	TB-20	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	TB-21	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-84A	Water	No	--	--	--	--	VOC – CCAL %D
660-20857	1/29/2008	SW-846	MW-82	Water	Yes	--	--	--	--	
660-20857	1/29/2008	SW-846	MW-74	Water	Yes	--	--	--	--	

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20857	1/29/2008	SW-846	DUP-19	Water	Yes	--	--	--	--	

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 6

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20907

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8124R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20907 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-71	660-20907-1	Water	1/30/2008	X				
MW-30	660-20907-10	Water	1/30/2008	X				
MW-48	660-20907-11	Water	1/30/2008	X				
MW-232	660-20907-12	Water	1/30/2008	X				
Dup-22	660-20907-13	Water	1/30/2008	X				
EW-110	660-20907-14	Water	1/30/2008	X				
EW-109	660-20907-15	Water	1/30/2008	X				
EW-106	660-20907-16	Water	1/30/2008	X				
EW-105	660-20907-17	Water	1/30/2008	X				
EW-104	660-20907-18	Water	1/30/2008	X				
EW-103	660-20907-19	Water	1/30/2008	X				
MW-149	660-20907-2	Water	1/30/2008	X				
EW-101	660-20907-20	Water	1/30/2008	X				
MW-166	660-20907-21	Water	1/30/2008	X				
MW-165	660-20907-22	Water	1/30/2008	X				
MW-164	660-20907-23	Water	1/30/2008	X				
Dup-25	660-20907-24	Water	1/30/2008	X				
MW-163	660-20907-25	Water	1/30/2008	X				
MW-162	660-20907-26	Water	1/30/2008	X				
MW-97	660-20907-27	Water	1/30/2008	X				
MW-87	660-20907-28	Water	1/30/2008	X				
MW-75	660-20907-29	Water	1/30/2008	X				
MW-129	660-20907-3	Water	1/30/2008	X				
MW-34	660-20907-30	Water	1/30/2008	X				
MW-35	660-20907-31	Water	1/30/2008	X				
MW-251	660-20907-32	Water	1/30/2008	X				
MW-69	660-20907-33	Water	1/30/2008	X				
MW-61	660-20907-34	Water	1/30/2008	X				
MW-196	660-20907-35	Water	1/30/2008	X				
MW-197	660-20907-36	Water	1/30/2008	X				
Private Well 48	660-20907-37	Water	1/30/2008	X				
Private Well 47	660-20907-38	Water	1/30/2008	X				
EB-22	660-20907-39	Water	1/30/2008	X				
MW-58	660-20907-4	Water	1/30/2008	X				
EB-25	660-20907-40	Water	1/30/2008	X				
EB-EW-1	660-20907-41	Water	1/30/2008	X				
TB-22	660-20907-42	Water	1/30/2008	X				
TB-25	660-20907-43	Water	1/30/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
TB-EW-1	660-20907-44	Water	1/30/2008	X				
EB-23	660-20907-45	Water	1/30/2008	X				
TB-23	660-20907-46	Water	1/30/2008	X				
Dup-23	660-20907-47	Water	1/30/2008	X				
MW-70	660-20907-48	Water	1/30/2008	X				
MW-47	660-20907-49	Water	1/30/2008	X				
MW-130	660-20907-5	Water	1/30/2008	X				
MW-81	660-20907-50	Water	1/30/2008	X				
TB-24	660-20907-51	Water	1/30/2008	X				
Dup-24	660-20907-52	Water	1/30/2008	X				
MW-57	660-20907-6	Water	1/30/2008	X				
MW-80	660-20907-7	Water	1/30/2008	X				
MW-72	660-20907-8	Water	1/30/2008	X				
MW-6	660-20907-9	Water	1/30/2008	X				

Notes:

1. Sample location Dup-22 is the field duplicate of parent sample location MW-166.
2. Sample location Dup-23 is the field duplicate of parent sample location MW-58.
3. Sample location Dup-24 is the field duplicate of parent sample location MW-232.
4. Sample location Dup-25 is the field duplicate of parent sample location MW-61.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

The analyses that exceeded the holding are presented in the following table.

Sample Locations	Holding Time	Criteria
MW-87 (SIM ID) DL	Analysis Completed	23 Days

Sample results reported from the diluted associated sample locations analyzed by analytical method SW-846 8260 were qualified, as specified in the table below. All other holding times were met.

Criteria	Qualification	
	Detected Analytes	Non-detect Analytes
Analysis completed less than two times holding time	J	UJ

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No compounds were detected in the associated blanks.

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
MW-72 Dup-22 EW-110 EW-109 EW-106 EW-105 EW-104 EW-103	CCV %D	1,3-Dichlorobenzene	-21.9%
		1,2,3-Trichlorobenzene	-25.1%
		1,2,4-Trichlorobenzene	-21.0%
		4-Chlorotoluene	-20.2%
		4-Isopropyltoluene	-20.6%
		Hexachlorobutadiene	-27.5%
		n-Butylbenzene	-23.3%
		Private Well 47 EB-22 Dup-23 MW-70 MW-47 MW-81 TB-24 Dup-24	CCV %D
1,3-Dichlorobenzene	-20.8%		
1,2,3-Trichlorobenzene	-29.4%		
1,2,4-Trichlorobenzene	-24.8%		
4-Chlorotoluene	-20.1%		
Bromomethane	67.4%		
Chloroethane	34.5%		
Hexachlorobutadiene	-28.4%		
n-Butylbenzene	-23.3%		
Naphthalene	-22.0%		

Sample Locations	Initial/Continuing	Compound	Criteria
EW-101 MW-166 MW-165 MW-164 Dup-25 MW-163 MW-162 MW-97 MW-87 MW-75 MW-34 MW-35 MW-251 MW-69 MW-61 MW-196 MW-197 Private Well 48	CCV %D	Bromomethane	53.6%
		Carbon disulfide	25.2%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

Sample locations associated with surrogates exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Surrogate	Recovery
EB-22	Toluene-d8	> UL
	4-Bromofluorobenzene	AC
	Dibromofluoromethane	AC
Dup-24	Toluene-d8	AC
	4-Bromofluorobenzene	> UL
	Dibromofluoromethane	AC

Upper control limit (UL)

Lower control limit (LL)

Diluted (D)

Acceptable (AC)

The criteria used to evaluate the surrogate recoveries are presented in the following table. In the case of a surrogate deviation, the sample results associated with the deviant fraction are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	No Action
	Detect	J
< LL but > 10%	Non-detect	J
	Detect	J
< 10%	Non-detect	R
	Detect	J
Surrogates diluted below the calibration curve due to the high concentration of a target compounds	Non-detect	No Action
	Detect	

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-166/Dup-22	All Compounds	ND	ND	AC
MW-58/Dup-23	All Compounds	ND	ND	AC
MW-232/Dup-24	Toluene	0.89 I	ND(0.51)	AC
	1,4-Dioxane (SIM ID)	7.6	7.6	0.0%
MW-61/Dup-25	1,4-Dioxane (SIM ID)	ND(1.0)	1.3	AC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
MW-80	1,1-Dichloroethene	160 E	170 D	170 D
EW-106	cis-1,2-Dichloroethene	580 E	430 D	580 EJ
	Trichloroethene	1500 E	740 D	1500 EJ
EW-104	Trichloroethene	1000 E	500 D	1000 EJ
MW-87	1,1-Dichloroethane	190 E	150 D	190 EJ
	1,1-Dichloroethene	970 E	720 D	970 EJ

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
	Trichloroethene	360 E	200 D	360 EJ
	1,4-Dioxane (SIM ID)	280 E	400 D	400 DJ
MW-81	Trichloroethene	890 E	430 D	890 EJ

Note: All diluted sample results were analyzed beyond the acceptable holding time criteria. Therefore sample results that exhibited a concentration greater than and/or within the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-71

Sdg Number: 660-20907

Lab Sample ID: 660-20907-1

Date Sampled: 01/30/2008 1340

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1122.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0029		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0029			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	9.5		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	2.0		0.45	1.0
cis-1,2-Dichloroethene	70		0.65	1.0
trans-1,2-Dichloroethene	3.2		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-71

Sdg Number: 660-20907

Lab Sample ID: 660-20907-1

Date Sampled: 01/30/2008 1340

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1122.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0029		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0029			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	32		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	95		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-149

Sdg Number: 660-20907

Lab Sample ID: 660-20907-2

Date Sampled: 01/30/2008 1549

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1123.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0051		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0051			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-149

Sdg Number: 660-20907

Lab Sample ID: 660-20907-2

Date Sampled: 01/30/2008 1549

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JB1123.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0051			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0051				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	95		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-129

Sdg Number: 660-20907

Lab Sample ID: 660-20907-3

Date Sampled: 01/30/2008 1431

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1124.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0114		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0114			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-129

Sdg Number: 660-20907

Lab Sample ID: 660-20907-3

Date Sampled: 01/30/2008 1431

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1124.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0114		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0114			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	115		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-58

Sdg Number: 660-20907

Lab Sample ID: 660-20907-4

Date Sampled: 01/30/2008 1436

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1125.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0136		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0136			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-58

Sdg Number: 660-20907

Lab Sample ID: 660-20907-4

Date Sampled: 01/30/2008 1436

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1125.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0136		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0136			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	113		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-130

Sdg Number: 660-20907

Lab Sample ID: 660-20907-5

Date Sampled: 01/30/2008 0922

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57754

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1126.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0159

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0159

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	10		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	28		0.45	1.0
cis-1,2-Dichloroethene	1.1		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-130

Sdg Number: 660-20907

Lab Sample ID: 660-20907-5

Date Sampled: 01/30/2008 0922

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JB1126.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0159			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0159				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	3.7		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	114		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-57

Sdg Number: 660-20907

Lab Sample ID: 660-20907-6

Date Sampled: 01/30/2008 0930

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JB1127.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0221			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0221				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-57

Sdg Number: 660-20907

Lab Sample ID: 660-20907-6

Date Sampled: 01/30/2008 0930

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1127.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0221		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0221			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	430		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	110		74 - 126	
Dibromofluoromethane	102		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-80

Sdg Number: 660-20907

Lab Sample ID: 660-20907-7

Date Sampled: 01/30/2008 1026

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JB1128.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/12/2008 0244		Final Weight/Volume: 5 mL
Date Prepared:	02/12/2008 0244		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	4.1		0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	17		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	6.2		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-80

Sdg Number: 660-20907

Lab Sample ID: 660-20907-7

Date Sampled: 01/30/2008 1026

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1128.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0244		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0244			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Tetrachloroethene	4.5		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	23		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.94	I	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	117		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-80

Sdg Number: 660-20907

Lab Sample ID: 660-20907-7

Date Sampled: 01/30/2008 1026

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57902

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1152.D

Dilution: 2.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1143

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1143

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	170	D	0.90	2.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-72

Sdg Number: 660-20907

Lab Sample ID: 660-20907-8

Date Sampled: 01/30/2008 1027

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1151.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1120		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1120			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-72

Sdg Number: 660-20907

Lab Sample ID: 660-20907-8

Date Sampled: 01/30/2008 1027

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1151.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1120		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1120			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.7		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2.5		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	111		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-6

Sdg Number: 660-20907

Lab Sample ID: 660-20907-9

Date Sampled: 01/30/2008 1118

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1130.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0329		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0329			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-6

Sdg Number: 660-20907

Lab Sample ID: 660-20907-9

Date Sampled: 01/30/2008 1118

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1130.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0329		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0329			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.0		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-30

Sdg Number: 660-20907

Lab Sample ID: 660-20907-10

Date Sampled: 01/30/2008 1620

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1131.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0351		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0351			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-30

Sdg Number: 660-20907

Lab Sample ID: 660-20907-10

Date Sampled: 01/30/2008 1620

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1131.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0351		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0351			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.1		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	18	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	101	77 - 122
4-Bromofluorobenzene	105	74 - 126
Dibromofluoromethane	101	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-48

Sdg Number: 660-20907

Lab Sample ID: 660-20907-11

Date Sampled: 01/30/2008 1206

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1132.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0414		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0414			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-48

Sdg Number: 660-20907

Lab Sample ID: 660-20907-11

Date Sampled: 01/30/2008 1206

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57754

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1132.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0414

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0414

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.88	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	111	77 - 122
4-Bromofluorobenzene	116	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-232

Sdg Number: 660-20907

Lab Sample ID: 660-20907-12

Date Sampled: 01/30/2008 1503

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1133.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0436		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0436			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-232

Sdg Number: 660-20907

Lab Sample ID: 660-20907-12

Date Sampled: 01/30/2008 1503

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57754	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1133.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0436		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0436			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.89	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	108	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Dup-22

Lab Sample ID: 660-20907-13

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JB1153.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/12/2008 1205		Final Weight/Volume: 5 mL
Date Prepared:	02/12/2008 1205		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-22

Sdg Number: 660-20907

Lab Sample ID: 660-20907-13

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JB1153.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/12/2008 1205		Final Weight/Volume: 5 mL
Date Prepared:	02/12/2008 1205		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	123		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-110

Sdg Number: 660-20907

Lab Sample ID: 660-20907-14

Date Sampled: 01/30/2008 1536

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1160.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1446		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1446			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	28	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	68	U	0.45	1.0
cis-1,2-Dichloroethene	5.4	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-110

Sdg Number: 660-20907

Lab Sample ID: 660-20907-14

Date Sampled: 01/30/2008 1536

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1160.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1446		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1446			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	4.0		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	73		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	10	I	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.0		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	113		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-109

Sdg Number: 660-20907

Lab Sample ID: 660-20907-15

Date Sampled: 01/30/2008 1538

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JB1161.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1508			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1508				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-109

Sdg Number: 660-20907

Lab Sample ID: 660-20907-15

Date Sampled: 01/30/2008 1538

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1161.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1508		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1508			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.1		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	113		74 - 126	
Dibromofluoromethane	108		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-106

Sdg Number: 660-20907

Lab Sample ID: 660-20907-16

Date Sampled: 01/30/2008 1208

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 2JB1162.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/12/2008 1530		Final Weight/Volume: 5 mL
Date Prepared:	02/12/2008 1530		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U J	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U J	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U J	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	24	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	43	U	0.45	1.0
cis-1,2-Dichloroethene	580	EJ L	0.65	1.0
trans-1,2-Dichloroethene	1.6	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U J	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-106

Sdg Number: 660-20907

Lab Sample ID: 660-20907-16

Date Sampled: 01/30/2008 1208

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1162.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1530		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1530			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	26		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1500	ES U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.57	I	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	112	74 - 126
Dibromofluoromethane	110	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-106

Sdg Number: 660-20907

Lab Sample ID: 660-20907-16

Date Sampled: 01/30/2008 1208

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58162

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2216.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 1604

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 1604

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	430	Q	13	20
Trichloroethene	740	Q	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-105

Sdg Number: 660-20907

Lab Sample ID: 660-20907-17

Date Sampled: 01/30/2008 1213

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1163.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1553		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1553			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	1.3	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.3	U	0.45	1.0
cis-1,2-Dichloroethene	1.4	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-105

Sdg Number: 660-20907

Lab Sample ID: 660-20907-17

Date Sampled: 01/30/2008 1213

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1163.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1553		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1553			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.79	I	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	19		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	114	74 - 126
Dibromofluoromethane	113	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-104

Sdg Number: 660-20907

Lab Sample ID: 660-20907-18

Date Sampled: 01/30/2008 1447

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1164.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1615		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1615			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	49		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	14		0.45	1.0
cis-1,2-Dichloroethene	71		0.65	1.0
trans-1,2-Dichloroethene	0.89	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-104

Lab Sample ID: 660-20907-18

Date Sampled: 01/30/2008 1447

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1164.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1615		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1615			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,1,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	4.9		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1000	ES	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	112	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-104

Lab Sample ID: 660-20907-18

Date Sampled: 01/30/2008 1447

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58162

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2217.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 1627

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 1627

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	500	Q	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-103

Sdg Number: 660-20907

Lab Sample ID: 660-20907-19

Date Sampled: 01/30/2008 1450

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57902	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	2JB1165.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1638			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1638				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	18	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	4.2	U	0.45	1.0
cis-1,2-Dichloroethene	14	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-103

Sdg Number: 660-20907

Lab Sample ID: 660-20907-19

Date Sampled: 01/30/2008 1450

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57902

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1165.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1638

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1638

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.0		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	31		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	117	74 - 126
Dibromofluoromethane	112	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-101

Lab Sample ID: 660-20907-20

Date Sampled: 01/30/2008 1251

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID: BVMG GC/MS
Preparation:	5030B		Lab File ID: 1GB1312.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 1425		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 1425		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	1.1		0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	3.7		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	3.2		0.45	1.0
cis-1,2-Dichloroethene	14		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-101

Sdg Number: 660-20907

Lab Sample ID: 660-20907-20

Date Sampled: 01/30/2008 1251

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1312.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1425		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1425			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	7.3		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	140		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	101	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-166

Sdg Number: 660-20907

Lab Sample ID: 660-20907-21

Date Sampled: 01/30/2008 1445

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B			Lab File ID:	1GB1313.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1451			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1451				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-166

Sdg Number: 660-20907

Lab Sample ID: 660-20907-21

Date Sampled: 01/30/2008 1445

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID: BVMG GC/MS
Preparation:	5030B		Lab File ID: 1GB1313.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 1451		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 1451		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	119	77 - 122
4-Bromofluorobenzene	107	74 - 126
Dibromofluoromethane	110	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-165

Sdg Number: 660-20907

Lab Sample ID: 660-20907-22

Date Sampled: 01/30/2008 1140

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1314.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1516		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1516			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-165

Sdg Number: 660-20907

Lab Sample ID: 660-20907-22

Date Sampled: 01/30/2008 1140

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B			Lab File ID:	1GB1314.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1516			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1516				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	111	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	109	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-164

Sdg Number: 660-20907

Lab Sample ID: 660-20907-23

Date Sampled: 01/30/2008 1140

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1315.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1540		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1540			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-164

Sdg Number: 660-20907

Lab Sample ID: 660-20907-23

Date Sampled: 01/30/2008 1140

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1315.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1540		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1540			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	117		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-25

Sdg Number: 660-20907

Lab Sample ID: 660-20907-24

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID: BVMG GC/MS
Preparation:	5030B		Lab File ID: 1GB1316.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 1605		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 1605		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-25

Sdg Number: 660-20907

Lab Sample ID: 660-20907-24

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1316.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1605		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1605			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	108		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-163

Sdg Number: 660-20907

Lab Sample ID: 660-20907-25

Date Sampled: 01/30/2008 1436

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1317.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1629		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1629			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-163

Sdg Number: 660-20907

Lab Sample ID: 660-20907-25

Date Sampled: 01/30/2008 1436

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B			Lab File ID:	1GB1317.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1629			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1629				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	115	77 - 122
4-Bromofluorobenzene	104	74 - 126
Dibromofluoromethane	110	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-162

Lab Sample ID: 660-20907-26

Date Sampled: 01/30/2008 1225

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1318.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1654		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1654			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-162

Sdg Number: 660-20907

Lab Sample ID: 660-20907-26

Date Sampled: 01/30/2008 1225

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58411

Instrument ID: BVMG GC/MS

Preparation: 5030B

Lab File ID: 1GB1318.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1654

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1654

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-97

Sdg Number: 660-20907

Lab Sample ID: 660-20907-27

Date Sampled: 01/30/2008 1235

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1319.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1718		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1718			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-97

Lab Sample ID: 660-20907-27

Date Sampled: 01/30/2008 1235

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1319.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1718		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1718			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	101		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-87

Sdg Number: 660-20907

Lab Sample ID: 660-20907-28

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1320.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1743		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1743			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	190	ES U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	970	ES U	0.45	1.0
cis-1,2-Dichloroethene	8.5		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-87

Sdg Number: 660-20907

Lab Sample ID: 660-20907-28

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B			Lab File ID:	1GB1320.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1743			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1743				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	61		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	360	EJ L	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-87

Sdg Number: 660-20907

Lab Sample ID: 660-20907-28

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58162

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2218.D

Dilution: 10

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 1651

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 1651

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethane	150	Q	5.2	10
1,1-Dichloroethene	720	Q	4.5	10
Trichloroethene	200	Q	5.0	10

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-75

Sdg Number: 660-20907

Lab Sample ID: 660-20907-29

Date Sampled: 01/30/2008 0945

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1321.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1808		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1808			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	16		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	5.1		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-75

Sdg Number: 660-20907

Lab Sample ID: 660-20907-29

Date Sampled: 01/30/2008 0945

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1321.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1808		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1808			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	115	77 - 122
4-Bromofluorobenzene	105	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-34

Sdg Number: 660-20907

Lab Sample ID: 660-20907-30

Date Sampled: 01/30/2008 0820

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B			Lab File ID:	1GB1322.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1832			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1832				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-34

Sdg Number: 660-20907

Lab Sample ID: 660-20907-30

Date Sampled: 01/30/2008 0820

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1322.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1832		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1832			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-35

Lab Sample ID: 660-20907-31

Date Sampled: 01/30/2008 0820

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58411

Instrument ID: BVMG GC/MS

Preparation: 5030B

Lab File ID: 1GB1323.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1856

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1856

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	3.7	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-35

Lab Sample ID: 660-20907-31

Date Sampled: 01/30/2008 0820

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58411

Instrument ID: BVMG GC/MS

Preparation: 5030B

Lab File ID: 1GB1323.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1856

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1856

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	15		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	11		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	120		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-251

Sdg Number: 660-20907

Lab Sample ID: 660-20907-32

Date Sampled: 01/30/2008 0957

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1324.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1921		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1921			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-251

Lab Sample ID: 660-20907-32

Date Sampled: 01/30/2008 0957

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58411

Instrument ID: BVMG GC/MS

Preparation: 5030B

Lab File ID: 1GB1324.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1921

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1921

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	116	77 - 122
4-Bromofluorobenzene	102	74 - 126
Dibromofluoromethane	110	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-69

Sdg Number: 660-20907

Lab Sample ID: 660-20907-33

Date Sampled: 01/30/2008 1102

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1325.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1945		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1945			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	1.1		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-69

Lab Sample ID: 660-20907-33

Date Sampled: 01/30/2008 1102

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58411

Instrument ID: BVMG GC/MS

Preparation: 5030B

Lab File ID: 1GB1325.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1945

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1945

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	104	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-61

Lab Sample ID: 660-20907-34

Date Sampled: 01/30/2008 0956

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1326.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2009		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2009			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-61

Sdg Number: 660-20907

Lab Sample ID: 660-20907-34

Date Sampled: 01/30/2008 0956

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1326.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2009		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2009			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	113		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	108		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-196

Lab Sample ID: 660-20907-35

Date Sampled: 01/30/2008 1455

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1327.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2034		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2034			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-196

Lab Sample ID: 660-20907-35

Date Sampled: 01/30/2008 1455

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58411

Instrument ID: BVMG GC/MS

Preparation: 5030B

Lab File ID: 1GB1327.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2034

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2034

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2.2		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-197

Sdg Number: 660-20907

Lab Sample ID: 660-20907-36

Date Sampled: 01/30/2008 1415

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B		Lab File ID:	1GB1328.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2058		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2058			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-197

Sdg Number: 660-20907

Lab Sample ID: 660-20907-36

Date Sampled: 01/30/2008 1415

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58411	Instrument ID: BVMG GC/MS
Preparation:	5030B		Lab File ID: 1GB1328.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 2058		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 2058		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	105		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Private Well 48

Sdg Number: 660-20907

Lab Sample ID: 660-20907-37

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58411 Instrument ID: BVMG GC/MS
 Preparation: 5030B Lab File ID: 1GB1329.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/13/2008 2122 Final Weight/Volume: 5 mL
 Date Prepared: 02/13/2008 2122

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Private Well 48

Lab Sample ID: 660-20907-37

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58411	Instrument ID:	BVMG GC/MS
Preparation:	5030B			Lab File ID:	1GB1329.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2122			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2122				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2.2		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	116	77 - 122
4-Bromofluorobenzene	97	74 - 126
Dibromofluoromethane	107	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Private Well 47

Sdg Number: 660-20907

Lab Sample ID: 660-20907-38

Date Sampled: 01/30/2008 0900

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1246.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1213

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1213

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Private Well 47

Lab Sample ID: 660-20907-38

Date Sampled: 01/30/2008 0900

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1246.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1213

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1213

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1.9		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	113	74 - 126
Dibromofluoromethane	106	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EB-22

Lab Sample ID: 660-20907-39

Date Sampled: 01/30/2008 1500

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1247.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1235

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1235

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U ✓	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U ✓	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U ✓	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U ✓	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U ✓	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U ✓	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EB-22

Lab Sample ID: 660-20907-39

Date Sampled: 01/30/2008 1500

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1247.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1235

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1235

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	125	J1	77 - 122
4-Bromofluorobenzene	124		74 - 126
Dibromofluoromethane	118		70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EB-25

Lab Sample ID: 660-20907-40

Date Sampled: 01/30/2008 1525

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1227.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 2310		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 2310			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EB-25

Lab Sample ID: 660-20907-40

Date Sampled: 01/30/2008 1525

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1227.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 2310		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 2310			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	94	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EB-EW-1

Sdg Number: 660-20907

Lab Sample ID: 660-20907-41

Date Sampled: 01/30/2008 1620

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1228.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 2333			Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 2333				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EB-EW-1

Lab Sample ID: 660-20907-41

Date Sampled: 01/30/2008 1620

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57611

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1228.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2333

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2333

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	108	77 - 122
4-Bromofluorobenzene	97	74 - 126
Dibromofluoromethane	96	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-22

Sdg Number: 660-20907

Lab Sample ID: 660-20907-42

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57611

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1229.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2356

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2356

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-22

Sdg Number: 660-20907

Lab Sample ID: 660-20907-42

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1229.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 2356		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 2356			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	98	77 - 122
4-Bromofluorobenzene	91	74 - 126
Dibromofluoromethane	88	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-25

Sdg Number: 660-20907

Lab Sample ID: 660-20907-43

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1230.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 0019			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 0019				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-25

Lab Sample ID: 660-20907-43

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57611

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1230.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0019

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0019

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	91	74 - 126
Dibromofluoromethane	90	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-EW-1

Lab Sample ID: 660-20907-44

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1231.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 0042		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 0042			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-EW-1

Sdg Number: 660-20907

Lab Sample ID: 660-20907-44

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57611

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1231.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0042

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0042

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	113		77 - 122	
4-Bromofluorobenzene	94		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EB-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-45

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1232.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 0105		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 0105			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EB-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-45

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57611	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1232.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 0105		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 0105			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	98	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-46
Client Matrix: WaterDate Sampled: 01/30/2008 0000
Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-57611 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB1233.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/13/2008 0128 Final Weight/Volume: 5 mL
 Date Prepared: 02/13/2008 0128

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-23

Lab Sample ID: 660-20907-46

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57611

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1233.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0128

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0128

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	98		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-47

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1248.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 1258		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 1258		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-47

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1248.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 1258		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 1258		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	112		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-70

Sdg Number: 660-20907

Lab Sample ID: 660-20907-48

Date Sampled: 01/30/2008 1257

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1249.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1320

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1320

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	0.58	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	17	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	4.6	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	I	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-70

Sdg Number: 660-20907

Lab Sample ID: 660-20907-48

Date Sampled: 01/30/2008 1257

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1249.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1320			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1320				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	20		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	310		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	114		74 - 126	
Dibromofluoromethane	111		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-47

Sdg Number: 660-20907

Lab Sample ID: 660-20907-49

Date Sampled: 01/30/2008 1705

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1250.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1342			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1342				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	2.2	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-47

Sdg Number: 660-20907

Lab Sample ID: 660-20907-49

Date Sampled: 01/30/2008 1705

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1250.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 1342		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 1342		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	104	77 - 122
4-Bromofluorobenzene	108	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-81

Lab Sample ID: 660-20907-50

Date Sampled: 01/30/2008 1106

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1251.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1405		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1405			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	8.1		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	78		0.45	1.0
cis-1,2-Dichloroethene	49		0.65	1.0
trans-1,2-Dichloroethene	0.44	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	9.9		4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-81

Sdg Number: 660-20907

Lab Sample ID: 660-20907-50

Date Sampled: 01/30/2008 1106

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1251.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1405		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1405			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	890	ES +	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	115		74 - 126	
Dibromofluoromethane	114		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-81

Sdg Number: 660-20907

Lab Sample ID: 660-20907-50

Date Sampled: 01/30/2008 1106

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2121.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2109

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2109

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	430	Q	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-24

Sdg Number: 660-20907

Lab Sample ID: 660-20907-51

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1252.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1427			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1427				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-24

Sdg Number: 660-20907

Lab Sample ID: 660-20907-51

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1252.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1427			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1427				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	111		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-24

Sdg Number: 660-20907

Lab Sample ID: 660-20907-52

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1253.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1449		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1449			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U J	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U J	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U J	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U J	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U J	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U J	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Dup-24

Lab Sample ID: 660-20907-52

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1253.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1449		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1449			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	119		77 - 122	
4-Bromofluorobenzene	128	J1	74 - 126	
Dibromofluoromethane	121		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-71

Lab Sample ID: 660-20907-1

Date Sampled: 01/30/2008 1340

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57643	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1149.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1035		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1035			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	5.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-149

Lab Sample ID: 660-20907-2

Date Sampled: 01/30/2008 1549

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1150.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1058

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1058

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-129

Lab Sample ID: 660-20907-3

Date Sampled: 01/30/2008 1431

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57755	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1124.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 0114		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 0114			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	4.2		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-58

Sdg Number: 660-20907

Lab Sample ID: 660-20907-4

Date Sampled: 01/30/2008 1436

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1125.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0136

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0136

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-130

Lab Sample ID: 660-20907-5

Date Sampled: 01/30/2008 0922

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1126.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0159

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0159

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	92		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-57

Sdg Number: 660-20907

Lab Sample ID: 660-20907-6

Date Sampled: 01/30/2008 0930

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1127.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0221

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0221

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-80

Lab Sample ID: 660-20907-7

Date Sampled: 01/30/2008 1026

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1128.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0244

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0244

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	95		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-72

Lab Sample ID: 660-20907-8

Date Sampled: 01/30/2008 1027

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1151.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1120

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1120

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-6

Sdg Number: 660-20907

Lab Sample ID: 660-20907-9

Date Sampled: 01/30/2008 1118

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1130.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0329

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0329

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-30

Lab Sample ID: 660-20907-10

Date Sampled: 01/30/2008 1620

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1131.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0351

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0351

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-48

Sdg Number: 660-20907

Lab Sample ID: 660-20907-11

Date Sampled: 01/30/2008 1206

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1132.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 0414

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 0414

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	6.1		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-232

Lab Sample ID: 660-20907-12

Date Sampled: 01/30/2008 1503

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/12/2008 0436
Date Prepared: 02/12/2008 0436

Analysis Batch: 660-57755

Instrument ID: BVMJ GC/MS

Lab File ID: 2JB1133.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Dup-22

Lab Sample ID: 660-20907-13

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1153.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1205

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1205

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-110

Lab Sample ID: 660-20907-14

Date Sampled: 01/30/2008 1536

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57643	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	2JB1160.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/12/2008 1446		Final Weight/Volume:	5 mL
Date Prepared:	02/12/2008 1446			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	27		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-109

Lab Sample ID: 660-20907-15

Date Sampled: 01/30/2008 1538

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1161.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1508

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1508

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-106

Sdg Number: 660-20907

Lab Sample ID: 660-20907-16

Date Sampled: 01/30/2008 1208

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1162.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1530

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1530

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	20		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-105

Lab Sample ID: 660-20907-17

Date Sampled: 01/30/2008 1213

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1163.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1553

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1553

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	3.9		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EW-104

Sdg Number: 660-20907

Lab Sample ID: 660-20907-18

Date Sampled: 01/30/2008 1447

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1164.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1615

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1615

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	25		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-103

Lab Sample ID: 660-20907-19

Date Sampled: 01/30/2008 1450

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57643

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 2JB1165.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 1638

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 1638

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EW-101

Lab Sample ID: 660-20907-20

Date Sampled: 01/30/2008 1251

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1214.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2256

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2256

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-166

Sdg Number: 660-20907

Lab Sample ID: 660-20907-21

Date Sampled: 01/30/2008 1445

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1215.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2319

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2319

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-165

Sdg Number: 660-20907

Lab Sample ID: 660-20907-22

Date Sampled: 01/30/2008 1140

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1216.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2341

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2341

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-164

Lab Sample ID: 660-20907-23

Date Sampled: 01/30/2008 1140

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1217.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0003

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0003

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: Dup-25

Sdg Number: 660-20907

Lab Sample ID: 660-20907-24

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1218.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0025

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0025

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.3		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-163

Lab Sample ID: 660-20907-25

Date Sampled: 01/30/2008 1436

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1219.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0047

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0047

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-162

Sdg Number: 660-20907

Lab Sample ID: 660-20907-26

Date Sampled: 01/30/2008 1225

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1220.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0109

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0109

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-97

Sdg Number: 660-20907

Lab Sample ID: 660-20907-27

Date Sampled: 01/30/2008 1235

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57608	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1221.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 0132		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 0132			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-87

Lab Sample ID: 660-20907-28

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1222.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0154

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0154

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	280 400	← DJ	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-87

Sdg Number: 660-20907

Lab Sample ID: 660-20907-28

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58129

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2218.D

Dilution: 10

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 1651

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 1651

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	400	Q	10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-75

Sdg Number: 660-20907

Lab Sample ID: 660-20907-29

Date Sampled: 01/30/2008 0945

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1223.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0216

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0216

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	33		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-34

Sdg Number: 660-20907

Lab Sample ID: 660-20907-30

Date Sampled: 01/30/2008 0820

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1224.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0238

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0238

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-35

Sdg Number: 660-20907

Lab Sample ID: 660-20907-31

Date Sampled: 01/30/2008 0820

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1225.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0300

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0300

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-251

Lab Sample ID: 660-20907-32

Date Sampled: 01/30/2008 0957

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1226.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0322

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0322

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-69

Sdg Number: 660-20907

Lab Sample ID: 660-20907-33

Date Sampled: 01/30/2008 1102

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1227.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0344

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0344

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.4		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-61

Lab Sample ID: 660-20907-34

Date Sampled: 01/30/2008 0956

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1228.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0407

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0407

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: MW-196

Sdg Number: 660-20907

Lab Sample ID: 660-20907-35

Date Sampled: 01/30/2008 1455

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1229.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0429

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0429

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-197

Lab Sample ID: 660-20907-36

Date Sampled: 01/30/2008 1415

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/13/2008 0451
Date Prepared: 02/13/2008 0451

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1230.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Private Well 48

Lab Sample ID: 660-20907-37

Date Sampled: 01/30/2008 0855

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57608

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1231.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0513

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0513

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Private Well 47

Lab Sample ID: 660-20907-38

Date Sampled: 01/30/2008 0900

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1246.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1213

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1213

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.7		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EB-22

Sdg Number: 660-20907

Lab Sample ID: 660-20907-39

Date Sampled: 01/30/2008 1500

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1247.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1235

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1235

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: EB-25

Lab Sample ID: 660-20907-40

Date Sampled: 01/30/2008 1525

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1227.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2310

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2310

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EB-EW-1

Sdg Number: 660-20907

Lab Sample ID: 660-20907-41

Date Sampled: 01/30/2008 1620

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/12/2008 2333
 Date Prepared: 02/12/2008 2333

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Lab File ID: 1HB1228.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-22

Lab Sample ID: 660-20907-42

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1229.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/12/2008 2356

Final Weight/Volume: 5 mL

Date Prepared: 02/12/2008 2356

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-25

Lab Sample ID: 660-20907-43

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Preparation: 5030B .

Lab File ID: 1HB1230.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0019

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0019

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-EW-1

Lab Sample ID: 660-20907-44

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/13/2008 0042
Date Prepared: 02/13/2008 0042

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Lab File ID: 1HB1231.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: EB-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-45

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1232.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0105

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0105

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Client Sample ID: TB-23

Sdg Number: 660-20907

Lab Sample ID: 660-20907-46

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57599

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1233.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 0128

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 0128

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Dup-23

Lab Sample ID: 660-20907-47

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1248.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1258

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1258

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-70

Lab Sample ID: 660-20907-48

Date Sampled: 01/30/2008 1257

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/13/2008 1320
Date Prepared: 02/13/2008 1320

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1249.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	3.1		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: **MW-47**

Lab Sample ID: 660-20907-49

Date Sampled: 01/30/2008 1705

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1250.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1342

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1342

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	39		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: MW-81

Lab Sample ID: 660-20907-50

Date Sampled: 01/30/2008 1106

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 20
Date Analyzed: 02/13/2008 1929
Date Prepared: 02/13/2008 1929

Analysis Batch: 660-57632

Instrument ID: BVMH GC/MS
Lab File ID: 1HB1325.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	150		20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: TB-24

Lab Sample ID: 660-20907-51

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1252.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1427

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1427

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20907-1

Sdg Number: 660-20907

Client Sample ID: Dup-24

Lab Sample ID: 660-20907-52

Date Sampled: 01/30/2008 0000

Client Matrix: Water

Date Received: 01/31/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/13/2008 1449
Date Prepared: 02/13/2008 1449

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1253.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.6		1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20907	1/30/2008	SW-846	MW-71	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-30	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-48	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-232	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	Dup-22	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EW-110	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EW-109	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EW-106	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EW-105	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EW-104	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EW-103	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-149	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	EW-101	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-166	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-165	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-164	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	Dup-25	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-163	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-162	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-97	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-87	Water	No	--	--	--	--	VOC – Holding Time
660-20907	1/30/2008	SW-846	MW-75	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-129	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-34	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-35	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-251	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-69	Water	Yes	--	--	--	--	

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20907	1/30/2008	SW-846	MW-61	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-196	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-197	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	Private Well 48	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	Private Well 47	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	EB-22	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-58	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	EB-25	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	EB-EW-1	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	TB-22	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	TB-25	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	TB-EW-1	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	EB-23	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	TB-23	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	Dup-23	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-70	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-47	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-130	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-81	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	TB-24	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	Dup-24	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-57	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-80	Water	Yes	--	--	--	--	
660-20907	1/30/2008	SW-846	MW-72	Water	No	--	--	--	--	VOC – CCAL %D
660-20907	1/30/2008	SW-846	MW-6	Water	Yes	--	--	--	--	

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 7

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20926

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8160R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20926 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-249	660-20926-1	Water	1/31/2008	X				
MW-237	660-20926-10	Water	1/31/2008	X				
MW-238	660-20926-11	Water	1/31/2008	X				
MW-213	660-20926-12	Water	1/31/2008	X				
MW-62	660-20926-13	Water	1/31/2008	X				
MW-92	660-20926-14	Water	1/31/2008	X				
MW-46	660-20926-15	Water	1/31/2008	X				
MW-63	660-20926-16	Water	1/31/2008	X				
MW-131	660-20926-17	Water	1/31/2008	X				
PZ-LSAS-1	660-20926-18	Water	1/31/2008	X				
PZ-LSAS-2	660-20926-19	Water	1/31/2008	X				
MW-250	660-20926-2	Water	1/31/2008	X				
PZ-LSAS-3	660-20926-20	Water	1/31/2008	X				
PZ-LSAS-4	660-20926-21	Water	1/31/2008	X				
PZ-LSAS-5	660-20926-22	Water	1/31/2008	X				
PZ-LSAS-6	660-20926-23	Water	1/31/2008	X				
PZ-LSAS-7	660-20926-24	Water	1/31/2008	X				
Dup-30	660-20926-25	Water	1/31/2008	X				
EB-30	660-20926-26	Water	1/31/2008	X				
TB-30	660-20926-27	Water	1/31/2008	X				
Dup-27	660-20926-28	Water	1/31/2008	X				
EB-27	660-20926-29	Water	1/31/2008	X				
MW-214	660-20926-3	Water	1/31/2008	X				
MW-40	660-20926-30	Water	1/31/2008	X				
MW-41	660-20926-31	Water	1/31/2008	X				
MW-33	660-20926-32	Water	1/31/2008	X				
EW-102	660-20926-33	Water	1/31/2008	X				
EW-107	660-20926-34	Water	1/31/2008	X				
EW-108	660-20926-35	Water	1/31/2008	X				
MW-10	660-20926-36	Water	1/31/2008	X				
MW-36	660-20926-37	Water	1/31/2008	X				
MW-37	660-20926-38	Water	1/31/2008	X				
MW-127	660-20926-39	Water	1/31/2008	X				
MW-215	660-20926-4	Water	1/31/2008	X				
MW-128	660-20926-40	Water	1/31/2008	X				
Dup-28	660-20926-41	Water	1/31/2008	X				
EB-28	660-20926-42	Water	1/31/2008	X				
TB-28	660-20926-43	Water	1/31/2008	X				

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
Dup-29	660-20926-44	Water	1/31/2008	X				
EB-29	660-20926-45	Water	1/31/2008	X				
TB-29	660-20926-46	Water	1/31/2008	X				
Private Well 85	660-20926-47	Water	1/31/2008	X				
MW-19	660-20926-48	Water	1/31/2008	X				
MW-252	660-20926-49	Water	1/31/2008	X				
MW-216	660-20926-5	Water	1/31/2008	X				
MW-217	660-20926-6	Water	1/31/2008	X				
MW-234	660-20926-7	Water	1/31/2008	X				
MW-235	660-20926-8	Water	1/31/2008	X				
MW-236	660-20926-9	Water	1/31/2008	X				

Notes:

1. Sample location Dup-27 is the field duplicate of parent sample location MW-46.
2. Sample location Dup-28 is the field duplicate of parent sample location MW-128.
3. Sample location Dup-29 is the field duplicate of parent sample location EW-107.
4. Sample location Dup-30 is the field duplicate of parent sample location MW-237.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

The analyses that exceeded the holding are presented in the following table.

Sample Locations	Holding Time	Criteria
PZ-LSAS-2 (SIM ID) DL PZ-LSAS-4 (SIM ID) DL Dup-30 (SIM ID)	Analysis Completed	18 Days
MW-37 DL MW-37 (SIM ID) DL MW-127 DL MW-127 (SIM ID) DL		19 Days
PZ-LSAS-5 DL PZ-LSAS-5 (SIM ID) PZ-LSAS-6 (SIM ID) PZ-LSAS-7 DL PZ-LSAS-7 (SIM ID) Dup-27 (SIM ID) EB-27 (SIM ID) MW-40 (SIM ID) MW-41 (SIM ID) MW-33 DL MW-33 (SIM ID) EW-102 (SIM ID) DL EW-107 (SIM ID)		21 Days
MW-213 EW-108 (SIM ID)		22 Days
PZ-LSAS-6 DL		25 Days

Sample results associated with sample locations analyzed by analytical method SW-846 8260 were qualified, as specified in the table below. All other holding times were met.

Criteria	Qualification	
	Detected Analytes	Non-detect Analytes
Analysis completed less than two times holding time	J	UJ

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

Target compounds were detected in the associated QA blanks. Sample results associated with blank contamination that were greater than the BAL and/or non-detect did not result in any qualification of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
MW-238 MW-128 Dup-28	Acetone	Detected sample results <RL and <BAL	"U" at the PQL
EW-108	Acetone	Detected sample results >RL and <BAL	"U" at detected sample concentration
MW-131 EW-102 MW-10	Trichloroethene	Detected sample results >RL and <BAL	"U" at detected sample concentration

RL = reporting limit

3. Mass Spectrometer Tuning

The initial 1,4-dioxane (SIM ID) analysis for sample locations MW-213, PZ-LSAS-4, Dup-30, Dup-27, EB-27 and EW-107 were analyzed outside of the 12-hour tune clock. These samples were reanalyzed outside of holding time and reported. The reported results were qualified as estimated.

The initial TIM analysis for sample location MW-213 was analyzed outside of the 12-hour tune clock. This sample was reanalyzed outside of holding time and reported. The reported results were qualified as estimated.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the

control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
MW-249 MW-250 MW-214 MW-215 MW-216 MW-217 MW-234 MW-235 MW-236 MW-237 MW-238	CCV %D	1,2-Dibromo-3-chloropropane	-26.0%
		1,3-Dichlorobenzene	-20.8%
		1,2,3-Trichlorobenzene	-24.8%
		1,2,4-Trichlorobenzene	-29.4%
		4-Chlorotoluene	-20.1%
		Bromomethane	67.4%
		Chloroethane	34.5%
		Hexachlorobutadiene	-28.4%
		n-Butylbenzene	-23.3%
		Naphthalene	-22.0%
		MW-213	CCV %D
Chloroethane	-21.9%		
Chloromethane	-30.0%		
Methylene chloride	-20.5%		

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J

Initial/Continuing	Criteria	Sample Result	Qualification
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

Sample locations associated with surrogates exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Surrogate	Recovery
MW-33	Toluene-d8	> UL
	4-Bromofluorobenzene	AC
	Dibromofluoromethane	AC

Upper control limit (UL)

Lower control limit (LL)

Diluted (D)

Acceptable (AC)

The criteria used to evaluate the surrogate recoveries are presented in the following table. In the case of a surrogate deviation, the sample results associated with the deviant fraction are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	No Action
	Detect	J
< LL but > 10%	Non-detect	J
	Detect	J
< 10%	Non-detect	R
	Detect	J
Surrogates diluted below the calibration curve due to the high concentration of a target compounds	Non-detect	No Action
	Detect	

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-46/Dup-27	All Compounds	ND	ND	AC
MW-128/Dup-28	1,1-Dichloroethane	2.3	2.6	AC
	1,1-Dichloroethene	3.6	4	AC
	Carbon disulfide	0.99 l	1.3	AC
	cis-1,2-Dichloroethene	2.5	2.5	AC
	Toluene	1.5	1.5	AC
	Trichloroethene	34	35	2.8%
	1,4-Dioxane (SIM ID)	7.2	7.6	5.4%
EW-107/Dup-29	Carbon disulfide	6.4	5.1	22.6%
	Toluene	0.52 l	ND(0.51)	AC
	Trichloroethene	6.1	4.2	AC
MW-237/Dup-30	All Compounds	ND	ND	AC

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
PZ-LSAS-1	Trichloroethene	310 E	280 D	310 EJ
	1,4-Dioxane (SIM ID)	660 E	470 D	660 EJ
PZ-LSAS-2	cis-1,2-Dichloroethene	680 E	540 D	680 EJ
	Trichloroethene	1000 E	880 D	1000 EJ
	1,4-Dioxane (SIM ID)	570 E	580 D	580 DJ
PZ-LSAS-3	Trichloroethene	2500 E	1900 D	2500 EJ
PZ-LSAS-4	cis-1,2-Dichloroethene	730 E	560 D	730 EJ
	Trichloroethene	4100 E	3800 D	4100 EJ
PZ-LSAS-5	cis-1,2-Dichloroethene	690 E	290 D	690 EJ
	Trichloroethene	2100 E	3400 D	3400 DJ
PZ-LSAS-6	cis-1,2-Dichloroethene	510 E	120 D	510 EJ
	Trichloroethene	2500 E	8900 D	8900 DJ
PZ-LSAS-7	cis-1,2-Dichloroethene	480 E	70 D	480 EJ
	Trichloroethene	1900 E	2300 D	2300 DJ
MW-40	Tetrachloroethene	160 E	150 D	160 EJ
	Trichloroethene	970 E	880 D	970 EJ
MW-41	Trichloroethene	250 E	190 D	250 EJ
MW-33	1,1-Dichloroethane	160 E	140 D	160 EJ
	1,1-Dichloroethene	570 E	580 D	580 DJ
	Trichloroethene	410 E	290 D	410 EJ
EW-108	cis-1,2-Dichloroethene	930 E	460 D	930 EJ
	Carbon disulfide	100 E	85 D	100 EJ
MW-36	Trichloroethene	190 E	130 D	190 EJ
MW-37	1,1-Dichloroethene	230 E	230 D	230 DJ
	cis-1,2-Dichloroethene	420 E	400 D	420 EJ

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
	Trichloroethene	4500 E	9400 D	9400 DJ
	1,4-Dioxane (SIM ID)	300 E	760 D	760 DJ
MW-127	1,1-Dichloroethene	240 E	270 D	270 DJ
	Trichloroethene	3200 E	3300 D	3300 DJ
	1,4-Dioxane (SIM ID)	200 E	300 D	300 DJ
EW-102	1,4-Dioxane (SIM ID)	120 E	120 D	120 DJ

Note: All diluted sample results were analyzed beyond the acceptable holding time criteria. Therefore sample results that exhibited a concentration greater than and/or within the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

Sample locations EB-30 and TB-30 were analyzed for 1,4-dioxane (SIM ID) by the laboratory but were not reported due to technical issues.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-249

Sdg Number: 660-20926

Lab Sample ID: 660-20926-1

Date Sampled: 01/31/2008 1428

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1254.D

Dilution: 1.0

Date Analyzed: 02/13/2008 1512

Initial Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1512

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	3.0		0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	4.4		0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-249

Sdg Number: 660-20926

Lab Sample ID: 660-20926-1

Date Sampled: 01/31/2008 1428

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1254.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1512		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1512			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	107		74 - 126	
Dibromofluoromethane	111		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-250

Sdg Number: 660-20926

Lab Sample ID: 660-20926-2

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1255.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1534			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1534				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-250

Sdg Number: 660-20926

Lab Sample ID: 660-20926-2

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1255.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1534			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1534				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	114		77 - 122	
4-Bromofluorobenzene	115		74 - 126	
Dibromofluoromethane	116		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-214

Sdg Number: 660-20926

Lab Sample ID: 660-20926-3

Date Sampled: 01/31/2008 1535

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1256.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1557		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1557			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	5.2	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-214

Sdg Number: 660-20926

Lab Sample ID: 660-20926-3

Date Sampled: 01/31/2008 1535

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1256.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1557		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1557			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	107		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-215

Sdg Number: 660-20926

Lab Sample ID: 660-20926-4

Date Sampled: 01/31/2008 1545

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1257.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1619

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1619

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	0.58	1.0
n-Butylbenzene	0.67	U	2.5	5.0
sec-Butylbenzene	0.63	U	0.67	1.0
tert-Butylbenzene	0.84	U	0.63	1.0
Carbon tetrachloride	0.42	U	0.84	1.0
Chlorobenzene	0.63	U	0.42	1.0
Chloroethane	2.5	U	0.63	1.0
Chloroform	0.90	U	2.5	5.0
Chloromethane	1.0	U	0.90	1.0
2-Chlorotoluene	0.65	U	1.0	4.0
4-Chlorotoluene	0.52	U	0.65	1.0
Chlorodibromomethane	0.34	U	0.52	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	0.34	1.0
Ethylene Dibromide	0.50	U	2.5	5.0
Dibromomethane	0.41	U	0.50	1.0
1,2-Dichlorobenzene	0.44	U	0.41	1.0
1,3-Dichlorobenzene	0.64	U	0.44	1.0
1,4-Dichlorobenzene	0.52	U	0.64	1.0
Dichlorodifluoromethane	2.5	U	0.52	1.0
1,1-Dichloroethane	0.52	U	2.5	5.0
1,2-Dichloroethane	0.57	U	0.52	1.0
1,1-Dichloroethene	0.45	U	0.57	1.0
cis-1,2-Dichloroethene	0.65	U	0.45	1.0
trans-1,2-Dichloroethene	0.44	U	0.65	1.0
1,2-Dichloropropane	0.52	U	0.44	1.0
1,3-Dichloropropane	0.39	U	0.52	1.0
2,2-Dichloropropane	0.36	U	0.39	1.0
1,1-Dichloropropene	0.31	U	0.36	1.0
cis-1,3-Dichloropropene	0.14	U	0.31	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.14	1.0
Hexachlorobutadiene	0.40	U	0.44	1.0
Isopropylbenzene	0.19	U	0.40	1.0
4-Isopropyltoluene	0.69	U	0.19	1.0
Methylene Chloride	4.0	U	0.69	1.0
Naphthalene	2.5	U	4.0	5.0
N-Propylbenzene	0.59	U	2.5	5.0
Styrene	0.98	U	0.59	1.0
1,1,1,2-Tetrachloroethane	0.63	U	0.98	2.0
			0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-215

Sdg Number: 660-20926

Lab Sample ID: 660-20926-4

Date Sampled: 01/31/2008 1545

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1257.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1619

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1619

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	112		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-216

Sdg Number: 660-20926

Lab Sample ID: 660-20926-5

Date Sampled: 01/31/2008 1525

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1258.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1641

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1641

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	0.58	1.0
n-Butylbenzene	0.67	U	2.5	5.0
sec-Butylbenzene	0.63	U	0.67	1.0
tert-Butylbenzene	0.63	U	0.63	1.0
Carbon tetrachloride	0.84	U	0.84	1.0
Chlorobenzene	0.42	U	0.42	1.0
Chloroethane	0.63	U	0.63	1.0
Chloroform	2.5	U	2.5	5.0
Chloromethane	0.90	U	0.90	1.0
2-Chlorotoluene	1.0	U	1.0	4.0
4-Chlorotoluene	0.65	U	0.65	1.0
Chlorodibromomethane	0.52	U	0.52	1.0
1,2-Dibromo-3-Chloropropane	0.34	U	0.34	1.0
Ethylene Dibromide	2.5	U	2.5	5.0
Dibromomethane	0.50	U	0.50	1.0
1,2-Dichlorobenzene	0.41	U	0.41	1.0
1,3-Dichlorobenzene	0.44	U	0.44	1.0
1,4-Dichlorobenzene	0.64	U	0.64	1.0
Dichlorodifluoromethane	0.52	U	0.52	1.0
1,1-Dichloroethane	2.5	U	2.5	5.0
1,2-Dichloroethane	0.52	U	0.52	1.0
1,1-Dichloroethene	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	0.45	U	0.45	1.0
trans-1,2-Dichloroethene	0.65	U	0.65	1.0
1,2-Dichloropropane	0.44	U	0.44	1.0
1,3-Dichloropropane	0.52	U	0.52	1.0
2,2-Dichloropropane	0.39	U	0.39	1.0
1,1-Dichloropropene	0.36	U	0.36	1.0
cis-1,3-Dichloropropene	0.31	U	0.31	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.14	U	0.14	1.0
Hexachlorobutadiene	0.44	U	0.44	1.0
Isopropylbenzene	0.40	U	0.40	1.0
4-Isopropyltoluene	0.19	U	0.19	1.0
Methylene Chloride	0.69	U	0.69	1.0
Naphthalene	4.0	U	4.0	5.0
N-Propylbenzene	2.5	U	2.5	5.0
Styrene	0.59	U	0.59	1.0
1,1,1,2-Tetrachloroethane	0.98	U	0.98	2.0
	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-216

Sdg Number: 660-20926

Lab Sample ID: 660-20926-5

Date Sampled: 01/31/2008 1525

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1258.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1641

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1641

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	106		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-217

Sdg Number: 660-20926

Lab Sample ID: 660-20926-6

Date Sampled: 01/31/2008 1625

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1259.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1704		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1704			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-217

Sdg Number: 660-20926

Lab Sample ID: 660-20926-6

Date Sampled: 01/31/2008 1625

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1259.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1704		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1704			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-234

Sdg Number: 660-20926

Lab Sample ID: 660-20926-7

Date Sampled: 01/31/2008 1100

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1260.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1726

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1726

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	5.7	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-234

Sdg Number: 660-20926

Lab Sample ID: 660-20926-7

Date Sampled: 01/31/2008 1100

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1260.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1726

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1726

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	107		74 - 126	
Dibromofluoromethane	110		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-235

Sdg Number: 660-20926

Lab Sample ID: 660-20926-8

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1261.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1748		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1748			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-235

Sdg Number: 660-20926

Lab Sample ID: 660-20926-8

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1261.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1748		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1748			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	108		74 - 126	
Dibromofluoromethane	112		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-236

Sdg Number: 660-20926

Lab Sample ID: 660-20926-9

Date Sampled: 01/31/2008 1108

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58005

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1262.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1811

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1811

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	0.58	1.0
n-Butylbenzene	0.67	U	2.5	5.0
sec-Butylbenzene	0.63	U	0.67	1.0
tert-Butylbenzene	0.84	U	0.63	1.0
Carbon tetrachloride	0.42	U	0.84	1.0
Chlorobenzene	0.63	U	0.42	1.0
Chloroethane	2.5	U	0.63	1.0
Chloroform	0.90	U	2.5	5.0
Chloromethane	1.0	U	0.90	1.0
2-Chlorotoluene	0.65	U	1.0	4.0
4-Chlorotoluene	0.52	U	0.65	1.0
Chlorodibromomethane	0.34	U	0.52	1.0
1,2-Dibromo-3-Chloropropane	0.34	U	0.34	1.0
Ethylene Dibromide	2.5	U	2.5	5.0
Dibromomethane	0.50	U	0.50	1.0
1,2-Dichlorobenzene	0.41	U	0.41	1.0
1,3-Dichlorobenzene	0.44	U	0.44	1.0
1,4-Dichlorobenzene	0.64	U	0.64	1.0
Dichlorodifluoromethane	0.52	U	0.52	1.0
1,1-Dichloroethane	2.5	U	2.5	5.0
1,2-Dichloroethane	0.52	U	0.52	1.0
1,1-Dichloroethene	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	0.45	U	0.45	1.0
trans-1,2-Dichloroethene	0.65	U	0.65	1.0
1,2-Dichloropropane	0.44	U	0.44	1.0
1,3-Dichloropropane	0.52	U	0.52	1.0
2,2-Dichloropropane	0.39	U	0.39	1.0
1,1-Dichloropropene	0.36	U	0.36	1.0
cis-1,3-Dichloropropene	0.31	U	0.31	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.14	U	0.14	1.0
Hexachlorobutadiene	0.44	U	0.44	1.0
Isopropylbenzene	0.40	U	0.40	1.0
4-Isopropyltoluene	0.19	U	0.19	1.0
Methylene Chloride	0.69	U	0.69	1.0
Naphthalene	4.0	U	4.0	5.0
N-Propylbenzene	2.5	U	2.5	5.0
Styrene	0.59	U	0.59	1.0
1,1,1,2-Tetrachloroethane	0.98	U	0.98	2.0
	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-236

Sdg Number: 660-20926

Lab Sample ID: 660-20926-9

Date Sampled: 01/31/2008 1108

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1262.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1811		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1811			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	103		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-237

Sdg Number: 660-20926

Lab Sample ID: 660-20926-10

Date Sampled: 01/31/2008 1100

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1263.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1833		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1833			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-237

Sdg Number: 660-20926

Lab Sample ID: 660-20926-10

Date Sampled: 01/31/2008 1100

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1263.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1833		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1833			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	115		77 - 122	
4-Bromofluorobenzene	112		74 - 126	
Dibromofluoromethane	117		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-238

Sdg Number: 660-20926

Lab Sample ID: 660-20926-11

Date Sampled: 01/31/2008 1116

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1264.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1855		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1855			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-238

Sdg Number: 660-20926

Lab Sample ID: 660-20926-11

Date Sampled: 01/31/2008 1116

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58005	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1264.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1855		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1855			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-213

Sdg Number: 660-20926

Lab Sample ID: 660-20926-12

Date Sampled: 01/31/2008 1546

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58162

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2219.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 1714

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 1714

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	UQ	0.50	1.0
Bromobenzene	0.58	UQ	0.58	1.0
Chlorobromomethane	0.58	UQ	0.58	1.0
Dichlorobromomethane	0.35	UQ	0.35	1.0
Bromoform	0.58	UQ	0.58	1.0
Bromomethane	2.5	UQ	2.5	5.0
n-Butylbenzene	0.67	UQ	0.67	1.0
sec-Butylbenzene	0.63	UQ	0.63	1.0
tert-Butylbenzene	0.84	UQ	0.84	1.0
Carbon tetrachloride	0.42	UQ	0.42	1.0
Chlorobenzene	0.63	UQ	0.63	1.0
Chloroethane	2.5	UQ	2.5	5.0
Chloroform	0.90	UQ	0.90	1.0
Chloromethane	1.0	UQ	1.0	4.0
2-Chlorotoluene	0.65	UQ	0.65	1.0
4-Chlorotoluene	0.52	UQ	0.52	1.0
Chlorodibromomethane	0.34	UQ	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	UQ	2.5	5.0
Ethylene Dibromide	0.50	UQ	0.50	1.0
Dibromomethane	0.41	UQ	0.41	1.0
1,2-Dichlorobenzene	0.44	UQ	0.44	1.0
1,3-Dichlorobenzene	0.64	UQ	0.64	1.0
1,4-Dichlorobenzene	0.52	UQ	0.52	1.0
Dichlorodifluoromethane	2.5	UQ	2.5	5.0
1,1-Dichloroethane	0.52	UQ	0.52	1.0
1,2-Dichloroethane	0.57	UQ	0.57	1.0
1,1-Dichloroethene	0.45	UQ	0.45	1.0
cis-1,2-Dichloroethene	0.65	UQ	0.65	1.0
trans-1,2-Dichloroethene	0.44	UQ	0.44	1.0
1,2-Dichloropropane	0.52	UQ	0.52	1.0
1,3-Dichloropropane	0.39	UQ	0.39	1.0
2,2-Dichloropropane	0.36	UQ	0.36	1.0
1,1-Dichloropropene	0.31	UQ	0.31	1.0
cis-1,3-Dichloropropene	0.14	UQ	0.14	1.0
trans-1,3-Dichloropropene	0.14	UQ	0.14	1.0
Ethylbenzene	0.44	UQ	0.44	1.0
Hexachlorobutadiene	0.40	UQ	0.40	1.0
Isopropylbenzene	0.19	UQ	0.19	1.0
4-Isopropyltoluene	0.69	UQ	0.69	1.0
Methylene Chloride	4.0	UQ	4.0	5.0
Naphthalene	2.5	UQ	2.5	5.0
N-Propylbenzene	0.59	UQ	0.59	1.0
Styrene	0.98	UQ	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	UQ	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-213

Sdg Number: 660-20926

Lab Sample ID: 660-20926-12

Date Sampled: 01/31/2008 1546

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58162

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2219.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 1714

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 1714

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	UQ	0.15	1.0
Tetrachloroethene	0.50	UQ	0.50	1.0
Toluene	0.51	UQ	0.51	1.0
1,2,3-Trichlorobenzene	0.77	UQ	0.77	1.0
1,2,4-Trichlorobenzene	0.58	UQ	0.58	1.0
1,1,1-Trichloroethane	0.46	UQ	0.46	1.0
1,1,2-Trichloroethane	0.47	UQ	0.47	1.0
Trichloroethene	0.50	UQ	0.50	1.0
Trichlorofluoromethane	2.5	UQ	2.5	5.0
1,2,3-Trichloropropane	0.18	UQ	0.18	1.0
1,2,4-Trimethylbenzene	0.86	UQ	0.86	1.0
1,3,5-Trimethylbenzene	0.54	UQ	0.54	1.0
Vinyl chloride	0.50	UQ	0.50	1.0
o-Xylene	0.50	UQ	0.50	1.0
Acetone	9.9	UQ	9.9	20
2-Butanone (MEK)	8.4	UQ	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	UQ	3.8	10
Carbon disulfide	0.85	UQ	0.85	1.0
2-Hexanone	4.4	UQ	4.4	10
Methyl tert-butyl ether	0.44	UQ	0.44	1.0
m-Xylene & p-Xylene	0.60	UQ	0.60	2.0
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Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	103		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	87		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-62

Sdg Number: 660-20926

Lab Sample ID: 660-20926-13

Date Sampled: 01/31/2008 1025

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1326.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1952		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1952			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-62

Sdg Number: 660-20926

Lab Sample ID: 660-20926-13

Date Sampled: 01/31/2008 1025

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1326.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1952		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1952			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	74		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-92

Sdg Number: 660-20926

Lab Sample ID: 660-20926-14

Date Sampled: 01/31/2008 0920

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1327.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2015		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2015			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-92

Sdg Number: 660-20926

Lab Sample ID: 660-20926-14

Date Sampled: 01/31/2008 0920

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1327.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 2015		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 2015		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	96		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-46

Sdg Number: 660-20926

Lab Sample ID: 660-20926-15

Date Sampled: 01/31/2008 1010

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57631

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1328.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2038

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2038

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-46

Sdg Number: 660-20926

Lab Sample ID: 660-20926-15

Date Sampled: 01/31/2008 1010

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1328.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2038		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2038			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-63

Sdg Number: 660-20926

Lab Sample ID: 660-20926-16

Date Sampled: 01/31/2008 0955

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57631

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1329.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2101

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2101

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-63

Sdg Number: 660-20926

Lab Sample ID: 660-20926-16

Date Sampled: 01/31/2008 0955

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1329.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2101		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2101			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-131

Sdg Number: 660-20926

Lab Sample ID: 660-20926-17

Date Sampled: 01/31/2008 0920

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57631

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1330.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2124

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2124

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-131

Sdg Number: 660-20926

Lab Sample ID: 660-20926-17

Date Sampled: 01/31/2008 0920

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1330.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2124			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2124				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	3.2	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	113		77 - 122	
4-Bromofluorobenzene	100		74 - 126	
Dibromofluoromethane	106		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-1

Sdg Number: 660-20926

Lab Sample ID: 660-20926-18

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1331.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2147		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2147			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	3.3		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	48		0.45	1.0
cis-1,2-Dichloroethene	120		0.65	1.0
trans-1,2-Dichloroethene	0.97	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-1

Sdg Number: 660-20926

Lab Sample ID: 660-20926-18

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57631

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1331.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2147

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2147

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	310	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	102		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-1

Sdg Number: 660-20926

Lab Sample ID: 660-20926-18

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58013

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1450.D

Dilution: 10

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 0437

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 0437

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	280	Q D	5.0	10

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-2

Sdg Number: 660-20926

Lab Sample ID: 660-20926-19

Date Sampled: 01/31/2008 1215

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57631

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1332.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2210

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2210

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	18		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	77		0.45	1.0
cis-1,2-Dichloroethene	680	ES +	0.65	1.0
trans-1,2-Dichloroethene	2.3		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: PZ-LSAS-2

Lab Sample ID: 660-20926-19

Date Sampled: 01/31/2008 1215

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1332.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/13/2008 2210		Final Weight/Volume: 5 mL
Date Prepared:	02/13/2008 2210		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.7		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1000	EST	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.2		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	120	77 - 122
4-Bromofluorobenzene	106	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-2

Sdg Number: 660-20926

Lab Sample ID: 660-20926-19

Date Sampled: 01/31/2008 1215

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57963	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1829.D
Dilution:	20			Initial Weight/Volume:	5 mL
Date Analyzed:	02/18/2008 2151	Run Type:	DL	Final Weight/Volume:	5 mL
Date Prepared:	02/18/2008 2151				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	540	Q D	13	20
Trichloroethene	880	Q D	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: PZ-LSAS-3

Lab Sample ID: 660-20926-20

Date Sampled: 01/31/2008 1320

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57631

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1333.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2233

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2233

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	11		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	37		0.45	1.0
cis-1,2-Dichloroethene	56		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-3

Sdg Number: 660-20926

Lab Sample ID: 660-20926-20

Date Sampled: 01/31/2008 1320

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1333.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2233		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2233			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	13		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2500	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.98	I	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	740		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	103		74 - 126	
Dibromofluoromethane	94		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-3

Sdg Number: 660-20926

Lab Sample ID: 660-20926-20

Date Sampled: 01/31/2008 1320

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57963

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1830.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/18/2008 2213

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/18/2008 2213

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	1900	Q D	25	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-4

Sdg Number: 660-20926

Lab Sample ID: 660-20926-21

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1334.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2256			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2256				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	8.8	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	100	U	0.45	1.0
cis-1,2-Dichloroethene	730	E) +	0.65	1.0
trans-1,2-Dichloroethene	1.7	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: PZ-LSAS-4

Lab Sample ID: 660-20926-21

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57631	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1334.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2256			Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2256				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	4100	EST	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.7		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	113	77 - 122
4-Bromofluorobenzene	107	74 - 126
Dibromofluoromethane	104	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: PZ-LSAS-4

Lab Sample ID: 660-20926-21

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57963

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1831.D

Dilution: 100

Initial Weight/Volume: 5 mL

Date Analyzed: 02/18/2008 2235

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/18/2008 2235

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	560	Q D	65	100
Trichloroethene	3800	Q D	50	100

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-5

Sdg Number: 660-20926

Lab Sample ID: 660-20926-22

Date Sampled: 01/31/2008 1325

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1421.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1856		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1856			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	18		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	110		0.45	1.0
cis-1,2-Dichloroethene	690	ES -	0.65	1.0
trans-1,2-Dichloroethene	2.1		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-5

Sdg Number: 660-20926

Lab Sample ID: 660-20926-22

Date Sampled: 01/31/2008 1325

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1421.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1856			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1856				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2100 3400	U DJ	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	4.6		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	114	77 - 122
4-Bromofluorobenzene	96	74 - 126
Dibromofluoromethane	105	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-5

Sdg Number: 660-20926

Lab Sample ID: 660-20926-22

Date Sampled: 01/31/2008 1325

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2123.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2155

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	290	Q-D	33	50
Trichloroethene	3400	Q-D	25	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-6

Sdg Number: 660-20926

Lab Sample ID: 660-20926-23

Date Sampled: 01/31/2008 1425

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1422.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1919		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1919			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	33		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	150		0.45	1.0
cis-1,2-Dichloroethene	510	ES +	0.65	1.0
trans-1,2-Dichloroethene	1.6		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-6

Sdg Number: 660-20926

Lab Sample ID: 660-20926-23

Date Sampled: 01/31/2008 1425

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57981

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1422.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1919

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1919

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	3.1		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	2500 8100	U DJ	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	4.8		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	109		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	99		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-6

Sdg Number: 660-20926

Lab Sample ID: 660-20926-23

Date Sampled: 01/31/2008 1425

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58229

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2532.D

Dilution: 100

Initial Weight/Volume: 5 mL

Date Analyzed: 02/25/2008 2357

Final Weight/Volume: 5 mL

Date Prepared: 02/25/2008 2357

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	120	Q D	65	100
Trichloroethene	8900	Q D	50	100

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-7

Sdg Number: 660-20926

Lab Sample ID: 660-20926-24

Date Sampled: 01/31/2008 1250

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1423.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1942			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1942				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	49	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	65	U	0.45	1.0
cis-1,2-Dichloroethene	480	U	0.65	1.0
trans-1,2-Dichloroethene	1.8	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-7

Sdg Number: 660-20926

Lab Sample ID: 660-20926-24

Date Sampled: 01/31/2008 1250

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1423.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1942		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1942			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	6.3		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.97	I	0.47	1.0
Trichloroethene	1900 ^{***} 2500	U DJ	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.9		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone-	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	116		77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-7

Sdg Number: 660-20926

Lab Sample ID: 660-20926-24

Date Sampled: 01/31/2008 1250

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2125.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2242

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2242

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	70	Q-D	33	50
Trichloroethene	2300	Q-D	25	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-30

Sdg Number: 660-20926

Lab Sample ID: 660-20926-25

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1424.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2005			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2005				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-30

Sdg Number: 660-20926

Lab Sample ID: 660-20926-25

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1424.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2005			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2005				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	90		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-30

Sdg Number: 660-20926

Lab Sample ID: 660-20926-26

Date Sampled: 01/31/2008 1554

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1425.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2028		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2028			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: EB-30

Lab Sample ID: 660-20926-26

Date Sampled: 01/31/2008 1554

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1425.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2028		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2028			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	114	77 - 122
4-Bromofluorobenzene	96	74 - 126
Dibromofluoromethane	100	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-30

Sdg Number: 660-20926

Lab Sample ID: 660-20926-27

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1426.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2051		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2051			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: TB-30

Lab Sample ID: 660-20926-27

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1426.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2051			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2051				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	90		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-27

Sdg Number: 660-20926

Lab Sample ID: 660-20926-28

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1427.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2114		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2114			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: Dup-27

Lab Sample ID: 660-20926-28

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57981

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1427.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2114

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2114

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	109	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	102	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-27

Sdg Number: 660-20926

Lab Sample ID: 660-20926-29

Date Sampled: 01/31/2008 1550

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID: BVMH GC/MS
Preparation:	5030B		Lab File ID: 1HB1428.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/14/2008 2137		Final Weight/Volume: 5 mL
Date Prepared:	02/14/2008 2137		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-27

Sdg Number: 660-20926

Lab Sample ID: 660-20926-29

Date Sampled: 01/31/2008 1550

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1428.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2137		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2137			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	107	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	100	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-40

Sdg Number: 660-20926

Lab Sample ID: 660-20926-30

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1429.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2200			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2200				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	30		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	25		0.45	1.0
cis-1,2-Dichloroethene	8.2		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-40

Sdg Number: 660-20926

Lab Sample ID: 660-20926-30

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1429.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2200			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2200				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	160	EJ U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	970	EJ U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.4		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	116	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	106	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-40

Sdg Number: 660-20926

Lab Sample ID: 660-20926-30

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58090	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB2126.D
Dilution:	20			Initial Weight/Volume:	5 mL
Date Analyzed:	02/21/2008 2305	Run Type:	DL	Final Weight/Volume:	5 mL
Date Prepared:	02/21/2008 2305				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Tetrachloroethene	150	Q D	10	20
Trichloroethene	880	Q D	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-41

Sdg Number: 660-20926

Lab Sample ID: 660-20926-31

Date Sampled: 01/31/2008 1111

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1430.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2223		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2223			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	22		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	68		0.45	1.0
cis-1,2-Dichloroethene	96		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-41

Sdg Number: 660-20926

Lab Sample ID: 660-20926-31

Date Sampled: 01/31/2008 1111

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1430.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2223		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2223			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.59	I	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	250	ES L	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.68	I	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	118	77 - 122
4-Bromofluorobenzene	95	74 - 126
Dibromofluoromethane	103	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-41

Sdg Number: 660-20926

Lab Sample ID: 660-20926-31

Date Sampled: 01/31/2008 1111

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2127.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2328

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2328

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	190	Q D	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: MW-33

Lab Sample ID: 660-20926-32

Date Sampled: 01/31/2008 1020

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1431.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2246			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2246				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	160	U EJ	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethane	570 580	U DJ	0.45	1.0
cis-1,2-Dichloroethene	22	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-33

Sdg Number: 660-20926

Lab Sample ID: 660-20926-32

Date Sampled: 01/31/2008 1020

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1431.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2246		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2246			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	140	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	410	U ES	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	4.4	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	125	J1	77 - 122	
4-Bromofluorobenzene	95		74 - 126	
Dibromofluoromethane	107		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-33

Sdg Number: 660-20926

Lab Sample ID: 660-20926-32

Date Sampled: 01/31/2008 1020

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2128.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2351

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2351

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethane	140	Q-D	10	20
1,1-Dichloroethene	580	Q-D	9.0	20
Trichloroethene	290	Q-D	10	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-102

Sdg Number: 660-20926

Lab Sample ID: 660-20926-33

Date Sampled: 01/31/2008 1652

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1432.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2309		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2309			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	6.3		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	55		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-102

Sdg Number: 660-20926

Lab Sample ID: 660-20926-33

Date Sampled: 01/31/2008 1652

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1432.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2309		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2309			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1.8	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	73		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	111		77 - 122	
4-Bromofluorobenzene	90		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-107

Sdg Number: 660-20926

Lab Sample ID: 660-20926-34

Date Sampled: 01/31/2008 1600

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B			Lab File ID:	1HB1433.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2332			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2332				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-107

Sdg Number: 660-20926

Lab Sample ID: 660-20926-34

Date Sampled: 01/31/2008 1600

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57981

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1433.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2332

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2332

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.52	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	6.1		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	6.4		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	111	77 - 122
4-Bromofluorobenzene	92	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-108

Sdg Number: 660-20926

Lab Sample ID: 660-20926-35

Date Sampled: 01/31/2008 1447

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1434.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2355		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2355			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	46	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	140	U	0.45	1.0
cis-1,2-Dichloroethene	930	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-108

Sdg Number: 660-20926

Lab Sample ID: 660-20926-35

Date Sampled: 01/31/2008 1447

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57981	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1434.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2355		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2355			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.90	I	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	78		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	2.5		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	57	U	9.9	20
2-Butanone (MEK)	30		8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	100	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	84		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: EW-108

Lab Sample ID: 660-20926-35

Date Sampled: 01/31/2008 1447

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2129.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 0014

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 0014

Analyte	Result (ug/L)	Qualifier	MDL	PQL
cis-1,2-Dichloroethene	460	Q D	13	20
Carbon disulfide	85	Q D	17	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-10

Sdg Number: 660-20926

Lab Sample ID: 660-20926-36

Date Sampled: 01/31/2008 1030

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1419.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1702		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1702			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	2.7		0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	5.3		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	1.0		0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-10

Sdg Number: 660-20926

Lab Sample ID: 660-20926-36

Date Sampled: 01/31/2008 1030

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1419.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1702			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1702				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.5		0.50	1.0
Toluene	1.5		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1.8	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	95		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	90		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-36

Sdg Number: 660-20926

Lab Sample ID: 660-20926-37

Date Sampled: 01/31/2008 1120

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1420.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1727			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1727				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	3.1		0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	20		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	54		0.45	1.0
cis-1,2-Dichloroethene	5.8		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-36

Sdg Number: 660-20926

Lab Sample ID: 660-20926-37

Date Sampled: 01/31/2008 1120

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1420.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1727			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1727				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	1.3		0.50	1.0
Toluene	1.6		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	190	ES	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	114	74 - 126
Dibromofluoromethane	101	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-36

Sdg Number: 660-20926

Lab Sample ID: 660-20926-37

Date Sampled: 01/31/2008 1120

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58090

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2130.D

Dilution: 5.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/22/2008 0037

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/22/2008 0037

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	130	Q D	2.5	5.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-37

Sdg Number: 660-20926

Lab Sample ID: 660-20926-38

Date Sampled: 01/31/2008 1038

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1421.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1749			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1749				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	49	U	0.52	1.0
1,2-Dichloroethane	0:57	U	0.57	1.0
1,1-Dichloroethene	230	U	0.45	1.0
cis-1,2-Dichloroethene	420	U	0.65	1.0
trans-1,2-Dichloroethene	1.8	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-37

Sdg Number: 660-20926

Lab Sample ID: 660-20926-38

Date Sampled: 01/31/2008 1038

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1421.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1749			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1749				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	24		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	4500 - 7400	U - DJ	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	12		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	115		74 - 126	
Dibromofluoromethane	102		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-37

Sdg Number: 660-20926

Lab Sample ID: 660-20926-38

Date Sampled: 01/31/2008 1038

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57968

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1923.D

Dilution: 100

Initial Weight/Volume: 5 mL

Date Analyzed: 02/19/2008 1809

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/19/2008 1809

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	230	Q-D	45	100
cis-1,2-Dichloroethene	400	Q-D	65	100
Trichloroethene	9400	Q-D	50	100

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-127

Sdg Number: 660-20926

Lab Sample ID: 660-20926-39

Date Sampled: 01/31/2008 1221

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1422.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1811			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1811				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	34	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	240 270	U	0.45	1.0
cis-1,2-Dichloroethene	100	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-127

Sdg Number: 660-20926

Lab Sample ID: 660-20926-39

Date Sampled: 01/31/2008 1221

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1422.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1811			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1811				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	3.3		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	3200 3300	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	2.4		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	101		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	95		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-127

Sdg Number: 660-20926

Lab Sample ID: 660-20926-39

Date Sampled: 01/31/2008 1221

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-57968

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1924.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/19/2008 1831

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/19/2008 1831

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	270	Q-D	23	50
Trichloroethene	3300	Q-D	25	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-128

Sdg Number: 660-20926

Lab Sample ID: 660-20926-40

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1423.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1834			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1834				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	2.3		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	3.6		0.45	1.0
cis-1,2-Dichloroethene	2.5		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-128

Sdg Number: 660-20926

Lab Sample ID: 660-20926-40

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1423.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1834			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1834				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.5		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	34		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	11.20	+ U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.99	I	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	108		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-41

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1424.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1856

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1856

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	2.6		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	4.0		0.45	1.0
cis-1,2-Dichloroethene	2.5		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-41

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1424.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1856

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1856

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.5		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	35		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	11.20	+ U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	1.3		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	112		77 - 122	
4-Bromofluorobenzene	116		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-42

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1425.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1918			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1918				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-42

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1425.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1918		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1918			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.73	I	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	31		9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	91		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-43

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1426.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1941		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1941			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-43

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1426.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1941		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1941			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	113		77 - 122	
4-Bromofluorobenzene	117		74 - 126	
Dibromofluoromethane	103		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-44

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1427.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2003

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2003

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-44

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1427.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2003		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2003			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	4.2		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	5.1		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	105		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-45

Date Sampled: 01/31/2008 1655

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1428.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2026

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2026

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-45

Date Sampled: 01/31/2008 1655

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1428.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2026

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2026

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	101	74 - 126
Dibromofluoromethane	95	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-46

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1429.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2048		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2048			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-46

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1429.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2048		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2048			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	110	77 - 122
4-Bromofluorobenzene	110	74 - 126
Dibromofluoromethane	98	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Private Well 85

Sdg Number: 660-20926

Lab Sample ID: 660-20926-47

Date Sampled: 01/31/2008 1505

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1430.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2110			Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2110				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.55	I	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Private Well 85

Sdg Number: 660-20926

Lab Sample ID: 660-20926-47

Date Sampled: 01/31/2008 1505

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1430.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2110		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2110			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	1.3		0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	110		74 - 126	
Dibromofluoromethane	102		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-19

Sdg Number: 660-20926

Lab Sample ID: 660-20926-48

Date Sampled: 01/31/2008 1643

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1431.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2133

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2133

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: MW-19

Lab Sample ID: 660-20926-48

Date Sampled: 01/31/2008 1643

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58025

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1431.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2133

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2133

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	112	77 - 122
4-Bromofluorobenzene	114	74 - 126
Dibromofluoromethane	108	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-252

Sdg Number: 660-20926

Lab Sample ID: 660-20926-49

Date Sampled: 01/31/2008 1617

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1432.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/14/2008 2155		Final Weight/Volume: 5 mL
Date Prepared:	02/14/2008 2155		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-252

Sdg Number: 660-20926

Lab Sample ID: 660-20926-49

Date Sampled: 01/31/2008 1617

Client Matrix: Water

Date Received: 02/01/2008 1000

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58025	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1432.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2155		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2155			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	104		77 - 122	
4-Bromofluorobenzene	104		74 - 126	
Dibromofluoromethane	98		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-249

Sdg Number: 660-20926

Lab Sample ID: 660-20926-1

Date Sampled: 01/31/2008 1428

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58006	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1254.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1512		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1512			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-250

Sdg Number: 660-20926

Lab Sample ID: 660-20926-2

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58006	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1255.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1534		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1534			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	15		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-214

Sdg Number: 660-20926

Lab Sample ID: 660-20926-3

Date Sampled: 01/31/2008 1535

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58006	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1256.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1557		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1557			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-215

Sdg Number: 660-20926

Lab Sample ID: 660-20926-4

Date Sampled: 01/31/2008 1545

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1257.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1619

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1619

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-216

Sdg Number: 660-20926

Lab Sample ID: 660-20926-5

Date Sampled: 01/31/2008 1525

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58006	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1258.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1641		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1641			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-217

Sdg Number: 660-20926

Lab Sample ID: 660-20926-6

Date Sampled: 01/31/2008 1625

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1259.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1704

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1704

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-234

Sdg Number: 660-20926

Lab Sample ID: 660-20926-7

Date Sampled: 01/31/2008 1100

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58006	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1260.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1726		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1726			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-235

Sdg Number: 660-20926

Lab Sample ID: 660-20926-8

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58006	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1261.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1748		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1748			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-236

Sdg Number: 660-20926

Lab Sample ID: 660-20926-9

Date Sampled: 01/31/2008 1108

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1262.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1811

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1811

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-237

Sdg Number: 660-20926

Lab Sample ID: 660-20926-10

Date Sampled: 01/31/2008 1100

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1263.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1833

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1833

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-238

Sdg Number: 660-20926

Lab Sample ID: 660-20926-11

Date Sampled: 01/31/2008 1116

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58006

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1264.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 1855

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 1855

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-213

Sdg Number: 660-20926

Lab Sample ID: 660-20926-12

Date Sampled: 01/31/2008 1546

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID Analysis Batch: 660-58005 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1265.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/13/2008 1918 Final Weight/Volume: 5 mL
 Date Prepared: 02/13/2008 1918

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Method: 8260C SIM/ID Analysis Batch: 660-58129 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB2219.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/22/2008 1714 Final Weight/Volume: 5 mL
 Date Prepared: 02/22/2008 1714

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U J	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-62

Sdg Number: 660-20926

Lab Sample ID: 660-20926-13

Date Sampled: 01/31/2008 1025

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57632	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1326.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 1952		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 1952			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-92

Sdg Number: 660-20926

Lab Sample ID: 660-20926-14

Date Sampled: 01/31/2008 0920

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57632	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1327.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2015		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2015			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: MW-46

Lab Sample ID: 660-20926-15

Date Sampled: 01/31/2008 1010

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57632

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1328.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2038

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2038

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-63

Sdg Number: 660-20926

Lab Sample ID: 660-20926-16

Date Sampled: 01/31/2008 0955

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57632

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1329.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2101

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2101

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	6.7		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-131

Sdg Number: 660-20926

Lab Sample ID: 660-20926-17

Date Sampled: 01/31/2008 0920

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57632	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1330.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/13/2008 2124		Final Weight/Volume:	5 mL
Date Prepared:	02/13/2008 2124			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	2.9		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-1

Sdg Number: 660-20926

Lab Sample ID: 660-20926-18

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57632

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1331.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/13/2008 2147

Final Weight/Volume: 5 mL

Date Prepared: 02/13/2008 2147

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	660	ES ±	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-1

Sdg Number: 660-20926

Lab Sample ID: 660-20926-18

Date Sampled: 01/31/2008 1410

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58015

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1450.D

Dilution: 10

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 0437

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 0437

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	470	Q D	10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-2

Sdg Number: 660-20926

Lab Sample ID: 660-20926-19

Date Sampled: 01/31/2008 1215

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/13/2008 2210
 Date Prepared: 02/13/2008 2210

Analysis Batch: 660-57632

Instrument ID: BVMH GC/MS

Lab File ID: 1HB1332.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	570	L	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: PZ-LSAS-2

Lab Sample ID: 660-20926-19

Date Sampled: 01/31/2008 1215

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57964	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1829.D
Dilution:	20		Initial Weight/Volume: 5 mL
Date Analyzed:	02/18/2008 2151	Run Type: DL	Final Weight/Volume: 5 mL
Date Prepared:	02/18/2008 2151		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	580	Q D3	20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-3

Sdg Number: 660-20926

Lab Sample ID: 660-20926-20

Date Sampled: 01/31/2008 1320

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1433.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2217

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2217

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	100		50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-4

Sdg Number: 660-20926

Lab Sample ID: 660-20926-21

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1434.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2240

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2240

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	580		50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-4

Sdg Number: 660-20926

Lab Sample ID: 660-20926-21

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57964	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1831.D
Dilution:	100		Initial Weight/Volume:	5 mL
Date Analyzed:	02/18/2008 2235	Run Type: DL	Final Weight/Volume:	5 mL
Date Prepared:	02/18/2008 2235			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	600	Q-DJ	100	100

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Sdg Number: 660-20926

Client Sample ID: PZ-LSAS-5

Lab Sample ID: 660-20926-22

Date Sampled: 01/31/2008 1325

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58089

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2123.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2155

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	380	Q J	50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-6

Sdg Number: 660-20926

Lab Sample ID: 660-20926-23

Date Sampled: 01/31/2008 1425

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58089

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2124.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2218

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2218

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	350	Q S	50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: PZ-LSAS-7

Sdg Number: 660-20926

Lab Sample ID: 660-20926-24

Date Sampled: 01/31/2008 1250

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58089	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB2125.D
Dilution:	50		Initial Weight/Volume:	5 mL
Date Analyzed:	02/21/2008 2242		Final Weight/Volume:	5 mL
Date Prepared:	02/21/2008 2242			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	50	U-Q	50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-30

Sdg Number: 660-20926

Lab Sample ID: 660-20926-25

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57964	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1832.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/18/2008 2257		Final Weight/Volume:	5 mL
Date Prepared:	02/18/2008 2257			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U E J	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-27

Sdg Number: 660-20926

Lab Sample ID: 660-20926-28

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB2036.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 0317

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 0317

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U-G	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-27

Sdg Number: 660-20926

Lab Sample ID: 660-20926-29

Date Sampled: 01/31/2008 1550

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB2037.D

Dilution: 1.0


Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 0339

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 0339

Analyte

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U 	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-40

Sdg Number: 660-20926

Lab Sample ID: 660-20926-30

Date Sampled: 01/31/2008 1150

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58089

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2126.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2305

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2305

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	20	UQ	20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-41

Sdg Number: 660-20926

Lab Sample ID: 660-20926-31

Date Sampled: 01/31/2008 1111

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58089	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB2127.D
Dilution:	20		Initial Weight/Volume:	5 mL
Date Analyzed:	02/21/2008 2328		Final Weight/Volume:	5 mL
Date Prepared:	02/21/2008 2328			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	450	Q S	20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-33

Sdg Number: 660-20926

Lab Sample ID: 660-20926-32

Date Sampled: 01/31/2008 1020

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58089

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB2128.D

Dilution: 20

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 2351

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 2351

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	590	Q-S	20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-102

Sdg Number: 660-20926

Lab Sample ID: 660-20926-33

Date Sampled: 01/31/2008 1652

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58063	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB2039.D
Dilution:	2.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/21/2008 0424		Final Weight/Volume:	5 mL
Date Prepared:	02/21/2008 0424			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	120	DJ-Q	2.0	2.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-107

Sdg Number: 660-20926

Lab Sample ID: 660-20926-34

Date Sampled: 01/31/2008 1600

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58063

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB2038.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/21/2008 0401

Final Weight/Volume: 5 mL

Date Prepared: 02/21/2008 0401

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U Q	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EW-108

Sdg Number: 660-20926

Lab Sample ID: 660-20926-35

Date Sampled: 01/31/2008 1447

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 20
 Date Analyzed: 02/22/2008 0014
 Date Prepared: 02/22/2008 0014

Analysis Batch: 660-58089

Instrument ID: BVMH GC/MS

Lab File ID: 1HB2129.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	540	Q S	20	20

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-10

Sdg Number: 660-20926

Lab Sample ID: 660-20926-36

Date Sampled: 01/31/2008 1030

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/14/2008 1702
 Date Prepared: 02/14/2008 1702

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1419.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-36

Sdg Number: 660-20926

Lab Sample ID: 660-20926-37

Date Sampled: 01/31/2008 1120

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58019	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1420.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1727		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1727			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	15		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-37

Sdg Number: 660-20926

Lab Sample ID: 660-20926-38

Date Sampled: 01/31/2008 1038

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1421.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1749

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1749

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	300	L	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-37

Sdg Number: 660-20926

Lab Sample ID: 660-20926-38

Date Sampled: 01/31/2008 1038

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57967	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1923.D
Dilution:	100		Initial Weight/Volume:	5 mL
Date Analyzed:	02/19/2008 1809		Final Weight/Volume:	5 mL
Date Prepared:	02/19/2008 1809			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	760	Q DS	100	100

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-127

Sdg Number: 660-20926

Lab Sample ID: 660-20926-39

Date Sampled: 01/31/2008 1221

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1422.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1811

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1811

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	200	L	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-127

Sdg Number: 660-20926

Lab Sample ID: 660-20926-39

Date Sampled: 01/31/2008 1221

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57967

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1924.D

Dilution: 50

Initial Weight/Volume: 5 mL

Date Analyzed: 02/19/2008 1831

Final Weight/Volume: 5 mL

Date Prepared: 02/19/2008 1831

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	300	Q-DJ	50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-128

Sdg Number: 660-20926

Lab Sample ID: 660-20926-40

Date Sampled: 01/31/2008 1220

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58019	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1423.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1834		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1834			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.2		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-41

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1424.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1856

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1856

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	7.6		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-42

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1425.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 1918

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 1918

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-28

Sdg Number: 660-20926

Lab Sample ID: 660-20926-43

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58019	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1426.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 1941		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 1941			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Dup-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-44

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1427.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2003

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2003

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: EB-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-45

Date Sampled: 01/31/2008 1655

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58019	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1428.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2026		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2026			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: TB-29

Sdg Number: 660-20926

Lab Sample ID: 660-20926-46

Date Sampled: 01/31/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58019	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1429.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/14/2008 2048		Final Weight/Volume:	5 mL
Date Prepared:	02/14/2008 2048			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: Private Well 85

Sdg Number: 660-20926

Lab Sample ID: 660-20926-47

Date Sampled: 01/31/2008 1505

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1430.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2110

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2110

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	22		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-19

Sdg Number: 660-20926

Lab Sample ID: 660-20926-48

Date Sampled: 01/31/2008 1643

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1431.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2133

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2133

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20926-1

Client Sample ID: MW-252

Sdg Number: 660-20926

Lab Sample ID: 660-20926-49

Date Sampled: 01/31/2008 1617

Client Matrix: Water

Date Received: 02/01/2008 1000

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58019

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1432.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/14/2008 2155

Final Weight/Volume: 5 mL

Date Prepared: 02/14/2008 2155

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20926	1/31/2008	SW-846	MW-249	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	MW-237	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	MW-238	Water	No	--	--	--	--	VOC – CCAL %D, Associated Blank
660-20926	1/31/2008	SW-846	MW-213	Water	No	--	--	--	--	VOC – CCAL %D, Holding Time
660-20926	1/31/2008	SW-846	MW-62	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-92	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-46	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-63	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-131	Water	No	--	--	--	--	VOC – Associated Blank
660-20926	1/31/2008	SW-846	PZ-LSAS-1	Water	No	--	--	--	--	VOC – Calibration Range
660-20926	1/31/2008	SW-846	PZ-LSAS-2	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	MW-250	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	PZ-LSAS-3	Water	No	--	--	--	--	VOC – Calibration Range
660-20926	1/31/2008	SW-846	PZ-LSAS-4	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	PZ-LSAS-5	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	PZ-LSAS-6	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	PZ-LSAS-7	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	Dup-30	Water	No	--	--	--	--	VOC – Holding Time
660-20926	1/31/2008	SW-846	EB-30	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	TB-30	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	Dup-27	Water	No	--	--	--	--	VOC – Holding Time
660-20926	1/31/2008	SW-846	EB-27	Water	No	--	--	--	--	VOC – Holding Time
660-20926	1/31/2008	SW-846	MW-214	Water	No	--	--	--	--	VOC – CCAL %D

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20926	1/31/2008	SW-846	MW-40	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	MW-41	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	MW-33	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time, Surrogate %Recovery
660-20926	1/31/2008	SW-846	EW-102	Water	No	--	--	--	--	VOC – Associated Blank, Holding Time
660-20926	1/31/2008	SW-846	EW-107	Water	No	--	--	--	--	VOC – Holding Time
660-20926	1/31/2008	SW-846	EW-108	Water	No	--	--	--	--	VOC – Associated Blank, Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	MW-10	Water	No	--	--	--	--	VOC – Associated Blank
660-20926	1/31/2008	SW-846	MW-36	Water	No	--	--	--	--	VOC – Calibration Range
660-20926	1/31/2008	SW-846	MW-37	Water	No	--	--	--	--	VOC – Calibration Range, Holding Time
660-20926	1/31/2008	SW-846	MW-127	Water	No	--	--	--	--	VOC – Holding Time
660-20926	1/31/2008	SW-846	MW-215	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	MW-128	Water	No	--	--	--	--	VOC – Associated Blank
660-20926	1/31/2008	SW-846	Dup-28	Water	No	--	--	--	--	VOC – Associated Blank
660-20926	1/31/2008	SW-846	EB-28	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	TB-28	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	Dup-29	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	EB-29	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	TB-29	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	Private Well 85	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-19	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-252	Water	Yes	--	--	--	--	
660-20926	1/31/2008	SW-846	MW-216	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	MW-217	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	MW-234	Water	No	--	--	--	--	VOC – CCAL %D

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20926	1/31/2008	SW-846	MW-235	Water	No	--	--	--	--	VOC – CCAL %D
660-20926	1/31/2008	SW-846	MW-236	Water	No	--	--	--	--	VOC – CCAL %D

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 8

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-20941

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8130R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-20941 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
MW-202	660-20941-1	Water	2/1/2008	X				
MW-39	660-20941-10	Water	2/1/2008	X				
MW-253	660-20941-11	Water	2/1/2008	X				
MW-254	660-20941-12	Water	2/1/2008	X				
MW-8D	660-20941-13	Water	2/1/2008	X				
MW-123	660-20941-14	Water	2/1/2008	X				
RW-1	660-20941-15	Water	2/1/2008	X				
Dup-31	660-20941-16	Water	2/1/2008	X				
EB-31	660-20941-17	Water	2/1/2008	X				
TB-31	660-20941-18	Water	2/1/2008	X				
Private Well 7	660-20941-19	Water	2/1/2008	X				
MW-134	660-20941-2	Water	2/1/2008	X				
Private Well 57	660-20941-20	Water	2/1/2008	X				
Private Well 59	660-20941-21	Water	2/1/2008	X				
Private Well 38	660-20941-22	Water	2/1/2008	X				
Private Well 102	660-20941-23	Water	2/1/2008	X				
TB-32	660-20941-24	Water	2/1/2008	X				
DW-1	660-20941-3	Water	2/1/2008	X				
MW-42	660-20941-4	Water	2/1/2008	X				
MW-43	660-20941-5	Water	2/1/2008	X				
IWI-1	660-20941-6	Water	2/1/2008	X				
IWI-2	660-20941-7	Water	2/1/2008	X				
EW-UAFG-1	660-20941-8	Water	2/1/2008	X				
MW-38	660-20941-9	Water	2/1/2008	X				

Notes:

1. Sample location Dup-31 is the field duplicate of parent sample location MW-43.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No compounds were detected in the associated blanks.

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
TB-32	CCV %D	1,2,3-Trichlorobenzene	-34.7%
		1,2,4-Trichlorobenzene	-30.6%
		2,2-Dichloropropane	21.6%
		4-Methyl-2-pentanone	21.2%
		Bromomethane	-50.5%
		Chloromethane	-33.8%
		Dichlorobromomethane	21.5%
		Naphthalene	-38.9%
		Vinyl chloride	-25.7%
MW-134 DW-1 MW-42 MW-43 IWI-1 EW-UAFG-1 MW-253 MW-123 RW-1 EB-31 TB-31 Private Well 7 Private Well 57	CCV %D	1,2,4-Trichlorobenzene	-20.5%
		Bromomethane	34.7%
		Naphthalene	-23.6%
MW-202 IWI-2 MW-38 MW-39 MW-254 MW-8D Dup-31 Private Well 59 Private Well 38 Private Well 102	CCV %D	2-Hexanone	32.6%
		4-Methyl-2-pentanone	29.3%
		Acetone	26.4%
		Chloroethane	32.9%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

1. RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
MW-43/Dup-31	1,1-Dichloroethane	36	31	AC
	1,1-Dichloroethene	150	140	6.8%
	cis-1,2-Dichloroethene	19	16	AC
	Trichloroethene	190	150	23.5%
	Vinyl chloride	0.64 l	ND(5.0)	AC
	1,4-Dioxane (SIM ID)	570	560	1.7%

NC = Not compliant.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The calculated RPDs between the parent sample and field duplicate were acceptable.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
MW-134	Trichloroethene	200 E	190 D	190 D
MW-42	Trichloroethene	2400 E	1800 D	1800 D
MW-43	1,1-Dichloroethene	150 E	150 D	150 D
	Trichloroethene	200 E	190 D	190 D

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
IWI-1	1,1-Dichloroethene	400 E	370 D	370 D
	cis-1,2-Dichloroethene	240 E	190 D	190 D
	Trichloroethene	5100 E	4900 D	4900 D
EW-UAFG-1	1,1-Dichloroethene	260 E	250 D	250 D
	cis-1,2-Dichloroethene	2200 E	3000 D	3000 D
MW-253	1,1-Dichloroethene	380 E	400 D	400 D
	Trichloroethene	5600 E	6200 D	6200 D
MW-254	Tetrachloroethene	15000 E	11000 D	11000 D

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-202

Sdg Number: 660-20941

Lab Sample ID: 660-20941-1

Date Sampled: 02/01/2008 0810

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1510.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1548		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1548			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-202

Sdg Number: 660-20941

Lab Sample ID: 660-20941-1

Date Sampled: 02/01/2008 0810

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1510.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1548		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1548			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	93		77 - 122	
4-Bromofluorobenzene	91		74 - 126	
Dibromofluoromethane	91		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-134

Sdg Number: 660-20941

Lab Sample ID: 660-20941-2

Date Sampled: 02/01/2008 1132

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1451.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0459			Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0459				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	5.4		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	23		0.45	1.0
cis-1,2-Dichloroethene	5.9		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-134

Sdg Number: 660-20941

Lab Sample ID: 660-20941-2

Date Sampled: 02/01/2008 1132

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1451.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0459		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0459			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	94		77 - 122	
4-Bromofluorobenzene	93		74 - 126	
Dibromofluoromethane	89		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-134

Sdg Number: 660-20941

Lab Sample ID: 660-20941-2

Date Sampled: 02/01/2008 1132

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58240

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1514.D

Dilution: 2.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1717

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1717

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	190	D	1.0	2.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: DW-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-3

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1452.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0522		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0522			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: DW-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-3

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1452.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/15/2008 0522		Final Weight/Volume: 5 mL
Date Prepared:	02/15/2008 0522		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	102	77 - 122
4-Bromofluorobenzene	100	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-42

Sdg Number: 660-20941

Lab Sample ID: 660-20941-4

Date Sampled: 02/01/2008 1056

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1453.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0544			Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0544				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	3.6		0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	59		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	89		0.45	1.0
cis-1,2-Dichloroethene	83		0.65	1.0
trans-1,2-Dichloroethene	0.59	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-42

Sdg Number: 660-20941

Lab Sample ID: 660-20941-4

Date Sampled: 02/01/2008 1056

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58013

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1453.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 0544

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 0544

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	14		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	3.6		0.47	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.2		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	100		77 - 122	
4-Bromofluorobenzene	97		74 - 126	
Dibromofluoromethane	97		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-42

Sdg Number: 660-20941

Lab Sample ID: 660-20941-4

Date Sampled: 02/01/2008 1056

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58240

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1519.D

Dilution: 40

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1910

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1910

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Trichloroethene	1800	D	20	40

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-43

Sdg Number: 660-20941

Lab Sample ID: 660-20941-5

Date Sampled: 02/01/2008 1103

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1454.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0606		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0606			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	36		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	19		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-43

Sdg Number: 660-20941

Lab Sample ID: 660-20941-5

Date Sampled: 02/01/2008 1103

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1454.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0606			Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0606				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.64	I	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	116	77 - 122
4-Bromofluorobenzene	112	74 - 126
Dibromofluoromethane	114	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-43

Sdg Number: 660-20941

Lab Sample ID: 660-20941-5

Date Sampled: 02/01/2008 1103

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58240

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1515.D

Dilution: 10

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1740

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1740

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	150	D	4.5	10
Trichloroethene	190	D	5.0	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-6

Date Sampled: 02/01/2008 0915

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1455.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0629		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0629			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	60		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
trans-1,2-Dichloroethene	0.90	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	6.8		0.50	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-6

Date Sampled: 02/01/2008 0915

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1455.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0629		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0629			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	21		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	102		77 - 122	
4-Bromofluorobenzene	99		74 - 126	
Dibromofluoromethane	99		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-6

Date Sampled: 02/01/2008 0915

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58240

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1520.D

Dilution: 100

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1932

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1932

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	370	D	45	100
cis-1,2-Dichloroethene	190	↓	65	100
Trichloroethene	4900		50	100

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-2

Sdg Number: 660-20941

Lab Sample ID: 660-20941-7

Date Sampled: 02/01/2008 1025

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1527.D
Dilution:	1.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2209			Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2209				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.64	I	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.83	I	0.45	1.0
cis-1,2-Dichloroethene	4.1		0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-2

Sdg Number: 660-20941

Lab Sample ID: 660-20941-7

Date Sampled: 02/01/2008 1025

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1527.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2209		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2209			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.99	I	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	84	77 - 122
4-Bromofluorobenzene	82	74 - 126
Dibromofluoromethane	83	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EW-UAFG-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-8

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1457.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0713		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0713			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	34		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
trans-1,2-Dichloroethene	11		0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EW-UAFG-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-8

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1457.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0713		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0713			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	19		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	3.6		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	97		77 - 122	
4-Bromofluorobenzene	96		74 - 126	
Dibromofluoromethane	101		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EW-UAFG-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-8

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1521.D
Dilution:	40		Initial Weight/Volume: 5 mL
Date Analyzed:	02/15/2008 1954	Run Type: DL	Final Weight/Volume: 5 mL
Date Prepared:	02/15/2008 1954		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	250	D	18	40
cis-1,2-Dichloroethene	3000	D	26	40

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-38

Sdg Number: 660-20941

Lab Sample ID: 660-20941-9

Date Sampled: 02/01/2008 0940

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58240 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1516.D
 Dilution: 10 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 1803 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 1803

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	5.0	U	5.0	10
Bromobenzene	5.8	U	5.8	10
Chlorobromomethane	5.8	U	5.8	10
Dichlorobromomethane	3.5	U	3.5	10
Bromoform	5.8	U	5.8	10
Bromomethane	25	U	25	50
n-Butylbenzene	6.7	U	6.7	10
sec-Butylbenzene	6.3	U	6.3	10
tert-Butylbenzene	8.4	U	8.4	10
Carbon tetrachloride	4.2	U	4.2	10
Chlorobenzene	6.3	U	6.3	10
Chloroethane	25	U	25	50
Chloroform	9.0	U	9.0	10
Chloromethane	10	U	10	40
2-Chlorotoluene	6.5	U	6.5	10
4-Chlorotoluene	5.2	U	5.2	10
Chlorodibromomethane	3.4	U	3.4	10
1,2-Dibromo-3-Chloropropane	25	U	25	50
Ethylene Dibromide	5.0	U	5.0	10
Dibromomethane	4.1	U	4.1	10
1,2-Dichlorobenzene	4.4	U	4.4	10
1,3-Dichlorobenzene	6.4	U	6.4	10
1,4-Dichlorobenzene	5.2	U	5.2	10
Dichlorodifluoromethane	25	U	25	50
1,1-Dichloroethane	20	U	5.2	10
1,2-Dichloroethane	5.7	U	5.7	10
1,1-Dichloroethene	27	U	4.5	10
cis-1,2-Dichloroethene	6.5	U	6.5	10
trans-1,2-Dichloroethene	4.4	U	4.4	10
1,2-Dichloropropane	5.2	U	5.2	10
1,3-Dichloropropane	3.9	U	3.9	10
2,2-Dichloropropane	3.6	U	3.6	10
1,1-Dichloropropene	3.1	U	3.1	10
cis-1,3-Dichloropropene	1.4	U	1.4	10
trans-1,3-Dichloropropene	1.4	U	1.4	10
Ethylbenzene	4.4	U	4.4	10
Hexachlorobutadiene	4.0	U	4.0	10
Isopropylbenzene	1.9	U	1.9	10
4-Isopropyltoluene	6.9	U	6.9	10
Methylene Chloride	40	U	40	50
Naphthalene	25	U	25	50
N-Propylbenzene	5.9	U	5.9	10
Styrene	9.8	U	9.8	20
1,1,1,2-Tetrachloroethane	6.3	U	6.3	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-38

Sdg Number: 660-20941

Lab Sample ID: 660-20941-9

Date Sampled: 02/01/2008 0940

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1516.D
Dilution:	10		Initial Weight/Volume: 5 mL
Date Analyzed:	02/15/2008 1803		Final Weight/Volume: 5 mL
Date Prepared:	02/15/2008 1803		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	1.5	U	1.5	10
Tetrachloroethene	450		5.0	10
Toluene	5.1	U	5.1	10
1,2,3-Trichlorobenzene	7.7	U	7.7	10
1,2,4-Trichlorobenzene	5.8	U	5.8	10
1,1,1-Trichloroethane	4.6	U	4.6	10
1,1,2-Trichloroethane	4.7	U	4.7	10
Trichloroethene	600		5.0	10
Trichlorofluoromethane	25	U	25	50
1,2,3-Trichloropropane	1.8	U	1.8	10
1,2,4-Trimethylbenzene	8.6	U	8.6	10
1,3,5-Trimethylbenzene	5.4	U	5.4	10
Vinyl chloride	5.0	U	5.0	10
o-Xylene	5.0	U	5.0	10
Acetone	99	U	99	200
2-Butanone (MEK)	84	U	84	100
4-Methyl-2-pentanone (MIBK)	38	U	38	100
Carbon disulfide	8.5	U	8.5	10
2-Hexanone	44	U	44	100
Methyl tert-butyl ether	4.4	U	4.4	10
m-Xylene & p-Xylene	6.0	U	6.0	20
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	107		77 - 122	
4-Bromofluorobenzene	109		74 - 126	
Dibromofluoromethane	100		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: MW-39

Lab Sample ID: 660-20941-10

Date Sampled: 02/01/2008 1016

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58240 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1517.D
 Dilution: 4.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 1825 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 1825

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	2.0	U	2.0	4.0
Bromobenzene	2.3	U	2.3	4.0
Chlorobromomethane	2.3	U	2.3	4.0
Dichlorobromomethane	1.4	U	1.4	4.0
Bromoform	2.3	U	2.3	4.0
Bromomethane	10	U	10	20
n-Butylbenzene	2.7	U	2.7	4.0
sec-Butylbenzene	2.5	U	2.5	4.0
tert-Butylbenzene	3.4	U	3.4	4.0
Carbon tetrachloride	1.7	U	1.7	4.0
Chlorobenzene	2.5	U	2.5	4.0
Chloroethane	10	U	10	20
Chloroform	3.6	U	3.6	4.0
Chloromethane	4.0	U	4.0	16
2-Chlorotoluene	2.6	U	2.6	4.0
4-Chlorotoluene	2.1	U	2.1	4.0
Chlorodibromomethane	1.4	U	1.4	4.0
1,2-Dibromo-3-Chloropropane	10	U	10	20
Ethylene Dibromide	2.0	U	2.0	4.0
Dibromomethane	1.6	U	1.6	4.0
1,2-Dichlorobenzene	1.8	U	1.8	4.0
1,3-Dichlorobenzene	2.6	U	2.6	4.0
1,4-Dichlorobenzene	2.1	U	2.1	4.0
Dichlorodifluoromethane	10	U	10	20
1,1-Dichloroethane	5.7		2.1	4.0
1,2-Dichloroethane	2.3	U	2.3	4.0
1,1-Dichloroethene	22		1.8	4.0
cis-1,2-Dichloroethene	44		2.6	4.0
trans-1,2-Dichloroethene	1.8	U	1.8	4.0
1,2-Dichloropropane	2.1	U	2.1	4.0
1,3-Dichloropropane	1.6	U	1.6	4.0
2,2-Dichloropropane	1.4	U	1.4	4.0
1,1-Dichloropropene	1.2	U	1.2	4.0
cis-1,3-Dichloropropene	0.56	U	0.56	4.0
trans-1,3-Dichloropropene	0.56	U	0.56	4.0
Ethylbenzene	1.8	U	1.8	4.0
Hexachlorobutadiene	1.6	U	1.6	4.0
Isopropylbenzene	0.76	U	0.76	4.0
4-Isopropyltoluene	2.8	U	2.8	4.0
Methylene Chloride	16	U	16	20
Naphthalene	10	U	10	20
N-Propylbenzene	2.4	U	2.4	4.0
Styrene	3.9	U	3.9	8.0
1,1,1,2-Tetrachloroethane	2.5	U	2.5	4.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-39

Sdg Number: 660-20941

Lab Sample ID: 660-20941-10

Date Sampled: 02/01/2008 1016

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch:	660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B			Lab File ID:	1JB1517.D
Dilution:	4.0			Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1825			Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1825				

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.60	U	0.60	4.0
Tetrachloroethene	2.0	U	2.0	4.0
Toluene	2.0	U	2.0	4.0
1,2,3-Trichlorobenzene	3.1	U	3.1	4.0
1,2,4-Trichlorobenzene	2.3	U	2.3	4.0
1,1,1-Trichloroethane	1.8	U	1.8	4.0
1,1,2-Trichloroethane	1.9	U	1.9	4.0
Trichloroethene	190		2.0	4.0
Trichlorofluoromethane	10	U	10	20
1,2,3-Trichloropropane	0.72	U	0.72	4.0
1,2,4-Trimethylbenzene	3.4	U	3.4	4.0
1,3,5-Trimethylbenzene	2.2	U	2.2	4.0
Vinyl chloride	2.0	U	2.0	4.0
o-Xylene	2.0	U	2.0	4.0
Acetone	40	U	40	80
2-Butanone (MEK)	34	U	34	40
4-Methyl-2-pentanone (MIBK)	15	U	15	40
Carbon disulfide	3.4	U	3.4	4.0
2-Hexanone	18	U	18	40
Methyl tert-butyl ether	1.8	U	1.8	4.0
m-Xylene & p-Xylene	2.4	U	2.4	8.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	121		77 - 122	
4-Bromofluorobenzene	121		74 - 126	
Dibromofluoromethane	117		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: MW-253

Lab Sample ID: 660-20941-11

Date Sampled: 02/01/2008 0930

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58013 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1460.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 0821 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 0821

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	75		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
cis-1,2-Dichloroethene	130		0.65	1.0
trans-1,2-Dichloroethene	0.51	I	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-253

Sdg Number: 660-20941

Lab Sample ID: 660-20941-11

Date Sampled: 02/01/2008 0930

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1460.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0821		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0821			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Tetrachloroethene	5.4		0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	1.1		0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	100	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	100	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-253

Sdg Number: 660-20941

Lab Sample ID: 660-20941-11

Date Sampled: 02/01/2008 0930

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58240

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1522.D

Dilution: 100

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 2017

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 2017

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1-Dichloroethene	400	D	45	100
Trichloroethene	6200	D	50	100

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-254

Sdg Number: 660-20941

Lab Sample ID: 660-20941-12

Date Sampled: 02/01/2008 0840

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1511.D
Dilution:	10		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1610		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1610			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	5.0	U	5.0	10
Bromobenzene	5.8	U	5.8	10
Chlorobromomethane	5.8	U	5.8	10
Dichlorobromomethane	3.5	U	3.5	10
Bromoform	5.8	U	5.8	10
Bromomethane	25	U	25	50
n-Butylbenzene	6.7	U	6.7	10
sec-Butylbenzene	6.3	U	6.3	10
tert-Butylbenzene	8.4	U	8.4	10
Carbon tetrachloride	4.2	U	4.2	10
Chlorobenzene	6.3	U	6.3	10
Chloroethane	25	U	25	50
Chloroform	9.0	U	9.0	10
Chloromethane	10	U	10	40
2-Chlorotoluene	6.5	U	6.5	10
4-Chlorotoluene	5.2	U	5.2	10
Chlorodibromomethane	3.4	U	3.4	10
1,2-Dibromo-3-Chloropropane	25	U	25	50
Ethylene Dibromide	5.0	U	5.0	10
Dibromomethane	4.1	U	4.1	10
1,2-Dichlorobenzene	4.4	U	4.4	10
1,3-Dichlorobenzene	6.4	U	6.4	10
1,4-Dichlorobenzene	5.2	U	5.2	10
Dichlorodifluoromethane	25	U	25	50
1,1-Dichloroethane	6.0	I	5.2	10
1,2-Dichloroethane	5.7	U	5.7	10
1,1-Dichloroethene	37		4.5	10
cis-1,2-Dichloroethene	6.5	U	6.5	10
trans-1,2-Dichloroethene	4.4	U	4.4	10
1,2-Dichloropropane	5.2	U	5.2	10
1,3-Dichloropropane	3.9	U	3.9	10
2,2-Dichloropropane	3.6	U	3.6	10
1,1-Dichloropropene	3.1	U	3.1	10
cis-1,3-Dichloropropene	1.4	U	1.4	10
trans-1,3-Dichloropropene	1.4	U	1.4	10
Ethylbenzene	4.4	U	4.4	10
Hexachlorobutadiene	4.0	U	4.0	10
Isopropylbenzene	1.9	U	1.9	10
4-Isopropyltoluene	6.9	U	6.9	10
Methylene Chloride	40	U	40	50
Naphthalene	25	U	25	50
N-Propylbenzene	5.9	U	5.9	10
Styrene	9.8	U	9.8	20
1,1,1,2-Tetrachloroethane	6.3	U	6.3	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-254

Sdg Number: 660-20941

Lab Sample ID: 660-20941-12

Date Sampled: 02/01/2008 0840

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1511.D
Dilution:	10		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1610		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1610			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	1.5	U	1.5	10
Toluene	5.1	U	5.1	10
1,2,3-Trichlorobenzene	7.7	U	7.7	10
1,2,4-Trichlorobenzene	5.8	U	5.8	10
1,1,1-Trichloroethane	4.6	U	4.6	10
1,1,2-Trichloroethane	4.7	U	4.7	10
Trichloroethene	760		5.0	10
Trichlorofluoromethane	25	U	25	50
1,2,3-Trichloropropane	1.8	U	1.8	10
1,2,4-Trimethylbenzene	8.6	U	8.6	10
1,3,5-Trimethylbenzene	5.4	U	5.4	10
Vinyl chloride	5.0	U	5.0	10
o-Xylene	5.0	U	5.0	10
Acetone	99	U	99	200
2-Butanone (MEK)	84	U	84	100
4-Methyl-2-pentanone (MIBK)	38	U	38	100
Carbon disulfide	8.5	U	8.5	10
2-Hexanone	44	U	44	100
Methyl tert-butyl ether	4.4	U	4.4	10
m-Xylene & p-Xylene	6.0	U	6.0	20
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	94		77 - 122	
4-Bromofluorobenzene	92		74 - 126	
Dibromofluoromethane	91		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: **MW-254**

Sdg Number: 660-20941

Lab Sample ID: 660-20941-12

Date Sampled: 02/01/2008 0840

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58264

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1530.D

Dilution: 500

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1953

Run Type: DL

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1953

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Tetrachloroethene	11000	D	250	500

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-8D

Sdg Number: 660-20941

Lab Sample ID: 660-20941-13

Date Sampled: 02/01/2008 0906

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58240 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1528.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 2231 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 2231

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: MW-8D

Lab Sample ID: 660-20941-13

Date Sampled: 02/01/2008 0906

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58240

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1528.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 2231

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 2231

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	89	77 - 122
4-Bromofluorobenzene	83	74 - 126
Dibromofluoromethane	88	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: MW-123

Lab Sample ID: 660-20941-14

Date Sampled: 02/01/2008 0921

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1463.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0928		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0928			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-123

Sdg Number: 660-20941

Lab Sample ID: 660-20941-14

Date Sampled: 02/01/2008 0921

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58013

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1463.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 0928

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 0928

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	98	77 - 122
4-Bromofluorobenzene	94	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: RW-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-15

Date Sampled: 02/01/2008 0910

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58013 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1464.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 0950 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 0950

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	2.0	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: RW-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-15

Date Sampled: 02/01/2008 0910

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1464.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0950		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0950			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	1.5		0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	109	77 - 122
4-Bromofluorobenzene	105	74 - 126
Dibromofluoromethane	107	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Dup-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-16

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58240 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1518.D
 Dilution: 10 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 1847 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 1847

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	5.0	U	5.0	10
Bromobenzene	5.8	U	5.8	10
Chlorobromomethane	5.8	U	5.8	10
Dichlorobromomethane	3.5	U	3.5	10
Bromoform	5.8	U	5.8	10
Bromomethane	25	U	25	50
n-Butylbenzene	6.7	U	6.7	10
sec-Butylbenzene	6.3	U	6.3	10
tert-Butylbenzene	8.4	U	8.4	10
Carbon tetrachloride	4.2	U	4.2	10
Chlorobenzene	6.3	U	6.3	10
Chloroethane	25	U	25	50
Chloroform	9.0	U	9.0	10
Chloromethane	10	U	10	40
2-Chlorotoluene	6.5	U	6.5	10
4-Chlorotoluene	5.2	U	5.2	10
Chlorodibromomethane	3.4	U	3.4	10
1,2-Dibromo-3-Chloropropane	25	U	25	50
Ethylene Dibromide	5.0	U	5.0	10
Dibromomethane	4.1	U	4.1	10
1,2-Dichlorobenzene	4.4	U	4.4	10
1,3-Dichlorobenzene	6.4	U	6.4	10
1,4-Dichlorobenzene	5.2	U	5.2	10
Dichlorodifluoromethane	25	U	25	50
1,1-Dichloroethane	31		5.2	10
1,2-Dichloroethane	5.7	U	5.7	10
1,1-Dichloroethene	140		4.5	10
cis-1,2-Dichloroethene	16		6.5	10
trans-1,2-Dichloroethene	4.4	U	4.4	10
1,2-Dichloropropane	5.2	U	5.2	10
1,3-Dichloropropane	3.9	U	3.9	10
2,2-Dichloropropane	3.6	U	3.6	10
1,1-Dichloropropene	3.1	U	3.1	10
cis-1,3-Dichloropropene	1.4	U	1.4	10
trans-1,3-Dichloropropene	1.4	U	1.4	10
Ethylbenzene	4.4	U	4.4	10
Hexachlorobutadiene	4.0	U	4.0	10
Isopropylbenzene	1.9	U	1.9	10
4-Isopropyltoluene	6.9	U	6.9	10
Methylene Chloride	40	U	40	50
Naphthalene	25	U	25	50
N-Propylbenzene	5.9	U	5.9	10
Styrene	9.8	U	9.8	20
1,1,1,2-Tetrachloroethane	6.3	U	6.3	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Dup-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-16

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1518.D
Dilution:	10		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1847		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1847			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	1.5	U	1.5	10
Tetrachloroethene	5.0	U	5.0	10
Toluene	5.1	U	5.1	10
1,2,3-Trichlorobenzene	7.7	U	7.7	10
1,2,4-Trichlorobenzene	5.8	U	5.8	10
1,1,1-Trichloroethane	4.6	U	4.6	10
1,1,2-Trichloroethane	4.7	U	4.7	10
Trichloroethene	150		5.0	10
Trichlorofluoromethane	25	U	25	50
1,2,3-Trichloropropane	1.8	U	1.8	10
1,2,4-Trimethylbenzene	8.6	U	8.6	10
1,3,5-Trimethylbenzene	5.4	U	5.4	10
Vinyl chloride	5.0	U	5.0	10
o-Xylene	5.0	U	5.0	10
Acetone	99	U	99	200
2-Butanone (MEK)	84	U	84	100
4-Methyl-2-pentanone (MIBK)	38	U	38	100
Carbon disulfide	8.5	U	8.5	10
2-Hexanone	44	U	44	100
Methyl tert-butyl ether	4.4	U	4.4	10
m-Xylene & p-Xylene	6.0	U	6.0	20

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	104	77 - 122
4-Bromofluorobenzene	107	74 - 126
Dibromofluoromethane	99	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EB-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-17

Date Sampled: 02/01/2008 1145

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1466.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1035		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1035			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EB-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-17

Date Sampled: 02/01/2008 1145

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1466.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1035		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1035			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	93	74 - 126
Dibromofluoromethane	95	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: TB-31

Lab Sample ID: 660-20941-18

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58013

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1467.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1057

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1057

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: TB-31

Lab Sample ID: 660-20941-18

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58013

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1467.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1057

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1057

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	105	77 - 122
4-Bromofluorobenzene	98	74 - 126
Dibromofluoromethane	102	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 7

Sdg Number: 660-20941

Lab Sample ID: 660-20941-19

Date Sampled: 02/01/2008 0854

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58013 Instrument ID: BVMJ GC/MS
 Preparation: 5030B Lab File ID: 1JB1468.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 1119 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 1119

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: Private Well 7

Lab Sample ID: 660-20941-19

Date Sampled: 02/01/2008 0854

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58013

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1468.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1119

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1119

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	99	77 - 122
4-Bromofluorobenzene	92	74 - 126
Dibromofluoromethane	95	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 57

Sdg Number: 660-20941

Lab Sample ID: 660-20941-20

Date Sampled: 02/01/2008 0820

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1469.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1142		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1142			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 57

Sdg Number: 660-20941

Lab Sample ID: 660-20941-20

Date Sampled: 02/01/2008 0820

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58013	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1469.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1142		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1142			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	110		77 - 122	
4-Bromofluorobenzene	105		74 - 126	
Dibromofluoromethane	109		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 59

Sdg Number: 660-20941

Lab Sample ID: 660-20941-21

Date Sampled: 02/01/2008 0908

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1526.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2146		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2146			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 59

Sdg Number: 660-20941

Lab Sample ID: 660-20941-21

Date Sampled: 02/01/2008 0908

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID: BVMJ GC/MS
Preparation:	5030B		Lab File ID: 1JB1526.D
Dilution:	1.0		Initial Weight/Volume: 5 mL
Date Analyzed:	02/15/2008 2146		Final Weight/Volume: 5 mL
Date Prepared:	02/15/2008 2146		

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	22	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0
Surrogate	%Rec		Acceptance Limits	
Toluene-d8 (Surr)	108		77 - 122	
4-Bromofluorobenzene	106		74 - 126	
Dibromofluoromethane	104		70 - 130	

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 38

Sdg Number: 660-20941

Lab Sample ID: 660-20941-22

Date Sampled: 02/01/2008 1055

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1525.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2124		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2124			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	3.4		0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	4.9		0.45	1.0
cis-1,2-Dichloroethene	0.93	I	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 38

Sdg Number: 660-20941

Lab Sample ID: 660-20941-22

Date Sampled: 02/01/2008 1055

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1525.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2124		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2124			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	97	77 - 122
4-Bromofluorobenzene	94	74 - 126
Dibromofluoromethane	95	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 102

Sdg Number: 660-20941

Lab Sample ID: 660-20941-23

Date Sampled: 02/01/2008 1115

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1524.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2101		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2101			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: Private Well 102

Lab Sample ID: 660-20941-23

Date Sampled: 02/01/2008 1115

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-58240	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1524.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2101		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2101			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	4.4		0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	101	77 - 122
4-Bromofluorobenzene	97	74 - 126
Dibromofluoromethane	100	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: TB-32

Lab Sample ID: 660-20941-24

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B

Analysis Batch: 660-58264

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1517.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1447

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1447

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: TB-32

Sdg Number: 660-20941

Lab Sample ID: 660-20941-24

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 660-58264 Instrument ID: BVMH GC/MS
 Preparation: 5030B Lab File ID: 1HB1517.D
 Dilution: 1.0 Initial Weight/Volume: 5 mL
 Date Analyzed: 02/15/2008 1447 Final Weight/Volume: 5 mL
 Date Prepared: 02/15/2008 1447

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	106	77 - 122
4-Bromofluorobenzene	83	74 - 126
Dibromofluoromethane	87	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-202

Sdg Number: 660-20941

Lab Sample ID: 660-20941-1

Date Sampled: 02/01/2008 0810

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/15/2008 1548
 Date Prepared: 02/15/2008 1548

Analysis Batch: 660-58239

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1510.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-134

Sdg Number: 660-20941

Lab Sample ID: 660-20941-2

Date Sampled: 02/01/2008 1132

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58239	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1514.D
Dilution:	2.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1717		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1717			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	28		2.0	2.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: DW-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-3

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58015	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1452.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0522		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0522			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Sdg Number: 660-20941

Client Sample ID: MW-42

Lab Sample ID: 660-20941-4

Date Sampled: 02/01/2008 1056

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58015

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1453.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 0544

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 0544

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	72		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-43

Sdg Number: 660-20941

Lab Sample ID: 660-20941-5

Date Sampled: 02/01/2008 1103

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58239	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1515.D
Dilution:	10		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1740		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1740			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	570		10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-6

Date Sampled: 02/01/2008 0915

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57848

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1529.D

Dilution: 100

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1930

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1930

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	300		100	100

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: IWI-2

Sdg Number: 660-20941

Lab Sample ID: 660-20941-7

Date Sampled: 02/01/2008 1025

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57848	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1521.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1627		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1627			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	19		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EW-UAFG-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-8

Date Sampled: 02/01/2008 1000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57848	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1527.D
Dilution:	50		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1844		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1844			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	200		50	50

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-38

Sdg Number: 660-20941

Lab Sample ID: 660-20941-9

Date Sampled: 02/01/2008 0940

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 1.0
Date Analyzed: 02/15/2008 0736
Date Prepared: 02/15/2008 0736

Analysis Batch: 660-58015

Instrument ID: BVMJ GC/MS
Lab File ID: 1JB1458.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	24		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-39

Sdg Number: 660-20941

Lab Sample ID: 660-20941-10

Date Sampled: 02/01/2008 1016

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58239	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1517.D
Dilution:	4.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1825		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1825			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	160		4.0	4.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-253

Sdg Number: 660-20941

Lab Sample ID: 660-20941-11

Date Sampled: 02/01/2008 0930

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 100
 Date Analyzed: 02/15/2008 2017
 Date Prepared: 02/15/2008 2017

Analysis Batch: 660-58239

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1522.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	350		100	100

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-254

Sdg Number: 660-20941

Lab Sample ID: 660-20941-12

Date Sampled: 02/01/2008 0840

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58239	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1511.D
Dilution:	10		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1610		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1610			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	10	U	10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-8D

Sdg Number: 660-20941

Lab Sample ID: 660-20941-13

Date Sampled: 02/01/2008 0906

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-58015

Instrument ID: BVMJ GC/MS

Preparation: 5030B

Lab File ID: 1JB1462.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 0905

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 0905

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: MW-123

Sdg Number: 660-20941

Lab Sample ID: 660-20941-14

Date Sampled: 02/01/2008 0921

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58015	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1463.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 0928		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 0928			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: RW-1

Sdg Number: 660-20941

Lab Sample ID: 660-20941-15

Date Sampled: 02/01/2008 0910

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57848	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1524.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1735		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1735			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Dup-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-16

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
Preparation: 5030B
Dilution: 10
Date Analyzed: 02/15/2008 1847
Date Prepared: 02/15/2008 1847

Analysis Batch: 660-58239

Instrument ID: BVMJ GC/MS
Lab File ID: 1JB1518.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	560		10	10

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: EB-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-17

Date Sampled: 02/01/2008 1145

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57848

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1525.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1758

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1758

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: TB-31

Sdg Number: 660-20941

Lab Sample ID: 660-20941-18

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58015	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1467.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1057		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1057			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 7

Sdg Number: 660-20941

Lab Sample ID: 660-20941-19

Date Sampled: 02/01/2008 0854

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/15/2008 1633
 Date Prepared: 02/15/2008 1633

Analysis Batch: 660-58239

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1512.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 57

Sdg Number: 660-20941

Lab Sample ID: 660-20941-20

Date Sampled: 02/01/2008 0820

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58239	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1513.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1655		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1655			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 59

Sdg Number: 660-20941

Lab Sample ID: 660-20941-21

Date Sampled: 02/01/2008 0908

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-58239	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1526.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 2146		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 2146			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 38

Sdg Number: 660-20941

Lab Sample ID: 660-20941-22

Date Sampled: 02/01/2008 1055

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method:	8260C SIM/ID	Analysis Batch: 660-57848	Instrument ID:	BVMH GC/MS
Preparation:	5030B		Lab File ID:	1HB1519.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/15/2008 1541		Final Weight/Volume:	5 mL
Date Prepared:	02/15/2008 1541			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	96		1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: Private Well 102

Sdg Number: 660-20941

Lab Sample ID: 660-20941-23

Date Sampled: 02/01/2008 1115

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57848

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1518.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1517

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1517

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-20941-1

Client Sample ID: TB-32

Sdg Number: 660-20941

Lab Sample ID: 660-20941-24

Date Sampled: 02/01/2008 0000

Client Matrix: Water

Date Received: 02/01/2008 1518

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID

Analysis Batch: 660-57848

Instrument ID: BVMH GC/MS

Preparation: 5030B

Lab File ID: 1HB1517.D

Dilution: 1.0

Initial Weight/Volume: 5 mL

Date Analyzed: 02/15/2008 1447

Final Weight/Volume: 5 mL

Date Prepared: 02/15/2008 1447

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-20941	2/1/2008	SW-846	MW-202	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	MW-39	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	MW-253	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	MW-254	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	MW-8D	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	MW-123	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	RW-1	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	Dup-31	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	EB-31	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	TB-31	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	Private Well 7	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	MW-134	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	Private Well 57	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	Private Well 59	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	Private Well 38	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	Private Well 102	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	TB-32	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	DW-1	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	MW-42	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	MW-43	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	IWI-1	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	IWI-2	Water	Yes	--	--	--	--	
660-20941	2/1/2008	SW-846	EW-UAFG-1	Water	No	--	--	--	--	VOC – CCAL %D
660-20941	2/1/2008	SW-846	MW-38	Water	Yes	--	--	--	--	

¹ Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

ARCADIS

Attachment 9

DATA REVIEW FOR
LOCKHEED MARTIN - TALLEVAST
TALLEVAST, FLORIDA

SDG #660-21028

VOLATILE ANALYSES

Analyses performed by:

TestAmerica Laboratories
Tampa, Florida

Review performed by:



Syracuse, New York
Report #8142R

Summary

The following is an assessment of the data package for sample delivery group (SDG) #660-21028 for sampling from the Lockheed Martin - Tallevast Site. Included with this assessment are the corrected sample results and the sample compliance report. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample Date	Analysis				
				VOC	SVOC	PCB	MET	MISC
PW-64	660-21028-1	Water	2/4/2008	X				
PW-65	660-21028-2	Water	2/4/2008	X				

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United States Environmental Protection Agency) USEPA SW-846 Method 8260B total ion monitoring (TIM) and 8260C selective ion monitoring (SIM) by isotope dilution. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999 and January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
- J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
- B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
- JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
- E The compound was quantitated above the calibration range.
- D Concentration is based on a diluted sample analysis.
- C Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).
- UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
- R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

Data Assessment

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation
SW-846 8260B/C	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	48 hours from collection to extraction and 14 days from extraction to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method, trip, and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure contamination of samples during shipment. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No compounds were detected in the associated blanks.

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to insure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds analyzed by 8260B associated with the initial calibration standards must exhibit a %RSD less than the control limit (15%) or a correlation coefficient greater than 0.99. The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

4.2 Continuing Calibration

All target compounds analyzed by 8260B associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (20%) with the exception of 1,4-Dioxane must exhibit a %D less than the control limit of (50%). The compounds analyzed by method 8260B TIM must exhibit a RRF value greater than the control limit (0.05). The analysis of the target compound 1,4-Dioxane performed by method 8260C SIM by isotope dilution must exhibit a RRF value greater than the control limit (0.005).

All compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
PW-64 PW-65	CCV %D	1,3-Dichloropropane	23.5%
		2-Hexanone	32.6%
		2-Butanone	23.0%
		4-Methyl-2-pentanone	29.9%
		Acetone	36.1%
		Bromomethane	29.2%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
Initial and Continuing Calibration	RRF <0.05	Non-detect	R
		Detect	J
	RRF <0.01 ¹	Non-detect	R
		Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
		Detect	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90%	Non-detect	R
		Detect	J

- RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e. ketones, 1,4-Dioxane, etc.)

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria insure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

All internal standard areas and retention times were within established limits.

7. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

A MS/MSD was not performed on a sample location within this SDG.

8. Laboratory Control Sample (LCS)/Laboratory Control Sample Duplicate (LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

All compounds associated with the LCS/LCSD analysis exhibited acceptable recoveries and RPD between the LCS/LCSD recoveries.

9. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate.

A field duplicate was not performed on a sample location within this SDG.

10. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

All identified compounds met the specified criteria.

11. System Performance and Overall Assessment

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

Analytical Data

Client: ARCADIS

Job Number: 660-21028-1

Client Sample ID: PW-64

Sdg Number: 660-21028

Lab Sample ID: 660-21028-1

Date Sampled: 02/04/2008 0955

Client Matrix: Water

Date Received: 02/07/2008 1100

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57855	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1627.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/16/2008 2053		Final Weight/Volume:	5 mL
Date Prepared:	02/16/2008 2053			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-21028-1

Sdg Number: 660-21028

Client Sample ID: PW-64

Lab Sample ID: 660-21028-1

Date Sampled: 02/04/2008 0955

Client Matrix: Water

Date Received: 02/07/2008 1100

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57855	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1627.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/16/2008 2053		Final Weight/Volume:	5 mL
Date Prepared:	02/16/2008 2053			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	103	77 - 122
4-Bromofluorobenzene	99	74 - 126
Dibromofluoromethane	103	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-21028-1

Client Sample ID: PW-65

Sdg Number: 660-21028

Lab Sample ID: 660-21028-2

Date Sampled: 02/04/2008 1118

Client Matrix: Water

Date Received: 02/07/2008 1100

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57855	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1628.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/16/2008 2116		Final Weight/Volume:	5 mL
Date Prepared:	02/16/2008 2116			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
Benzene	0.50	U	0.50	1.0
Bromobenzene	0.58	U	0.58	1.0
Chlorobromomethane	0.58	U	0.58	1.0
Dichlorobromomethane	0.35	U	0.35	1.0
Bromoform	0.58	U	0.58	1.0
Bromomethane	2.5	U	2.5	5.0
n-Butylbenzene	0.67	U	0.67	1.0
sec-Butylbenzene	0.63	U	0.63	1.0
tert-Butylbenzene	0.84	U	0.84	1.0
Carbon tetrachloride	0.42	U	0.42	1.0
Chlorobenzene	0.63	U	0.63	1.0
Chloroethane	2.5	U	2.5	5.0
Chloroform	0.90	U	0.90	1.0
Chloromethane	1.0	U	1.0	4.0
2-Chlorotoluene	0.65	U	0.65	1.0
4-Chlorotoluene	0.52	U	0.52	1.0
Chlorodibromomethane	0.34	U	0.34	1.0
1,2-Dibromo-3-Chloropropane	2.5	U	2.5	5.0
Ethylene Dibromide	0.50	U	0.50	1.0
Dibromomethane	0.41	U	0.41	1.0
1,2-Dichlorobenzene	0.44	U	0.44	1.0
1,3-Dichlorobenzene	0.64	U	0.64	1.0
1,4-Dichlorobenzene	0.52	U	0.52	1.0
Dichlorodifluoromethane	2.5	U	2.5	5.0
1,1-Dichloroethane	0.52	U	0.52	1.0
1,2-Dichloroethane	0.57	U	0.57	1.0
1,1-Dichloroethene	0.45	U	0.45	1.0
cis-1,2-Dichloroethene	0.65	U	0.65	1.0
trans-1,2-Dichloroethene	0.44	U	0.44	1.0
1,2-Dichloropropane	0.52	U	0.52	1.0
1,3-Dichloropropane	0.39	U	0.39	1.0
2,2-Dichloropropane	0.36	U	0.36	1.0
1,1-Dichloropropene	0.31	U	0.31	1.0
cis-1,3-Dichloropropene	0.14	U	0.14	1.0
trans-1,3-Dichloropropene	0.14	U	0.14	1.0
Ethylbenzene	0.44	U	0.44	1.0
Hexachlorobutadiene	0.40	U	0.40	1.0
Isopropylbenzene	0.19	U	0.19	1.0
4-Isopropyltoluene	0.69	U	0.69	1.0
Methylene Chloride	4.0	U	4.0	5.0
Naphthalene	2.5	U	2.5	5.0
N-Propylbenzene	0.59	U	0.59	1.0
Styrene	0.98	U	0.98	2.0
1,1,1,2-Tetrachloroethane	0.63	U	0.63	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-21028-1

Sdg Number: 660-21028

Client Sample ID: PW-65

Lab Sample ID: 660-21028-2

Date Sampled: 02/04/2008 1118

Client Matrix: Water

Date Received: 02/07/2008 1100

8260B Volatile Organic Compounds by GC/MS

Method:	8260B	Analysis Batch: 660-57855	Instrument ID:	BVMJ GC/MS
Preparation:	5030B		Lab File ID:	1JB1628.D
Dilution:	1.0		Initial Weight/Volume:	5 mL
Date Analyzed:	02/16/2008 2116		Final Weight/Volume:	5 mL
Date Prepared:	02/16/2008 2116			

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,1,2,2-Tetrachloroethane	0.15	U	0.15	1.0
Tetrachloroethene	0.50	U	0.50	1.0
Toluene	0.51	U	0.51	1.0
1,2,3-Trichlorobenzene	0.77	U	0.77	1.0
1,2,4-Trichlorobenzene	0.58	U	0.58	1.0
1,1,1-Trichloroethane	0.46	U	0.46	1.0
1,1,2-Trichloroethane	0.47	U	0.47	1.0
Trichloroethene	0.50	U	0.50	1.0
Trichlorofluoromethane	2.5	U	2.5	5.0
1,2,3-Trichloropropane	0.18	U	0.18	1.0
1,2,4-Trimethylbenzene	0.86	U	0.86	1.0
1,3,5-Trimethylbenzene	0.54	U	0.54	1.0
Vinyl chloride	0.50	U	0.50	1.0
o-Xylene	0.50	U	0.50	1.0
Acetone	9.9	U	9.9	20
2-Butanone (MEK)	8.4	U	8.4	10
4-Methyl-2-pentanone (MIBK)	3.8	U	3.8	10
Carbon disulfide	0.85	U	0.85	1.0
2-Hexanone	4.4	U	4.4	10
Methyl tert-butyl ether	0.44	U	0.44	1.0
m-Xylene & p-Xylene	0.60	U	0.60	2.0

Surrogate	%Rec	Acceptance Limits
Toluene-d8 (Surr)	103	77 - 122
4-Bromofluorobenzene	101	74 - 126
Dibromofluoromethane	103	70 - 130

Analytical Data

Client: ARCADIS

Job Number: 660-21028-1

Client Sample ID: PW-64

Sdg Number: 660-21028

Lab Sample ID: 660-21028-1

Date Sampled: 02/04/2008 0955

Client Matrix: Water

Date Received: 02/07/2008 1100

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/16/2008 2053
 Date Prepared: 02/16/2008 2053

Analysis Batch: 660-57857

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1627.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

Analytical Data

Client: ARCADIS

Job Number: 660-21028-1

Client Sample ID: PW-65

Sdg Number: 660-21028

Lab Sample ID: 660-21028-2

Date Sampled: 02/04/2008 1118

Client Matrix: Water

Date Received: 02/07/2008 1100

8260C SIM/ID 1,4-Dioxane by SIM/Isotopic Dilution

Method: 8260C SIM/ID
 Preparation: 5030B
 Dilution: 1.0
 Date Analyzed: 02/16/2008 2116
 Date Prepared: 02/16/2008 2116

Analysis Batch: 660-57857

Instrument ID: BVMJ GC/MS

Lab File ID: 1JB1628.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Result (ug/L)	Qualifier	MDL	PQL
1,4-Dioxane	1.0	U	1.0	1.0

SAMPLE COMPLIANCE REPORT

SAMPLE COMPLIANCE REPORT

Sample Delivery Group	Sampling Date	Protocol	Sample ID	Matrix	Compliance ¹					Noncompliance
					VOC	SVOC	PCB/PEST/HERB	MET	MISC	
660-21028	2/4/2008	SW-846	PW-64	Water	Yes	--	--	--	--	
660-21028	2/4/2008	SW-846	PW-65	Water	Yes	--	--	--	--	

1 Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

Appendix D

Historical Data

APPENDIX D-1
HISTORICAL SUMMARY OF MONITORING WELL GROUNDWATER ELEVATIONS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Well ID	Zone	Top of Inner Casing (ft-msl)	October 17-19 2006		December 28 2006		February 20 2007		April 23 2007		December 3-6 2007		December 7 2007		January 21-22 2008		January 25-31 2008	
			Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation
DW-1	AF Gravel	31.00	11.40	19.60	13.87	17.13	13.81	17.19	NM	NM	13.15	17.85	13.25	17.75	14.54	16.46	NM	NM
EW-102	LSAS	30.52	NM	NM	NM	NM	NM	NM	NM	NM	8.08	22.44	8.20	22.32	NM	NM	11.82	18.70
EW-103	USAS	29.86	NM	NM	NM	NM	NM	NM	NM	NM	3.73	26.13	3.77	26.09	NM	NM	9.30	20.56
EW-104	LSAS	29.77	NM	NM	NM	NM	NM	NM	NM	NM	5.13	24.64	5.18	24.59	NM	NM	16.50	13.27
EW-105	USAS	30.40	NM	NM	NM	NM	NM	NM	NM	NM	4.12	26.28	4.16	26.24	NM	NM	15.31	15.09
EW-106	LSAS	30.33	NM	NM	NM	NM	NM	NM	NM	NM	5.57	24.76	5.64	24.69	NM	NM	22.47	7.86
EW-107	USAS	29.64	NM	NM	NM	NM	NM	NM	NM	NM	3.32	26.32	3.39	26.25	NM	NM	3.48	26.16
EW-109	USAS	30.10	NM	NM	NM	NM	NM	NM	NM	NM	3.30	26.80	3.38	26.72	NM	NM	3.48	26.62
EW-110	LSAS	30.12	NM	NM	NM	NM	NM	NM	NM	NM	3.60	26.52	3.69	26.43	NM	NM	14.80	15.32
EW-UAFG-1	AF Gravel	31.66	NM	NM	NM	NM	NM	NM	NM	NM	16.09	15.57	15.91	15.75	NM	NM	16.84	14.82
EXL-1 (EW-108)	LSAS	30.09	NM	NM	NM	NM	NM	NM	NM	NM	8.28	21.81	8.38	21.71	NM	NM	16.50	13.59
EXU-1 (EW-101)	USAS	30.31	NM	NM	NM	NM	NM	NM	NM	NM	3.99	26.32	4.04	26.27	NM	NM	15.92	14.39
IWI-1	AF Gravel	31.71	NM	NM	16.52	15.19	16.89	14.82	NM	NM	16.25	15.46	16.28	15.43	16.45	15.26	NM	NM
IWI-2	Zone 3-4 Clay	31.62	NM	NM	18.40	13.22	16.94	14.68	NM	NM	18.68	12.94	18.79	12.83	19.97	11.65	NM	NM
MW-2	USAS	29.85	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
MW-3	USAS	30.52	2.49	28.03	3.65	26.87	4.01	26.51	4.91	25.61	4.11	26.41	4.16	26.36	4.32	26.20	NM	NM
MW-4	USAS	31.50	3.50	28.00	4.68	26.82	5.12	26.38	5.84	25.66	5.10	26.40	5.17	26.33	5.50	26.00	NM	NM
MW-5	USAS	32.17	4.22	27.95	4.91	27.26	5.30	26.87	6.18	25.99	5.44	26.73	5.53	26.64	5.72	26.45	NM	NM
MW-6	USAS	31.92	4.12	27.80	4.88	27.04	5.29	26.63	6.02	25.90	5.41	26.51	5.47	26.45	5.78	26.14	NM	NM
MW-7D	USAS	31.30	2.89	28.41	3.43	27.87	4.17	27.13	4.91	26.39	4.12	27.18	4.22	27.08	4.21	27.09	NM	NM
MW-7S	USAS	31.50	3.10	28.40	3.63	27.87	4.15	27.35	5.16	26.34	4.31	27.19	NM	NM	4.41	27.09	NM	NM
MW-8D	USAS	30.96	2.58	28.38	3.40	27.56	3.91	27.05	5.23	25.73	4.01	26.95	4.09	26.87	4.06	26.90	NM	NM
MW-8S	USAS	30.99	2.55	28.44	3.44	27.55	3.90	27.09	4.92	26.07	4.06	26.93	NM	NM	4.13	26.86	NM	NM
MW-9D	USAS	30.21	2.79	27.42	3.61	26.60	4.01	26.20	4.66	25.55	4.31	25.90	4.35	25.86	4.60	25.61	NM	NM
MW-9S	USAS	30.24	2.79	27.45	3.65	26.59	4.05	26.19	4.96	25.28	4.33	25.91	NM	NM	4.65	25.59	NM	NM
MW-10	USAS	31.74	3.79	27.95	5.35	26.39	5.71	26.03	6.40	25.34	5.40	26.34	5.46	26.28	5.65	26.09	11.25	20.49
MW-11	USAS	31.87	3.63	28.24	4.82	27.05	5.14	26.73	5.95	25.92	5.20	26.67	5.25	26.62	5.46	26.41	NM	NM
MW-12	USAS	31.04	NM	NM	4.49	26.55	4.94	26.10	5.80	25.24	4.90	26.14	4.93	26.11	5.17	25.87	NM	NM
MW-13D	USAS	30.85	3.48	27.37	4.30	26.55	4.95	25.90	5.60	25.25	4.96	25.89	5.01	25.84	5.30	25.55	NM	NM
MW-13S	USAS	30.66	NM	NM	4.10	26.56	4.77	25.89	NM	NM	4.91	25.75	4.81	25.85	5.12	25.54	NM	NM
MW-14D	USAS	29.75	2.75	27.00	3.99	25.76	3.98	25.77	5.52	24.23	4.06	25.69	4.12	25.63	4.36	25.39	NM	NM
MW-14S	USAS	29.74	2.74	27.00	3.32	26.42	3.95	25.79	4.65	25.09	4.04	25.70	NM	NM	4.35	25.39	NM	NM
MW-15D	USAS	30.20	3.07	27.13	3.76	26.44	4.31	25.89	4.90	25.30	4.48	25.72	4.54	25.66	4.87	25.33	NM	NM
MW-15S	USAS	30.09	2.93	27.16	3.63	26.46	4.18	25.91	5.00	25.09	4.36	25.73	NM	NM	4.74	25.35	NM	NM
MW-16D	USAS	27.26	2.69	24.57	2.36	24.90	2.72	24.54	NM	NM	3.04	24.22	NM	NM	3.16	24.10	NM	NM
MW-16S	USAS	27.26	NM	NM	2.49	24.77	2.85	24.41	NM	NM	3.15	24.11	NM	NM	3.24	24.02	NM	NM
MW-17D	USAS	30.23	3.18	27.05	3.79	26.44	4.40	25.83	5.21	25.02	4.57	25.66	4.63	25.60	4.91	25.32	NM	NM
MW-17S	USAS	30.09	3.00	27.09	3.63	26.46	5.28	24.81	5.01	25.08	4.40	25.69	NM	NM	4.74	25.35	NM	NM
MW-18D	USAS	28.04	NM	NM	2.75	25.29	3.12	24.92	NM	NM	NM	NM	NM	NM	NM	NM	3.04	25.00
MW-18S	USAS	28.00	NM	NM	2.67	25.33	3.04	24.96	NM	NM	NM	NM	NM	NM	NM	NM	2.97	25.03
MW-19	Lower AF	31.25	15.97	15.28	17.65	13.60	16.98	14.27	NM	NM	18.61	12.64	18.69	12.56	19.25	12.00	NM	NM
MW-20	USAS	30.29	NM	NM	4.88	25.41	NM	NM	NM	NM	5.76	24.53	NM	NM	5.87	24.42	NM	NM
MW-21	S&P Sand	28.88	NM	NM	16.63	12.25	15.24	13.64	NM	NM	NM	NM	NM	NM	NM	NM	17.95	10.93
MW-22	Lower AF	28.71	NM	NM	19.19	9.52	16.08	12.63	NM	NM	16.16	12.55	NM	NM	16.94	11.77	NM	NM
MW-23	S&P Sand	28.70	NM	NM	15.68	13.02	14.93	13.77	NM	NM	15.54	13.16	NM	NM	17.05	11.65	NM	NM
MW-24	USAS	30.01	NM	NM	NM	NM	5.74	24.27	NM	NM	5.87	24.14	NM	NM	5.82	24.19	NM	NM
MW-25	USAS	29.58	NM	NM	3.64	25.94	4.18	25.40	NM	NM	4.16	25.42	NM	NM	4.29	25.29	NM	NM
MW-26	USAS	26.76	NM	NM	3.49	23.27	3.57	23.19	NM	NM	NM	NM	NM	NM	NM	NM	3.58	23.18
MW-27	USAS	27.06	2.78	24.28	2.63	24.43	2.82	24.24	NM	NM	NM	NM	NM	NM	NM	NM	2.90	24.16
MW-28	USAS	27.81	NM	NM	2.65	25.16	3.00	24.81	NM	NM	NM	NM	NM	NM	NM	NM	3.00	24.81
MW-29	USAS	27.73	NM	NM	2.61	25.12	2.96	24.77	NM	NM	NM	NM	NM	NM	NM	NM	2.87	24.86
MW-30	USAS	29.24	3.08	26.16	NM	NM	3.68	25.56	4.50	24.74	3.99	25.25	NM	NM	4.26	24.98	NM	NM
MW-31	Lower AF	28.49	NM	NM	15.24	13.25	14.35	14.14	NM	NM	NM	NM	NM	NM	NM	NM	16.75	11.74

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APPENDIX D-1
 HISTORICAL SUMMARY OF MONITORING WELL GROUNDWATER ELEVATIONS

2008 GROUNDWATER MONITORING REPORT
 FORMER AMERICAN BERYLLIUM COMPANY SITE
 TALLEVAST, FLORIDA

Well ID	Zone	Top of Inner Casing (ft-msl)	October 17-19 2006		December 28 2006		February 20 2007		April 23 2007		December 3-6 2007		December 7 2007		January 21-22 2008		January 25-31 2008	
			Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)
MW-32	USAS	31.00	3.04	27.96	4.39	26.61	4.75	26.25	5.59	25.41	4.67	26.33	4.68	26.32	4.82	26.18	NM	NM
MW-33	LSAS	31.00	8.42	22.58	11.16	19.84	11.81	19.19	13.13	17.87	9.27	21.73	9.37	21.63	10.40	20.60	NM	NM
MW-34	S&P Sand	29.99	NM	NM	NM	NM	NM	NM	NM	NM	17.09	12.90	NM	NM	18.64	11.35	NM	NM
MW-35	USAS	29.88	NM	NM	NM	NM	NM	NM	3.75	26.13	2.83	27.05	2.93	26.95	3.02	26.86	NM	NM
MW-36	USAS	31.71	4.00	27.71	6.33	25.38	7.18	24.53	7.74	23.97	5.58	26.13	5.64	26.07	5.89	25.82	NM	NM
MW-37	LSAS	31.60	9.10	22.50	11.91	19.69	12.48	19.12	13.58	18.02	9.90	21.70	10.00	21.60	11.09	20.51	NM	NM
MW-38	USAS	31.15	3.39	27.76	NM	NM	5.51	25.64	6.40	24.75	4.99	26.16	5.06	26.09	5.29	25.86	NM	NM
MW-39	LSAS	31.18	7.75	23.43	NM	NM	10.83	20.35	12.94	18.24	8.83	22.35	8.94	22.24	9.96	21.22	NM	NM
MW-40	USAS	31.32	3.55	27.77	4.82	26.50	5.17	26.15	6.03	25.29	5.17	26.15	5.24	26.08	5.46	25.86	NM	NM
MW-41	LSAS	31.22	8.70	22.52	11.42	19.80	12.11	19.11	13.35	17.87	9.49	21.73	9.60	21.62	9.67	21.55	NM	NM
MW-42	USAS	31.49	3.82	27.67	5.28	26.21	5.85	25.64	6.67	24.82	5.43	26.06	5.50	25.99	5.74	25.75	NM	NM
MW-43	LSAS	31.48	8.97	22.51	11.69	19.79	12.21	19.27	13.62	17.86	9.76	21.72	9.91	21.57	10.96	20.52	NM	NM
MW-44	S&P Sand	30.88	15.00	15.88	18.00	12.88	16.84	14.04	NM	NM	NM	NM	NM	NM	NM	19.13	11.75	
MW-45	S&P Sand	30.58	14.72	15.86	17.85	12.73	16.40	14.18	NM	NM	17.89	12.69	NM	NM	19.45	11.13	NM	NM
MW-46	Lower AF	27.33	NM	NM	13.51	13.82	12.70	14.63	NM	NM	14.58	12.75	NM	NM	15.23	12.10	NM	NM
MW-47	USAS	29.42	NM	NM	4.06	25.36	4.34	25.08	NM	NM	4.68	24.74	NM	NM	5.06	24.36	NM	NM
MW-48	LSAS	30.40	8.83	21.57	12.00	18.40	11.76	18.64	12.94	17.46	NM	NM	NM	11.09	19.31	NM	NM	
MW-49	S&P Sand	29.37	NM	NM	16.84	12.53	16.03	13.34	NM	NM	17.57	11.80	NM	NM	18.41	10.96	NM	NM
MW-50	Lower AF	27.56	NM	NM	14.06	13.50	13.35	14.21	NM	NM	NM	NM	NM	NM	NM	NM	15.71	11.85
MW-51	Lower AF	26.89	NM	NM	13.28	13.61	12.23	14.66	NM	NM	NM	NM	NM	NM	NM	NM	14.90	11.99
MW-52	S&P Sand	27.11	NM	NM	14.19	12.92	13.58	13.53	NM	NM	NM	NM	NM	NM	NM	NM	15.62	11.49
MW-53	S&P Sand	27.77	NM	NM	14.82	12.95	14.01	13.76	NM	NM	NM	NM	NM	NM	NM	NM	16.24	11.53
MW-54	S&P Sand	26.88	NM	NM	13.86	13.02	13.19	13.69	NM	NM	NM	NM	NM	NM	NM	NM	15.30	11.58
MW-55	AF Gravel	30.03	NM	NM	NM	NM	18.25	11.78	NM	NM	16.58	13.45	NM	NM	17.42	12.61	NM	NM
MW-56	S&P Sand	27.28	11.57	15.71	14.84	12.44	13.63	13.65	NM	NM	14.60	12.68	NM	NM	16.18	11.10	NM	NM
MW-57	S&P Sand	30.35	14.53	15.82	17.55	12.80	16.29	14.06	NM	NM	17.34	13.01	NM	NM	19.29	11.06	NM	NM
MW-58	S&P Sand	31.26	15.45	15.81	18.41	12.85	16.90	14.36	NM	NM	18.68	12.58	NM	NM	19.95	11.31	NM	NM
MW-59	S&P Sand	28.48	NM	NM	17.04	11.44	13.91	14.57	NM	NM	NM	NM	NM	NM	NM	NM	16.89	11.59
MW-60	S&P Sand	28.33	NM	NM	17.59	10.74	14.88	13.45	NM	NM	NM	NM	NM	NM	NM	NM	17.11	11.22
MW-61	S&P Sand	27.50	NM	NM	20.12	7.38	16.44	11.06	NM	NM	17.10	10.40	NM	NM	17.45	10.05	NM	NM
MW-62	USAS	27.35	NM	NM	3.04	24.31	3.24	24.11	NM	NM	3.42	23.93	NM	NM	3.75	23.60	NM	NM
MW-63	USAS	27.37	2.88	24.49	2.70	24.67	3.03	24.34	NM	NM	3.33	24.04	NM	NM	3.54	23.83	NM	NM
MW-64	USAS	27.38	NM	NM	2.61	24.77	2.91	24.47	NM	NM	NM	NM	NM	NM	NM	NM	2.76	24.62
MW-65	USAS	28.76	NM	NM	NM	NM	3.65	25.11	NM	NM	NM	NM	NM	NM	NM	NM	3.35	25.41
MW-66	USAS	29.20	NM	NM	3.02	26.18	3.63	25.57	NM	NM	3.71	25.49	NM	NM	4.10	25.10	NM	NM
MW-67	USAS	30.79	3.70	27.09	4.56	26.23	5.10	25.69	NM	NM	5.36	25.43	NM	NM	NM	NM	5.11	25.68
MW-68	LSAS	28.60	NM	NM	11.61	16.99	11.19	17.41	NM	NM	NM	NM	NM	NM	NM	NM	10.90	17.70
MW-69	USAS	26.91	NM	NM	4.86	22.05	4.33	22.58	NM	NM	5.38	21.53	NM	NM	4.67	22.24	NM	NM
MW-70	USAS	31.89	4.03	27.86	5.19	26.70	5.60	26.29	6.37	25.52	5.78	26.11	5.85	26.04	6.16	25.73	NM	NM
MW-71	USAS	31.23	3.69	27.54	4.94	26.29	7.27	23.96	6.45	24.78	5.25	25.98	5.32	25.91	5.60	25.63	NM	NM
MW-72	USAS	30.97	3.13	27.84	3.61	27.36	4.03	26.94	4.63	26.34	4.17	26.80	4.26	26.71	4.39	26.58	NM	NM
MW-73	USAS	26.03	2.58	23.45	NM	NM	3.26	22.77	NM	NM	3.04	22.99	NM	NM	3.20	22.83	NM	NM
MW-74	USAS	27.90	1.84	26.06	2.38	25.52	2.81	25.09	NM	NM	2.67	25.23	NM	NM	2.82	25.08	NM	NM
MW-75	USAS	31.38	NM	NM	NM	NM	NM	NM	NM	NM	5.94	25.44	NM	NM	5.76	25.62	NM	NM
MW-76	USAS	30.84	3.30	27.54	4.40	26.44	5.29	25.55	5.69	25.15	4.82	26.02	4.90	25.94	5.13	25.71	NM	NM
MW-77	LSAS	29.73	4.06	25.67	4.55	25.18	5.03	24.70	6.00	23.73	5.27	24.46	5.33	24.40	6.19	23.54	NM	NM
MW-78	LSAS	30.23	5.06	25.17	6.80	23.43	7.43	22.80	8.25	21.98	NM	NM	NM	NM	NM	NM	8.02	22.21
MW-79	LSAS	30.11	7.54	22.57	10.62	19.49	11.02	19.09	12.15	17.96	8.29	21.82	8.35	21.76	9.26	20.85	NM	NM
MW-80	LSAS	30.99	8.01	22.98	9.84	21.15	10.17	20.82	11.35	19.64	8.65	22.34	8.83	22.16	9.41	21.58	NM	NM
MW-81	LSAS	31.01	9.69	21.32	12.61	18.40	12.77	18.24	14.39	16.62	10.51	20.50	10.52	20.49	11.71	19.30	NM	NM
MW-82	LSAS	27.24	2.52	24.72	3.97	23.27	4.51	22.73	NM	NM	3.85	23.39	NM	NM	4.70	22.54	NM	NM
MW-83	AF Gravel	25.51	10.42	15.09	NM	NM	12.30	13.21	NM	NM	12.57	12.94	NM	NM	13.81	11.70	NM	NM
MW-84	LSAS	31.15	8.21	22.94	9.93	21.22	10.14	21.01	7.31	23.84	8.90	22.25	9.07	22.08	9.58	21.57	NM	NM

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**APPENDIX D-1
HISTORICAL SUMMARY OF MONITORING WELL GROUNDWATER ELEVATIONS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Well ID	Zone	Top of Inner Casing (ft-msl)	October 17-19 2006		December 28 2006		February 20 2007		April 23 2007		December 3-6 2007		December 7 2007		January 21-22 2008		January 25-31 2008	
			Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation
MW-85	LSAS	29.55	4.36	25.19	5.51	24.04	6.83	22.72	NM	NM	5.35	24.20	NM	NM	5.58	23.97	NM	NM
MW-86	LSAS	28.77	9.45	19.32	11.52	17.25	11.11	17.66	NM	NM	NM	NM	NM	NM	NM	NM	11.25	17.52
MW-87	LSAS	30.26	NM	NM	NM	NM	NM	NM	5.90	24.36	4.65	25.61	4.77	25.49	4.56	25.70	NM	NM
MW-88	Zone 1 Clay	27.28	8.33	18.95	12.35	14.93	10.95	16.33	NM	NM	11.41	15.87	NM	NM	13.50	13.78	NM	NM
MW-89	USAS	29.50	NM	NM	3.41	26.09	3.84	25.66	NM	NM	4.00	25.50	NM	NM	4.43	25.07	NM	NM
MW-90	USAS	27.95	NM	NM	2.74	25.21	3.02	24.93	NM	NM	3.25	24.70	NM	NM	3.66	24.29	NM	NM
MW-91	LSAS	27.66	7.29	20.37	9.01	18.65	9.21	18.45	NM	NM	NM	NM	NM	NM	NM	NM	7.95	19.71
MW-92	LSAS	27.35	9.83	17.52	11.73	15.62	11.20	16.15	NM	NM	11.40	15.95	NM	NM	12.54	14.81	NM	NM
MW-93	LSAS	27.73	10.36	17.37	11.80	15.93	11.35	16.38	NM	NM	11.62	16.11	NM	NM	NM	NM	12.54	15.19
MW-94	USAS	25.40	2.70	22.70	NM	NM	2.50	22.90	NM	NM	3.11	22.29	NM	NM	3.09	22.31	NM	NM
MW-95	USAS	24.85	NM	NM	NM	NM	2.19	22.66	NM	NM	2.91	21.94	NM	NM	2.92	21.93	NM	NM
MW-96	Zone 3-4 Clay	25.14	NM	NM	12.81	12.33	12.29	12.85	NM	NM	13.47	11.67	NM	NM	14.14	11.00	NM	NM
MW-97	Zone 3-4 Clay	25.29	NM	NM	13.05	12.24	12.85	12.44	NM	NM	14.11	11.18	NM	NM	14.75	10.54	NM	NM
MW-98	LSAS	25.75	2.99	22.76	NM	NM	3.74	22.01	NM	NM	3.61	22.14	NM	NM	3.83	21.92	NM	NM
MW-99	Zone 1 Clay	25.26	NM	NM	10.07	15.19	9.26	16.00	NM	NM	9.03	16.23	NM	NM	11.60	13.66	NM	NM
MW-100	USAS	25.23	NM	NM	5.04	20.19	5.15	20.08	NM	NM	3.56	21.67	NM	NM	5.13	20.10	NM	NM
MW-101	LSAS	30.09	NM	NM	NM	NM	10.15	19.94	NM	NM	7.04	23.05	NM	NM	7.33	22.76	NM	NM
MW-102	AF Gravel	26.10	NM	NM	12.50	13.60	10.67	15.43	NM	NM	12.57	13.53	NM	NM	13.31	12.79	NM	NM
MW-103	USAS	26.19	NM	NM	4.07	22.12	4.45	21.74	NM	NM	4.34	21.85	NM	NM	4.61	21.58	NM	NM
MW-104	USAS	26.39	NM	NM	2.42	23.97	2.62	23.77	NM	NM	NM	NM	NM	NM	NM	NM	2.74	23.65
MW-105	LSAS	26.41	3.75	22.66	5.41	21.00	6.81	19.60	NM	NM	NM	NM	NM	NM	NM	NM	4.93	21.48
MW-106	LSAS	28.31	5.90	22.41	6.72	21.59	7.60	20.71	NM	NM	6.36	21.95	NM	NM	7.04	21.27	NM	NM
MW-107	USAS	24.72	NM	NM	NM	NM	2.99	21.73	NM	NM	3.62	21.10	NM	NM	3.75	20.97	NM	NM
MW-108	USAS	24.36	NM	NM	NM	NM	3.39	20.97	NM	NM	3.10	21.26	NM	NM	3.68	20.68	NM	NM
MW-109	USAS	28.45	NM	NM	4.52	23.93	4.40	24.05	NM	NM	4.96	23.49	NM	NM	NM	NM	4.45	24.00
MW-110	USAS	28.80	NM	NM	3.95	24.85	3.88	24.92	NM	NM	NM	NM	NM	NM	NM	NM	3.80	25.00
MW-111	USAS	26.37	NM	NM	2.61	23.76	2.88	23.49	NM	NM	NM	NM	NM	NM	NM	NM	2.82	23.55
MW-112	Zone 1 Clay	28.28	NM	NM	19.46	8.82	14.68	13.60	NM	NM	13.70	14.58	NM	NM	12.78	15.50	NM	NM
MW-113	LSAS	26.31	NM	NM	10.62	15.69	10.08	16.23	NM	NM	NM	NM	NM	NM	NM	NM	11.52	14.79
MW-114	USAS	24.79	2.32	22.47	NM	NM	2.63	22.16	NM	NM	3.00	21.79	NM	NM	3.11	21.68	NM	NM
MW-115	USAS	30.21	NM	NM	6.74	23.47	7.01	23.20	NM	NM	7.02	23.19	NM	NM	7.01	23.20	NM	NM
MW-116	USAS	21.84	NM	NM	NM	NM	2.02	19.82	NM	NM	2.70	19.14	NM	NM	2.19	19.65	NM	NM
MW-117	LSAS	21.56	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	6.60	14.96
MW-118	USAS	21.47	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	5.95	15.52
MW-119	LSAS	21.18	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	9.33	11.85
MW-120	USAS	21.18	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	2.85	18.33
MW-121	USAS	21.35	NM	NM	7.92	13.43	8.06	13.29	NM	NM	7.93	13.42	NM	NM	8.02	13.33	NM	NM
MW-122	USAS	20.06	NM	NM	6.49	13.57	6.36	13.70	NM	NM	6.12	13.94	NM	NM	6.53	13.53	NM	NM
MW-123	Floridan	30.90	NM	NM	16.19	14.71	13.55	17.35	18.80	12.10	16.84	14.06	17.37	13.53	NM	NM	18.38	12.52
MW-124	Zone 2 Clay	28.97	NM	NM	15.62	13.35	35.55	-6.58	NM	NM	16.44	12.53	NM	NM	13.42	15.55	NM	NM
MW-125	Venice Clay	29.52	NM	NM	6.93	22.59	7.90	21.62	NM	NM	8.21	21.31	NM	NM	7.93	21.59	NM	NM
MW-126	USAS	28.32	NM	NM	7.50	20.82	8.01	20.31	NM	NM	6.95	21.37	NM	NM	7.48	20.84	NM	NM
MW-127	AF Gravel	31.74	14.02	17.72	16.39	15.35	16.08	15.66	18.40	13.34	16.21	15.53	16.02	15.72	16.26	15.48	NM	NM
MW-128	S&P Sand	31.59	15.59	16.00	18.63	12.96	17.31	14.28	20.02	11.57	18.70	12.89	18.80	12.79	20.08	11.51	NM	NM
MW-129	AF Gravel	31.41	14.22	17.19	16.45	14.96	15.55	15.86	NM	NM	16.25	15.16	13.50	17.91	17.09	14.32	NM	NM
MW-130	AF Gravel	30.37	13.30	17.07	15.20	15.17	14.30	16.07	NM	NM	14.87	15.50	14.92	15.45	15.92	14.45	NM	NM
MW-131	AF Gravel	27.33	11.90	15.43	14.15	13.18	13.83	13.50	NM	NM	14.35	12.98	NM	NM	15.56	11.77	NM	NM
MW-132	AF Gravel	30.07	13.25	16.82	15.61	14.46	15.31	14.76	NM	NM	15.53	14.54	15.73	14.34	16.26	13.81	NM	NM
MW-133	AF Gravel	27.68	11.92	15.76	14.22	13.46	14.02	13.66	NM	NM	NM	NM	NM	NM	NM	NM	15.11	12.57
MW-134	AF Gravel	31.10	13.91	17.19	16.28	14.82	15.96	15.14	NM	NM	15.95	15.15	16.35	14.75	16.83	14.27	NM	NM
MW-135	AF Gravel	27.64	12.22	15.42	NM	NM	14.25	13.39	NM	NM	14.67	12.97	NM	NM	NM	NM	15.58	12.06
MW-136	AF Gravel	25.23	NM	NM	NM	NM	NM	NM	NM	NM	14.01	11.22	NM	NM	15.01	10.22	NM	NM
MW-137	USAS	25.49	NM	NM	2.31	23.18	2.74	22.75	NM	NM	2.80	22.69	NM	NM	2.95	22.54	NM	NM

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APPENDIX D-1
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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Well ID	Zone	Top of Inner Casing (ft-msl)	October 17-19 2006		December 28 2006		February 20 2007		April 23 2007		December 3-6 2007		December 7 2007		January 21-22 2008		January 25-31 2008	
			Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation
MW-138	LSAS	25.49	NM	NM	8.95	16.54	8.99	16.50	NM	NM	9.70	15.79	NM	NM	10.01	15.48	NM	NM
MW-139	S&P Sand	25.46	NM	NM	14.78	10.68	13.11	12.35	NM	NM	14.35	11.11	NM	NM	15.26	10.20	NM	NM
MW-140	Lower AF	25.52	NM	NM	12.23	13.29	11.89	13.63	NM	NM	13.40	12.12	NM	NM	13.98	11.54	NM	NM
MW-141	USAS	25.55	NM	NM	2.09	23.46	2.48	23.07	NM	NM	2.82	22.73	NM	NM	2.75	22.80	NM	NM
MW-142	LSAS	25.64	2.30	23.34	2.14	23.50	2.57	23.07	NM	NM	2.88	22.76	NM	NM	3.85	21.79	NM	NM
MW-143	AF Gravel	25.60	10.56	15.04	12.63	12.97	11.91	13.69	NM	NM	12.90	12.70	NM	NM	14.00	11.60	NM	NM
MW-144	S&P Sand	25.51	11.31	14.20	13.40	12.11	12.81	12.70	NM	NM	14.00	11.51	NM	NM	14.97	10.54	NM	NM
MW-145	Lower AF	25.45	NM	NM	14.11	11.34	11.09	14.36	NM	NM	12.50	12.95	NM	NM	13.21	12.24	NM	NM
MW-146	USAS	26.06	2.70	23.36	12.32	13.74	2.61	23.45	NM	NM	3.35	22.71	NM	NM	3.19	22.87	NM	NM
MW-147	LSAS	25.94	2.54	23.40	12.78	13.16	2.55	23.39	NM	NM	3.21	22.73	NM	NM	3.11	22.83	NM	NM
MW-148	AF Gravel	25.90	10.44	15.46	13.91	11.99	12.18	13.72	NM	NM	13.36	12.54	NM	NM	14.15	11.75	NM	NM
MW-149	S&P Sand	26.20	12.21	13.99	2.90	23.30	13.02	13.18	NM	NM	14.33	11.87	NM	NM	15.05	11.15	NM	NM
MW-150	Lower AF	26.09	10.44	15.65	1.32	24.77	11.40	14.69	NM	NM	13.40	12.69	NM	NM	14.10	11.99	NM	NM
MW-151	USAS	22.44	3.90	18.54	3.97	18.47	4.31	18.13	NM	NM	4.03	18.41	NM	NM	4.36	18.08	NM	NM
MW-152	LSAS	22.21	4.14	18.07	4.98	17.23	4.96	17.25	NM	NM	4.97	17.24	NM	NM	5.41	16.80	NM	NM
MW-153	AF Gravel	22.50	NM	NM	9.41	13.09	8.25	14.25	NM	NM	8.35	14.15	NM	NM	10.54	11.96	NM	NM
MW-154	S&P Sand	22.25	NM	NM	10.57	11.68	8.53	13.72	NM	NM	11.22	11.03	NM	NM	11.55	10.70	NM	NM
MW-155	Lower AF	22.34	NM	NM	9.25	13.09	8.81	13.53	NM	NM	10.24	12.10	NM	NM	10.87	11.47	NM	NM
MW-156	USAS	24.81	NM	NM	2.51	22.30	2.55	22.26	NM	NM	3.47	21.34	NM	NM	2.44	22.37	NM	NM
MW-157	LSAS	24.74	NM	NM	8.26	16.48	8.25	16.49	NM	NM	10.02	14.72	NM	NM	9.69	15.05	NM	NM
MW-158	AF Gravel	24.78	NM	NM	11.62	13.16	11.40	13.38	NM	NM	11.81	12.97	NM	NM	13.08	11.70	NM	NM
MW-159	S&P Sand	24.68	NM	NM	NM	NM	11.44	13.24	NM	NM	12.59	12.09	NM	NM	13.45	11.23	NM	NM
MW-160	Lower AF	24.72	NM	NM	10.81	13.91	17.90	6.82	NM	NM	11.72	13.00	NM	NM	12.68	12.04	NM	NM
MW-161	Floridan	24.91	NM	NM	10.54	14.37	7.77	17.14	13.90	11.01	11.42	13.49	NM	NM	12.56	12.35	NM	NM
MW-162	USAS	25.38	NM	NM	2.01	23.37	2.41	22.97	NM	NM	2.55	22.83	NM	NM	2.75	22.63	NM	NM
MW-163	LSAS	25.60	8.76	16.84	8.88	16.72	9.01	16.59	NM	NM	9.70	15.90	NM	NM	9.90	15.70	NM	NM
MW-164	AF Gravel	25.59	12.22	13.37	14.27	11.32	13.81	11.78	NM	NM	14.40	11.19	NM	NM	15.74	9.85	NM	NM
MW-165	S&P Sand	25.35	17.66	7.69	14.52	10.83	14.02	11.33	NM	NM	14.54	10.81	NM	NM	15.13	10.22	NM	NM
MW-166	Lower AF	25.69	NM	NM	12.25	13.44	11.66	14.03	NM	NM	13.45	12.24	NM	NM	14.02	11.67	NM	NM
MW-167	USAS	27.05	3.48	23.57	3.74	23.31	3.38	23.67	NM	NM	3.54	23.51	NM	NM	3.73	23.32	NM	NM
MW-168	LSAS	27.41	5.21	22.20	5.77	21.64	7.73	19.68	NM	NM	6.25	21.16	NM	NM	6.55	20.86	NM	NM
MW-169	AF Gravel	27.48	13.97	13.51	16.48	11.00	15.35	12.13	NM	NM	14.12	13.36	NM	NM	12.68	14.80	NM	NM
MW-170	Lower AF	27.50	14.02	13.48	17.11	10.39	14.52	12.98	NM	NM	14.91	12.59	NM	NM	16.52	10.98	NM	NM
MW-171	LSAS	21.49	NM	NM	10.27	11.22	9.50	11.99	NM	NM	9.90	11.59	NM	NM	8.22	13.27	NM	NM
MW-172	AF Gravel	21.53	NM	NM	11.59	9.94	11.08	10.45	NM	NM	11.25	10.28	NM	NM	12.33	9.20	NM	NM
MW-173	S&P Sand	21.42	NM	NM	11.87	9.55	10.83	10.59	NM	NM	12.39	9.03	NM	NM	12.96	8.46	NM	NM
MW-174	Lower AF	21.39	NM	NM	13.33	8.06	9.22	12.17	NM	NM	10.69	10.70	NM	NM	10.77	10.62	NM	NM
MW-175	AF Gravel	27.80	NM	NM	18.24	9.56	16.93	10.87	NM	NM	21.39	6.41	NM	NM	18.84	8.96	NM	NM
MW-176	S&P Sand	29.01	NM	NM	19.79	9.22	18.19	10.82	NM	NM	19.55	9.46	NM	NM	21.11	7.90	NM	NM
MW-177	Lower AF	29.28	NM	NM	16.50	12.78	15.57	13.71	NM	NM	17.29	11.99	NM	NM	18.27	11.01	NM	NM
MW-178	LSAS	21.82	NM	NM	13.03	8.79	12.08	9.74	NM	NM	13.30	8.52	NM	NM	12.80	9.02	NM	NM
MW-179	AF Gravel	21.87	NM	NM	12.84	9.03	11.82	10.05	NM	NM	13.87	8.00	NM	NM	13.54	8.33	NM	NM
MW-180	S&P Sand	21.97	NM	NM	15.13	6.84	14.05	7.92	NM	NM	15.08	6.89	NM	NM	15.79	6.18	NM	NM
MW-181	Lower AF	22.09	NM	NM	12.52	9.57	11.83	10.26	NM	NM	12.31	9.78	NM	NM	12.31	9.78	NM	NM
MW-182	S&P Sand	27.19	11.24	15.95	16.38	10.81	30.89	-3.70	NM	NM	14.65	12.54	NM	NM	14.36	12.83	NM	NM
MW-183	USAS	24.20	3.79	20.41	2.98	21.22	2.97	21.23	NM	NM	3.86	20.34	NM	NM	3.52	20.68	3.54	20.66
MW-184	LSAS	24.22	8.15	16.07	8.61	15.61	8.54	15.68	NM	NM	9.24	14.98	NM	NM	9.75	14.47	10.72	13.50
MW-185	AF Gravel	24.30	9.50	14.80	11.67	12.63	10.71	13.59	NM	NM	11.95	12.35	NM	NM	13.25	11.05	13.07	11.23
MW-186	S&P Sand	24.37	10.09	14.28	11.96	12.41	11.24	13.13	NM	NM	12.64	11.73	NM	NM	13.80	10.57	13.50	10.87
MW-187	Lower AF	24.34	8.88	15.46	10.10	14.24	9.39	14.95	NM	NM	11.41	12.93	NM	NM	12.00	12.34	NM	NM
MW-188	USAS	21.97	3.68	18.29	3.41	18.56	3.52	18.45	NM	NM	4.30	17.67	NM	NM	4.21	17.76	NM	NM
MW-189	LSAS	22.21	6.04	16.17	6.96	15.25	6.75	15.46	NM	NM	7.48	14.73	NM	NM	8.08	14.13	NM	NM
MW-190	AF Gravel	22.11	8.02	14.09	11.32	10.79	8.90	13.21	NM	NM	10.74	11.37	NM	NM	12.42	9.69	NM	NM

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APPENDIX D-1
HISTORICAL SUMMARY OF MONITORING WELL GROUNDWATER ELEVATIONS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Well ID	Zone	Top of Inner Casing (ft-msl)	October 17-19 2006		December 28 2006		February 20 2007		April 23 2007		December 3-6 2007		December 7 2007		January 21-22 2008		January 25-31 2008	
			Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation	Depth To Water (ft-toc)	Water Elevation
MW-191	S&P Sand	21.97	7.99	13.98	10.03	11.94	9.06	12.91	NM	NM	10.62	11.35	NM	NM	12.09	9.88	NM	NM
MW-192	Lower AF	21.85	8.33	13.52	13.12	8.73	10.68	11.17	NM	NM	8.69	13.16	NM	NM	8.52	13.33	NM	NM
MW-193	AF Gravel	21.77	NM	NM	10.44	11.33	8.57	13.20	NM	NM	12.78	8.99	NM	NM	11.93	9.84	12.39	9.38
MW-194	S&P Sand	21.70	NM	NM	9.52	12.18	8.12	13.58	NM	NM	10.06	11.64	NM	NM	11.15	10.55	12.32	9.38
MW-195	Lower AF	22.08	NM	NM	9.54	12.54	8.38	13.70	NM	NM	10.22	11.86	NM	NM	10.20	11.88	10.27	11.81
MW-196	AF Gravel	26.67	NM	NM	13.08	13.59	14.84	11.83	NM	NM	17.30	9.37	NM	NM	17.29	9.38	NM	NM
MW-197	AF Gravel	28.99	NM	NM	19.86	9.13	17.83	11.16	NM	NM	20.67	8.32	NM	NM	21.39	7.60	NM	NM
MW-198	USAS	20.55	NM	NM	3.98	16.57	3.19	17.36	NM	NM	3.46	17.09	NM	NM	3.27	17.28	NM	NM
MW-199	LSAS	20.42	NM	NM	7.82	12.60	6.88	13.54	NM	NM	8.23	12.19	NM	NM	9.26	11.16	NM	NM
MW-200	AF Gravel	20.62	NM	NM	8.51	12.11	7.13	13.49	NM	NM	8.94	11.68	NM	NM	10.10	10.52	NM	NM
MW-201	S&P Sand	20.54	NM	NM	8.55	11.99	7.43	13.11	NM	NM	8.88	11.66	NM	NM	10.20	10.34	NM	NM
MW-202	Lower AF	20.62	NM	NM	7.05	13.57	5.74	14.88	NM	NM	8.17	12.45	NM	NM	9.02	11.60	NM	NM
MW-204	USAS	21.14	4.03	17.11	4.03	17.11	3.83	17.31	NM	NM	4.07	17.07	NM	NM	4.27	16.87	NM	NM
MW-205	LSAS	21.21	7.00	14.21	8.43	12.78	7.27	13.94	NM	NM	8.49	12.72	NM	NM	9.39	11.82	NM	NM
MW-206	AF Gravel	21.24	8.04	13.20	9.92	11.32	7.83	13.41	NM	NM	10.94	10.30	NM	NM	11.41	9.83	NM	NM
MW-207	Lower AF	21.57	6.95	14.62	3.21	18.36	7.23	14.34	NM	NM	8.10	13.47	NM	NM	9.74	11.83	NM	NM
MW-208	USAS	15.43	NM	NM	3.03	12.40	2.88	12.55	NM	NM	3.75	11.68	NM	NM	3.60	11.83	NM	NM
MW-209	LSAS	15.24	NM	NM	5.70	9.54	4.15	11.09	NM	NM	6.80	8.44	NM	NM	7.15	8.09	NM	NM
MW-210	AF Gravel	15.52	NM	NM	6.11	9.41	3.20	12.32	NM	NM	6.34	9.18	NM	NM	8.06	7.46	NM	NM
MW-211	S&P Sand	15.39	NM	NM	5.81	9.58	2.83	12.56	NM	NM	4.29	11.10	NM	NM	7.73	7.66	NM	NM
MW-212	Lower AF	15.66	NM	NM	9.59	5.97	4.51	11.05	NM	NM	6.75	8.81	NM	NM	5.39	10.17	NM	NM
MW-213	USAS	25.28	NM	NM	NM	NM	2.06	23.22	NM	NM	NM	NM	NM	NM	NM	NM	2.27	23.01
MW-214	LSAS	25.19	NM	NM	NM	NM	8.91	16.28	NM	NM	NM	NM	NM	NM	NM	NM	9.92	15.27
MW-215	AF Gravel	25.16	NM	NM	NM	NM	11.57	13.59	NM	NM	NM	NM	NM	NM	NM	NM	13.48	11.68
MW-216	S&P Sand	25.20	NM	NM	NM	NM	12.57	12.63	NM	NM	NM	NM	NM	NM	NM	NM	14.64	10.56
MW-217	Lower AF	25.14	NM	NM	NM	NM	10.82	14.32	NM	NM	NM	NM	NM	NM	NM	NM	13.04	12.10
MW-218	Floridan	26.03	NM	NM	NM	NM	NM	7.59	18.44	NM	NM	NM	NM	NM	NM	NM	NM	NM
MW-219	USAS	21.91	NM	NM	3.83	18.08	4.15	17.76	NM	NM	3.34	18.57	NM	NM	4.00	17.91	NM	NM
MW-220	LSAS	22.04	NM	NM	10.71	11.33	5.35	16.69	NM	NM	6.83	15.21	NM	NM	5.20	16.84	NM	NM
MW-221	AF Gravel	22.24	NM	NM	10.03	12.21	7.71	14.53	NM	NM	7.60	14.64	NM	NM	9.29	12.95	NM	NM
MW-222	S&P Sand	22.23	NM	NM	13.28	8.95	7.71	14.52	NM	NM	9.02	13.21	NM	NM	10.61	11.62	NM	NM
MW-223	Hardstreak Clay	17.11	NM	NM	5.30	11.81	5.34	11.77	NM	NM	5.96	11.15	NM	NM	5.63	11.48	NM	NM
MW-224	Venice Clay	17.22	NM	NM	5.35	11.87	5.44	11.78	NM	NM	5.70	11.52	NM	NM	5.92	11.30	NM	NM
MW-225	Venice Clay	17.13	NM	NM	4.45	12.68	4.46	12.67	NM	NM	5.22	11.91	NM	NM	5.48	11.65	NM	NM
MW-226	AF Gravel	17.14	NM	NM	3.28	13.86	1.45	15.69	NM	NM	5.92	11.22	NM	NM	5.42	11.72	NM	NM
MW-227	S&P Sand	17.40	NM	NM	10.98	6.42	4.10	13.30	NM	NM	4.67	12.73	NM	NM	7.19	10.21	NM	NM
MW-228	AF Gravel	20.73	NM	NM	NM	NM	NM	NM	NM	NM	11.13	9.60	NM	NM	13.13	7.60	NM	NM
MW-229	USAS	30.14	NM	NM	6.35	23.79	6.72	23.42	NM	NM	6.69	23.45	NM	NM	6.87	23.27	NM	NM
MW-230	LSAS	30.02	NM	NM	9.20	20.82	7.05	22.97	NM	NM	6.90	23.12	NM	NM	7.21	22.81	NM	NM
MW-231	AF Gravel	29.97	NM	NM	19.65	10.32	18.41	11.56	NM	NM	20.77	9.20	NM	NM	20.66	9.31	NM	NM
MW-232	AF Gravel	29.51	11.92	17.59	14.00	15.51	13.42	16.09	NM	NM	13.62	15.89	NM	NM	14.24	15.27	NM	NM
MW-233	AF Gravel	30.49	NM	NM	15.35	15.14	14.53	15.96	NM	NM	15.29	15.20	NM	NM	16.21	14.28	NM	NM
MW-234	USAS	24.68	NM	NM	NM	NM	2.73	21.95	NM	NM	NM	NM	NM	NM	NM	NM	2.57	22.11
MW-235	LSAS	24.71	NM	NM	NM	NM	9.13	15.58	NM	NM	NM	NM	NM	NM	NM	NM	10.33	14.38
MW-236	AF Gravel	24.77	NM	NM	NM	NM	11.95	12.82	NM	NM	NM	NM	NM	NM	NM	NM	15.56	9.21
MW-237	S&P Sand	24.76	NM	NM	NM	NM	12.15	12.61	NM	NM	NM	NM	NM	NM	NM	NM	14.48	10.28
MW-238	Lower AF	24.54	NM	NM	NM	NM	11.21	13.33	NM	NM	NM	NM	NM	NM	NM	NM	12.86	11.68
MW-239	AF Gravel	28.43	NM	NM	14.75	13.68	14.36	14.07	NM	NM	15.43	13.00	NM	NM	15.69	12.74	NM	NM
MW-240	S&P Sand	27.58	NM	NM	20.83	6.75	30.34	-2.76	NM	NM	21.41	6.17	NM	NM	14.83	12.75	NM	NM
MW-241	Lower AF	17.28	NM	NM	5.14	12.14	4.80	12.48	NM	NM	4.95	12.33	NM	NM	5.96	11.32	NM	NM
MW-242	USAS	22.60	NM	NM	3.19	19.41	3.17	19.43	NM	NM	3.34	19.26	NM	NM	3.58	19.02	NM	NM
MW-243	LSAS	22.62	NM	NM	11.98	10.64	11.36	11.26	NM	NM	11.55	11.07	NM	NM	12.14	10.48	NM	NM
MW-244	AF Gravel	22.66	NM	NM	13.19	9.47	12.42	10.24	NM	NM	12.45	10.21	NM	NM	13.08	9.58	NM	NM

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			Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)	Depth To Water (ft-toc)	Water Elevation (ft-msl)
MW-245	Hardstreak Clay	18.92	NM	NM	4.67	14.25	4.76	14.16	NM	NM	4.67	14.25	NM	NM	5.35	13.57	NM	NM
MW-246	LSAS	18.96	NM	NM	13.05	5.91	11.42	7.54	NM	NM	8.89	10.07	NM	NM	9.49	9.47	NM	NM
MW-247	AF Gravel	19.01	NM	NM	12.07	6.94	11.67	7.34	NM	NM	13.30	5.71	NM	NM	12.80	6.21	NM	NM
MW-248	AF Gravel	26.57	NM	NM	13.20	13.37	13.04	13.53	NM	NM	NM	NM	NM	NM	NM	NM	14.23	12.34
MW-249	AF Gravel	22.60	NM	NM	NM	NM	9.08	13.52	NM	NM	NM	NM	NM	NM	NM	NM	11.26	11.34
MW-250	AF Gravel	24.83	NM	NM	NM	NM	11.17	13.66	NM	NM	NM	NM	NM	NM	NM	NM	12.67	12.16
MW-251	Floridan	27.37	NM	NM	NM	NM	NM	NM	15.12	12.25	13.34	14.03	NM	NM	14.28	13.09	NM	NM
MW-252	S&P Sand	31.56	NM	NM	NM	NM	NM	NM	NM	NM	18.81	12.75	NM	NM	20.17	11.39	NM	NM
MW-253	AF Gravel	31.48	NM	NM	NM	NM	NM	NM	NM	NM	15.93	15.55	NM	NM	16.42	15.06	NM	NM
MW-254 (MW-BT-1)	USAS	31.39	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	5.23	26.16	NM	NM
PZ-LSAS-1	LSAS	31.12	NM	NM	NM	NM	NM	NM	NM	NM	5.46	25.66	5.53	25.59	5.98	25.14	NM	NM
PZ-LSAS-2	LSAS	31.44	NM	NM	NM	NM	NM	NM	NM	NM	5.76	25.68	5.82	25.62	6.24	25.20	NM	NM
PZ-LSAS-3	LSAS	32.16	NM	NM	NM	NM	NM	NM	NM	NM	6.46	25.70	6.39	25.77	6.71	25.45	NM	NM
PZ-LSAS-4	LSAS	31.60	NM	NM	NM	NM	NM	NM	NM	NM	6.19	25.41	6.13	25.47	6.65	24.95	NM	NM
PZ-LSAS-5	LSAS	31.61	NM	NM	NM	NM	NM	NM	NM	NM	6.18	25.43	6.09	25.52	6.59	25.02	NM	NM
PZ-LSAS-6	LSAS	32.73	NM	NM	NM	NM	NM	NM	NM	NM	7.40	25.33	7.25	25.48	7.83	24.90	NM	NM
PZ-LSAS-7	LSAS	31.90	NM	NM	NM	NM	NM	NM	NM	NM	6.20	25.70	6.12	25.78	6.66	25.24	NM	NM
RW-1	USAS	30.68	NM	NM	5.33	25.35	5.71	24.97	NM	NM	5.49	25.19	5.53	25.15	5.76	24.92	NM	NM
RW-2	USAS	29.98	NM	NM	4.58	25.40	4.96	25.02	NM	NM	4.95	25.03	5.02	24.96	5.25	24.73	NM	NM
Staff Gauge-1	Unassigned	23.29	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	0.65	22.64	NM	NM
Staff Gauge-1 (old)	Unassigned	23.17	NM	NM	NM	NM	1.28	24.45	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
Staff Gauge-2	Unassigned	25.62	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	4.26	21.36	0.76	24.86	NM	NM
Staff Gauge-2 (old)	Unassigned	25.50	NM	NM	NM	NM	1.15	26.65	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
Staff Gauge-3	Unassigned	14.41	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	0.98	13.43	NM	NM
Staff Gauge-3 (old)	Unassigned	14.16	NM	NM	NM	NM	1.12	15.28	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
Staff Gauge-4	Unassigned	21.18	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	0.72	20.46	NM	NM
Staff Gauge-4 (old)	Unassigned	21.06	NM	NM	NM	NM	0.84	21.90	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
Staff Gauge-5	Unassigned	23.91	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	0.70	23.21	NM	NM
Staff Gauge-5 (old)	Unassigned	23.77	NM	NM	NM	NM	1.19	24.96	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
Staff Gauge-6 (old)	Unassigned	19.82	NM	NM	NM	NM	1.05	20.87	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
Staff Gauge-8	Unassigned	23.38	NM	NM	NM	NM	NM	NM	NM	NM	4.26	21.36	NM	NM	NM	NM	NM	NM
Staff Gauge-9	Unassigned	21.93	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	0.72	21.21
Stilling Well-1	Unassigned	30.83	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	4.46	26.37	NM	NM
Stilling Well-2	Unassigned	14.55	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	1.76	12.79	NM	NM
Stilling Well-3	Unassigned	26.04	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	2.88	20.32	NM	NM
Stilling Well-4	Unassigned	26.96	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	4.30	22.66
TW-84-A	USAS	32.10	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	6.85	25.25	NM	NM
TW-84-B	USAS	32.07	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	7.18	24.89	NM	NM

Notes:

AF Gravel = Arcadia Formation Gravels
ft msl = feet above mean sea level
ft toc = feet below top of casing
LSAS = Lower Shallow Aquifer System
Lower AF = Lower Arcadia Formation
NI = Not installed
NM = Not Measured
S&P Sand = Salt & Pepper sands
USAS = Upper Surficial Aquifer System

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	DW-1 AF Gravel 12/14/01	DW-1 AF Gravel 03/10/03	DW-1 AF Gravel 12/20/04	DW-1 AF Gravel 06/16/05	DW-1 AF Gravel 12/21/05	DW-1 AF Gravel 12/05/06	DW-1 AF Gravel 02/01/08	DW-1 AF Gravel 06/17/04
Semivolatile Organics								
1,4-Dioxane	ug/L	NA	NA	1 U	NA	NA	NA	NA
Semivolatile Organics (8270C)								
1,4-Dioxane	ug/L	NA	NA	NA	2.1 U	1 U	NA	NA
Volatile Organics (8260B) - SIM								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	1 U	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	1 U	1 U	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1 U	1 U	0.65 U	0.5 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	1 U	1 U	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	1 U	1.9	0.28 U	0.5 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	IWI-1 AF Gravel 10/10/05	IWI-1 AF Gravel 02/01/08	IWI-2 Zone 3-4 Clay 10/10/05	IWI-2 Zone 3-4 Clay 02/01/08	MW-2 USAS 02/13/01	MW-2 USAS 06/22/05	MW-2 USAS 01/25/06	MW-2 USAS 09/11/07	
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	61.2	NA	4.1 J	NA	NA	2 U	1 U	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	300	NA	19	NA	NA	NA	7.7
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	43	60	0.54 J	0.64 I	18	1.6	0.52 U	4.7
1,1-Dichloroethene	ug/L	165	370 D	0.53 J	0.83 I	3.4	0.98 J	0.45 U	1.5
cis-1,2-Dichloroethene	ug/L	92.5	190 D	0.5 U	4.1	11	9.4	2.5	6.4
Tetrachloroethene	ug/L	4.6	6.8	0.5 U	0.5 U	130	45.2	7.8 J	53
Trichloroethene	ug/L	3,310	4,900 D	0.65 U	0.99 I	60	32.4	7.9	64

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-3 USAS 02/13/01	MW-3 USAS 03/10/03	MW-3 USAS 01/05/05	MW-3 USAS 06/14/05	MW-3 USAS 01/25/06	MW-3 USAS 10/11/06	MW-3 USAS 12/06/06	MW-3 USAS 09/11/07	MW-3 USAS 01/25/08	MW-3 USAS 06/17/04
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	1 U	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	1.9 U	1 U	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	R [R]	R	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	2 UJ [2 UJ]	0.76 U	0.69 UJ	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	1 U	0.52 U	0.5 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	1 U	1 U	0.45 U	0.5 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1 U	1 U	0.65 U	0.5 U	0.65 U	0.75 U [0.75 U]	0.75 UJ	0.75 U	0.65 U
Tetrachloroethene	ug/L	1 U	1 U	0.34 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	1 U	1 U	0.28 U	0.5 U	0.28 U	0.71 U [0.71 U]	0.71 U	0.71 U	0.5 U

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-4 USAS 02/13/01	MW-4 USAS 03/10/03	MW-4 USAS 06/22/05	MW-4 USAS 01/27/06	MW-4 USAS 10/11/06	MW-4 USAS 12/07/06	MW-4 USAS 09/11/07	MW-4 USAS 01/25/08	MW-4 USAS 12/22/04	MW-4 USAS 06/17/04
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	1 U	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	2 U	1 U	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	1 J	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	1.8 I	1 U	NA
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	2	1 U [1 U]	1.1	6.6	90	180	38	6.1	0.52 U 0.42 I
1,1-Dichloroethene	ug/L	1 U	1 U [1 U]	0.5 U	0.8 I	28	79	16	1.7	0.45 U 0.2 U
cis-1,2-Dichloroethene	ug/L	1 U	1 U [1 U]	0.5 U	0.65 U	1.9	0.8 I	4.8	0.65 U	0.65 U NA
Tetrachloroethene	ug/L	1 U	1 U [1 U]	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U	0.34 U 0.2 U
Trichloroethene	ug/L	5.6	2.1 [3.1]	0.5 U	1.1	3	11	27	5.9	0.28 U 1.1

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-5 USAS 03/10/03	MW-5 USAS 12/21/04	MW-5 USAS 06/23/05	MW-5 USAS 01/24/06	MW-5 USAS 03/21/06	MW-5 USAS 10/10/06	MW-5 USAS 12/06/06	MW-5 USAS 09/11/07	MW-5 USAS 01/24/08	MW-5 USAS 06/16/04
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	1.9 U	1 U	1 U	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	R	R	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	0.52 U	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	1 U	0.45 U	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1 U	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.75 UJ	0.75 U	0.65 U
Tetrachloroethene	ug/L	1 U	0.34 U	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	1 U	0.28 U	0.5 U	0.28 U	0.28 U	0.71 U	0.71 U	0.71 U	0.5 U

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:		MW-6 USAS 03/10/03	MW-6 USAS 12/21/04	MW-6 USAS 06/21/05	MW-6 USAS 06/23/05	MW-6 USAS 01/24/06	MW-6 USAS 10/11/06	MW-6 USAS 12/07/06	MW-6 USAS 09/11/07	MW-6 USAS 01/30/08	MW-6 USAS 06/17/04	MW-6 USAS 12/22/04
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	3.2 J
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	NA	NA	2 U	1 U	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	R	R	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	2 UJ	0.76 U	1 I	1 U	NA	NA
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	0.52 U	0.5 U	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.2 U	0.52 U
1,1-Dichloroethene	ug/L	1 U	0.45 U	0.5 U	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	0.2 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1 U	0.65 U	0.5 U	0.5 U	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U	NA	0.65 U
Tetrachloroethene	ug/L	1 U	0.34 U	0.5 U	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U	0.2 U	0.34 U
Trichloroethene	ug/L	1 U	0.28 U	0.5 U	0.5 U	0.28 U	0.71 U	0.71 U	0.71 U	0.5 U	0.5 U	0.28 U

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-7D USAS 03/06/03	MW-7D USAS 12/19/04	MW-7D USAS 06/15/05	MW-7D USAS 01/24/06	MW-7D USAS 10/09/06	MW-7D USAS 12/05/06	MW-7D USAS 09/10/07	MW-7D USAS 01/24/08	MW-7D USAS 06/17/04	MW-7S USAS 03/06/03	MW-7S USAS 01/05/05	MW-7S USAS 06/15/05	MW-7S USAS 01/24/06	MW-7S USAS 10/09/06	MW-7S USAS 12/05/06	MW-7S USAS 09/10/07	MW-7S USAS 01/24/08	MW-7S USAS 06/17/04	
Semivolatile Organics																			
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	1 U [1 U]	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																			
1,4-Dioxane	ug/L	NA	NA	2 U	1 U	NA	NA	NA	NA	NA	NA	1.9 U	1 U	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	R [53 J]	R [R]	NA	NA	NA	NA	NA	NA	R	R	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ [2 UJ]	0.76 U [0.76 U]	0.69 UJ	1 U	NA	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	NA
Volatile Organics (8260B)																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	0.52 U	0.52 J	0.52 U	0.6 U [0.6 U]	0.6 U [0.6 U]	0.6 U	0.52 U	0.2 U	1 U	0.52 U [0.52 U]	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.2 U
1,1-Dichloroethene	ug/L	1 U	0.45 U	1.3	0.45 U	0.83 U [0.83 U]	0.83 U [0.83 U]	0.83 U	0.45 U	0.2 U	1 U	0.45 U [0.45 U]	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	0.3 U
cis-1,2-Dichloroethene	ug/L	1 U	0.65 U	0.5 U	0.65 U	0.75 U [0.75 U]	0.75 U [0.75 U]	0.75 U	0.65 U	NA	1 U	0.65 U [0.65 U]	0.5 U	0.65 U	0.75 U	0.75 UJ	0.75 U	0.65 U	NA
Tetrachloroethene	ug/L	1 U	0.34 U	4	0.34 U	0.65 U [0.65 U]	0.65 U [0.65 U]	0.65 U	0.5 U	0.46 U	1 U	0.34 U [0.34 U]	0.5 U	0.34 U	0.77 U	0.65 U	0.65 U	0.5 U	0.5 U
Trichloroethene	ug/L	1 U	0.28 U	3.1	0.28 U	0.71 U [0.71 U]	0.71 U [0.71 U]	0.71 U	0.5 U	0.5 U	1 U	0.28 U [0.28 U]	0.5 U	0.28 U	0.71 U	0.71 U	0.71 U	0.5 U	0.5 U

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-8D USAS 03/06/03	MW-8D USAS 12/19/04	MW-8D USAS 06/14/05	MW-8D USAS 01/25/06	MW-8D USAS 03/24/06	MW-8D USAS 12/05/06	MW-8D USAS 09/11/07	MW-8D USAS 02/01/08	MW-8D USAS 06/15/04	MW-8S USAS 03/06/03	MW-8S USAS 12/19/04	MW-8S USAS 06/14/05	MW-8S USAS 01/25/06	MW-8S USAS 03/24/06	MW-8S USAS 12/05/06	MW-8S USAS 09/11/07	MW-8S USAS 01/25/08	MW-8S USAS 06/15/04	
Semivolatile Organics																			
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	1 U	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																			
1,4-Dioxane	ug/L	NA	NA	1.9 U	1 U	1 U	NA	NA	NA	NA	NA	1.9 U	1 U	1 U	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	R	NA	NA	NA	NA	NA	NA	NA	NA	R	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	0.76 U	0.69 U	1 U	NA	NA	NA	NA	NA	NA	0.76 U	0.69 U	1 U	NA
Volatile Organics (8260B)																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	0.52 U	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U	0.45 I	1 U	0.52 U	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U	0.2 U
1,1-Dichloroethene	ug/L	1 U	0.45 U	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.45 U	0.24 I	1 U	0.45 U	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.45 U	0.2 U
cis-1,2-Dichloroethene	ug/L	1 U	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.75 U	0.65 U	NA	1 U	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.75 U	0.65 U	NA
Tetrachloroethene	ug/L	1 U	0.34 U	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U	0.68 I	1 U	0.34 U	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U	0.5 U
Trichloroethene	ug/L	1 U	0.35 J	0.5 U	0.53 I	0.28 U	0.71 U	0.71 U	0.5 U	0.5 U	1 U	0.28 U	0.5 U	0.28 U	0.28 U	0.71 U	0.71 U	0.5 U	0.5 U

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-9D USAS 03/07/03	MW-9D USAS 12/20/04	MW-9D USAS 06/15/05	MW-9D USAS 01/20/06	MW-9D USAS 10/11/06	MW-9D USAS 12/13/06	MW-9D USAS 09/12/07	MW-9D USAS 01/24/08	MW-9D USAS 06/15/04	MW-9S USAS 03/07/03	MW-9S USAS 12/20/04	MW-9S USAS 06/15/05	MW-9S USAS 01/20/06	MW-9S USAS 10/11/06	MW-9S USAS 12/13/06	MW-9S USAS 09/12/07	MW-9S USAS 01/24/08	MW-9S USAS 06/15/04	
Semivolatile Organics																			
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)																			
1,4-Dioxane	ug/L	NA	NA	1.9 U	1 U	NA	NA	NA	NA	NA	NA	2 U	1 U	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	R	NA	NA	NA	NA	NA	NA	NA	34 J	R	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	NA	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	NA
Volatile Organics (8260B)																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	0.52 U	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.2 U	1 U	0.52 U	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.2 U
1,1-Dichloroethene	ug/L	1 U	0.45 U	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	0.2 U	1 U	0.45 U	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	0.2 U
cis-1,2-Dichloroethene	ug/L	1 U	0.65 U	0.5 U	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U	NA	1 U	0.65 U	0.5 U	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U	NA
Tetrachloroethene	ug/L	1 U	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U	0.5 U	1 U	0.34 U	0.62 J	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U	0.5 U
Trichloroethene	ug/L	1 U	0.28 J	0.5 U	0.28 UJ	0.71 U	0.71 U	0.71 U	0.5 U	0.5 U	1 U	0.28 U	0.5 U	0.28 UJ	0.71 U	0.71 U	0.71 U	0.5 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-10 USAS 03/10/03	MW-10 USAS 01/04/05	MW-10 USAS 06/14/05	MW-10 USAS 01/24/06	MW-10 USAS 10/11/06	MW-10 USAS 12/05/06	MW-10 USAS 09/10/07	MW-10 USAS 01/31/08	MW-10 USAS 06/16/04	MW-11 USAS 03/10/03	MW-11 USAS 12/20/04	MW-11 USAS 06/22/05	MW-11 USAS 01/27/06	MW-11 USAS 10/11/06	MW-11 USAS 12/07/06	MW-11 USAS 12/11/06	MW-11 USAS 09/11/07	MW-11 USAS 01/24/08	MW-11 USAS 06/17/04	
Semivolatile Organics																				
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	1 U [1.6 J]	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																				
1,4-Dioxane	ug/L	NA	NA	1.9 U	1 U	NA	NA	NA	NA	NA	NA	6.2	3 I	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	R	NA	NA	NA	NA	NA	NA	20 J	3.4 J	0.99 J	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	0.69 UJ	1 U	NA	NA	NA	NA	40 UJ	2.2	0.76 U	0.69 U	1 U	NA	
Volatile Organics (8260B)																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	90	7.8	0.84 J	0.52 U	0.6 U	0.6 U	0.6 U	5.3	260	14 [12]	6.7 [5.3]	92.8 J [53.9 J]	72	310 D	36	0.6 U	9.2	9.1	3.2
1,1-Dichloroethene	ug/L	140	0.46 J	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	1	130	7.9 [6.8]	0.45 U [0.45 U]	49.3 J [28.4 J]	150	940 D	130	0.83 U	1.6	1.5	2.3
cis-1,2-Dichloroethene	ug/L	48	0.65 U	0.5 U	0.65 U	0.75 U	0.75 UJ	0.75 U	0.65 U	NA	6.2 [5.2]	3.8 [3.2]	68.4 J [39.4 J]	12	50 D	1	0.75 UJ	1	0.65 U	NA
Tetrachloroethene	ug/L	13	12	5.5	0.34 U	0.65 U	0.65 U	0.65 U	1.5	22	4.4 [4]	3.7 [4.1]	14.4 [5 U]	4.3 J	7.7 D	3	0.65 U	1.2	1.4	4.8
Trichloroethene	ug/L	820 L	27	3.9	0.7 I	0.71 U	0.71 U	0.71 U	1.8 U	550	140 [190]	100 [100]	937 J [544 J]	200	640 D	79	0.71 U	33	17	71

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS	MW-12 USAS
	03/10/03	12/21/04	06/22/05	01/25/06	10/05/06	12/05/06	09/11/07	01/24/08	06/17/04	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	10	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	14.3 [16.6]	11	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	22 J	16 J	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	33 J	16	0.69 U [0.69 U]	23	NA
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	23	4.7	12.6 [12.7]	11	38 D	18	5.1 [4.2]	10	21
1,1-Dichloroethene	ug/L	30	0.45 U	15.8 [15.4]	12	24 D	12	0.83 U [0.83 U]	6.9	24
cis-1,2-Dichloroethene	ug/L	8.1	2.2	5.6 [5.4]	2.8	5.5 D	3.8 U	11 [8.1]	5.1	NA
Tetrachloroethene	ug/L	290	110	199 [200]	120	300 D	120	55 [39]	58	300
Trichloroethene	ug/L	710	150	364 [375]	310	470 D	210	13 [11]	190 D	600

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-13D USAS 03/07/03	MW-13D USAS 12/22/04	MW-13D USAS 06/14/05	MW-13D USAS 01/19/06	MW-13D USAS 10/10/06	MW-13D USAS 12/15/06	MW-13D USAS 09/13/07	MW-13D USAS 01/29/08	MW-13D USAS 06/17/04	MW-13S USAS 03/07/03	MW-13S USAS 12/22/04	MW-13S USAS 06/14/05	MW-13S USAS 01/19/06	MW-13S USAS 01/29/08	MW-13S USAS 06/17/04
Semivolatile Organics															
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	1 U [1 U]	NA	NA	NA	NA
Semivolatile Organics (8270C)															
1,4-Dioxane	ug/L	NA	NA	2.1 U	1 U	NA	NA	NA	NA	NA	NA	1.9 U	1 U	NA	NA
Volatile Organics (8260B) - SIM															
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	R	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution															
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	NA	NA	NA	NA	4.5	NA
Volatile Organics (8260B)															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U [1 U]	0.52 U	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.2 U	1 U	0.52 U [0.52 U]	0.5 U	0.52 U	0.52 U
1,1-Dichloroethene	ug/L	1 U [1 U]	0.45 U	1.1	0.71	1.1	0.95 I	0.83 U	0.44 U	0.42 I	1 U	0.45 U [0.45 U]	0.5 U	0.45 U	0.44 U
cis-1,2-Dichloroethene	ug/L	1 U [1 U]	0.65 U	0.5 U	0.65 U	0.75 U	0.75 UJ	0.75 U	0.64 U	NA	1 U	0.65 U [0.65 U]	0.5 U	0.65 U	0.64 U
Tetrachloroethene	ug/L	1 U [1 U]	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U	0.2 U	1 U	0.34 U [0.34 U]	0.5 U	0.34 U	0.5 U
Trichloroethene	ug/L	1 U [1 U]	0.28 U	0.5 U	0.28 I	0.71 U	0.71 UJ	0.71 U	0.5 U	0.5 U	1 U	0.28 U [0.28 U]	0.5 U	0.28 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-14D USAS 03/06/03	MW-14D USAS 12/20/04	MW-14D USAS 06/14/05	MW-14D USAS 01/19/06	MW-14D USAS 10/11/06	MW-14D USAS 12/13/06	MW-14D USAS 09/13/07	MW-14D USAS 01/24/08	MW-14D USAS 06/16/04	MW-14S USAS 03/06/03	MW-14S USAS 12/20/04	MW-14S USAS 06/20/05	MW-14S USAS 01/19/06	MW-14S USAS 10/11/06	MW-14S USAS 06/16/04	MW-14S USAS 12/13/06	MW-14S USAS 09/13/07	MW-14S USAS 01/24/08	
Semivolatile Organics																			
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	1 U	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																			
1,4-Dioxane	ug/L	NA	NA	2 U	1 U [1 U]	NA	NA	NA	NA	NA	NA	2 U	1 U	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	R	NA	NA	NA	NA	NA	NA	R	NA	0.94 U [R]	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	NA	NA	NA	NA	2 UJ	NA	0.76 U [0.76 U]	0.69 U	1 U	
Volatile Organics (8260B)																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	1 U	0.52 U	0.5 U	0.52 U [0.52 U]	0.6 U	0.6 U	0.6 U	0.52 U	0.2 U	1 U	0.52 U	0.5 U	0.52 U	0.6 U	0.2 U	0.6 U [0.6 U]	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	1 U	0.45 U	0.5 U	0.45 U [0.45 U]	0.83 U	0.83 U	0.83 U	0.45 U	0.2 U	1 U	0.45 U	0.5 U	0.45 U	0.83 U	0.2 U	0.83 U [0.83 U]	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1 U	0.65 U	0.5 U	0.65 U [0.65 U]	0.75 U	0.75 UJ	0.75 U	0.65 U	NA	1 U	0.65 U	0.5 U	0.65 U	0.75 U	NA	0.75 U [0.75 U]	0.75 U	0.65 U
Tetrachloroethene	ug/L	1 U	0.34 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.65 U	0.65 U	0.5 U	0.5 U	1 U	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U	0.65 U [0.65 U]	0.65 U	0.5 U
Trichloroethene	ug/L	1 U	0.28 U	0.5 U	0.28 U [0.28 U]	0.71 U	0.71 UJ	0.71 U	0.5 U	0.5 U	1 U	0.28 U	0.5 U	0.28 U	0.71 U	0.5 U	0.71 U [0.71 U]	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-15D USAS 03/06/03	MW-15D USAS 12/21/04	MW-15D USAS 06/20/05	MW-15D USAS 01/25/06	MW-15D USAS 10/16/06	MW-15D USAS 12/18/06	MW-15D USAS 01/18/07	MW-15D USAS 04/25/07	MW-15D USAS 07/10/07	MW-15D USAS 09/12/07	MW-15D USAS 01/24/08	MW-15S USAS 03/06/03	MW-15S USAS 01/05/05	MW-15S USAS 06/20/05	MW-15S USAS 01/25/06	MW-15S USAS 10/16/06	MW-15S USAS 12/18/06	MW-15S USAS 09/12/07	MW-15S USAS 01/24/08	
Semivolatile Organics																				
1,4-Dioxane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 UJ	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																				
1,4-Dioxane	ug/L	NA	NA	2.2 J	1 U	NA	NA	NA	NA	NA	NA	NA	NA	2 U [2 U]	1 U	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																				
1,4-Dioxane	ug/L	NA	NA	NA	7.1 J	4.7 J [5.2 J]	NA	NA	NA	NA	NA	NA	NA	NA	NA	R	R	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																				
1,4-Dioxane	ug/L	NA	NA	NA	6.4 J	4.6 [8.2]	6.9	8.7	7.3 J	10	7.5	NA	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	
Volatile Organics (8260B)																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	1 U [1.7]	1.3	2.4	2.4	2.8	4.7 [4.8]	2.8	3.7	6.1	2.6	4.9	1 U	0.52 U	0.5 U [0.5 U]	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	1 U [1 U]	0.82 J	1.1	0.88 I	0.83 U	1.4 [1.4]	0.83 U	1.1	1.4	0.83 U	0.45 U	1 U	0.45 U	0.5 U [0.5 U]	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1 U [1 U]	0.65 U	0.77 J	0.8 I	1.9	2.2 [2]	2	2.5	3.2	1.4	3	1 U	0.65 U	0.5 U [0.5 U]	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	1 U [1 U]	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U [0.65 U]	0.65 U	0.65 U	0.65 U	0.65 U	0.5 U	1 U	0.34 U	0.5 U [0.5 U]	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	1 U [1 U]	0.28 U	0.62 J	1.4	3.1	2.9 [2.6]	2.5	3.4	4.2	2.4	4	1 U	0.28 U	0.5 U [0.5 U]	0.28 U	0.71 U	0.71 U	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-16D USAS 03/07/03	MW-16D USAS 06/16/05	MW-16D USAS 01/18/06	MW-16D USAS 10/13/06	MW-16D USAS 12/11/06	MW-16D USAS 01/29/08	MW-16D USAS 06/16/04	MW-16S USAS 03/07/03	MW-16S USAS 01/04/05	MW-16S USAS 06/16/05	MW-16S USAS 01/18/06	MW-16S USAS 01/29/08	MW-16S USAS 06/16/04
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	1 U	NA	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	NA	2 U	1 U	NA	NA	NA	NA	NA	2 U	1 U	NA	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	5.3 J	5.8 J	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	5.1 J	6	6	NA	NA	NA	NA	1 U	NA
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U	0.2 U	1 U	0.52 U	0.5 U	0.52 U	0.2 U
1,1-Dichloroethene	ug/L	1 U	0.5 U	0.45 U	0.83 U	0.83 U	0.44 U	0.2 U	1 U	0.45 U	0.5 U	0.45 U	0.2 U
cis-1,2-Dichloroethene	ug/L	1 U	0.5 U	0.65 U	0.75 U	0.75 UJ	0.64 U	NA	1 U	0.65 U	0.5 U	0.65 U	NA
Tetrachloroethene	ug/L	1 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U	0.5 U	1 U	0.34 U	0.5 U	0.34 U	0.5 U
Trichloroethene	ug/L	1 U	0.5 U	0.28 U	0.71 U	0.71 U	0.5 U	0.5 U	1 U	0.28 U	0.5 U	0.28 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-17D USAS 12/21/04	MW-17D USAS 06/15/05	MW-17D USAS 01/19/06	MW-17D USAS 10/11/06	MW-17D USAS 12/13/06	MW-17D USAS 09/12/07	MW-17D USAS 01/23/08	MW-17D USAS 06/16/04	MW-17S USAS 01/04/05	MW-17S USAS 06/15/05	MW-17S USAS 01/19/06	MW-17S USAS 10/11/06	MW-17S USAS 09/12/07	MW-17S USAS 01/23/08	MW-17S USAS 06/16/04
Semivolatile Organics															
1,4-Dioxane	ug/L	1 U	NA	NA	NA	NA	NA	NA	1 U	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)															
1,4-Dioxane	ug/L	NA	2 U	1 U	NA	NA	NA	NA	NA	2 U	1 U	NA	NA	NA	NA
Volatile Organics (8260B) - SIM															
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	NA	NA	R	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution															
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.87 I	0.69 U [0.69 U	1 U	NA	NA	NA	2 UJ	0.69 U	1 U	NA
Volatile Organics (8260B)															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U [0.6 U	0.52 U	0.2 U	0.52 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U [0.83 U	0.45 U	0.2 U	0.45 U	0.5 U	0.45 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U	0.65 U	0.75 U	0.75 UJ	0.75 U [0.75 U	0.65 U	NA	0.65 U	0.5 U	0.65 U	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U [0.65 U	0.5 U	0.5 U	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 J	0.5 U	0.36 I	0.71 U	0.71 UJ	0.71 U [0.71 U	0.5 U	0.5 U	0.28 U	0.5 U	0.28 U	0.71 U	0.71 U	0.5 U

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TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-18D USAS 12/21/04	MW-18D USAS 06/17/05	MW-18D USAS 01/17/06	MW-18D USAS 02/28/06	MW-18D USAS 03/23/06	MW-18D USAS 04/01/06	MW-18D USAS 12/18/06	MW-18D USAS 01/28/08	MW-18D USAS 06/16/04
Semivolatile Organics									
1,4-Dioxane	ug/L	1 U	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	5.6	4 I	1 U	6.4 I	1 U	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	7.8 J	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	11 J	2.6	NA
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	3.1	3.9	1.6	NA	1.9	1.7	2.3	0.74 I
1,1-Dichloroethene	ug/L	1.8	2.4	0.95 I	NA	1.2	0.98 I	1.8	0.45 U
cis-1,2-Dichloroethene	ug/L	8.9	9	3.1	NA	3.7	3.5	5.3	2.1
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	NA	0.34 U	0.34 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	1.3	7.1	5.1	NA	5.9	4.6	3	0.52 J

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-18S	MW-18S	MW-18S	MW-18S	MW-18S	MW-18S
Zone:	USAS	USAS	USAS	USAS	USAS	USAS
Date Collected:	12/21/04	06/17/05	01/17/06	12/18/06	01/28/08	06/16/04
Semivolatile Organics						
1,4-Dioxane	ug/L	1 U	NA	NA	NA	NA
Semivolatile Organics (8270C)						
1,4-Dioxane	ug/L	NA	1.9 U	1.7 I	NA	NA
Volatile Organics (8260B) - SIM						
1,4-Dioxane	ug/L	NA	NA	NA	2.9 J	NA
Volatile Organics (8260) - SIM Isotope Dilution						
1,4-Dioxane	ug/L	NA	NA	NA	3.2	1 U
Volatile Organics (8260B)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1.6	0.5 U	1.9	1.5	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	1.2	0.5 U	0.33 I	0.71 U	0.5 UJ

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-19 Lower AF 01/12/05	MW-19 Lower AF 06/24/05	MW-19 Lower AF 01/31/06	MW-19 Lower AF 03/21/06	MW-19 Lower AF 12/19/06	MW-19 Lower AF 07/11/07	MW-19 Lower AF 09/11/07	MW-19 Lower AF 01/31/08	MW-20 USAS 01/17/05	MW-20 USAS 06/21/05	MW-20 USAS 01/26/06	MW-20 USAS 12/19/06	MW-20 USAS 01/24/08
Semivolatile Organics													
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	NA	NA	13	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	NA	1.9 U	1 U [1 U]	1 U	NA	NA	NA	NA	NA	4.4 J	1 U	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	NA	NA	NA	NA	NA	9.4 J	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	NA	0.76 U	0.69 U	0.69 U	1 U	NA	NA	NA	9.6
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	1.2 [1.2]	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	4.8	7	7.7	11
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	2.4 [2.6]	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	2.8	4.1	5.5	9
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U [0.65 U]	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U	0.34 I	0.75 J	0.7 I	0.92 J
Tetrachloroethene	ug/L	0.85 I	0.91 J	0.34 U [0.34 U]	2.4	0.65 U	0.86 I	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.7 I	0.5 U	2.7 [2.6]	1.1	0.71 U	3.8	0.71 U	0.5 U	0.38 I	0.61 J	1.2	2.4 J
													1.3

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TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-21 S&P Sand 01/28/08	MW-21 S&P Sand 01/12/05	MW-21 S&P Sand 06/14/05	MW-21 S&P Sand 02/20/06	MW-21 S&P Sand 12/18/06	MW-22 Lower AF 01/11/05	MW-22 Lower AF 06/24/05	MW-22 Lower AF 02/02/06	MW-22 Lower AF 12/19/06	MW-22 Lower AF 01/25/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	2.5 U	NA	NA	NA	2.5 U	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	2 U	1 U	NA	NA	2 U	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	1.7 J	NA	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	1 U [2.6]	NA	NA	NA	0.76 U	NA	NA	NA	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U [0.52 U]	0.18 U	0.5 U	0.52 U	0.6 U	0.18 U	0.5 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U [0.45 U]	0.36 U	0.5 U	0.45 U	0.83 U	0.36 U	0.5 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U [0.65 U]	0.18 U	0.5 U	0.65 U	0.75 U	0.18 U	0.5 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.5 U [0.5 U]	0.18 U	0.5 U	0.34 U	0.65 U	0.18 U	1.4	1.4	0.82 I
Trichloroethene	ug/L	0.5 U [0.5 U]	0.11 U	0.5 U	0.28 U	0.71 U	0.11 U	0.5 U	0.28 U	0.71 U

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-23 S&P Sand 01/11/05	MW-23 S&P Sand 06/21/05	MW-23 S&P Sand 01/21/06	MW-23 S&P Sand 03/17/06	MW-23 S&P Sand 03/21/06	MW-23 S&P Sand 03/27/06	MW-23 S&P Sand 12/19/06	MW-23 S&P Sand 01/25/08
Semivolatile Organics								
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)								
1,4-Dioxane	ug/L	NA	2.4 J	1 U [1 U]	2.1 I	1 U	1 U	NA
Volatile Organics (8260B) - SIM								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	3.7 J
Volatile Organics (8260) - SIM Isotope Dilution								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	2.4
Volatile Organics (8260B)								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U [0.52 U]	0.52 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethane	ug/L	0.36 U	0.5 U	0.45 U [0.45 U]	0.45 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U [0.65 U]	0.65 U	0.65 U	0.65 U	0.75 UJ
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U [0.34 U]	0.34 U	0.34 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U [0.28 U]	0.28 U	0.28 U	0.28 U	0.71 UJ

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TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-24 USAS 01/11/05	MW-24 USAS 06/15/05	MW-24 USAS 01/23/06	MW-24 USAS 10/11/06	MW-24 USAS 12/13/06	MW-24 USAS 01/29/08	MW-25 USAS 01/06/05	MW-25 USAS 06/21/05	MW-25 USAS 01/20/06	MW-25 USAS 02/28/06	MW-25 USAS 03/19/06	MW-25 USAS 04/04/06	MW-25 USAS 12/13/06	MW-25 USAS 01/24/08	
Semivolatile Organics															
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	9.8	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)															
1,4-Dioxane	ug/L	NA	1.9 U	1 U	NA	NA	NA	7.7 [7.7]	4 I	6.8 IJ	10	1 UJ	NA	NA	
Volatile Organics (8260B) - SIM															
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	NA	NA	NA	9.2 J [11 J]	NA	
Volatile Organics (8260) - SIM Isotope Dilution															
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	1 U	NA	NA	NA	NA	NA	12 [16]	7.7	
Volatile Organics (8260B)															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U	39	42.6 [42.4]	37	NA	38	35	27 [26]	14
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.83 U	0.44 U	13	18.8 [17.8]	16	NA	17	17	16 [16]	7.5
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U	0.75 U	0.64 U	3	6.8 [6.8]	5.1	NA	5.2	5.4	3.5 [3.6]	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.5 U [0.5 U]	0.34 U	NA	0.34 U	0.34 U	65 U [0.65]	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U	0.71 U	0.5 U	2.1	3.6 [3.6]	4.4 J	NA	3.8	3.9	3.3 [3.3]	2.3

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TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-26 USAS 01/13/05	MW-26 USAS 06/17/05	MW-26 USAS 01/19/06	MW-26 USAS 12/11/06	MW-26 USAS 01/28/08	MW-27 USAS 01/06/05	MW-27 USAS 06/17/05	MW-27 USAS 02/21/06	MW-27 USAS 06/28/06	MW-27 USAS 10/10/06	MW-27 USAS 12/18/06	MW-27 USAS 01/28/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	390	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	NA	2 U	1 U	NA	NA	457	440	530 D	NA	NA	NA	
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	1.2 J	NA	NA	NA	NA	370 DJ	530 J	NA	
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	NA	790 DJ	760	69	
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U	140	266	260	140	230 D	290	69
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U	290	459	450	330	410 D	600	120
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 UJ	0.65 U	10	34.6	37	17	74 D	59	11
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.5 UJ	290	779	790	370	2,000 D	960 D	79 J
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 UJ	0.5 UJ	42	94.3	110	66	120 D	91	23 J

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FORMER AMERICAN BERYLLIUM COMPANY SITE
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Location ID:	MW-28	MW-28	MW-28	MW-28	MW-28	MW-28	MW-28	MW-29	MW-29	MW-29	MW-29	
Zone:	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	
Date Collected:	01/03/05	06/20/05	01/18/06	03/19/06	12/18/06	01/28/08	01/06/05	06/17/05	01/17/06	12/18/06	01/28/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	40	NA	NA	NA	NA	NA	28	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	41.3	35	18	NA	NA	NA	34.4	71	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	49 J	NA	NA	NA	NA	130 J	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	80	28	NA	NA	NA	120	91
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	3.2	8.5	6.8	6.1	5.3	3.1	8.5	12	11	13	10
1,1-Dichloroethene	ug/L	0.71 J	4.7	3.3	2.9	3.7	1.4	5.3	7.3	7.6	10	7.5
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.65 U	6.6	11.7	11	12	12
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 UJ	0.34 U	0.5 U	0.34 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	0.88 J	1.7	1.4	1	0.88 I	0.5 UJ	37	51.6	69	75	91 J

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Location ID: Zone: Date Collected:	MW-30 USAS 01/11/05	MW-30 USAS 06/15/05	MW-30 USAS 01/20/06	MW-30 USAS 10/11/06	MW-30 USAS 12/19/06	MW-30 USAS 09/13/07	MW-30 USAS 01/30/08	MW-31 Lower AF 01/12/05	MW-31 Lower AF 06/15/05	MW-31 Lower AF 01/20/06	MW-31 Lower AF 12/18/06	MW-31 Lower AF 01/28/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	NA	2.1 U	1 U	NA	NA	NA	NA	2 U	1 U	NA	NA	
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	NA	R	NA	
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	0.69 U	1 U	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.48 I	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.62 I	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.23 I	0.5 U	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U	0.18 U	0.5 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U	0.18 U	1.2	1.6	1.2	0.7 I
Trichloroethene	ug/L	3.6	0.5 U	0.28 UJ	0.71 U	0.71 U	0.71 U	0.5 U	0.11 U	0.5 U	0.28 UJ	0.71 U	0.5 U

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Location ID: Zone: Date Collected:	MW-32 USAS 12/30/04	MW-32 USAS 06/27/05	MW-32 USAS 01/31/06	MW-32 USAS 10/05/06	MW-32 USAS 12/05/06	MW-32 USAS 09/10/07	MW-32 USAS 01/24/08	MW-33 LSAS 12/28/04	MW-33 LSAS 06/22/05	MW-33 LSAS 01/25/06	MW-33 LSAS 10/05/06	MW-33 LSAS 12/05/06	MW-33 LSAS 01/17/07	MW-33 LSAS 04/24/07	MW-33 LSAS 07/10/07	MW-33 LSAS 09/10/07	MW-33 LSAS 01/31/08	
Semivolatile Organics																		
1,4-Dioxane	ug/L	6.5	NA	NA	NA	NA	NA	580	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																		
1,4-Dioxane	ug/L	NA	3.1 J	7.7 I	NA	NA	NA	NA	418	400	NA	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																		
1,4-Dioxane	ug/L	NA	NA	NA	R	6.6 J [8.4 J]	NA	NA	NA	NA	NA	530 J	NA	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																		
1,4-Dioxane	ug/L	NA	NA	NA	9.9 J	6.3 [8.5]	3.5 J	2.6	NA	NA	NA	930 DJ	670	750	570	750	900	590 J
Volatile Organics (8260B)																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	650 DJ	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	87	76.6	86	60	56 [53]	9.8	9.9	42	60.4	110	130 D	150	61 J	69	260	140	160 EJ
1,1-Dichloroethene	ug/L	120	87.6	98	65	72 [110]	7.8	8.6	140	163	380	440 D	670	170	240	1,400 D	550	580 DJ
cis-1,2-Dichloroethene	ug/L	5.9	3.8	7.4	9.8	3.4 [3.8 U]	2.1	0.65 U	3.2	17.4	17	25 D	6.3	15 J	17	16	9.1 I	22 J
Tetrachloroethene	ug/L	2.4	8.1	9.3	120	72 [87]	8	3.5	0.34 U	9.1	20	63 D	59	60 J	60	43	24	140 J
Trichloroethene	ug/L	65	71.4	59	200	120 [140]	13	8.3	17	106	150	330 D	300	230	330	310	200	410 EJ

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-34 S&P Sand 12/19/04	MW-34 S&P Sand 01/13/05	MW-34 S&P Sand 06/22/05	MW-34 S&P Sand 02/16/06	MW-34 S&P Sand 10/13/06	MW-34 S&P Sand 12/13/06	MW-34 S&P Sand 01/30/08	MW-35 USAS 01/13/05	MW-35 USAS 06/22/05	MW-35 USAS 02/16/06	MW-35 USAS 10/13/06	MW-35 USAS 12/13/06	MW-35 USAS 01/18/07	MW-35 USAS 04/25/07	MW-35 USAS 07/10/07	MW-35 USAS 09/13/07	MW-35 USAS 01/30/08	
Semivolatile Organics																		
1,4-Dioxane	ug/L	NA	2.5 U	NA	NA	NA	NA	13	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																		
1,4-Dioxane	ug/L	NA	NA	2 U	8 UJ	NA	NA	NA	1.9 U	16 J	NA	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	R	NA	NA	NA	NA	3.2 J	12	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	1 U	NA	NA	NA	3.5 J	12	2	1.2 I	1.3 J	0.69 U	1 U
Volatile Organics (8260B)																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	660	0.18 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U	20	0.99 J	12	3.1	15	2.3	0.61 I	1.5	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	1,700	0.36 U	0.5 U	0.45 U	0.83 U	0.83 U	0.45 U	48	2.1	24	7.2	39	5.1	2.7	9.4	2.6	3.7
cis-1,2-Dichloroethene	ug/L	15	0.18 U	0.5 U	0.65 U	0.75 U	0.75 U	0.65 U	1.1	0.5 U	0.65 U	0.75 U	0.75 U	0.75 U	0.75 U	1.2	0.75 U	0.65 U
Tetrachloroethene	ug/L	43	0.18 U	0.5 U	0.34 UJ	0.65 U	0.65 U	0.5 U	41	13	18 J	18	44	29	12	26	27	15
Trichloroethene	ug/L	230	0.28 I	0.5 U	0.28 U	0.71 U	0.71 U	0.5 U	58	6.3	52	19	52	15	6.9	16	9.6	11

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-36 USAS 01/11/05	MW-36 USAS 06/23/05	MW-36 USAS 01/24/06	MW-36 USAS 10/04/06	MW-36 USAS 12/05/06	MW-36 USAS 01/17/07	MW-36 USAS 04/25/07	MW-36 USAS 07/10/07	MW-36 USAS 09/10/07	MW-36 USAS 01/31/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	34	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	18.9	3.8 I	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	R	2.7 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	4.1 J	0.76 U	0.76 U	4	1.4 J	2.8 J	15
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	180	113	41	10 D	6	2 J	0.8 I	2.4	3.4	20
1,1-Dichloroethene	ug/L	440	242	96	23 D	15	3.8	2.4	8.5	8.9	54
cis-1,2-Dichloroethene	ug/L	210	115	39	18 D	3.7 J	4.4 J	1.6	2.2	1.6	5.8
Tetrachloroethene	ug/L	19 I	25 U	3.7	6.5 UD	3.2	2.3 J	1.2	0.82 I	0.86 I	1.3
Trichloroethene	ug/L	2,600	1,400	490	250 D	130	55	55	53	52	190 EJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-37 LSAS 01/07/05	MW-37 LSAS 06/14/05	MW-37 LSAS 01/27/06	MW-37 LSAS 10/04/06	MW-37 LSAS 12/05/06	MW-37 LSAS 01/17/07	MW-37 LSAS 04/25/07	MW-37 LSAS 07/10/07	MW-37 LSAS 09/10/07	MW-37 LSAS 01/31/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	190	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	232	330	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	900 DJ	580 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	520 DJ	560	680 J [720]	1,100	690 J	780	760 DJ
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	2.4	25 U	20	60 ID	40 I	120 U [55 J]	60 U	60 U	44 I	49
1,1-Dichloroethene	ug/L	74	51.6	280	360 D	350	320 [330 DJ]	300	810	370	230 DJ
cis-1,2-Dichloroethene	ug/L	440	217	3,300	1,300 D	590	530 [640 DJ]	380	490	240	420 EJ
Tetrachloroethene	ug/L	0.34 U	25 U	0.37 IJ	65 UD	32 U	130 U [2.4 J]	65 U	65 U	32 U	24
Trichloroethene	ug/L	2,000	2,310	6,100	7,700 D	13,000 D	8,100 J [8,300 DJ]	7,500	8,600	4,000	9,400 DJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-38	MW-38	MW-38	MW-38	MW-38	MW-38	MW-38	MW-38	MW-38	MW-38	MW-38	
Zone:	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	USAS	
Date Collected:	01/11/05	06/14/05	01/25/06	03/22/06	10/04/06	12/05/06	01/17/07	04/24/07	07/10/07	09/10/07	02/01/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	40	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	30.9	15	18	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	21 J	28 J	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	22 J	25 J	36	37	25 J [28]	36 J	24
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	61 I	123	83	59	36 D	34	32 J	41	44 [58]	22	20
1,1-Dichloroethene	ug/L	140	253	130	100	52 D	61	44	68	99 [110]	32	27
cis-1,2-Dichloroethene	ug/L	52 I	10.7 J	8.9	9.5	7.8 D	3.8 UJ	6.4 J	7.5 U	7.5 U [5.8]	7.5 U	6.5 U
Tetrachloroethene	ug/L	250	347	53	130	93 D	130	510 J	780	900 [830 D]	720	450
Trichloroethene	ug/L	2,200	1,160	460	620	580 D	570	630	940	1,100 [1,100 D]	750	600

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-39 LSAS 01/07/05	MW-39 LSAS 06/14/05	MW-39 LSAS 01/25/06	MW-39 LSAS 10/11/06	MW-39 LSAS 12/05/06	MW-39 LSAS 01/17/07	MW-39 LSAS 04/24/07	MW-39 LSAS 07/10/07	MW-39 LSAS 09/10/07	MW-39 LSAS 02/01/08	MW-40 USAS 01/11/05	MW-40 USAS 06/14/05	MW-40 USAS 01/24/06	MW-40 USAS 10/05/06	MW-40 USAS 12/05/06	MW-40 USAS 09/11/07	MW-40 USAS 01/31/08
Semivolatile Organics																	
1,4-Dioxane	ug/L	150	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	9.4	NA	NA	NA	NA
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	NA	179 J	40	NA	NA	NA	NA	NA	NA	NA	NA	9.8	15	NA	NA	NA
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	40 J	6.4 J	NA	NA	NA	NA	NA	NA	NA	NA	25 J	29 J	NA
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	23 J	8	26	3.4	9 J	0.69 UJ	160	NA	NA	NA	32 J	22	36
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	10 U	18	7.9 D	7.5	2.4 J	0.6 U	1.1	0.6 U	5.7	6.1 I	20.2	32	57 D	54	24
1,1-Dichloroethene	ug/L	26	34.2	11	11 D	8.1	3.4	0.89 I	1.5	0.83 U	22	11	23.1	39	57 D	66	20
cis-1,2-Dichloroethene	ug/L	130	94.5	13	15 D	1 J	10 J	1.4	3.6	0.75 U	44	2.4 I	2.6 J	3.9	8.3 D	6.8	4.6 I
Tetrachloroethene	ug/L	0.34 U	15.7 J	24	50 D	50	22 J	2.7	1.1	0.65 U	2 U	29	70.1	89	130 D	140	95
Trichloroethene	ug/L	360	526	210	210 D	100	110	13	29	2.6	190	180	137	220	470 D	500	580
																	970 EJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-41 LSAS 12/28/04	MW-41 LSAS 06/14/05	MW-41 LSAS 01/24/06	MW-41 LSAS 10/05/06	MW-41 LSAS 12/05/06	MW-41 LSAS 09/11/07	MW-41 LSAS 01/31/08	MW-42 USAS 12/29/04	MW-42 USAS 06/23/05	MW-42 USAS 01/31/06	MW-42 USAS 10/04/06	MW-42 USAS 12/05/06	MW-42 USAS 01/17/07	MW-42 USAS 04/24/07	MW-42 USAS 07/11/07	MW-42 USAS 09/10/07	MW-42 USAS 02/01/08
Semivolatile Organics																	
1,4-Dioxane	ug/L	180	NA	NA	NA	NA	NA	83	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	NA	185	250	NA	NA	NA	NA	45.7	97	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	310 J	NA	NA	NA	NA	70 DJ	86 J	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	540 DJ	230	610	450 J	NA	NA	100 DJ	96	96	140	84	80 J [80]	72
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	460 J	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	10	7.9	7.9	17	19	17	22	41	50 U	42	67 D	87	39 J	55	92	49 [52]
1,1-Dichloroethene	ug/L	48	43.4	44	54	81	57	68	70	50.4 J	95	110 D	170	48	77	170	81 [74]
cis-1,2-Dichloroethene	ug/L	64	69.9	65	85	72	72	96	140	133	130	180 D	68	85 J	74	120	90 [110]
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.7 I	1.1	0.65 U	0.59 I	6.5	50 U	10	16 UD	20 I	13 J	25	13 U	13 U [13]
Trichloroethene	ug/L	81	71.8	130	190	200	180	250 EJ	3,700	3,160	4,600	3,700 D	3,000	1,700	1,900	2,700	2,000 [2,000 D]

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-43 LSAS 12/28/04	MW-43 LSAS 06/14/05	MW-43 LSAS 09/07/05	MW-43 LSAS 01/24/06	MW-43 LSAS 10/11/06	MW-43 LSAS 12/05/06	MW-43 LSAS 01/17/07	MW-43 LSAS 04/24/07	MW-43 LSAS 07/11/07	MW-43 LSAS 09/10/07	MW-43 LSAS 02/01/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	140	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	169	121 J	270	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	290 DJ	500 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	570 J [660 J]	590	500	800	600	740	570 [560]
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	2.8	12	7.9	31	55 D [56]	53	26 J	32	67	37	36 [31]
1,1-Dichloroethene	ug/L	40	67.8	55.7	110 J	250 D [180 D]	270	78	120	430	150	150 D [140]
cis-1,2-Dichloroethene	ug/L	12	17.1	28.3	19	42 D [47]	6.8 J	21 J	16	18	14	19 [16]
Tetrachloroethene	ug/L	0.34 U	1 U	0.5 U	0.34 U	3.2 UD [0.78 I]	3.2 U	3.2 U	3.2 U	3.2 U	3.2 U	0.5 U [5 U]
Trichloroethene	ug/L	26	69.6	72.3	110 J	310 D [280 D]	220	220	220	220	150	190 D [150]

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-44 S&P Sand 01/12/05	MW-44 S&P Sand 06/24/05	MW-44 S&P Sand 01/31/06	MW-44 S&P Sand 06/29/06	MW-44 S&P Sand 10/12/06	MW-44 S&P Sand 12/13/06	MW-44 S&P Sand 07/10/07	MW-44 S&P Sand 09/13/07	MW-44 S&P Sand 01/28/08	MW-45 S&P Sand 01/12/05	MW-45 S&P Sand 07/14/05	MW-45 S&P Sand 01/26/06	MW-45 S&P Sand 10/10/06	MW-45 S&P Sand 12/15/06	MW-45 S&P Sand 01/29/08	
Semivolatile Organics																
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																
1,4-Dioxane	ug/L	NA	1.9 U	1.1 U [1.1 U]	3 I	NA	NA	NA	NA	NA	2.1 U	1 U	NA	NA	NA	
Volatile Organics (8260B) - SIM																
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	1.2 J	NA	NA	NA	NA	NA	NA	R	R	NA
Volatile Organics (8260) - SIM Isotope Dilution																
1,4-Dioxane	ug/L	NA	NA	NA	NA	2 UJ	0.76 U	1.2 J	1.5 I	4.8	NA	NA	NA	2 UJ	0.76 U	1 U
Volatile Organics (8260B)																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	9.3	7.8	9.5 [8.9]	2.7	2.5	3.3	2.4	1.2	5.1	0.18 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	12	8	7.1 [7.2]	4.5	4	6.3	3.6	1.3	2.9	0.36 U	0.5 U	0.45 U	0.83 U	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.73 I	0.5 U	0.65 U [0.65 U]	0.65 U	0.75 U	0.75 U	0.75 U	0.75 U	0.65 U	0.18 U	0.5 U	0.65 U	0.75 U	0.75 UJ	0.64 U
Tetrachloroethene	ug/L	0.55 I	0.5 U	0.34 U [0.34 U]	5.7	0.65 U	0.65 U	0.65 U	0.65 U	0.5 UJ	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.52 I	0.5 U	0.42 I [0.57 I]	1	0.71 U	0.71 U	0.71 U	0.71 U	0.5 UJ	0.11 U	0.5 U	0.28 U	0.71 U	0.71 UJ	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-46 Lower AF 01/20/05	MW-46 Lower AF 06/16/05	MW-46 Lower AF 02/18/06	MW-46 Lower AF 12/12/06	MW-46 Lower AF 01/31/08	MW-47 USAS 01/06/05	MW-47 USAS 06/17/05	MW-47 USAS 01/26/06	MW-47 USAS 12/18/06	MW-47 USAS 01/30/08	MW-48 LSAS 06/14/05	MW-48 LSAS 01/23/06	MW-48 LSAS 03/21/06	MW-48 LSAS 10/10/06	MW-48 LSAS 12/18/06	MW-48 LSAS 09/13/07	MW-48 LSAS 01/30/08		
Semivolatile Organics																			
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	5.4	NA	NA	NA	NA	4 J	NA	NA	NA	NA	NA		
Semivolatile Organics (8270C)																			
1,4-Dioxane	ug/L	NA	2.1 U	1 U	NA	NA	8.5	3.7 [2.4]	NA	NA	NA	2.3 J [2.1 J]	1 U	1.1 U	NA	NA	NA		
Volatile Organics (8260B) - SIM																			
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	17 J	NA	NA	NA	NA	NA	R	3.9 J	NA		
Volatile Organics (8260) - SIM Isotope Dilution																			
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U [1 UJ]	NA	NA	NA	27	39	NA	NA	NA	NA	3.6 J	4.2	7.8	6.1
Volatile Organics (8260B)																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA		
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U [0.52 U]	0.52 U	1.7	1.7 [1.5]	2	2.2	0.52 U	0.5 U [0.5 U]	0.52 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U [0.45 U]	0.45 U	0.5 U	0.45 U [0.45 U]	0.83 U	0.45 U	0.45 U	0.5 U [0.5 U]	0.45 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U	0.65 U [0.65 U]	0.65 U	0.5 U	0.65 U [0.65 U]	0.75 U	0.65 U	0.65 U	0.5 U [0.5 U]	0.65 U	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U [0.5 U]	0.34 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.5 U	0.34 U	0.5 U [0.5 U]	0.34 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 UJ	0.5 U [0.5 U]	0.28 U	0.5 U	0.28 U [0.28 U]	0.71 U	0.5 U	0.28 U	0.5 U [0.5 U]	0.28 U	0.28 U	0.71 U	0.71 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-49 S&P Sand 01/12/05	MW-49 S&P Sand 06/17/05	MW-49 S&P Sand 01/30/06	MW-49 S&P Sand 12/13/06	MW-49 S&P Sand 01/23/08	MW-50 Lower AF 01/19/05	MW-50 Lower AF 06/17/05	MW-50 Lower AF 01/18/06	MW-50 Lower AF 12/12/06	MW-50 Lower AF 01/28/08
Semivolatile Organics										
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	2.5 U	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	2.1 U	1 U	NA	NA	NA	2.1 U	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	NA	0.86 I
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 UJ	0.65 U	0.18 U	0.5 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 UJ	0.5 U	0.11 U	0.5 U	0.28 U	0.71 UJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-51 Lower AF 01/20/05	MW-51 Lower AF 06/15/05	MW-51 Lower AF 01/19/06	MW-51 Lower AF 12/18/06	MW-51 Lower AF 01/28/08	MW-52 S&P Sand 01/13/05	MW-52 S&P Sand 06/20/05	MW-52 S&P Sand 01/18/06	MW-52 S&P Sand 03/24/06	MW-52 S&P Sand 12/18/06	MW-52 S&P Sand 01/28/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	2.5 U	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	2 U	1 U	NA	NA	2 U	1 U	1.1 U	NA	NA	
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	1.8 J	NA	
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	NA	0.76 U	1 U	
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U	0.65 U	0.18 U	0.5 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.5 UJ	0.18 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U	0.5 UJ	0.11 U	0.5 U	0.28 U	0.28 U	0.71 U	0.5 UJ

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-53 S&P Sand 01/12/05	MW-53 S&P Sand 06/20/05	MW-53 S&P Sand 01/18/06	MW-53 S&P Sand 03/24/06	MW-53 S&P Sand 12/18/06	MW-53 S&P Sand 01/28/08	MW-54 S&P Sand 01/12/05	MW-54 S&P Sand 06/20/05	MW-54 S&P Sand 02/21/06	MW-54 S&P Sand 12/18/06	MW-54 S&P Sand 01/28/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	1.9 U	1 U [1 U]	1.1 U	NA	NA	2 U	1 U	NA	NA	
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	1.1 J	NA	NA	NA	R	NA	
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	0.76 U	1 U [1 U]	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U [0.52 U]	0.52 U	0.6 U	0.52 U [0.52 U]	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U [0.45 U]	0.45 U	0.83 U	0.45 U [0.45 U]	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U [0.65 U]	0.65 U	0.75 U	0.65 U [0.65 U]	0.18 U	0.5 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U [0.34 U]	0.34 U	0.65 U	0.5 UJ [0.5 U]	0.18 U	0.5 U	0.34 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U [0.28 U]	0.28 U	0.71 U	0.5 UJ [0.5 U]	0.41 I	0.5 U	0.28 U	0.71 U	0.5 UJ

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-55 AF Gravel 01/12/05	MW-55 AF Gravel 06/16/05	MW-55 AF Gravel 07/19/05	MW-55 AF Gravel 01/23/06	MW-55 AF Gravel 02/28/06	MW-55 AF Gravel 03/22/06	MW-55 AF Gravel 04/04/06	MW-55 AF Gravel 10/11/06	MW-55 AF Gravel 12/13/06	MW-55 AF Gravel 01/29/08
Semivolatile Organics										
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	5.4	9.3	8.7 I	10	5.5 I	6.9 IJ	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	16 J	5.7 J
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	15 J	3
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	2.2	1.9	1.8	NA	3.6	2.6	1.4	0.67 I
1,1-Dichloroethene	ug/L	0.36 U	1.2	0.95 J	1	NA	1.9	1.3	0.85 I	0.83 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 J	0.63 J	0.65 U	NA	1	0.88 I	0.75 U	0.75 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.5 U	0.34 U	NA	0.34 U	0.34 U	0.65 U	0.65 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.5 U	0.28 U	NA	0.28 U	0.28 U	0.71 U	0.71 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-56 S&P Sand 01/18/05	MW-56 S&P Sand 06/16/05	MW-56 S&P Sand 01/21/06	MW-56 S&P Sand 10/12/06	MW-56 S&P Sand 12/07/06	MW-56 S&P Sand 01/29/08	MW-57 S&P Sand 01/12/05	MW-57 S&P Sand 06/27/05	MW-57 S&P Sand 01/30/06	MW-57 S&P Sand 03/27/06	MW-57 S&P Sand 10/09/06	MW-57 S&P Sand 12/06/06	MW-57 S&P Sand 07/11/07	MW-57 S&P Sand 09/11/07	MW-57 S&P Sand 01/30/08	
Semivolatile Organics																
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																
1,4-Dioxane	ug/L	NA	1.9 U	1 U	NA	NA	NA	2 U	17	1 U [1 UJ]	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																
1,4-Dioxane	ug/L	NA	NA	NA	R [R]	R	NA	NA	NA	NA	NA	R	R [R]	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ [2 U]	0.76 U	1 U	NA	NA	NA	NA	2 UJ	0.76 U [0.76 U]	0.69 U	0.69 UJ	1 U
Volatile Organics (8260B)																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.52 U	0.18 U	0.5 U	11	0.52 U [0.52 U]	0.6 U	0.6 UJ [0.6 U]	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.44 U	0.36 U	0.5 U	27	0.45 U [0.45 U]	0.83 U	0.83 UJ [0.83 U]	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U [0.75 U]	0.75 U	0.64 U	0.18 U	0.5 U	0.69 U	0.65 U [0.65 U]	0.75 UJ	0.75 UJ [0.75 U]	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.46 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.34 U [0.34 U]	0.65 U	0.65 UJ [0.65 U]	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U [0.71 U]	0.71 U	0.5 U	0.11 U	0.5 U	3.4	0.28 U [0.28 U]	0.71 U	0.71 UJ [0.71 U]	0.71 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-58 S&P Sand 01/13/05	MW-58 S&P Sand 06/23/05	MW-58 S&P Sand 01/23/06	MW-58 S&P Sand 10/10/06	MW-58 S&P Sand 12/06/06	MW-58 S&P Sand 01/30/08	MW-59 S&P Sand 01/12/05	MW-59 S&P Sand 06/15/05	MW-59 S&P Sand 01/20/06	MW-59 S&P Sand 12/18/06	MW-59 S&P Sand 01/28/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	2 U	1 U	NA	NA	NA	2 U	1 U	NA	NA	
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	R [R]	R	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ [2 UJ]	0.76 U	1 U [1 U]	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.52 U [0.52 U]	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.45 U [0.45 U]	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U [0.75 U]	0.75 U	0.65 U [0.65 U]	0.18 U	0.5 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.23 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.5 U [0.5 U]	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.2 U	0.5 U	0.28 U	0.71 U [0.71 U]	0.71 U	0.5 U [0.5 U]	0.11 U	0.5 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-60 S&P Sand 01/13/05	MW-60 S&P Sand 06/15/05	MW-60 S&P Sand 01/23/06	MW-60 S&P Sand 12/18/06	MW-60 S&P Sand 01/28/08	MW-61 S&P Sand 01/20/05	MW-61 S&P Sand 06/15/05	MW-61 S&P Sand 01/20/06	MW-61 S&P Sand 12/11/06	MW-61 S&P Sand 01/30/08	MW-62 USAS 01/20/05	MW-62 USAS 06/16/05	MW-62 USAS 01/26/06	MW-62 USAS 12/12/06	MW-62 USAS 01/31/08	
Semivolatile Organics																
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	2.5 U	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	
Semivolatile Organics (8270C)																
1,4-Dioxane	ug/L	NA	1.9 U	1.1 U	NA	NA	2 U	1 U	NA	NA	NA	1.9 U	1 U	NA	NA	
Volatile Organics (8260B) - SIM																
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	R	NA	NA	NA	R [R]	NA	
Volatile Organics (8260) - SIM Isotope Dilution																
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	2.5	NA	NA	NA	0.76 U	1 U [1.3]	NA	NA	NA	0.76 U [0.76 U]	1 U
Volatile Organics (8260B)																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U [0.52 U]	0.18 U	0.5 U	0.52 U	0.6 U [0.6 U]	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U [0.45 U]	0.36 U	0.5 U	0.45 U	0.83 U [0.83 U]	0.45 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U	0.65 U	0.18 U	0.5 U	0.65 U	0.75 UJ	0.65 U [0.65 U]	0.18 U	0.5 U	0.65 U	0.75 UJ [0.75 UJ]	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U [0.5 U]	0.18 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U	0.5 U	0.11 U	0.5 U	0.28 UJ	0.71 U	0.5 U [0.5 U]	0.11 U	0.5 U	0.29 U	0.71 UJ [0.71 UJ]	0.5 U

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-63 USAS 01/10/05	MW-63 USAS 06/16/05	MW-63 USAS 01/18/06	MW-63 USAS 10/12/06	MW-63 USAS 12/11/06	MW-63 USAS 01/31/08	MW-64 USAS 01/06/05	MW-64 USAS 06/17/05	MW-64 USAS 01/18/06	MW-64 USAS 02/28/06	MW-64 USAS 04/01/06	MW-64 USAS 12/18/06	MW-64 USAS 01/28/08	MW-65 USAS 01/07/05	MW-65 USAS 06/20/05	MW-65 USAS 01/20/06	MW-65 USAS 12/14/06	MW-65 USAS 01/28/08
Semivolatile Organics																		
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	15	NA	NA	NA	NA	NA	NA	1 U	NA	NA	NA	NA
Semivolatile Organics (8270C)																		
1,4-Dioxane	ug/L	NA	2.1 U	1 U	NA	NA	NA	16.4	23	17 J	24	NA	NA	NA	2 U	1 U	NA	NA
Volatile Organics (8260B) - SIM																		
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	NA	NA	44 J	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution																		
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	6.7	NA	NA	NA	NA	71	19	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U	0.94	4.8	4.4	NA	4.3	4.9	1.7	0.52 U	0.5 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	3	2.9	NA	3.1	4.4	0.79 I	0.45 U	0.5 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.71 I	0.5 U	0.65 U	0.75 U	0.75 UJ	0.65 U	0.65 U	4	4.1	NA	4.3	4.9	2	0.65 U	0.5 U	0.65 U	0.75 UJ
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	NA	0.34 U	0.65 U	0.5 UJ	0.34 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.53 I	0.5 U	0.28 U	0.71 U	0.71 UJ	0.5 U	5.5	16.1	17	NA	16	18	10 J	0.28 U	0.5 U	0.28 UJ	0.71 UJ

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-66 USAS 01/20/05	MW-66 USAS 06/16/05	MW-66 USAS 01/30/06	MW-66 USAS 12/13/06	MW-66 USAS 01/23/08	MW-67 USAS 01/20/05	MW-67 USAS 06/14/05	MW-67 USAS 01/20/06	MW-67 USAS 03/02/06	MW-67 USAS 04/03/06	MW-67 USAS 06/23/06	MW-67 USAS 10/11/06	MW-67 USAS 12/15/06	MW-67 USAS 01/28/08
Semivolatile Organics														
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	44	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)														
1,4-Dioxane	ug/L	NA	1.9 U	1 U	NA	NA	52	42	64	23	51	NA	NA	NA
Volatile Organics (8260B) - SIM														
1,4-Dioxane	ug/L	NA	NA	NA	1.1 J	NA	NA	NA	NA	NA	NA	30 J	96 J	NA
Volatile Organics (8260) - SIM Isotope Dilution														
1,4-Dioxane	ug/L	NA	NA	NA	0.76 UJ	1 U	NA	NA	NA	NA	NA	100 J	82	54
Volatile Organics (8260B)														
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U	74	79.7	55	NA	67	68	69	68
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U	120	88.4	92	NA	96	130	120	140
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 UJ	0.65 U	4.6 I	3	2.4	NA	3.1	2.3	3.2	2.2 J
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U	0.9 U	0.5 U	0.34 U	NA	0.34 U	0.5 U	0.65 U	0.65 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 UJ	0.5 U	14	11.1	14 J	NA	13	7	15	9.5 J

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-68 LSAS 01/07/05	MW-68 LSAS 06/15/05	MW-68 LSAS 01/20/06	MW-68 LSAS 03/21/06	MW-68 LSAS 12/18/06	MW-68 LSAS 01/28/08	MW-69 USAS 01/11/05	MW-69 USAS 06/20/05	MW-69 USAS 01/20/06	MW-69 USAS 12/11/06	MW-69 USAS 01/30/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	2.7 J	NA	NA	NA	NA	5.5	NA	NA	NA	NA	
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	2 U	1 U [1 U]	1.1 U	NA	NA	1.9 U	1 U	NA	NA	
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	3.3 J	NA	NA	NA	5.6 J	NA	
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	4.3	16	NA	NA	5.7	7.4	
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.52 U	0.5 U	0.52 U [0.52 U]	0.52 U	0.6 U	0.52 U	2.1	0.5 U	0.52 U	1.2	1.1
1,1-Dichloroethene	ug/L	0.45 U	0.5 U	0.86 I [0.72 I]	0.45 U	0.83 U	1.8	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U	0.75 I [0.69 I]	0.65 U	0.8 I	1.6	0.18 U	0.5 U	0.65 U	0.75 UJ	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U [0.34 U]	0.34 U	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	2.4	0.5 U	3.1 J [0.28 UJ]	2.3	1.8	5.2	0.43 I	0.5 U	0.28 UJ	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-70 USAS 01/11/05	MW-70 USAS 06/23/05	MW-70 USAS 02/02/06	MW-70 USAS 10/11/06	MW-70 USAS 12/07/06	MW-70 USAS 01/17/07	MW-70 USAS 04/24/07	MW-70 USAS 07/11/07	MW-70 USAS 09/11/07	MW-70 USAS 01/30/08	MW-71 USAS 01/11/05	MW-71 USAS 06/20/05	MW-71 USAS 01/21/06	MW-71 USAS 10/11/06	MW-71 USAS 12/14/06	MW-71 USAS 01/17/07	MW-71 USAS 04/25/07	MW-71 USAS 07/11/07	MW-71 USAS 09/12/07	MW-71 USAS 01/30/08	
Semivolatile Organics																					
1,4-Dioxane	ug/L	3.4 I	NA	NA	NA	NA	NA	NA	NA	NA	20	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																					
1,4-Dioxane	ug/L	NA	2.1 U	1 U	NA	NA	NA	NA	NA	NA	NA	5.3	1.1 U	NA	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																					
1,4-Dioxane	ug/L	NA	NA	NA	13 J	2.4 J [2.5 J]	NA	NA	NA	NA	NA	NA	NA	10 J	11 J	NA	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																					
1,4-Dioxane	ug/L	NA	NA	NA	4.6 J	1.4 I [3.2]	1.4 I	4.6	0.69 U	0.69 UJ	3.1	NA	NA	NA	13 J	7.6	8.4	8.2	6.2 J	5	5.6
Volatile Organics (8260B)																					
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	1	0.94 J	1.5	2.7	4.5 [4.5]	2.7 J	9.5	23	2.8	17	11	7.1	7.6	14	13	5.6 J	5.4	4.8	4.1	9.5
1,1-Dichloroethene	ug/L	1.8	0.5 U	3.2	5.5	5.1 [5]	2	3.4	8.1	0.83 U	4.6	9.3 I	5	6.5	4.3	4.9	1.2	1.9	2	1.4	2
cis-1,2-Dichloroethene	ug/L	0.36 I	0.5 U	0.65 U	0.75 U	0.75 U [0.75 U]	0.75 UJ	0.75 U	0.75 U	0.75 U	0.65 I	31	36.1	56	39	31 J	23 J	45	89	84	70
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U [0.65 U]	0.65 U	0.65 U	0.65 U	0.65 U	0.5 U	2.6 I	1 U	0.34 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	
Trichloroethene	ug/L	0.97 I	0.5 U	3.3	13	15 [15]	9.3	17	28	4	20	360	192	220	180	100 J	47	54	36	31	32

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-72 USAS 01/11/05	MW-72 USAS 06/22/05	MW-72 USAS 06/23/05	MW-72 USAS 01/24/06	MW-72 USAS 03/27/06	MW-72 USAS 10/09/06	MW-72 USAS 12/06/06	MW-72 USAS 01/17/07	MW-72 USAS 04/24/07	MW-72 USAS 07/11/07	MW-72 USAS 09/10/07	MW-72 USAS 01/30/08	MW-72 USAS 01/06/05	MW-73 USAS 06/22/05	MW-73 USAS 02/02/06	MW-73 USAS 10/12/06	MW-73 USAS 12/13/06	MW-73 USAS 01/25/08	
Semivolatile Organics																			
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	300	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)																			
1,4-Dioxane	ug/L	NA	2 U	2 U	1 U [1 U]	1 U	NA	NA	NA	NA	NA	NA	NA	NA	42.7	9.2 I	NA	NA	NA
Volatile Organics (8260B) - SIM																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	5.3 J	R	NA	NA	NA	NA	NA	NA	NA	NA	140 J	110 J	NA
Volatile Organics (8260) - SIM Isotope Dilution																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	2 UJ	0.76 U	0.76 U [0.76 U]	1.1 I	3.1	0.72 I	1 U	NA	NA	NA	180 J	110	84
Volatile Organics (8260B)																			
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.32 I	NA	0.5 U	0.52 U [0.52 U]	0.52 U	0.6 U	0.6 U	0.6 UJ [0.6 U]	0.6 U	0.6 U	0.6 U	0.52 U	210	64.9	82	44	54	53
1,1-Dichloroethene	ug/L	0.65 I	NA	0.5 U	0.45 U [0.45 U]	0.45 U	0.83 U	0.83 U	0.83 U [0.83 U]	1.2	0.83 U	0.83 U	0.45 U	370	98.5	130	89	110	91
cis-1,2-Dichloroethene	ug/L	0.5 I	NA	0.5 U	0.65 U [0.65 U]	0.65 U	0.75 U	0.75 UJ	0.75 UJ [0.75 U]	0.75 U	0.75 U	0.75 U	0.65 U	0.65 U	0.5 U	3.3 U	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	3.6	NA	3.3	2.8 [3.7]	3.6	4.4	4.8	4.1 J [4.1]	3.1	2.9	1.9	1.7	29	12.1	18	15	9.2	7.9
Trichloroethene	ug/L	4.5	NA	3.7	2.8 [3.8]	2.9	3.6	4	2.4 [2.4]	2.8	2.9	1.9	2.5	29	15.5	27	14	8.4	11

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-74 USAS 01/17/05	MW-74 USAS 06/22/05	MW-74 USAS 01/21/06	MW-74 USAS 02/28/06	MW-74 USAS 04/04/06	MW-74 USAS 06/21/06	MW-74 USAS 10/12/06	MW-74 USAS 12/14/06	MW-74 USAS 01/29/08	MW-75 USAS 01/07/05	MW-75 USAS 06/22/05	MW-75 USAS 02/22/06	MW-75 USAS 10/13/06	MW-75 USAS 12/13/06	MW-75 USAS 01/30/08	
Semivolatile Organics																
1,4-Dioxane	ug/L	230	NA	NA	NA	NA	NA	NA	NA	NA	41	NA	NA	NA	NA	
Semivolatile Organics (8270C)																
1,4-Dioxane	ug/L	NA	117 J	110	110 J	130	110	NA	NA	NA	NA	39.1	170 J	NA	NA	
Volatile Organics (8260B) - SIM																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	170 J	NA	NA	NA	NA	93 J	NA	
Volatile Organics (8260) - SIM Isotope Dilution																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	350 J	200	130	NA	NA	NA	260 J	110	33
Volatile Organics (8260B)																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	240 J	NA	NA	NA	NA	NA	NA	310 J	NA	NA
1,1-Dichloroethane	ug/L	300	234	210	NA	130	160	160	140	82	37	45.6	180	150	150	16
1,1-Dichloroethene	ug/L	150	119	120	NA	88	85	85	90	37	4.5	7.6	44	41	45	5.1
cis-1,2-Dichloroethene	ug/L	11	8.8	4.8	NA	5.3	4.4	5.3	3.8 J	3.3 U	0.65 U	2.9	9.5	9.3	5.6	0.65 U
Tetrachloroethene	ug/L	5 IV	0.5 U	0.68 U	NA	0.34 U	1 U	1.4	0.74 I	2.4 U	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	11 V	5	5.5	NA	5	4.1	6.2	3.2 J	2.4 U	0.29 J	0.5 U	0.49 I	0.74 I	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-76 USAS 01/11/05	MW-76 USAS 06/21/05	MW-76 USAS 01/25/06	MW-76 USAS 10/11/06	MW-76 USAS 12/13/06	MW-76 USAS 01/18/07	MW-76 USAS 04/24/07	MW-76 USAS 07/10/07	MW-76 USAS 09/11/07	MW-76 USAS 01/24/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	32	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	9	4.2 I	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	13 J	23 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	47 J	22	18	9.6	7.6 J	6.8 J	9.9
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	24 I	10.7	15	18 D	12	3.1 J	4.7	3.3	1.1	3.5
1,1-Dichloroethene	ug/L	23 I	7.9	13	15 D	13	2.4 J	3.9	4.6	1.2	3.1
cis-1,2-Dichloroethene	ug/L	13 I	4.2 J	3.8	7.2 D	3.8 I	2.3 J	2	0.93 I	0.75 U	0.65 U
Tetrachloroethene	ug/L	8.7 I	5.8	3.8 J	4.8 ID	5.4	4 J	5	4.8	6.6	3.1
Trichloroethene	ug/L	620	232	260	430 D	200	61 J	41	51	15	51

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-77 LSAS 01/12/05	MW-77 LSAS 06/21/05	MW-77 LSAS 01/19/06	MW-77 LSAS 03/02/06	MW-77 LSAS 04/03/06	MW-77 LSAS 10/16/06	MW-77 LSAS 12/18/06	MW-77 LSAS 01/18/07	MW-77 LSAS 04/24/07	MW-77 LSAS 07/10/07	MW-77 LSAS 09/12/07	MW-77 LSAS 01/24/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	120	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	NA	65.3	72	1 U	95	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	170 DJ	89 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	270 J	170	190	280	160 J	200	150 [220]
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	12	50.2	41	NA	36	15	50	23	33	45	30	39 [45]
1,1-Dichloroethene	ug/L	9.9	28.5	26	NA	20	13	26	8.7	14	34	17	18 [20]
cis-1,2-Dichloroethene	ug/L	1.8	2.3	2	NA	2	1.8	2.6	2.2	3.7	3.9	2.6	3.5 [3.8]
Tetrachloroethene	ug/L	0.18 U	1.7	1.6	NA	0.61 l	0.97 l	1.4	1.3	1.4	0.93 l	1.4	1.3 [1.4]
Trichloroethene	ug/L	1	9	12	NA	10	9.3	16	13	20	29	26	24 [26]

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-78 LSAS 01/10/05	MW-78 LSAS 06/24/05	MW-78 LSAS 01/31/06	MW-78 LSAS 10/12/06	MW-78 LSAS 12/13/06	MW-78 LSAS 01/18/07	MW-78 LSAS 04/24/07	MW-78 LSAS 07/10/07	MW-78 LSAS 09/11/07	MW-78 LSAS 01/28/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	290	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	86.5	140 [170]	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	86 J	100 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	210 J	140	200	130	140 J	170	99
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	320	382	400 [420]	490 D	480	160 J	170	490	300	340 D
1,1-Dichloroethene	ug/L	140	199	220 [220]	340 D	340	92 J	180	440	220	320 D
cis-1,2-Dichloroethene	ug/L	34	93.5	140 [140]	190 D	190	75 J	120	190	130	110
Tetrachloroethene	ug/L	2.4 I	2.5 U	3.8 [3.6]	4.8 ID	6.1	3.8 J	5.8	4.6 I	6.9	6.4 J
Trichloroethene	ug/L	13	29.9	55 [54]	76 D	67	33 J	59	68	70	48 J

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-79 LSAS 01/10/05	MW-79 LSAS 06/15/05	MW-79 LSAS 01/31/06	MW-79 LSAS 10/11/06	MW-79 LSAS 12/13/06	MW-79 LSAS 01/18/07	MW-79 LSAS 04/25/07	MW-79 LSAS 07/10/07	MW-79 LSAS 09/12/07	MW-79 LSAS 01/23/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	22	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	44.6 [44.8]	65 J	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	30 J	61 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	76 J	56	51	64	57 J [97 J]	57 [64]	52
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	1.8 U	5 U [5 U]	1.5	3 UD	3 U	3 U	3.6 I	5.5 [6 U]	3.1 I [3.7]	4
1,1-Dichloroethene	ug/L	5.9 I	13.4 [12.9]	5.6	6.5 D	8.2	4.2 U	8.5	10 [16]	4.7 I [4.4]	14
cis-1,2-Dichloroethene	ug/L	67	420 [417]	220	350 D	320	220	190	340 D [280]	220 [200]	140
Tetrachloroethene	ug/L	1.8 U	5 U [5 U]	0.34 U	3.2 UD	3.2 U	3.2 U	3.2 U	0.65 U [6.5 U]	3.2 U [0.65 U]	0.5 U
Trichloroethene	ug/L	310	827 [988]	220	190 D	180	140	130	120 [130]	81 [87]	100

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-80 LSAS 01/10/05	MW-80 LSAS 06/22/05	MW-80 LSAS 01/24/06	MW-80 LSAS 10/09/06	MW-80 LSAS 12/06/06	MW-80 LSAS 01/17/07	MW-80 LSAS 04/24/07	MW-80 LSAS 07/11/07	MW-80 LSAS 09/10/07	MW-80 LSAS 01/30/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	5	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	4.7 J	3.7 I [2.4 I]	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	R [R]	23 J	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	47 J [35 J]	31	48	64	61	72	95
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	29	51.8	21 [22]	21 [18]	23	11 J	12	21	13	17
1,1-Dichloroethene	ug/L	280	485	250 [240]	170 [160]	240 D	81 J	110	410 D	140	170 D
cis-1,2-Dichloroethene	ug/L	7.4 I	10.5	3.5 [3.6]	6.8 [6]	1.7 J	3.9 J	4.7	5.2	3.9	6.2
Tetrachloroethene	ug/L	4.6 I	10.8	6.8 [4.8]	7 [6.3]	9.9	5.7 J	4.6	8	5.1	4.5
Trichloroethene	ug/L	28	44.7	23 [22]	26 [23]	36	15 J	15	23	17	23

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-81 LSAS 01/11/05	MW-81 LSAS 06/23/05	MW-81 LSAS 02/01/06	MW-81 LSAS 10/10/06	MW-81 LSAS 12/15/06	MW-81 LSAS 01/18/07	MW-81 LSAS 04/25/07	MW-81 LSAS 07/10/07	MW-81 LSAS 09/12/07	MW-81 LSAS 01/30/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	41	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	30.1	27	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	20 J	68 J [79 J]	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	55 J	79 [79]	78	53	27 J	60	150
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	2.3	2.1	6.6	1.5	1.7 [2.5]	2.1 J	0.6 U	0.6 U	0.6 U	8.1
1,1-Dichloroethene	ug/L	9.6	2.3	46	2.3	3.5 J [7.2 J]	9.3	0.92 I	1.1	1.5	78
cis-1,2-Dichloroethene	ug/L	2.4	0.5 U	19	0.75 UJ	0.75 UJ [1.2 J]	6.8 J	0.75 U	0.75 U	0.75 U	49
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U [0.65 U]	0.65 U	0.65 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	22	2.1	240	0.89 I	0.71 J [15 J]	88	5	1.7	7.9	890 EJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-82 LSAS 01/17/05	MW-82 LSAS 06/16/05	MW-82 LSAS 01/21/06	MW-82 LSAS 03/01/06	MW-82 LSAS 04/04/06	MW-82 LSAS 06/22/06	MW-82 LSAS 10/12/06	MW-82 LSAS 12/14/06	MW-82 LSAS 01/29/08	MW-83 AF Gravel 01/20/05	MW-83 AF Gravel 06/21/05	MW-83 AF Gravel 01/19/06	MW-83 AF Gravel 03/22/06	MW-83 AF Gravel 10/18/06	MW-83 AF Gravel 12/14/06	MW-83 AF Gravel 01/24/08	
Semivolatile Organics																	
1,4-Dioxane	ug/L	180	NA	NA	NA	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	NA	131	280	180 J	200	220 D	NA	NA	NA	2 U	1 U	1 U	NA	NA	NA	
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	430 J	NA	NA	NA	NA	NA	R [R]	R	NA	
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	680 J	560	580	NA	NA	NA	NA	2 UJ [2 UJ]	0.76 U	1 U	
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	440 J	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	25	57.6	100	NA	54	26	96	110	73	0.18 U	0.5 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	5.6	15.3	40	NA	26	13	45	52	17	0.36 U	0.5 U	0.45 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	1	1.9	3.7	NA	2.3	0.87 I	4.3	4.4 J	6.4 U	0.18 U	0.5 U	0.65 U	0.65 U	0.75 U [0.75 U]	0.75 UJ	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	NA	0.34 U	0.5 U	0.65 U	0.65 U	5 U	0.18 U	0.5 U	0.34 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	NA	0.28 U	0.5 U	0.71 U	0.71 UJ	5 U	0.11 U	0.5 U	0.28 U	0.28 U	0.71 U [0.71 U]	0.71 UJ	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-84 LSAS 01/17/05	MW-84 LSAS 06/14/05	MW-84 LSAS 06/22/05	MW-84 LSAS 01/24/06	MW-84 LSAS 03/20/06	MW-84 LSAS 10/10/06	MW-84 LSAS 12/06/06	MW-84 LSAS 01/17/07	MW-84 LSAS 04/24/07	MW-84 LSAS 07/11/07	MW-84 LSAS 09/11/07	MW-84 LSAS 01/25/08	MW-84 LSAS 01/17/05	MW-85 LSAS 06/21/05	MW-85 LSAS 01/20/06	MW-85 LSAS 02/28/06	MW-85 LSAS 04/03/06	MW-85 LSAS 10/12/06	MW-85 LSAS 12/13/06	MW-85 LSAS 01/24/08	
Semivolatile Organics																					
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	230	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																					
1,4-Dioxane	ug/L	NA	42.9	2 U	1 U	1 U	NA	NA	NA	NA	NA	NA	NA	27.6	150	140	190	NA	NA	NA	
Volatile Organics (8260B) - SIM																					
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	1.1 J	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	300 J	NA	
Volatile Organics (8260) - SIM Isotope Dilution																					
1,4-Dioxane	ug/L	NA	NA	NA	NA	2.4 J	0.76 U	1.5 I	2.3	3.8	2.1 J [2.2]	1 U	NA	NA	NA	NA	NA	490 J	330	590	
Volatile Organics (8260B)																					
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	430 J	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	3.8	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.6 UJ	0.6 U	0.6 U	0.6 U [0.6 U]	0.52 U	82	46.5	53	NA	88	67	91	98
1,1-Dichloroethene	ug/L	0.36 U	20.7	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.83 U	0.83 U	0.83 U	0.83 U [0.83 U]	0.45 U	68	32.3	36	NA	76	66	94	92
cis-1,2-Dichloroethene	ug/L	0.18 U	9.7	0.5 U	0.65 U	0.65 U	0.75 U	0.75 UJ	0.75 UJ	0.75 U	0.75 U	0.75 U [0.75 U]	0.65 U	49	24.6	22	NA	34	35	37	34
Tetrachloroethene	ug/L	0.18 U	1 U	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U [0.65 U]	0.5 U	1.1 IV	0.5 U	0.34 U	NA	0.34 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.11 U	109	0.5 U	0.28 U	0.28 U	0.71 U	0.71 U	0.71 U	0.71 U	0.71 U	0.71 U [0.71 U]	0.5 U	6.9 V	2.1	4.1 J	NA	9.4	9.2	6.2	7.6

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-86 LSAS 01/18/05	MW-86 LSAS 06/14/05	MW-86 LSAS 02/20/06	MW-86 LSAS 10/11/06	MW-86 LSAS 12/18/06	MW-86 LSAS 01/28/08	MW-87 LSAS 01/19/05	MW-87 LSAS 06/23/05	MW-87 LSAS 02/22/06	MW-87 LSAS 10/13/06	MW-87 LSAS 12/13/06	MW-87 LSAS 01/18/07	MW-87 LSAS 04/25/07	MW-87 LSAS 07/10/07	MW-87 LSAS 09/13/07	MW-87 LSAS 01/30/08	
Semivolatile Organics																	
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	400	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	NA	6.8	13	NA	NA	NA	184	200 J	NA	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	6.7 J	12 J	NA	NA	NA	NA	380 J	NA	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	16 J	18	13	NA	NA	NA	500 J [500 DJ]	520	310	630	340 J	270	400 DJ
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	560 DJ [420 J]	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.18 U	4.9	5.5	2.8	3.1	1.6	460	142	220	130 D [130]	240	140	110	170	110	190 EJ
1,1-Dichloroethene	ug/L	0.36 U	4.9	5.1	3.2	4	1.8	1,700	443	870	510 D [700 D]	990	560	490	1,100	470	970 EJ
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U	0.75 U	0.65 U	32.1	10.4	14	8.7 D [8.6]	9	7.5 U	5.4	7.5 U	7.5 U	8.5
Tetrachloroethene	ug/L	0.46 IV	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U	150	62.9	83	110 D [99]	91	84	68	76	67	61
Trichloroethene	ug/L	0.64 IV	0.5 U	0.28 U	0.71 U	0.71 U	0.5 U	1,000	381	400	460 D [480 D]	380	250	280	380	320	360 EJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-88 Zone 1 Clay 01/20/05	MW-88 Zone 1 Clay 06/16/05	MW-88 Zone 1 Clay 01/21/06	MW-88 Zone 1 Clay 10/12/06	MW-88 Zone 1 Clay 12/07/06	MW-88 Zone 1 Clay 01/29/08	MW-89 USAS 01/13/05	MW-89 USAS 06/16/05	MW-89 USAS 01/27/06	MW-89 USAS 12/13/06	MW-89 USAS 01/23/08	MW-90 USAS 01/19/05	MW-90 USAS 06/17/05	MW-90 USAS 01/25/06	MW-90 USAS 12/14/06	MW-90 USAS 01/23/08
Semivolatile Organics																
1,4-Dioxane	ug/L	2.5 U	NA	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	2.5 U	NA	NA	NA
Semivolatile Organics (8270C)																
1,4-Dioxane	ug/L	NA	1.9 U	1 U	NA	NA	NA	2.1 U	1 U [1 U]	NA	NA	NA	2 U	1 U [1 U]	NA	NA
Volatile Organics (8260B) - SIM																
1,4-Dioxane	ug/L	NA	NA	NA	R [R]	R	NA	NA	NA	NA	3.6 J	NA	NA	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution																
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ [2 UJ]	0.76 U	1 U	NA	NA	NA	2.6	2.9	NA	NA	NA	0.76 U
Volatile Organics (8260B)																
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.52 U	0.71 I	0.67 J	0.6 I [0.52 U]	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U [0.52 U]	0.6 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.44 U	0.36 U	0.5 U	0.45 U [0.45 U]	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U [0.45 U]	0.83 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U [0.75 U]	0.75 U	0.64 U	0.18 U	0.5 U	0.65 U [0.65 U]	0.75 UJ	0.65 U	0.18 U	0.5 U	0.65 U [0.65 U]	0.75 UJ
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U [0.34 U]	0.65 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U [0.71 U]	0.71 U	0.5 U	0.2 I	0.5 U	0.28 U [0.28 U]	0.71 UJ	0.5 U	0.11 U	0.5 U	0.28 U [0.28 U]	0.71 UJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-91 LSAS 01/19/05	MW-91 LSAS 06/17/05	MW-91 LSAS 01/30/06	MW-91 LSAS 10/10/06	MW-91 LSAS 12/18/06	MW-91 LSAS 01/28/08	MW-92 LSAS 01/19/05	MW-92 LSAS 06/16/05	MW-92 LSAS 01/18/06	MW-92 LSAS 10/12/06	MW-92 LSAS 12/12/06	MW-92 LSAS 01/31/08	MW-93 LSAS 01/20/05	MW-93 LSAS 06/16/05	MW-93 LSAS 01/18/06	MW-93 LSAS 06/27/06	MW-93 LSAS 10/12/06	MW-93 LSAS 12/12/06	MW-93 LSAS 01/28/08	
Semivolatile Organics																				
1,4-Dioxane	ug/L	51	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	NA	2.5 U	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																				
1,4-Dioxane	ug/L	NA	30.2	36	NA	NA	NA	2.1 U [1.9 U]	1 U	NA	NA	NA	NA	1.9 U	1 U	1.7 I	NA	NA	NA	
Volatile Organics (8260B) - SIM																				
1,4-Dioxane	ug/L	NA	NA	NA	29 J	58 J	NA	NA	NA	NA	R	R	NA	NA	NA	NA	NA	61 J	R	NA
Volatile Organics (8260) - SIM Isotope Dilution																				
1,4-Dioxane	ug/L	NA	NA	NA	130 DJ	89	85	NA	NA	NA	2 UJ	0.76 U	1 U	NA	NA	NA	NA	2 UJ	0.76 U	1 U
Volatile Organics (8260B)																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	2.3	2.8	3.4	5.8	5.5	3.8	0.18 U	0.5 U [0.5 U]	0.52 U	0.6 U	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	3.2	2.7	3.6	4.6	5.3	3.1	0.36 U	0.5 U [0.5 U]	0.45 U	0.83 U	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	5.4	5.9	4.4	9.8	7.2	7	0.18 U	0.5 U [0.5 U]	0.65 U	0.75 U	0.75 U	0.65 U	0.18 U	0.5 U	0.65 U	0.65 U	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 UJ	0.18 U	0.5 U [0.5 U]	0.34 U	0.65 U	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.5 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	3.5	5	6.9	22	14	16 J	0.11 U	0.5 U [0.5 U]	0.28 U	0.71 U	0.71 UJ	0.5 U	0.11 U	0.5 U	0.28 U	0.5 U	0.71 U	0.71 UJ	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-94 USAS 01/20/05	MW-94 USAS 06/21/05	MW-94 USAS 01/26/06	MW-94 USAS 10/18/06	MW-94 USAS 12/14/06	MW-94 USAS 01/24/08	MW-95 USAS 01/20/05	MW-95 USAS 06/21/05	MW-95 USAS 06/21/05	MW-95 USAS 01/26/06	MW-95 USAS 12/19/06	MW-95 USAS 01/24/08	MW-96 Zone 3-4 Clay 02/16/05	MW-96 Zone 3-4 Clay 06/17/05	MW-96 Zone 3-4 Clay 01/26/06	MW-96 Zone 3-4 Clay 12/15/06	MW-96 Zone 3-4 Clay 01/29/08	
Semivolatile Organics																		
1,4-Dioxane	ug/L	3.8 I	NA	NA	NA	NA	11	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)																		
1,4-Dioxane	ug/L	NA	2.9 J	1 U	NA	NA	NA	2 U	2 U	1 U	NA	NA	2.5 U	2 U	1 U	NA	NA	NA
Volatile Organics (8260B) - SIM																		
1,4-Dioxane	ug/L	NA	NA	NA	10 J	9.7 J	NA	NA	NA	NA	NA	2.7 J	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution																		
1,4-Dioxane	ug/L	NA	NA	NA	13 J	11	19	NA	NA	NA	NA	3	35	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U	0.18 U	0.5 U	0.5 U	0.52 U	0.6 U	0.52 U	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.83 U	0.45 U	0.36 U	0.5 U	0.5 U	0.45 U	0.83 U	0.45 U	0.36 U	0.5 U	0.45 U	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.18 U	0.5 U	0.65 U	0.75 U	0.75 UJ	0.65 U	0.18 U	0.5 U	0.5 U	0.65 U	0.75 UJ	0.65 U	NA	0.5 U	0.65 U	0.75 U	0.64 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 U	0.18 U	0.5 U	0.5 U	0.34 U	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U	0.71 UJ	0.5 U	0.16 I	0.5 U	0.5 U	0.28 U	0.71 UJ	0.5 U	0.11 U	0.5 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-97 Zone 3-4 Clay 02/16/05	MW-97 Zone 3-4 Clay 06/21/05	MW-97 Zone 3-4 Clay 02/16/06	MW-97 Zone 3-4 Clay 12/12/06	MW-97 Zone 3-4 Clay 01/30/08	MW-98 LSAS 02/15/05	MW-98 LSAS 06/16/05	MW-98 LSAS 02/02/06	MW-98 LSAS 10/12/06	MW-98 LSAS 12/13/06	MW-98 LSAS 01/25/08
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	2.5 U	2 U	R	NA	NA	470	205	420 Q	NA	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	650 J	NA
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	NA	1,200 J	900
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	940 DJ	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.52 U	190	149	200	170 D	35
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.45 U	390	332	440	420 D	86
cis-1,2-Dichloroethene	ug/L	NA	0.5 U	0.65 U	0.75 U	0.65 U	NA	5 U	4.7	5.6 D	0.75 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 UJ	0.65 U	0.5 U	5.6	8.9 J	13	21 D	2.4
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 UJ	0.5 U	65 D	27.2	73	72 D	7.7

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-99 Zone 1 Clay 02/15/05	MW-99 Zone 1 Clay 06/21/05	MW-99 Zone 1 Clay 06/21/05	MW-99 Zone 1 Clay 01/23/06	MW-99 Zone 1 Clay 02/28/06	MW-99 Zone 1 Clay 03/15/06	MW-99 Zone 1 Clay 04/01/06	MW-99 Zone 1 Clay 12/14/06	MW-99 Zone 1 Clay 01/25/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	2.5 U	2 U	2 U	3.9 I	1 U	1.7 I	1 U	NA	
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	1.2 J	
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	0.76 U	1 U [1 U]
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.45 I	0.5 U	NA	4.6	NA	0.52 U	0.52 U	0.6 U	0.52 U [0.52 U]
1,1-Dichloroethene	ug/L	0.67 I	0.5 U	NA	0.63 I	NA	0.45 U	0.45 U	0.83 U	0.45 U [0.45 U]
cis-1,2-Dichloroethene	ug/L	NA	0.5 U	NA	0.65 U	NA	0.65 U	0.65 U	0.75 UJ	0.65 U [0.65 U]
Tetrachloroethene	ug/L	0.18 U	0.5 U	NA	0.34 U	NA	0.34 U	0.34 U	0.65 U	0.5 U [0.5 U]
Trichloroethene	ug/L	0.11 I	0.5 U	NA	0.28 U	NA	0.28 U	0.28 U	0.71 UJ	0.5 U [0.5 U]

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-100 USAS 02/15/05	MW-100 USAS 06/22/05	MW-100 USAS 01/23/06	MW-100 USAS 01/31/06	MW-100 USAS 02/28/06	MW-100 USAS 03/15/06	MW-100 USAS 04/01/06	MW-100 USAS 10/12/06	MW-100 USAS 12/14/06	MW-100 USAS 01/25/08	MW-101 LSAS 02/14/05	MW-101 LSAS 06/16/05	MW-101 LSAS 01/23/06	MW-101 LSAS 10/11/06	MW-101 LSAS 12/13/06	MW-101 LSAS 01/29/08	
Semivolatile Organics																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	14	3.2 J	2.9 I	8.7 I [14]	7.8 I	12	8.2 I	NA	NA	NA	2.5 U	2 U	1 U	NA	NA	
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	R	8.9 J	NA	NA	NA	NA	20 J	1.3 J	NA
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	32 J	12	28	NA	NA	NA	19 J	0.76 U	2.5
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	3.1	4	4.5	4.3 [4.6]	NA	4.9	4.4	7.4	3.8	7.5	0.18 U	0.5 U	0.91 I	2	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.86 I	1.1	1.6	0.45 U [1.4]	NA	4.6	1.1	2.2	1.4	2.2	0.36 U	0.5 U	0.55 I	1.1	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	NA	0.5 U	0.65 U	0.65 U [0.65 U]	NA	0.65 U	0.65 U	0.75 U	0.75 UJ	0.65 U	NA	0.5 U	0.65 U	0.89 I	0.75 U	0.64 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.34 U [0.34 U]	NA	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.28 U [0.28 U]	NA	0.28 U	0.28 U	0.71 U	0.71 UJ	0.5 U	0.11 U	0.5 U	0.28 U	0.71 U	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-102 AF Gravel 02/15/05	MW-102 AF Gravel 06/22/05	MW-102 AF Gravel 02/01/06	MW-102 AF Gravel 12/19/06	MW-102 AF Gravel 01/25/08	MW-103 USAS 02/15/05	MW-103 USAS 06/21/05	MW-103 USAS 02/01/06	MW-103 USAS 12/19/06	MW-103 USAS 01/25/08	MW-104 USAS 02/14/05	MW-104 USAS 06/21/05	MW-104 USAS 02/21/06	MW-104 USAS 10/11/06	MW-104 USAS 12/18/06	MW-104 USAS 01/28/08	
Semivolatile Organics																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	2.5 U	2 U	1 U [1 U]	NA	NA	2.5 U	2 U [2 UR]	1 U [1 U]	NA	NA	110	93.6	140	NA	NA	
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	0.96 J	NA	NA	NA	NA	1.9 J	NA	NA	NA	NA	170 DJ	140 J [160 J]	NA
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	NA	1.5 I	1 U	NA	NA	NA	230 J	220 [170 J]	91
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U [0.52 U]	0.6 U	0.52 U	0.18 U	0.5 U [0.5 U]	0.52 U [0.52 U]	0.6 U	0.52 U	25	32.7	38	37	32 [36]	23
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U [0.45 U]	0.83 U	0.45 U	0.36 U	0.5 U [0.5 U]	0.45 U [0.45 U]	0.83 U	0.45 U	14	24.6	32	38	38 [41]	25
cis-1,2-Dichloroethene	ug/L	NA	0.5 U	0.65 U [0.65 U]	0.75 U	0.65 U	NA	0.5 U [0.5 U]	0.65 U [0.65 U]	0.75 U	0.65 U	NA	0.5 U	0.65 U	0.75 U	0.75 U [0.75 U]	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.5 U	0.18 U	0.5 U [0.5 U]	0.34 U [0.34 U]	0.65 U	0.5 U	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U [0.65 U]	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U [0.28 U]	0.71 U	0.5 U	0.11 U	0.5 U [0.5 U]	0.28 U [0.28 U]	0.71 U	0.5 U	1.3	1.2	1.8	1.6	1.1 [1.1]	0.85 J

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-105 LSAS 02/14/05	MW-105 LSAS 06/23/05	MW-105 LSAS 02/21/06	MW-105 LSAS 10/11/06	MW-105 LSAS 12/18/06	MW-105 LSAS 01/28/08	MW-106 LSAS 03/21/05	MW-106 LSAS 06/16/05	MW-106 LSAS 01/23/06	MW-106 LSAS 10/11/06	MW-106 LSAS 12/07/06	MW-106 LSAS 01/24/08	MW-107 USAS 04/06/05	MW-107 USAS 06/23/05	MW-107 USAS 03/01/06	MW-107 USAS 12/19/06	MW-107 USAS 01/24/08	
Semivolatile Organics																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)																		
1,4-Dioxane	ug/L	2.5 U	1.9 U	1.1 U	NA	NA	NA	2.2 U	2 U	1 U	NA	NA	NA	2.3 U	2 U	1 U	NA	NA
Volatile Organics (8260B) - SIM																		
1,4-Dioxane	ug/L	NA	NA	NA	R	0.95 J	NA	NA	NA	NA	R	R [R]	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution																		
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	1 U	NA	NA	NA	2 UJ	0.76 U [0.76 U]	1 U	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)																		
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.18 U	0.5 U	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.36 U	0.5 U	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	NA	0.5 U	0.65 U	0.75 U	0.75 U	0.65 U	0.65 U	0.5 U	0.65 U	0.75 U	0.75 U [0.75 U]	0.65 U	0.65 U	0.5 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.18 U	0.5 U	0.34 U	0.65 U	0.65 U	0.5 UJ	0.34 U	0.5 U	0.34 U	0.65 U	0.65 U [0.65 U]	0.5 U	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.11 U	0.5 U	0.28 U	0.71 U	0.71 U	0.5 UJ	0.28 U	0.5 U	0.28 U	0.71 U	0.71 U [0.71 U]	0.5 U	0.28 U	0.5 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-108 USAS 03/21/05	MW-108 USAS 06/15/05	MW-108 USAS 01/20/06	MW-108 USAS 03/20/06	MW-108 USAS 12/15/06	MW-108 USAS 01/29/08	MW-109 USAS 03/21/05	MW-109 USAS 06/15/05	MW-109 USAS 01/31/06	MW-109 USAS 03/24/06	MW-109 USAS 10/10/06	MW-109 USAS 12/18/06	MW-109 USAS 01/28/08	MW-110 USAS 03/21/05	MW-110 USAS 06/14/05	MW-110 USAS 02/20/06	MW-110 USAS 03/24/06	MW-110 USAS 12/18/06	MW-110 USAS 01/28/08	
Semivolatile Organics																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																				
1,4-Dioxane	ug/L	9.4 J [4.5 J]	4 J [4.1 J]	1 U	2.7 I	NA	NA	12	7.2	2.8 I	9.1 I	NA	NA	NA	33	26.6	37	69	NA	NA
Volatile Organics (8260B) - SIM																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	9.9 J	NA	NA	NA	NA	NA	R	8.5 J	NA	NA	NA	NA	NA	50 J	NA
Volatile Organics (8260) - SIM Isotope Dilution																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	11	12	NA	NA	NA	NA	18 J	10	15	NA	NA	NA	NA	80	62
Volatile Organics (8260B)																				
1,4-Dioxane	ug/L	NA	NA	NA	NA	25 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.88 J [0.97 J]	1.4 [1.3]	1.4	1.3	2.2	1.3	0.83 J	1.2	1.2	1.1	1.2	1.1	0.77 I	10	19.3	24	19	19	15
1,1-Dichloroethene	ug/L	0.45 U [0.45 U]	0.5 U [0.5 U]	0.45 U	0.45 U	0.83 U	0.44 U	0.45 U	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.45 U	6.4	20.5	22	19	19	15
cis-1,2-Dichloroethene	ug/L	0.65 U [0.65 U]	0.5 U [0.5 U]	0.65 U	0.65 U	0.75 U	0.64 U	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.75 U	0.65 U	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U [0.34 U]	0.5 U [0.5 U]	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U [0.28 U]	0.5 U [0.5 U]	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.5 U	0.28 U	0.28 U	0.71 U	0.71 U	0.5 U	0.28 U	0.5 U	0.28 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-111	MW-111	MW-111	MW-111	MW-112	MW-112	MW-112	MW-112	MW-112	MW-112	MW-113	MW-113	MW-113	MW-113	MW-113
Zone:	USAS	USAS	USAS	USAS	Zone 1 Clay	Zone 1 Clay	Zone 1 Clay	Zone 1 Clay	Zone 1 Clay	Zone 1 Clay	LSAS	LSAS	LSAS	LSAS	LSAS
Date Collected:	03/21/05	06/16/05	01/19/06	01/28/08	03/22/05	06/28/05	01/21/06	12/07/06	01/24/08	03/21/05	06/16/05	01/19/06	12/11/06	01/28/08	
Semivolatile Organics															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)															
1,4-Dioxane	ug/L	2.2 U	2 U	1 U	NA	NA	2 U	1 U	NA	NA	2.2 U	2 U	1 U	NA	NA
Volatile Organics (8260B) - SIM															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	R	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution															
1,4-Dioxane	ug/L	NA	NA	NA	1 U	NA	NA	NA	0.76 U	1 U	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)															
1,4-Dioxane	ug/L	NA	NA	NA	NA	2.2 U	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U	0.52 U	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.5 U	0.45 U	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U	0.65 U	0.65 U	0.65 U	0.5 U	0.65 U	0.75 U	0.65 U	0.65 U	0.5 U	0.65 U	0.75 UJ	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.5 UJ	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	0.65 U	0.5 UJ
Trichloroethene	ug/L	0.28 U	0.5 U	0.28 U	0.5 UJ	0.28 U	0.5 U	0.28 U	0.71 U	0.5 U	0.28 U	0.5 U	0.28 U	0.71 U	0.5 UJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-114 USAS 04/06/05	MW-114 USAS 06/21/05	MW-114 USAS 01/25/06	MW-114 USAS 10/19/06	MW-114 USAS 12/14/06	MW-114 USAS 01/24/08	MW-115 USAS 05/27/05	MW-115 USAS 06/15/05	MW-115 USAS 01/28/06	MW-115 USAS 12/08/06	MW-115 USAS 01/29/08
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	2.3 U	2.1 U [2 U]	1 U	NA	NA	NA	1 U	2 U	1 U	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	NA	R [R]
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	1.1	NA	NA	NA	0.76 U [0.76 U]
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U [0.5 U]	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U [0.6 U]
1,1-Dichloroethene	ug/L	0.45 U	0.5 U [0.5 U]	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U [0.83 U]
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U [0.5 U]	0.65 U	0.75 U	0.75 UJ	0.65 U	0.65 U	0.5 U	0.65 U	0.75 U [0.75 U]
Tetrachloroethene	ug/L	0.34 U	0.5 U [0.5 U]	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	0.65 U [0.65 U]
Trichloroethene	ug/L	0.28 U	0.5 U [0.5 U]	0.28 U	0.71 U	0.71 UJ	0.5 U	0.28 U	0.8 J	0.28 U	0.71 U [0.71 U]

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-116 USAS 05/27/05	MW-116 USAS 06/15/05	MW-116 USAS 01/27/06	MW-116 USAS 12/14/06	MW-116 USAS 01/25/08	MW-117 LSAS 05/26/05	MW-117 LSAS 06/30/05	MW-117 LSAS 04/03/06	MW-117 LSAS 12/27/06	MW-117 LSAS 01/25/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	1.9 U	1 U	NA	NA	1 U	2 U	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	NA	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.5 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.5 U	0.28 U	0.71 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-118 USAS 05/26/05	MW-118 USAS 06/30/05	MW-118 USAS 03/31/06	MW-118 USAS 12/27/06	MW-118 USAS 01/25/08	MW-119 LSAS 05/26/05	MW-119 LSAS 06/30/05	MW-119 LSAS 03/31/06	MW-119 LSAS 04/22/06	MW-119 LSAS 12/27/06	MW-119 LSAS 01/25/08	MW-120 USAS 05/26/05	MW-120 USAS 06/30/05	MW-120 USAS 03/31/06	MW-120 USAS 12/27/06	MW-120 USAS 01/25/08	
Semivolatile Organics																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)																	
1,4-Dioxane	ug/L	1 U	2 U	1 U	NA	NA	1 U	2 U	63	1.1 U [1.1 U]	NA	NA	1.5 J	2 U	1 U	NA	NA
Volatile Organics (8260B) - SIM																	
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	NA	R [R]	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution																	
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	2.3	NA	NA	NA	NA	0.76 U [0.76 U]	1 U	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)																	
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U	0.52 U	0.5 U	0.52 U	NA	0.6 U [0.6 U]	0.52 U	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U	0.45 U	0.5 U	0.45 U	NA	0.83 U [0.83 U]	0.45 U	0.45 U	0.5 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U	0.65 U	0.75 U	0.65 U	0.65 U	0.5 U	0.65 U	NA	0.75 U [0.75 U]	0.65 U	0.65 U	0.5 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.5 U	0.34 U	NA	0.65 U [0.65 U]	0.5 U	0.34 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.5 U	0.28 U	0.71 U	0.5 U	0.28 U	0.5 U	0.28 U	NA	0.71 U [0.71 U]	0.5 U	0.28 U	0.5 U	0.28 U	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-121 USAS 05/27/05	MW-121 USAS 06/15/05	MW-121 USAS 02/18/06	MW-121 USAS 03/21/06	MW-121 USAS 12/11/06	MW-121 USAS 01/23/08	MW-122 USAS 05/27/05	MW-122 USAS 06/15/05	MW-122 USAS 01/23/06	MW-122 USAS 03/21/06	MW-122 USAS 12/18/06	MW-122 USAS 01/25/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	1 U	2 U [1.9 U]	1 U [1.1 U]	1 U	NA	NA	1.1 U [1 U]	2 U	1 U	1 U	NA	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	NA	R	NA	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	NA	0.76 U	1 U	NA	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.5 U [0.5 U]	0.52 U [0.52 U]	0.52 U	0.6 U	0.52 U	0.52 U [0.52 U]	0.5 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.5 U [0.5 U]	0.45 U [0.45 U]	0.45 U	0.83 U	0.45 U	0.45 U [0.45 U]	0.5 U	0.45 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.5 U [0.5 U]	0.65 U [0.65 U]	0.65 U	0.75 UJ	0.65 U	0.65 U [0.65 U]	0.5 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.5 U [0.5 U]	0.34 U [0.34 U]	0.34 U	0.65 U	0.5 U	0.34 U [0.34 U]	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.5 U [0.5 U]	0.28 U [0.28 U]	0.28 U	0.71 U	0.5 U	0.28 U [0.28 U]	0.5 U	0.28 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-123 Floridan 06/28/05	MW-123 Floridan 07/21/05	MW-123 Floridan 02/20/06	MW-123 Floridan 03/03/06	MW-123 Floridan 03/08/06	MW-123 Floridan 03/10/06	MW-123 Floridan 03/14/06	MW-123 Floridan 03/16/06	MW-123 Floridan 03/21/06	MW-123 Floridan 12/19/06	MW-123 Floridan 02/01/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	2 U	1.9 U	1 U	1 U	1 U	29 I	1 U	1 U	1 U	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	0.76 U	1 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.5 U	3	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.52 J	7	1.7	0.45 U	0.45 U	0.45 U	0.45 U	0.45 U	0.45 U	1.4	0.45 U
cis-1,2-Dichloroethene	ug/L	0.5 U	0.5 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.5 U	0.5 U	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.5 U	0.51 J	8.2	0.43 I	0.28 U	1.6	0.28 U	0.28 U	0.28 U	2.2	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-124	MW-124	MW-124	MW-124	MW-125	MW-125	MW-125	MW-125
Zone:	Zone 2 Clay	Zone 2 Clay	Zone 2 Clay	Zone 2 Clay	Venice Clay	Venice Clay	Venice Clay	Venice Clay
Date Collected:	02/15/06	12/07/06	01/24/08	07/22/05	01/28/06	12/08/06	01/29/08	07/25/05
Semivolatile Organics								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)								
1,4-Dioxane	ug/L	1 U	NA	NA	2 U	1 U	NA	2.1 U
Volatile Organics (8260B) - SIM								
1,4-Dioxane	ug/L	NA	R	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution								
1,4-Dioxane	ug/L	NA	0.76 U	2.3	NA	NA	0.76 U	1 U
Volatile Organics (8260B)								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.5 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.5 U	0.45 UJ	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.65 U	0.5 U	0.65 U	0.75 U	0.64 U
Tetrachloroethene	ug/L	0.34 UJ	0.65 U	0.5 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.5 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-126 USAS 01/21/06	MW-126 USAS 10/11/06	MW-126 USAS 12/07/06	MW-126 USAS 01/24/08	MW-126 USAS 07/22/05	MW-127 AF Gravel 10/08/05	MW-127 AF Gravel 01/24/06	MW-127 AF Gravel 10/11/06	MW-127 AF Gravel 12/05/06	MW-127 AF Gravel 01/17/07	MW-127 AF Gravel 04/25/07	MW-127 AF Gravel 07/11/07	MW-127 AF Gravel 09/10/07	MW-127 AF Gravel 01/31/08	
Semivolatile Organics															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)															
1,4-Dioxane	ug/L	1 U	NA	NA	NA	2 U	73.6 [59.8]	150	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM															
1,4-Dioxane	ug/L	NA	R	R	NA	NA	NA	110 J	330 J	NA	NA	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution															
1,4-Dioxane	ug/L	NA	2 UJ	0.76 U	1 U	NA	NA	200 J	270	270	450	290	310	300 DJ	
Volatile Organics (8260B)															
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.6 U	0.52 U	0.5 U	52.2 J [51.4]	59	36 ID	44 I	29 J	40	38	29	34
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.83 U	0.45 U	0.5 U	275 J [177 J]	550	250 D	340	160	220	570	240	270 DJ
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.75 U	0.65 U	0.5 U	195 J [125 J]	140	100 D	38 UJ	140 J	130	140	94	100
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.65 U	0.5 U	0.5 U	50 U [9]	6	32 UD	32 U	16 U	16 U	16 U	16 U	3.3
Trichloroethene	ug/L	0.28 U	0.71 U	0.71 U	0.5 U	0.5 U	5,800 J [3,440 J]	5,500	3,800 D	3,600	2,400	2,800	3,600	2,800	3,300 DJ

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-128 S&P Sand 10/08/05	MW-128 S&P Sand 01/24/06	MW-128 S&P Sand 10/11/06	MW-128 S&P Sand 12/05/06	MW-128 S&P Sand 01/17/07	MW-128 S&P Sand 04/25/07	MW-128 S&P Sand 07/11/07	MW-128 S&P Sand 09/10/07	MW-128 S&P Sand 01/31/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	2.1 U	1 U	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	R	2.2 J	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	2 UJ	1.4 I	1.4 I	5	4.8	2.7 J	7.2 [7.6]
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	1.7	2	0.83 I	0.74 I	0.6 UJ	0.6 U	0.6 U	0.6 U	2.3 [2.6]
1,1-Dichloroethene	ug/L	1.9	1	0.83 U	1.3	0.83 U	0.83 U	0.83 U	0.83 U	3.6 [4]
cis-1,2-Dichloroethene	ug/L	1.2	0.98 I	2	0.91 I	1.4 J	1.6	1.5	1.1	2.5 [2.5]
Tetrachloroethene	ug/L	0.56 J	0.47 I	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.5 U [0.5 U]
Trichloroethene	ug/L	32.8	16	12	18	14	14	16	17	34 [35]

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-129 AF Gravel 10/10/05	MW-129 AF Gravel 01/23/06	MW-129 AF Gravel 10/10/06	MW-129 AF Gravel 12/06/06	MW-129 AF Gravel 01/30/08	MW-130 AF Gravel 10/07/05	MW-130 AF Gravel 03/14/06	MW-130 AF Gravel 10/09/06	MW-130 AF Gravel 12/06/06	MW-130 AF Gravel 01/30/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	2 U	1 U	NA	NA	NA	11.6	37	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	R [R]	2.3 J	NA	NA	NA	31 J	44 J
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	4 J [4.3 J]	2 I	4.2	NA	NA	110 DJ	70
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.5 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.52 U	5.7	11	15	12
1,1-Dichloroethene	ug/L	0.5 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.45 U	9.1	27	39	38
cis-1,2-Dichloroethene	ug/L	0.5 U	0.65 U	0.75 U [0.75 U]	0.75 U	0.65 U	0.53 J	1.1	1.5	0.75 U
Tetrachloroethene	ug/L	0.5 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.5 U	0.5 U	0.34 U	0.65 U	0.65 U
Trichloroethene	ug/L	1.5	0.28 U	0.71 U [0.71 U]	0.71 U	0.5 U	1.7	3.8	5.6	4.9

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-131 AF Gravel 10/06/05	MW-131 AF Gravel 01/23/06	MW-131 AF Gravel 06/26/06	MW-131 AF Gravel 10/12/06	MW-131 AF Gravel 12/11/06	MW-131 AF Gravel 01/31/08	MW-132 AF Gravel 10/07/05	MW-132 AF Gravel 01/19/06	MW-132 AF Gravel 03/22/06	MW-132 AF Gravel 10/11/06	MW-132 AF Gravel 12/13/06	MW-132 AF Gravel 01/23/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	2.4 J	1 U	21	NA	NA	NA	2.1 U	1 U	1.1 U	NA	NA	
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	NA	R	1.3 J	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	3 J	0.76 U	2.9	NA	NA	NA	2 UJ	0.76 U	2.8 [2.1]
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.5 U	0.52 U	0.98 I	0.6 U	0.6 U	0.52 U	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U [0.52 U]
1,1-Dichloroethene	ug/L	0.61 J	0.81 I	1.7	0.83 U	0.83 U	0.45 U	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	2.6 [0.9 I]
cis-1,2-Dichloroethene	ug/L	3.3	3.4	6.5	1.1	0.75 UJ	0.65 U	0.5 U	0.65 U	0.65 U	0.75 U	0.75 UJ	0.65 U [0.65 U]
Tetrachloroethene	ug/L	0.5 U	0.34 U	0.5 U	0.65 U	0.65 U	0.5 U	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U [0.5 U]
Trichloroethene	ug/L	13.9	18	29	4.6	0.71 UJ	3.2 U	0.5 U	0.28 U	0.28 U	0.71 U	0.71 UJ	3.7 [1.7]

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-133 AF Gravel 10/06/05	MW-133 AF Gravel 01/30/06	MW-133 AF Gravel 03/19/06	MW-133 AF Gravel 10/10/06	MW-133 AF Gravel 12/18/06	MW-133 AF Gravel 01/28/08	MW-134 AF Gravel 10/07/05	MW-134 AF Gravel 01/24/06	MW-134 AF Gravel 03/23/06	MW-134 AF Gravel 10/05/06	MW-134 AF Gravel 12/05/06	MW-134 AF Gravel 02/01/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	2 U	1 U	1 U	NA	NA	2.1 U	1 U	1 U	NA	NA	NA	
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	25 J	1.5 J	NA	NA	NA	NA	R	2.3 J	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	1 U [1 U]	NA	NA	NA	4.1 J	0.93 I	28
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.5 U	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U [0.52 U]	0.5 U	0.52 U	0.52 U	1.4	0.6 U	5.4
1,1-Dichloroethene	ug/L	0.5 U	0.45 U	0.45 U	0.83 U	0.83 U	0.45 U [0.45 U]	0.5 U	0.45 U	0.45 U	1.3 J	2.7	23
cis-1,2-Dichloroethene	ug/L	0.5 U	0.65 U	0.65 U	0.75 U	0.75 U	0.65 U [0.65 U]	0.5 U	0.65 U	0.65 U	0.75 U	0.75 U	5.9
Tetrachloroethene	ug/L	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.5 UJ [0.5 U]	0.5 U	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.5 U	0.28 U	0.28 U	0.71 U	0.71 U	0.5 UJ [0.5 U]	2.6 U	0.45 I	1	10	6.6	190 D

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-135 AF Gravel 10/06/05	MW-135 AF Gravel 01/26/06	MW-135 AF Gravel 02/28/06	MW-135 AF Gravel 03/22/06	MW-135 AF Gravel 04/06/06	MW-135 AF Gravel 10/12/06	MW-135 AF Gravel 12/13/06	MW-135 AF Gravel 12/15/06	MW-135 AF Gravel 01/28/08	MW-136 AF Gravel 01/26/06	MW-136 AF Gravel 01/29/08
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	7.3	11	11	24	20	NA	NA	NA	NA	1 U
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	41 J	29 J	R	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	35 J	34	0.76 U	29 [33]	NA
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.86 J	1.1	NA	1.1	1.2	0.88 I	0.91 I	0.6 U	0.72 I [0.98 I]	0.52 U
1,1-Dichloroethene	ug/L	1.1	1.8	NA	2	1.5	1.8	2.2	0.83 U	1.2 [2.1]	0.45 U
cis-1,2-Dichloroethene	ug/L	38.2	55	NA	67	71	92	130	0.75 UJ	63 [93]	0.65 U
Tetrachloroethene	ug/L	0.5 U	0.34 U	NA	0.34 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U [0.5 U]	0.34 U
Trichloroethene	ug/L	95.2	130	NA	140	110	150	43	0.71 UJ	80 [120]	0.28 U

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-137 USAS 01/26/06	MW-137 USAS 12/11/06	MW-137 USAS 01/29/08	MW-138 LSAS 01/26/06	MW-138 LSAS 12/11/06	MW-138 LSAS 01/29/08	MW-139 S&P Sand 01/26/06	MW-139 S&P Sand 12/15/06	MW-139 S&P Sand 01/29/08	MW-140 Lower AF 01/26/06	MW-140 Lower AF 12/11/06	MW-140 Lower AF 01/29/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	NA	NA	1 U	NA	NA	1 U	NA	
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	R [R]	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U [0.76 U]	1 U	NA	0.76 U	1 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U [0.83 U]	0.44 U	0.45 U	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.64 U	0.65 U	0.75 UJ	0.64 U	0.65 U	0.75 U [0.75 U]	0.64 U	0.65 U	0.75 UJ	0.64 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 U [0.71 U]	0.5 U	0.28 U	0.71 UJ	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-141	MW-141	MW-141	MW-142	MW-142	MW-142	MW-143	MW-143	MW-143	MW-143	MW-143	
Zone:	USAS	USAS	USAS	LSAS	LSAS	LSAS	AF Gravel	AF Gravel	AF Gravel	AF Gravel	AF Gravel	
Date Collected:	01/26/06	12/11/06	01/29/08	01/26/06	12/11/06	01/29/08	01/26/06	03/24/06	10/19/06	12/11/06	01/29/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	NA	NA	1 U	1 U	NA	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	NA	R [R]	R	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	NA	2 UJ [2 UJ]	0.76 U	1 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U	0.45 U	0.45 U	0.83 U [0.83 U]	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.64 U	0.65 U	0.75 UJ	0.64 U	0.65 U	0.65 U	0.75 U [0.75 U]	0.75 UJ	0.64 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U [0.65 U]	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.28 U	0.71 U [0.71 U]	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-144 S&P Sand 01/26/06	MW-144 S&P Sand 03/24/06	MW-144 S&P Sand 10/13/06	MW-144 S&P Sand 12/11/06	MW-144 S&P Sand 01/29/08	MW-145 Lower AF 01/26/06	MW-145 Lower AF 03/22/06	MW-145 Lower AF 12/15/06	MW-145 Lower AF 01/29/08	MW-146 USAS 02/02/06	MW-146 USAS 12/11/06	MW-146 USAS 01/29/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	1 U	1 U	NA	NA	NA	1 U	3.2 I	NA	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	R	R	NA	NA	NA	R	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	2 UJ	0.76 U	1 U	NA	NA	0.76 U	1 U	NA	0.76 U	1 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.83 U	0.44 U	0.45 U	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 U	0.75 UJ	0.64 U	0.65 U	0.65 U	0.75 U	0.64 U	0.65 U	0.75 UJ	0.64 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U	0.71 U	0.5 U	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-147 LSAS 02/02/06	MW-147 LSAS 12/11/06	MW-147 LSAS 01/29/08	MW-148 AF Gravel 02/02/06	MW-148 AF Gravel 10/19/06	MW-148 AF Gravel 12/15/06	MW-148 AF Gravel 01/29/08	MW-149 S&P Sand 02/02/06	MW-149 S&P Sand 12/11/06	MW-149 S&P Sand 01/30/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	NA	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	R	NA	NA	R	R	NA	NA	R [R]
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	2 UJ	0.76 U	1 U	NA	0.76 U [0.76 U]
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 UJ	0.6 U	0.52 U	0.52 UJ	0.6 U [0.6 U]
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.44 U	0.45 U	0.83 UJ	0.83 U	0.44 U	0.45 UJ	0.83 U [0.83 U]
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.64 U	0.65 U	0.75 UJ	0.75 U	0.64 U	0.65 UJ	0.75 UJ [0.75 UJ]
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 UJ	0.34 U	0.65 UJ	0.65 U	0.5 UJ	0.34 UJ	0.65 U [0.65 U]
Trichloroethene	ug/L	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.71 U	0.5 U	0.28 UJ	0.71 UJ [0.71 UJ]

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-150 Lower AF 02/02/06	MW-150 Lower AF 03/01/06	MW-150 Lower AF 04/04/06	MW-150 Lower AF 12/11/06	MW-150 Lower AF 01/29/08	MW-151 USAS 01/27/06	MW-151 USAS 02/16/06	MW-151 USAS 10/09/06	MW-151 USAS 12/11/06	MW-151 USAS 01/23/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	1 UJ	1 UJ	NA	NA	1 U	3.3 IJ	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	R	2.3 J
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	2 J	0.9 I
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	NA	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	NA	0.45 U	0.83 U	0.44 U	0.45 U	0.45 U	0.83 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	NA	0.65 U	0.75 UJ	0.64 U	0.65 U	0.65 U	0.75 U	0.75 UJ
Tetrachloroethene	ug/L	0.34 U	NA	0.34 U	0.65 U	0.5 UJ	0.34 U	0.34 UJ	0.65 U	0.65 U
Trichloroethene	ug/L	0.28 U	NA	0.28 U	0.71 UJ	0.5 U	0.28 U	0.28 U	0.71 U	0.71 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-152 LSAS 01/27/06	MW-152 LSAS 02/16/06	MW-152 LSAS 03/23/06	MW-152 LSAS 10/09/06	MW-152 LSAS 12/11/06	MW-152 LSAS 01/23/08	MW-153 AF Gravel 01/27/06	MW-153 AF Gravel 02/16/06	MW-153 AF Gravel 12/11/06	MW-153 AF Gravel 01/24/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	1 U	21 J	1 U	NA	NA	NA	1 U	R	NA	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	NA	R	R	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	NA	2 UJ	0.76 U	1 U	NA	NA	0.76 U	1 U [1 U]
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U [0.52 U]
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U [0.45 U]
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.65 U	0.75 U	0.75 UJ	0.65 U	0.65 U	0.65 U	0.75 UJ	0.65 U [0.65 U]
Tetrachloroethene	ug/L	0.34 U	0.34 UJ	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.34 UJ	0.65 U	0.5 U [0.5 U]
Trichloroethene	ug/L	0.28 U	0.28 U	0.28 U	2.4	0.71 U	0.5 U	0.28 U	0.28 U	0.71 U	0.5 U [0.5 U]

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-154 S&P Sand 01/27/06	MW-154 S&P Sand 02/16/06	MW-154 S&P Sand 12/15/06	MW-154 S&P Sand 01/23/08	MW-155 Lower AF 01/27/06	MW-155 Lower AF 02/16/06	MW-155 Lower AF 12/19/06	MW-155 Lower AF 01/24/08	MW-156 USAS 02/01/06	MW-156 USAS 12/14/06	MW-156 USAS 01/24/08	MW-157 LSAS 02/01/06	MW-157 LSAS 12/14/06	MW-157 LSAS 01/24/08
Semivolatile Organics														
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)														
1,4-Dioxane	ug/L	1 U	R	NA	NA	1 U	R	NA	NA	1 U	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM														
1,4-Dioxane	ug/L	NA	NA	R [R]	NA	NA	NA	R [R]	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution														
1,4-Dioxane	ug/L	NA	NA	0.76 U [0.76 U]	1 U	NA	NA	0.76 U [0.76 U]	1 U	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)														
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.45 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 U [0.75 U]	0.65 U	0.65 U	0.65 U	0.75 UJ [0.75 UJ]	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 UJ
Tetrachloroethene	ug/L	0.34 U	0.34 UJ	0.65 U [0.65 U]	0.5 U	0.34 U	0.34 UJ	0.65 U [0.65 U]	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U [0.71 U]	0.5 U	0.28 U	0.28 U	0.71 UJ [0.71 UJ]	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-158 AF Gravel 02/01/06	MW-158 AF Gravel 03/23/06	MW-158 AF Gravel 10/10/06	MW-158 AF Gravel 12/14/06	MW-158 AF Gravel 01/24/08	MW-159 S&P Sand 02/01/06	MW-159 S&P Sand 12/14/06	MW-159 S&P Sand 01/24/08	MW-160 Lower AF 02/01/06	MW-160 Lower AF 12/14/06	MW-160 Lower AF 01/24/08
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	56	42	NA	NA	NA	1 U	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	NA	43 J	78 J	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	NA	120 DJ	92	84	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	4.3	3.7	5.2	1.8	4.7	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	4.3	3.6	5.5	1.8	4.6	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.68 I	0.65 I	1.3	0.75 UJ	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 UJ
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U	0.71 UJ	0.5 U	0.28 I	0.71 UJ	0.5 U	0.28 U	0.71 UJ

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-161 Floridan 02/01/06	MW-161 Floridan 03/08/06	MW-161 Floridan 12/14/06	MW-161 Floridan 01/24/08	MW-162 USAS 02/16/06	MW-162 USAS 12/12/06	MW-162 USAS 01/30/08	MW-163 LSAS 02/16/06	MW-163 LSAS 12/12/06	MW-163 LSAS 01/30/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	1 U	NA	NA	R [R]	NA	NA	R [2.3 U]	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U [0.52 U]	0.6 U	0.52 U	0.52 U [0.52 U]	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U [0.45 U]	0.83 U	0.45 U	0.45 U [0.45 U]	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U [0.65 U]	0.75 UJ	0.65 U	0.65 U [0.65 U]	0.75 UJ
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.5 U	0.34 UJ [0.34 UJ]	0.65 U	0.5 U	0.34 UJ [0.34 UJ]	0.65 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 UJ	0.5 U	0.28 U [0.28 U]	0.71 U	0.5 U	0.28 U [0.28 U]	0.71 UJ

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-164 AF Gravel 02/16/06	MW-164 AF Gravel 10/12/06	MW-164 AF Gravel 12/12/06	MW-164 AF Gravel 01/30/08	MW-165 S&P Sand 03/02/06	MW-165 S&P Sand 07/10/06	MW-165 S&P Sand 10/12/06	MW-165 S&P Sand 12/11/06	MW-165 S&P Sand 12/12/06	MW-165 S&P Sand 01/30/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	R	NA	NA	NA	120 J [1 UJ]	1 U	NA	NA	NA	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	R	R	NA	NA	NA	R	1.1 J	R	NA
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	2 UJ	0.76 U	1 U	NA	NA	2 UJ	0.76 U	0.76 U	1 U
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U [0.52 U]	NA	0.6 UJ	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U [0.45 U]	NA	0.83 UJ	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.75 U	0.65 U	0.65 U [0.65 U]	NA	0.75 UJ	0.75 UJ	0.75 UJ	0.65 U
Tetrachloroethene	ug/L	0.34 UJ	0.65 U	0.65 U	0.5 U	0.34 U [0.34 U]	NA	0.65 UJ	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.71 UJ	0.5 U	0.28 U [0.28 U]	NA	0.71 UJ	0.71 U	0.71 UJ	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-166 Lower AF 02/16/06	MW-166 Lower AF 12/13/06	MW-166 Lower AF 01/30/08	MW-167 USAS 02/15/06	MW-167 USAS 12/07/06	MW-167 USAS 01/23/08	MW-168 LSAS 02/15/06	MW-168 LSAS 10/12/06	MW-168 LSAS 12/07/06	MW-168 LSAS 01/23/08	MW-169 AF Gravel 02/15/06	MW-169 AF Gravel 12/07/06	MW-169 AF Gravel 01/23/08
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	R	NA	NA	1 U	NA	NA	1 U	NA	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	R	NA	NA	2.6 J	NA	NA	R	R [R]	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	0.76 U	1 U [1 U]	NA	0.76 U	1 U	NA	2 UJ	0.76 U [0.76 U]	1 U	NA	0.76 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U [0.52 U]	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U [0.45 U]	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.65 U [0.65 U]	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.75 U [0.75 U]	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 UJ	0.65 U	0.5 U [0.5 U]	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.65 U [0.65 U]	0.5 U	0.34 UJ	0.65 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U [0.5 U]	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U	0.71 U [0.71 U]	0.5 U	0.28 U	0.71 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-170 Lower AF 02/15/06	MW-170 Lower AF 12/07/06	MW-170 Lower AF 01/23/08	MW-171 LSAS 02/18/06	MW-171 LSAS 12/11/06	MW-171 LSAS 01/23/08	MW-172 AF Gravel 02/18/06	MW-172 AF Gravel 12/11/06	MW-172 AF Gravel 01/23/08	MW-173 S&P Sand 02/18/06	MW-173 S&P Sand 12/11/06	MW-173 S&P Sand 01/23/08
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	1 U	NA	NA	R [1 U]	NA	NA	R [1 U]	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	R	NA	NA	0.94 U
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U [0.52 U]	0.6 U	0.52 U	0.52 U [0.52 U]	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.45 U [0.45 U]	0.83 U	0.45 U	0.45 U [0.45 U]	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.65 U	0.65 U [0.65 U]	0.75 UJ	0.65 U	0.65 U [0.65 U]	0.75 UJ	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U [0.28 U]	0.71 U	0.5 U	0.28 U [0.28 U]	0.71 U	0.5 U	0.28 U	0.71 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-174 Lower AF 02/18/06	MW-174 Lower AF 12/11/06	MW-174 Lower AF 01/25/08	MW-175 AF Gravel 01/28/06	MW-175 AF Gravel 03/22/06	MW-175 AF Gravel 12/11/06	MW-175 AF Gravel 12/13/06	MW-175 AF Gravel 01/29/08
Semivolatile Organics								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)								
1,4-Dioxane	ug/L	1 U	NA	NA	1 U [1 U]	1 U	NA	NA
Volatile Organics (8260B) - SIM								
1,4-Dioxane	ug/L	NA	R	NA	NA	NA	R	R
Volatile Organics (8260) - SIM Isotope Dilution								
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	NA	0.76 U	0.76 U
Volatile Organics (8260B)								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U [0.52 U]	0.52 U	0.6 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.45 U [0.45 U]	0.45 U	0.83 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.65 U	0.65 U [0.65 U]	0.65 U	0.75 UJ	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U [0.34 U]	0.34 U	0.65 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U [0.28 U]	0.28 U	0.71 U	0.71 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-176 S&P Sand 01/28/06	MW-176 S&P Sand 12/07/06	MW-176 S&P Sand 01/29/08	MW-177 Lower AF 01/28/06	MW-177 Lower AF 12/07/06	MW-177 Lower AF 01/29/08	MW-178 LSAS 01/27/06	MW-178 LSAS 12/14/06	MW-178 LSAS 01/25/08	MW-179 AF Gravel 01/27/06	MW-179 AF Gravel 12/14/06	MW-179 AF Gravel 01/25/08
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	1 U	NA	NA	1.1 U	NA	NA	1 U	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.64 U	0.65 U	0.75 U	0.64 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 UJ
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-180 S&P Sand 01/27/06	MW-180 S&P Sand 12/14/06	MW-180 S&P Sand 01/25/08	MW-181 Lower AF 01/27/06	MW-181 Lower AF 12/14/06	MW-181 Lower AF 01/25/08	MW-182 S&P Sand 02/15/06	MW-182 S&P Sand 12/07/06	MW-182 S&P Sand 01/23/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	NA	NA	1.1 U [1.1 U]	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	R	NA	NA	R [R]	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U [0.76 U]	1 U	NA	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U [0.52 U]	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U [0.83 U]	0.45 U	0.45 U [0.45 U]	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 UJ [0.75 UJ]	0.65 U	0.65 U [0.65 U]	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U [0.65 U]	0.5 U	0.34 UJ [0.34 UJ]	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ [0.71 UJ]	0.5 U	0.28 U [0.28 U]	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-183 USAS 02/17/06	MW-183 USAS 10/09/06	MW-183 USAS 12/06/06	MW-183 USAS 01/24/08	MW-184 LSAS 02/17/06	MW-184 LSAS 10/09/06	MW-184 LSAS 12/06/06	MW-184 LSAS 01/24/08	MW-185 AF Gravel 02/17/06	MW-185 AF Gravel 10/09/06	MW-185 AF Gravel 12/06/06	MW-185 AF Gravel 01/24/08
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	1 U	NA	NA	NA	1 U [1 U]	NA	NA	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	R	R	NA	NA	R	R	NA	NA	R	R
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	2 UJ	0.76 U	1 U	NA	2 UJ	0.76 U	1 U	NA	2 UJ	0.76 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U [0.52 U]	0.6 U	0.6 U	0.52 U	0.52 U	0.6 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U [0.45 U]	0.83 U	0.83 U	0.45 U	0.45 U	0.83 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.75 U	0.65 U	0.65 U [0.65 U]	0.75 U	0.75 UJ	0.65 U	0.65 U	0.75 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U [0.34 U]	0.65 U	0.65 U	0.5 U	0.34 U	0.65 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.75 U	0.71 U	0.5 U	0.28 U [0.28 U]	0.71 U	0.71 U	0.5 U	0.28 U	0.71 U	0.71 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-186 S&P Sand 02/17/06	MW-186 S&P Sand 10/10/06	MW-186 S&P Sand 12/06/06	MW-186 S&P Sand 01/24/08	MW-187 Lower AF 02/17/06	MW-187 Lower AF 12/06/06	MW-187 Lower AF 01/24/08	MW-188 USAS 12/06/06	MW-188 USAS 01/23/08	MW-188 USAS 03/07/06	MW-188 USAS 03/17/06	MW-188 USAS 03/20/06
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	1 U	NA	NA	NA	1 U	NA	NA	NA	1 U	1 U	1 U
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	R	R	NA	NA	R	NA	R	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	2 UJ	0.76 U	1 U	NA	0.76 U	1 U	0.76 U	1 U	NA	NA
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.83 U	0.45 U	0.45 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.75 UJ	0.65 U	0.65 U	0.75 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.65 U	0.5 U	0.34 U	0.34 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.71 U	0.5 U	0.28 U	0.71 U	0.5 U	0.71 U	0.5 U	0.28 U	0.28 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-189	MW-189	MW-189	MW-189	MW-189	MW-190	MW-190	MW-190	MW-190	MW-190	MW-190
Zone:	LSAS	LSAS	LSAS	LSAS	LSAS	AF Gravel	AF Gravel	AF Gravel	AF Gravel	AF Gravel	AF Gravel
Date Collected:	12/06/06	01/23/08	03/08/06	03/17/06	03/20/06	03/07/06	03/17/06	03/20/06	03/20/06	12/06/06	01/25/08
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	R	NA	NA	NA	NA	NA	NA	NA	R [R]	NA
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	0.76 U	1 U [1 U]	NA	NA	NA	NA	NA	NA	0.76 U [0.76 U]	1 U
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.6 U	0.52 U [0.52 U]	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.52 U
1,1-Dichloroethene	ug/L	0.83 U	0.45 U [0.45 U]	0.45 U	0.45 U	0.45 U	0.45 U	0.45 U	0.45 U	0.83 U [0.83 U]	0.45 U
cis-1,2-Dichloroethene	ug/L	0.75 U	0.65 U [0.65 U]	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U	0.75 U [0.75 U]	0.65 U
Tetrachloroethene	ug/L	0.65 U	0.5 U [0.5 U]	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U	0.65 U [0.65 U]	0.5 U
Trichloroethene	ug/L	0.71 U	0.5 U [0.5 U]	0.28 U	0.28 U	0.28 U	0.28 U	0.28 U	0.28 U	0.71 U [0.71 U]	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-191	MW-191	MW-191	MW-191	MW-191	MW-192	MW-192	MW-192	MW-192	
Zone:	S&P Sand	S&P Sand	S&P Sand	S&P Sand	S&P Sand	Lower AF	Lower AF	Lower AF	Lower AF	
Date Collected:	03/08/06	03/17/06	03/20/06	12/06/06	01/23/08	03/07/06	03/18/06	12/06/06	01/23/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	1 U	1 U	NA	NA	1 U	1 U	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	0.76 U	1 U	NA	NA	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.28 U	0.79 I	0.5 U

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2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-193 AF Gravel 03/05/06	MW-193 AF Gravel 03/16/06	MW-193 AF Gravel 12/06/06	MW-193 AF Gravel 01/24/08	MW-194 S&P Sand 03/06/06	MW-194 S&P Sand 03/16/06	MW-194 S&P Sand 12/06/06	MW-194 S&P Sand 01/24/08
Semivolatile Organics								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)								
1,4-Dioxane	ug/L	1 U	1 U	NA	NA	1 U	1 U	NA
Volatile Organics (8260B) - SIM								
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution								
1,4-Dioxane	ug/L	NA	NA	0.76 U	1 U	NA	NA	0.76 U
Volatile Organics (8260B)								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.28 U	0.71 U

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FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-195 Lower AF 03/07/06	MW-195 Lower AF 03/18/06	MW-195 Lower AF 12/06/06	MW-195 Lower AF 01/24/08	MW-196 AF Gravel 03/15/06	MW-196 AF Gravel 03/23/06	MW-196 AF Gravel 12/15/06	MW-196 AF Gravel 01/30/08	MW-197 AF Gravel 03/15/06	MW-197 AF Gravel 12/12/06	MW-197 AF Gravel 01/30/08	MW-198 USAS 01/28/08	MW-199 LSAS 01/28/08	MW-200 AF Gravel 01/28/08
Semivolatile Organics														
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)														
1,4-Dioxane	ug/L	1 U	1 U	NA	NA	1 U	1 U	NA	NA	1 UJ	NA	NA	NA	NA
Volatile Organics (8260B) - SIM														
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	NA	R	NA	NA	R	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution														
1,4-Dioxane	ug/L	NA	NA	0.76 U	1 U [1 U]	NA	NA	0.76 U	1 U	NA	0.76 U	1 U	1 U	1 U
Volatile Organics (8260B)														
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.52 U [0.52 U]	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.45 U [0.45 U]	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 UJ	0.65 U [0.65 U]	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.5 U [0.5 U]	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.5 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U	0.5 U [0.5 U]	7.2	4.9	3.7	2.2	0.28 U	0.71 UJ	0.5 U	0.5 U	0.5 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-201 S&P Sand 01/28/08	MW-202 Lower AF 02/01/08	MW-204 USAS 01/23/08	MW-205 LSAS 01/23/08	MW-206 AF Gravel 01/29/08	MW-207 Lower AF 01/29/08	MW-208 USAS 03/30/06	MW-208 USAS 12/07/06	MW-208 USAS 01/23/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	1 U	1 U	1 U	1 U	1 U	1 U	NA	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.45 U	0.45 U	0.44 U	0.44 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.65 U	0.65 U	0.64 U	0.64 U	11	5.8	11
Tetrachloroethene	ug/L	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-209	MW-209	MW-209	MW-210	MW-210	MW-210	MW-210	MW-211	MW-211	MW-211
Zone:	LSAS	LSAS	LSAS	AF Gravel	AF Gravel	AF Gravel	AF Gravel	S&P Sand	S&P Sand	S&P Sand
Date Collected:	03/30/06	12/07/06	01/23/08	03/25/06	04/03/06	12/07/06	01/23/08	03/30/06	12/07/06	01/23/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	1.1 U	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	R	NA	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.65 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U

APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-212	MW-212	MW-212	MW-213	MW-213	MW-213	MW-214	MW-214	MW-214	
Zone:	Lower AF	Lower AF	Lower AF	USAS	USAS	USAS	LSAS	LSAS	LSAS	
Date Collected:	03/25/06	12/07/06	01/23/08	03/31/06	12/12/06	01/31/08	03/31/06	12/12/06	01/31/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1.1 U	NA	NA	1 U	NA	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 UJ	NA	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 UJ	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 UJ	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.65 UJ	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 UJ	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U	0.71 UJ	0.5 UJ	0.28 U	0.71 UJ	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-215 AF Gravel 03/31/06	MW-215 AF Gravel 11/13/06	MW-215 AF Gravel 12/12/06	MW-215 AF Gravel 01/31/08	MW-216 S&P Sand 03/30/06	MW-216 S&P Sand 12/12/06	MW-216 S&P Sand 01/31/08	MW-217 Lower AF 03/30/06	MW-217 Lower AF 12/12/06	MW-217 Lower AF 01/31/08	MW-218 Floridan 04/06/06	MW-218 Floridan 12/12/06
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	4.1 I	NA	NA	NA	1 U	NA	NA	1 U	NA	NA	1.1 U
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	R	NA	NA	R	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	2 U	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U	1 U	NA
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.75 U	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-219	MW-219	MW-219	MW-219	MW-220	MW-220	MW-220	MW-220	MW-221	MW-221	MW-221	MW-221	
Zone:	USAS	USAS	USAS	USAS	LSAS	LSAS	LSAS	LSAS	AF Gravel	AF Gravel	AF Gravel	AF Gravel	
Date Collected:	03/20/06	03/25/06	12/13/06	01/25/08	03/20/06	03/25/06	12/13/06	01/25/08	03/20/06	03/26/06	12/13/06	01/25/08	
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	1 U	1 U	NA	NA	1 U	1 U	NA	NA	1 U	1 U	NA	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	NA	R [R]	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	0.76 U	1 U	NA	NA	0.76 U [0.76 U]	1 U	NA	NA	0.76 U	1 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.45 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.65 U	0.75 U [0.75 U]	0.65 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U [0.65 U]	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.28 U	0.71 U [0.71 U]	NA	0.28 U	0.28 U	0.71 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-222 S&P Sand 03/20/06	MW-222 S&P Sand 03/26/06	MW-222 S&P Sand 12/13/06	MW-222 S&P Sand 01/25/08	MW-223 Hardstreak Clay 03/26/06	MW-223 Hardstreak Clay 12/07/06	MW-223 Hardstreak Clay 01/25/08
Semivolatile Organics							
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)							
1,4-Dioxane	ug/L	1 U	1.1 U	NA	NA	1.1 U	NA
Volatile Organics (8260B) - SIM							
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution							
1,4-Dioxane	ug/L	NA	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)							
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-224 Venice Clay 03/26/06	MW-224 Venice Clay 12/08/06	MW-224 Venice Clay 01/25/08	MW-225 Venice Clay 03/26/06	MW-225 Venice Clay 12/07/06	MW-225 Venice Clay 01/25/08	MW-226 AF Gravel 03/26/06	MW-226 AF Gravel 04/04/06	MW-226 AF Gravel 12/07/06	MW-226 AF Gravel 01/25/08	MW-227 S&P Sand 03/26/06	MW-227 S&P Sand 12/08/06	MW-227 S&P Sand 01/25/08
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	NA	NA	1 U	1 UJ	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	1.1	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.45 U	0.45 U	0.96 I	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U	0.5 U	0.28 U	0.28 U	0.71 U	0.5 U	0.28 U	0.71 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID:	MW-228	MW-228	MW-229	MW-229	MW-229	MW-230	MW-230	MW-230	MW-231	MW-231	MW-231	
Zone:	AF Gravel	AF Gravel	USAS	USAS	USAS	LSAS	LSAS	LSAS	AF Gravel	AF Gravel	AF Gravel	
Date Collected:	04/04/06	01/23/08	04/02/06	12/14/06	01/29/08	04/02/06	12/14/06	01/29/08	04/02/06	12/14/06	01/29/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	1 UJ	NA	1 U	NA	NA	1 U	NA	NA	1 U [1 U]	NA	NA
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	R [R]	NA	NA	R	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	NA	1 U [1 U]	NA	0.76 U [0.76 U]	1 U	NA	0.76 U	1 U	NA	0.76 U	1 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U [0.52 U]	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U [0.52 U]	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U [0.45 U]	0.45 U	0.83 U [0.83 U]	0.44 U	0.45 U	0.83 U	0.44 U	0.45 U [0.45 U]	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U [0.65 U]	0.65 U	0.75 UJ [0.75 UJ]	0.64 U	0.65 U	0.75 UJ	0.64 U	0.65 U [0.65 U]	0.75 UJ	0.64 U
Tetrachloroethene	ug/L	0.34 U	0.5 U [0.5 U]	0.34 U	0.65 U [0.65 U]	0.5 U	0.34 U	0.65 U	0.5 UJ	0.34 U [0.34 U]	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.5 U [0.5 U]	0.28 U	0.71 UJ [0.71 UJ]	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U [0.28 U]	0.71 UJ	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Location ID: Zone: Date Collected:	MW-232 AF Gravel 04/04/06	MW-232 AF Gravel 10/10/06	MW-232 AF Gravel 12/18/06	MW-232 AF Gravel 01/30/08	MW-233 AF Gravel 04/04/06	MW-233 AF Gravel 12/13/06	MW-233 AF Gravel 01/24/08	MW-234 USAS 04/02/06	MW-234 USAS 12/12/06	MW-234 USAS 01/31/08	MW-235 LSAS 04/02/06	MW-235 LSAS 12/12/06	MW-235 LSAS 01/31/08
Semivolatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	1 UJ [12 J]	NA	NA	NA	1 UJ	NA	NA	1 U	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	R	5.2 J [4.5 J]	NA	NA	R	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	6.1 J	5.2 [2.7]	7.6 [7.6]	NA	0.76 U	1 U	NA	0.76 U	1 U	NA	0.76 U
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U [0.52 U]	0.6 U	0.6 U [0.6 U]	0.52 U [0.52 U]	0.56 I	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U [0.45 U]	0.83 U	0.83 U [0.83 U]	0.45 U [0.45 U]	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U [0.65 U]	0.75 U	0.75 U [0.75 U]	0.65 U [0.65 U]	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U [0.34 U]	0.65 U	0.65 U [0.65 U]	0.5 U [0.5 U]	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.28 U [0.28 U]	0.71 U	0.71 U [0.71 U]	0.5 U [0.5 U]	0.28 U	0.71 U	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	MW-236 AF Gravel 04/02/06	MW-236 AF Gravel 12/12/06	MW-236 AF Gravel 01/31/08	MW-237 S&P Sand 04/07/06	MW-237 S&P Sand 12/12/06	MW-237 S&P Sand 01/31/08	MW-238 Lower AF 04/07/06	MW-238 Lower AF 12/12/06	MW-238 Lower AF 01/31/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1.1 U	NA	NA	1 U	NA	NA	1.1 U	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	R [R]	NA	NA	R	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	0.76 U [0.76 U]	1 U	NA	0.76 U	1 U [1 UJ]	NA	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.6 U	0.52 U [0.52 U]	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.83 U	0.45 U [0.45 U]	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U [0.75 U]	0.65 U	0.65 U	0.75 U	0.65 U [0.65 U]	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.34 U	0.65 U [0.65 U]	0.5 U	0.34 U	0.65 U	0.5 U [0.5 U]	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 UJ [0.71 UJ]	0.5 U	0.28 U	0.71 UJ	0.5 U [0.5 U]	0.28 U	0.71 UJ	0.5 U

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	MW-239 AF Gravel 04/05/06	MW-239 AF Gravel 12/14/06	MW-239 AF Gravel 01/23/08	MW-240 S&P Sand 04/05/06	MW-240 S&P Sand 12/07/06	MW-240 S&P Sand 01/23/08	MW-241 Lower AF 12/19/06	MW-241 Lower AF 01/25/08	MW-242 USAS 04/06/06	MW-242 USAS 12/12/06	MW-242 USAS 01/29/08
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	1 U	NA	NA	1 U	NA	NA	NA	NA	1.1 U	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	4.6 J [5 J]	NA	NA	R	NA	R	NA	NA	R [R]
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	3.6 J [7.1 J]	6.7	NA	0.76 U	1 U	0.76 U	1 U	NA	0.76 U [0.76 U]
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	R
1,1-Dichloroethane	ug/L	0.52 U	0.6 U [0.6 U]	0.52 U	0.52 U	0.6 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U [0.6 U]
1,1-Dichloroethene	ug/L	0.45 U	0.83 U [0.83 U]	0.45 U	0.45 U	0.83 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U [0.83 U]
cis-1,2-Dichloroethene	ug/L	0.65 U	0.98 J [1 J]	2.1	0.65 U	0.75 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 U [0.75 UJ]
Tetrachloroethene	ug/L	0.34 U	0.65 U [0.65 U]	0.5 U	0.34 U	0.65 U	0.5 U	0.65 U	0.5 U	0.34 U	0.65 U [0.65 U]
Trichloroethene	ug/L	0.41 I	0.92 J [0.86 J]	2.9	0.28 U	0.71 U	0.5 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ [0.71 UJ]

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	MW-243 LSAS 04/06/06	MW-243 LSAS 12/12/06	MW-243 LSAS 01/29/08	MW-244 AF Gravel 04/06/06	MW-244 AF Gravel 12/12/06	MW-244 AF Gravel 01/29/08	MW-245 Hardstreak Clay 04/06/06	MW-245 Hardstreak Clay 12/14/06	MW-245 Hardstreak Clay 01/29/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	1 U	NA	NA	1.1 U	NA	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.8 I	1 U	NA	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 U	0.64 U	0.65 U	0.75 U	0.64 U	0.65 U	0.75 UJ	0.64 U
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID:	MW-246	MW-246	MW-246	MW-247	MW-247	MW-247	MW-248	MW-248	MW-248	MW-248	
Zone:	LSAS	LSAS	LSAS	AF Gravel	AF Gravel	AF Gravel	AF Gravel	AF Gravel	AF Gravel	AF Gravel	
Date Collected:	04/06/06	12/14/06	01/29/08	04/06/06	12/14/06	01/29/08	04/07/06	04/11/06	12/18/06	01/28/08	
Semivolatile Organics											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)											
1,4-Dioxane	ug/L	1 U	NA	NA	1.1 U	NA	NA	1.2 U	1 U	NA	NA
Volatile Organics (8260B) - SIM											
1,4-Dioxane	ug/L	NA	R	NA	NA	R	NA	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution											
1,4-Dioxane	ug/L	NA	0.76 U	1 U	NA	0.76 U	1 U [1 U]	NA	NA	0.76 U	12 J [1 UJ]
Volatile Organics (8260B)											
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U [0.52 U]	0.52 U	0.52 U	0.6 U	1.1 [0.52 U]
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.44 U	0.45 U	0.83 U	0.44 U [0.44 U]	0.45 U	0.45 U	0.83 U	0.68 U [0.45 U]
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.64 U	0.65 U	0.75 UJ	0.64 U [0.64 U]	0.65 U	0.65 U	0.75 U	0.65 U [0.65 U]
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U [0.5 U]	0.34 U	0.34 U	0.65 U	0.5 U [0.5 U]
Trichloroethene	ug/L	0.28 U	0.71 UJ	0.5 U	0.28 U	0.71 UJ	0.5 U [0.5 U]	0.28 U	0.28 U	0.71 U	0.5 U [0.5 U]

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	MW-249 AF Gravel 02/16/07	MW-249 AF Gravel 01/31/08	MW-250 AF Gravel 02/16/07	MW-250 AF Gravel 01/31/08	MW-251 Floridan 04/23/07	MW-251 Floridan 01/30/08	MW-252 S&P Sand 01/31/08	MW-253 AF Gravel 02/01/08	MW-254 (MW-BT-1) USAS 02/01/08	
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	0.76 U	1 U	0.76 U [4.4]	15	0.69 U	1 U	1 U	350	10 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.6 U	0.52 U	0.6 U [0.6 U]	0.52 U	0.6 U	0.52 U	0.52 U	75	61
1,1-Dichloroethene	ug/L	0.83 U	0.45 U	0.83 U [0.83 U]	0.45 U	0.83 U	0.45 U	0.45 U	400 D	37
cis-1,2-Dichloroethene	ug/L	0.75 U	0.65 U	0.75 U [0.75 U]	0.65 U	0.75 U	0.65 U	0.65 U	130	6.5 U
Tetrachloroethene	ug/L	0.65 U	0.5 U	0.65 U [0.65 U]	0.5 U	0.65 U	0.5 U	0.5 U	5.4	11,000 D
Trichloroethene	ug/L	0.71 U	0.5 U	0.71 U [0.71 U]	0.5 U	0.71 U	0.5 U	0.5 U	6,200 D	760

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	PZ-LSAS-1 LSAS 01/31/08	PZ-LSAS-2 LSAS 01/31/08	PZ-LSAS-3 LSAS 01/31/08	PZ-LSAS-4 LSAS 01/31/08	PZ-LSAS-5 LSAS 01/31/08	PZ-LSAS-6 LSAS 01/31/08	PZ-LSAS-7 LSAS 01/31/08	RW-1 USAS 12/05/06	RW-1 USAS 02/01/08	RW-2 USAS 12/05/06	RW-2 USAS 01/24/08	
Semivolatile Organics												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Volatile Organics (8260B) - SIM												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	R	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution												
1,4-Dioxane	ug/L	660 EJ	580 DJ	100	600 DJ	380 J	350 J	50 UJ	0.76 U	1 U	0.76 U	1 U
Volatile Organics (8260B)												
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	3.3	18	11	8.8	18	33	49	0.6 U	2	6.9	4.1
1,1-Dichloroethene	ug/L	48	77	37	100	110	150	65	0.83 U	0.45 U	1.2	0.56 I
cis-1,2-Dichloroethene	ug/L	120	680 EJ	56	730 EJ	690 EJ	510 EJ	480 EJ	0.75 UJ	0.65 U	4.5	15
Tetrachloroethene	ug/L	0.5 U	1.7	13	0.5 U	0.5 U	3.1	6.3	0.65 U	0.5 U	72	21
Trichloroethene	ug/L	310 EJ	1,000 EJ	2,500 EJ	4,100 EJ	3,400 DJ	8,900 DJ	2,300 DJ	0.71 U	1.5	17	9.7

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID:	TW-84-A	TW-84-A	TW-84-B	TW-84-B
Zone:	USAS	USAS	USAS	USAS
Date Collected:	10/12/07	01/29/08	10/12/07	01/29/08
Semivolatile Organics				
1,4-Dioxane	ug/L	NA	NA	NA
Semivolatile Organics (8270C)				
1,4-Dioxane	ug/L	NA	NA	NA
Volatile Organics (8260B) - SIM				
1,4-Dioxane	ug/L	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution				
1,4-Dioxane	ug/L	NA	1 U	1 U
Volatile Organics (8260B)				
1,4-Dioxane	ug/L	1 U	NA	1 U
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.52 U
1,1-Dichloroethene	ug/L	0.45 U	0.44 U	0.44 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.64 U	0.64 U
Tetrachloroethene	ug/L	0.5 U	0.5 U	0.5 U
Trichloroethene	ug/L	0.5 U	0.5 U	0.5 U

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EXU-1 (EW-101) USAS 01/30/08	EXU-1 (EW-101) USAS 02/05/08	EXU-1 (EW-101) USAS 02/12/08	EXU-1 (EW-101) USAS 02/19/08	EXU-1 (EW-101) USAS 02/26/08
Semivolatile Organics						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution						
1,4-Dioxane	ug/L	1 U	15	27	44	32
Volatile Organics (8260B)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	3.7	11	19	25	25
1,1-Dichloroethene	ug/L	3.2	9.7	27	37	41
cis-1,2-Dichloroethene	ug/L	14	16	35	47	42
Tetrachloroethene	ug/L	7.3	4.4	8.3	6.8	7.9
Trichloroethene	ug/L	140	530	940	1,400	1,500

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-102 LSAS 01/31/08	EW-102 LSAS 02/05/08	EW-102 LSAS 02/12/08	EW-102 LSAS 02/19/08	EW-102 LSAS 02/26/08
Semivolatile Organics						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution						
1,4-Dioxane	ug/L	120 DJ	84	43	48	66
Volatile Organics (8260B)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	6.3	4.3	3.7	4.9	7.1
1,1-Dichloroethene	ug/L	55	58	57	67	78
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U
Tetrachloroethene	ug/L	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
Trichloroethene	ug/L	1.8 U	0.5 U	0.5 U	0.5 U	0.57 I

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-103 USAS 09/20/06	EW-103 USAS 11/21/06	EW-103 USAS 12/06/06	EW-103 USAS 12/13/06	EW-103 USAS 01/15/07	EW-103 USAS 02/07/07	EW-103 USAS 03/22/07	EW-103 USAS 04/18/07
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	2.3 l	1.9 l	10 U	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	4.6	10	11 J	0.76 U	5.1
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	41	50	37	53	41	49	15	61
1,1-Dichloroethene	ug/L	9.1	4.5 U	4.5 U	19	14	17	3.1	12
cis-1,2-Dichloroethene	ug/L	42	22	21	14	18	34	30	21
Tetrachloroethene	ug/L	3.1	5 U	5 U	6	6 J	6.5 U	1.5	6.5 U
Trichloroethene	ug/L	950 D	480	450	750 D	870	610	110	470

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-103 USAS 05/21/07	EW-103 USAS 06/20/07	EW-103 USAS 07/16/07	EW-103 USAS 08/16/07	EW-103 USAS 09/26/07	EW-103 USAS 10/24/07	EW-103 USAS 11/08/07	EW-103 USAS 01/30/08
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	9.3	8	5.5	5.9	5.2	3.8	4.5	1 U
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	R	NA	NA
1,1-Dichloroethane	ug/L	53	80	64	35	38	39	34	18
1,1-Dichloroethene	ug/L	12	19	13	8.4	6.4	9.4	12	4.2
cis-1,2-Dichloroethene	ug/L	23	27	13	9.6	11	14	14	14
Tetrachloroethene	ug/L	7.1	6.6	3.2 U	3.2 U	2.4 U	3.2 U	4.2 I	1
Trichloroethene	ug/L	680	690 J	420	330	330	340	470	31

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-104 LSAS 09/20/06	EW-104 LSAS 11/21/06	EW-104 LSAS 12/06/06	EW-104 LSAS 12/13/06	EW-104 LSAS 01/15/07	EW-104 LSAS 02/07/07	EW-104 LSAS 03/22/07	EW-104 LSAS 04/18/07
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	110	120	150	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	300 J	360	300	250	260
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	54	150	49 I	70	48	54	35 I	65
1,1-Dichloroethene	ug/L	130	71	130	200 D	190	180	150	170
cis-1,2-Dichloroethene	ug/L	170	180	160	120	98	140	120	180
Tetrachloroethene	ug/L	13 U	25 U	25 U	33	28 J	16 U	32 U	19 I
Trichloroethene	ug/L	3,900 D	3,600	3,600	3,900 D	3,900	2,600	2,800	3,500

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-104 LSAS 05/21/07	EW-104 LSAS 06/20/07	EW-104 LSAS 07/16/07	EW-104 LSAS 08/16/07	EW-104 LSAS 09/26/07	EW-104 LSAS 10/24/07	EW-104 LSAS 11/08/07	EW-104 LSAS 01/30/08
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	260	250	200	290	340	250	280	25
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	R	NA	NA
1,1-Dichloroethane	ug/L	51	64	56	50	50	52	39	49
1,1-Dichloroethene	ug/L	140	200	230	280	140	140	170	14
cis-1,2-Dichloroethene	ug/L	92	110	100	87	81	99	80	71
Tetrachloroethene	ug/L	32 U	27	13 U	13 U	12 U	13 I	19 I	4.9
Trichloroethene	ug/L	2,300	3,100	3,100	2,700	2,100	2,100	2,600	1,000 EJ

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-105 USAS 09/20/06	EW-105 USAS 11/21/06	EW-105 USAS 12/06/06	EW-105 USAS 12/13/06	EW-105 USAS 01/15/07	EW-105 USAS 02/07/07	EW-105 USAS 03/22/07	EW-105 USAS 04/18/07
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	1 U	1.3 I	10 U	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	22 J	55	47 J	56	63
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	4.1	7.4	6	11	11 J	20	14 I	25 I
1,1-Dichloroethene	ug/L	9.3	6.6	0.9 U	39	39 J	55	48	65
cis-1,2-Dichloroethene	ug/L	13	15	24	23	44 J	100	63	88
Tetrachloroethene	ug/L	1.7 I	7.7	2 I	5	6.5 UJ	6.5 U	13 U	16 U
Trichloroethene	ug/L	240	180	450 D	1,000 D	2,000 J	2,900 DJ	1,900	2,500

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-105 USAS 05/21/07	EW-105 USAS 06/20/07	EW-105 USAS 07/16/07	EW-105 USAS 08/16/07	EW-105 USAS 09/26/07	EW-105 USAS 10/24/07	EW-105 USAS 11/08/07	EW-105 USAS 01/30/08
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	50	61 J	17	32	46 J	33	50	3.9
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	R	NA	NA
1,1-Dichloroethane	ug/L	25	31	12 U	12 U	21	16	16	1.3
1,1-Dichloroethene	ug/L	63	70	27	36	34	28	42	1.3
cis-1,2-Dichloroethene	ug/L	56	51	15 U	15 U	30	26	24	1.4
Tetrachloroethene	ug/L	27	46	13 U	16 I	79	35	110	0.79 I
Trichloroethene	ug/L	2,200	2,000	710	930	1,200 D	800	1,200	19

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-106 LSAS 09/20/06	EW-106 LSAS 11/21/06	EW-106 LSAS 12/06/06	EW-106 LSAS 12/13/06	EW-106 LSAS 01/15/07	EW-106 LSAS 02/07/07	EW-106 LSAS 03/22/07	EW-106 LSAS 04/18/07	EW-106 LSAS 05/21/07
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	130	150	130	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	540 J	630	620 J	580	510	490
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	77 I	52 U	52 U	96	60 U	110	64 I	100	92 I
1,1-Dichloroethene	ug/L	330	180	210	560 D	480	460	350	410	360
cis-1,2-Dichloroethene	ug/L	1,200	870	1,100	100 D	380	320	250	390	220
Tetrachloroethene	ug/L	50 U	50 U	50 U	9.3	65 U	65 U	65 U	32 U	65 U
Trichloroethene	ug/L	8,900	8,100	12,000	12,000 D	12,000	6,400	7,000	8,600	8,200

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-106 LSAS 06/20/07	EW-106 LSAS 07/16/07	EW-106 LSAS 08/16/07	EW-106 LSAS 09/26/07	EW-106 LSAS 10/24/07	EW-106 LSAS 08/16/07	EW-106 LSAS 10/24/07	EW-106 LSAS 11/08/07	EW-106 LSAS 01/30/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	540	460	650	650	610	NA	NA	610	20
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	R	NA	R	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	90 I	110	74	72 I	93	NA	NA	62	24
1,1-Dichloroethene	ug/L	430	760	650	350	NA	NA	420	380	43
cis-1,2-Dichloroethene	ug/L	180	240	120	110	NA	NA	150	160	580 EJ
Tetrachloroethene	ug/L	65 U	32 U	32 U	47 U	NA	NA	130	310	26
Trichloroethene	ug/L	7,500	10,000	9,500	9,900	NA	NA	7,800	11,000 D	1,500 EJ

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-107 USAS 01/31/08	EW-107 USAS 02/05/08	EW-107 USAS 02/12/08	EW-107 USAS 02/19/08	EXL-1 (EW-108) LSAS 01/31/08	EXL-1 (EW-108) LSAS 02/05/08	EXL-1 (EW-108) LSAS 02/12/08	EXL-1 (EW-108) LSAS 02/19/08	EXL-1 (EW-108) LSAS 02/26/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	1 UJ [1 U]	3.5	3.2	9.8	540 J	920	680 V	690	660
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	52 U [0.52]	0.97 I	2	3.4	46	180	130	120	96
1,1-Dichloroethene	ug/L	45 U [0.45]	1.2	2.8	5.5	140	990	670	630	470
cis-1,2-Dichloroethene	ug/L	65 U [0.65]	0.65 U	0.65 U	2	930 EJ	39	34	32	29
Tetrachloroethene	ug/L	0.5 U [0.5 U]	0.99 I	4.5	6.4	0.5 U	6	14	13	20 I
Trichloroethene	ug/L	6.1 [4.2]	28	64	140	78	730	470	530	610

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-109 USAS 09/20/06	EW-109 USAS 11/21/06	EW-109 USAS 12/06/06	EW-109 USAS 12/13/06	EW-109 USAS 01/15/07	EW-109 USAS 02/07/07	EW-109 USAS 03/22/07	EW-109 USAS 04/18/07
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	13	11	14	0.76 U	0.87 I
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	6.4	16	18	12	12	14	2.9	0.6 U
1,1-Dichloroethene	ug/L	9.7	24	30	32	25	33	5.5	0.83 U
cis-1,2-Dichloroethene	ug/L	1.4	2.9	3.2	1.3	1.7	1.7	0.75 U	0.75 U
Tetrachloroethene	ug/L	0.95 I	4.6	1.4	1.1	2.1 J	0.66 I	0.85 I	0.65 U
Trichloroethene	ug/L	3.1	7.9	7.4	35	6.8	5.8	1.8	0.71 U

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-109 USAS 05/21/07	EW-109 USAS 06/20/07	EW-109 USAS 07/16/07	EW-109 USAS 08/16/07	EW-109 USAS 09/26/07	EW-109 USAS 10/24/07	EW-109 USAS 11/08/07	EW-109 USAS 01/30/08
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	1.7 I	1.5 I	0.69 U	1.2 I	0.84 I	0.69 U	0.69 U	1 U
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	R	NA	NA
1,1-Dichloroethane	ug/L	3.4	0.82 I	0.77 I	0.98 I	0.28 U	0.7 I	0.28 U	0.52 U
1,1-Dichloroethene	ug/L	4.6	0.99 I	1.3	3.1	0.41 I	1.4	0.4 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.75 U	0.75 U	0.75 U	0.75 U	0.18 U	0.75 U	0.18 U	0.65 U
Tetrachloroethene	ug/L	1.3	0.65 U	0.65 U	0.65 U	0.47 U	0.65 U	0.47 U	0.5 U
Trichloroethene	ug/L	1.7	0.71 U	0.71 U	0.71 U	0.26 U	0.71 U	0.26 U	0.5 U

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-110 LSAS 09/20/06	EW-110 LSAS 11/21/06	EW-110 LSAS 12/06/06	EW-110 LSAS 12/13/06	EW-110 LSAS 01/15/07	EW-110 LSAS 02/07/07	EW-110 LSAS 03/22/07	EW-110 LSAS 04/18/07
Semivolatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	360	380	260	190	230
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	310 D	260	340	340 D	270	290	120	270
1,1-Dichloroethene	ug/L	720 D	650	870	1,300 D	1,100	910	470	850
cis-1,2-Dichloroethene	ug/L	36	24	22	18	16 I	22	15	20
Tetrachloroethene	ug/L	11	5 U	10	32	27 J	14 I	19	15
Trichloroethene	ug/L	120	79	89	140	210	73	68	110

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Location ID: Zone: Date Collected:	Units	EW-110 LSAS 05/21/07	EW-110 LSAS 06/20/07	EW-110 LSAS 07/16/07	EW-110 LSAS 08/16/07	EW-110 LSAS 09/26/07	EW-110 LSAS 10/24/07	EW-110 LSAS 11/08/07	EW-110 LSAS 01/30/08	EW-UAFG-1 AF Gravel 02/01/08
Semivolatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	270	200 J	120	180	200	190	290	27	200
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	R	NA	R	NA	NA	NA
1,1-Dichloroethane	ug/L	250	160	280	200	150	150	120	28	34
1,1-Dichloroethene	ug/L	760	510	990 D	980	390	430	570	68	250 D
cis-1,2-Dichloroethene	ug/L	16	10	11	8.6	8.6 I	12	11	5.4	3,000 D
Tetrachloroethene	ug/L	28	17	9.8	9.6	8.9 I	13	21	4	0.5 U
Trichloroethene	ug/L	130	70	94	84	69	83	120	73	19

**APPENDIX D-2
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Notes:

AF Gravel = Arcadian Formation Gravels.

LSAS = Lower Shallow Aquifer System.

Lower AF = Lower Arcadia Formation.

S&P Sand = Salt & Pepper sands.

USAS = Upper Surficial Aquifer System.

ug/L = micrograms per liter.

D = The value is the result of a secondary dilution.

E = Sample result is greater than calibration range

I = Detected but below reporting limit. Result is an estimated concentration.

J = Estimated value.

L = Estimated value, biased low.

Q = Sample held beyond accepted holding time.

R = Rejected.

U = The analyte was analyzed for, but not detected.

UJ = The analyte was analyzed for, but not detected. The reporting limit is an estimated value.

V = Indicates the analyte was detected in both the sample and the associated method blank.

[] = Duplicate sample result.

APPENDIX D-3
PRIVATE WELL DETAILS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map Well ID	Property Address	Owner's Name	Alt ID ¹	Well Usage ²	Closure Status ³	Closure Date	Total Depth (feet) ⁴	Casing Diameter (inches)	Zone
1	7620 15TH ST E	NA	NA	Former Monitoring	Closed	Unknown	NA	NA	AF Gravel
2	7300 15TH ST E	Nathan Benderson	NA	Former Monitoring	Closed	Unknown	NA	NA	AF Gravel
3	1306 ROME AVE	Theodor Vreuls	79	Former Irrigation	Closed	4/1/2006	460	NA	AF Gravel
4	1309 HARDIN AVE	Stephen Crowe	70	Former Potable	Closed	5/3/2006	139	NA	AF Gravel
5	1227/1231 HARDIN AVE	NA	69	Former Monitoring	Closed	Unknown	NA	NA	AF Gravel
6	7580 15TH ST E	Revco Properties LLC	NA	Current Monitoring	Open	NA	15	2"	USAS
7	7561/7571 15TH ST E	Dr. W. G. Thomas (Animal Hospital)	7/8	Current Potable	Open	NA	150	NA	AF Gravel
8	7602 16TH ST CT E	Pinkney Hall	24	Former Irrigation	Closed	Unknown	NA	NA	AF Gravel
9	7604/7608 16TH ST E	Willie Robinson	16/17	Former Potable	Closed	Unknown	74.5	NA	LSAS
10	7609 16TH ST E	Robert Smith	18	Former Potable	Closed	10/13/1992	18	NA	USAS
11	7616 16TH ST E	Marvin Washington	20	Former Potable	Closed	7/21/2006	98	NA	AF Gravel
12	7620 16TH ST E	Dorothy Mazon	21	Former Irrigation	Closed	7/25/2006	22	NA	USAS
13	7624 16TH ST E	Eugene Sloan	23	Former Irrigation	Closed	7/25/2006	91	NA	LSAS
14,16	1507 TALLEVAST RD	Frank Williams	83/100	Former Irrigation	Closed	3/6/2006	85	NA	LSAS
15	7671 15TH ST E	Quick Stop, Inc.	NA	Current Monitoring	Open	NA	20	2"	USAS
17	7621 16TH ST E	Beatrice Graydon	22	Former Potable	Closed	7/21/2006	233	NA	AF Gravel
18	1615 TALLEVAST RD	Mount Tabor Missionary	86	Former Irrigation	Closed	7/21/2006	119	NA	AF Gravel*
19	7605 17TH ST E	Flora Jean Bryant	25	Former Irrigation	Closed	7/24/2006	65	NA	LSAS*
20	1709 76TH AVE DR E	Flora Jean Bryant	61	Former Irrigation	Closed	Unknown	NA	NA	AF Gravel
21	1710/1714 76TH AVE DR E	Fred Bryant	62	Former Irrigation	Closed	7/24/2006	65	NA	LSAS
22	7619 17TH ST E	Ernestine Rowe	26	Former Irrigation	Closed	7/24/2006	26	NA	USAS
23	7623 17TH ST E	James Moore	27	Former Irrigation	Closed	7/13/2006	27	NA	USAS
24	1712 TALLEVAST RD	Wyman Pitts	87	Former Irrigation	Closed	7/18/2006	32	NA	USAS
25	7515 18TH ST E	Bennie Tooley	32	Former Potable	Closed	7/20/2006	23	NA	USAS
26	7519 18TH ST E	Mrs. Essie Sims	33	Former Potable	Closed	7/26/2006	171	NA	AF Gravel
27	7609 18TH ST E	Joel Carsten McNair	34/99	Former Potable	Closed	3/3/2006	154	NA	AF Gravel
28	7611 18TH ST E	Joel Carsten McNair	36	Former Potable	Closed	4/10/2006	116	NA	AF Gravel
29	7615 18TH ST E	Sylvester Brown	37	Former Potable	Closed	2/16/2006	65	NA	LSAS
30	7619 18TH ST E	Annie Mae Bryant	38	Former Potable	Closed	4/10/2006	105	NA	AF Gravel
31	1811 TALLEVAST RD	Joseph Bryant	90	Former Potable	Closed	7/18/2006	77	NA	LSAS
32	7624 19TH ST E	Willy McKnight	39	Former Potable	Closed	7/20/2006	114	NA	AF Gravel
33	7600 19TH ST E	Michael Wesley	40	Former Irrigation	Closed	7/25/2006	140	NA	AF Gravel
34	7603 19TH ST E	Calvin Sloan / Louise Sloan	41	Former Potable	Closed	5/11/2006	130	NA	AF Gravel*
35	1911/1913 TALLEVAST RD	Est. of John Stanley	93	Former Potable	Closed	1/16/2007	66	NA	LSAS
36	1955 TALLEVAST RD	Cliff Heathington	94	Former Irrigation	Closed	7/26/2006	33	NA	LSAS
37	2003 TALLEVAST RD	Lillian Granderson	95	Former Potable	Closed	12/10/2006	114	NA	AF Gravel
38	2105 TALLEVAST RD	Heidi Boothe	96	Current Irrigation	Open	NA	NA	NA	AF Gravel
39	1812 TALLEVAST RD	C Billy Ward	91	Former Irrigation	Closed	7/25/2006	74	NA	LSAS*
40	1808 TALLEVAST RD	Arthur Bryant / Mary Montgomery	89	Former Irrigation	Closed	11/28/2006	75	NA	LSAS
41	1864 TALLEVAST RD	Clifford Ward	92	Former Potable	Closed	7/20/2006	127	NA	LSAS
42	1804 TALLEVAST RD	Delores Williams	88	Former Potable	Closed	7/26/2006	213	NA	AF Gravel
43	7715 17TH ST CT E	Lewis Pryor	29	Former Irrigation	Closed	7/24/2006	110	NA	LSAS
44	7716 17TH ST CT E	Lizzie Davis	30	Former Potable	Closed	7/17/2006	30	NA	LSAS
45	7813 17TH ST CT E	Tim H Wallace	31	Former Irrigation	Closed	7/24/2006	15	NA	USAS*
46	7819 17TH ST E	Guirlene Duvignaud	28	Former Irrigation	Closed	7/24/2006	15	NA	USAS
47	7851 15TH ST E	Sarasota-Manatee Airport Authority (SRQ)	12	Current Irrigation	Open	NA	480	NA	Floridan
48	7741 15TH ST E	Sarasota-Manatee Airport Authority (SRQ)	11	Former Irrigation	Closed	11/16/2000	407	NA	Floridan*
49	7715 15TH ST E	Whogas Inc	10	Current Irrigation	Open	NA	460	4"	AF Gravel*
50	1234 CLYDE JONES RD	Sarasota-Manatee Airport Authority (SRQ)	68	Former Monitoring	Closed	Unknown	NA	NA	USAS
51	7850 15TH ST E	NA	NA	Former Monitoring	Closed	Unknown	NA	NA	AF Gravel
52	7881 15TH ST E	Sarasota-Manatee Airport Authority (SRQ)	NA	Current Monitoring	Open	NA	12	2"	USAS
53	8011 15TH ST E	Innovation Marine	13/73	Current Irrigation	Open	NA	375	NA	Floridan
54	8061 15TH ST E	NA	NA	Former Monitoring	Closed	Unknown	NA	NA	AF Gravel
55	7350 26TH CT E	NA	52	Current Monitoring	Open	NA	15	NA	USAS
56	7455 16TH ST E	Mader Electric Inc	15	Current Irrigation	Open	NA	NA	NA	AF Gravel
57	7500 26TH CT E	NA	53	Current Irrigation	Open	NA	440	NA	Floridan
58	7501 15TH ST	Goodwill Ind-Manasota, Inc	6	Former Irrigation	Closed	Unknown	NA	NA	AF Gravel
59	7524 COMMERCE PLACE	Chris Wentzell	65	Current Irrigation	Open	NA	NA	NA	AF Gravel
60	7575 COMMERCE CT	Commerce Court LLC	66	Current Irrigation	Open	NA	NA	NA	AF Gravel
61	7602 17TH ST E	Mount Tabor Missionary	NA	Former Irrigation	Closed	7/27/2006	120	NA	AF Gravel*
62	7603 18TH ST E	Joel Carsten McNair	35	Former Potable	Closed	Unknown	NA	NA	AF Gravel
63	1012 PONDEROSA PINE LANE	Janette Marie Girard	78	Current Irrigation	Open	NA	NA	NA	AF Gravel
64	1107 TALLEVAST RD	NA	81	Current Irrigation	Open	NA	NA	NA	AF Gravel
65	1201 TALLEVAST RD	NA	82	Current Irrigation	Open	NA	NA	NA	AF Gravel
66	1305 HARDIN AVE	NA	70	Former Potable	Closed	7/17/2006	120	NA	AF Gravel
67	1375 HARDIN AVE	Eric Woolley	71	Former Potable	Closed	Unknown	NA	NA	AF Gravel
68	1401 COMMERCE BLVD	Studio 21 Lighting Inc	67	Former Irrigation	Closed	Unknown	NA	NA	AF Gravel
69	1403 HARDIN AVE	Revco Properties LLC	72	Former Potable	Closed	Unknown	NA	NA	AF Gravel
70	1600 TALLEVAST RD	NA	NA	Former Irrigation	Closed	12/1/2006	178	NA	AF Gravel
71	1607 TALLEVAST RD	Vanessa Peterson	85	Former Irrigation	Closed	Unknown	NA	NA	AF Gravel
72	1611 TALLEVAST RD	NA	NA	Former Irrigation	Closed	11/29/2006	90	NA	AFG
73	1619 TALLEVAST RD	Clifford Ward	NA	Former Potable	Closed	7/21/2006	119	NA	AF Gravel
74	1701 BIOTECH WAY	HQ Properties, Inc	64	Current Irrigation	Open	NA	NA	NA	AF Gravel
75	1715 TALLEVAST RD	Ulysses Ward	NA	Former Irrigation	Closed	12/1/2006	84.2	NA	LSAS
76	1790 TALLEVAST RD	Dexter Baldwin	NA	Former Irrigation	Closed	12/13/2006	49	NA	LSAS
77	1807 TALLEVAST RD	Bryant Chapel	NA	Former Irrigation	Closed	11/28/2006	41	NA	USAS
78	1905 72ND DR E	Glasshouse LLC	55	Current Irrigation	Open	NA	120	NA	AF Gravel
79	1915 72ND DR E	Richard Bouck	56	Current Irrigation	Open	NA	125	NA	AF Gravel
80	2217 72ND AVE E	Rucha LLC	58	Current Irrigation	Open	NA	125	NA	AF Gravel
81	2227 72ND AVE E	Keith Goldstein	59	Current Irrigation	Open	NA	120	NA	AF Gravel
82	2305 72ND AVE E	Gap Properties of SW Florida-1 Inc	60	Current Irrigation	Open	NA	125	NA	AF Gravel

APPENDIX D-3
PRIVATE WELL DETAILS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map Well ID	Property Address	Owner's Name	Alt ID ¹	Well Usage ²	Closure Status ³	Closure Date	Total Depth (feet) ⁴	Casing Diameter (inches)	Zone
83	2337 72ND DR E	Clara Gingerich	57	Current Irrigation	Open	NA	130	NA	AF Gravel
84	2400 TALLEVAST RD	Walter Schmidt (owner not listed on permits)	NA	Current Irrigation	Open	NA	NA	NA	AF Gravel
85	2411 TALLEVAST RD	Covered Bride Holdings II LLC	98	Current Irrigation	Open	NA	NA	NA	AF Gravel
86	7061 15TH ST E	Genmar Ind Inc	1	Current Irrigation	Open	NA	NA	NA	AF Gravel
87	7116 24TH CT E	NA	49	Current Irrigation	Open	NA	120	NA	AF Gravel
88	7126 24TH CT E	NA	51	Current Irrigation	Open	NA	120	NA	AF Gravel
89	7175 21ST ST E	Florida Loaders Enterprises, Inc.	42	Current Irrigation	Open	NA	109	NA	AF Gravel*
90	7198 21ST ST E	NA	NA	Current Irrigation	Open	NA	NA	NA	AF Gravel
91	7205/7207 21ST ST E	McLeod Land Services inc	44	Current Irrigation	Open	NA	NA	NA	AF Gravel
92	7212 24TH CT E	Osborne Enterprises LLC	48	Current Irrigation	Open	NA	120	NA	AF Gravel
93	7216 21ST ST E	Florida Loaders Enterprises, Inc.	46	Current Irrigation	Open	NA	125	NA	AF Gravel
94	7220 24TH CT E	In Tandem Inc	50	Current Irrigation	Open	NA	120	NA	AF Gravel
95	7230 15TH ST E	Schmitt Family Partnership	4	Current Irrigation	Open	NA	NA	NA	AF Gravel
96	7245 21ST ST E	K E B Holdings LLC	47	Current Irrigation	Open	NA	NA	NA	AF Gravel
97	7260 15TH ST E	Mader Electric Inc	5	Current Irrigation	Open	NA	NA	NA	AF Gravel
98	7519 18TH ST E (East Well)	Essie Mae Sims	NA	Current Potable	Open	NA	NA	NA	AF Gravel
99	7606 16TH ST CT E	Robert Smith	NA	Former Irrigation	Closed	11/29/2006	150.5	NA	AF Gravel
100	7630 MATOAKA RD	NA	101	Current Irrigation	Open	NA	NA	NA	AF Gravel
101	7845 27TH ST E	Lucien Levy	54	Current Irrigation	Open	NA	NA	NA	AF Gravel
102	8005 15TH ST E	NA	NA	Current Irrigation	Open	NA	NA	NA	AF Gravel
103	8155 27TH ST E	NA	102	Current Irrigation	Open	NA	NA	NA	AF Gravel
104	8161 15TH ST E	CC Property Acquisition LLC	14	Current Irrigation	Open	NA	NA	NA	USAS
105	904 PONDEROSA PINE LANE	Richard Cashman	75	Current Irrigation	Open	NA	NA	NA	AF Gravel
106	905 PONDEROSA PINE LANE	Richard Cashman	76	Current Irrigation	Open	NA	NA	NA	AF Gravel
107	911 SOUTHERN PINE LANE	NA	80	Current Irrigation	Open	NA	NA	NA	AF Gravel
108	916 PONDEROSA PINE LANE	John Beghtel	77	Former Irrigation	Closed	Unknown	NA	NA	AF Gravel
109	7603 19TH ST E Well #2	Calvin Sloan / Louise Sloan ⁵	NA	Former Potable	Closed	5/9/2006	112	NA	AF Gravel*
110	7619 18TH ST E Well #2	Frank Bryant ⁵	NA	Former Potable	Closed	4/28/2006	320	NA	Floridan*
111	1710/1714 76TH AVE DR E Well #2	Fred Bryant ⁵	NA	Former Irrigation	Closed	7/18/2006	144	NA	AF Gravel*
112	7609 16TH ST E Well #2	Calvin Bryant ⁵	NA	Current Potable	Open	NA	140	NA	LSAS
113	1613 76TH AVE DR E	Earlene Walton Baker ⁵	NA	Former Potable	Closed	2/27/2006	86	NA	LSAS
114	7604 19TH ST E	Alley Hornes/ Michael Wesley ⁵	NA	Potable	Unknown	Unknown	50	NA	LSAS*
115	7828 17TH ST CT E	Michelle Bland ⁵	NA	Unknown	Closed	11/1/2002	40	NA	USAS*
116	7205 15TH ST E	Benderson Development Co. ⁵	NA	Former Irrigation	Closed	6/29/1999	780	NA	Floridan*
117	7205 15TH ST E Well #2	Bradenton Commerce Center ⁵	NA	Current Potable	Open	NA	235	NA	Lower AF*
118	7616 17TH ST CT E	Mr. Robinson ⁵	NA	Unknown	Closed	5/16/1973	NA	NA	
119	7201 15TH ST E	S.E.S. Groves (Midway Groves) ⁵	NA	Former Irrigation	Closed	5/7/1994	764	NA	Floridan*
120	2224 72ND AVE E	CD Development LLP ⁵	NA	Current Irrigation	Open	NA	125	NA	AF Gravel*
121	7178 21ST ST E	N.M.B., LLC ⁵	NA	Current Irrigation	Open	NA	120	NA	AF Gravel*

Notes:

NA - not available.

AF Gravel = Arcadian Formation Gravels.

LSAS = Lower Shallow Aquifer System.

Lower AF = Lower Arcadia Formation.

USAS = Upper Surficial Aquifer System.

* Aquifer zone assignments reflect consistency with other nearby wells that have similar well construction details.

¹ID from the May 2006 Private Well Survey Report.

²Current Irrigation - an open well that is primarily used for watering landscape plants; Former Irrigation - a closed irrigation well;

Current Monitoring - an open monitoring well installed for the purpose of water level measurements and/or collecting water quality samples;

Former Monitoring - a closed monitoring well; Current Potable - an open well that may have been used as a source of drinking and bathing water;

Former Potable - a closed potable well.

³Based on available information from the Lockheed Martin well closure program

⁴Reported total well depths were taken from Southwest Florida Water Management District (SWFWMD) well closure reports, where available.

⁵The owner's name listed is the owner at the time of well permitting.

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	3 1306 ROME AVE AF Gravel 07/11/05	4 1309 HARDIN AVE AF Gravel 04/11/05	4 1309 HARDIN AVE AF Gravel 05/11/05	4 1309 HARDIN AVE AF Gravel 07/12/05	4 1309 HARDIN AVE AF Gravel 03/19/06
Volatile Organics					
1,4-Dioxane	ug/L	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA
Volatile Organics (524)					
1,1-Dichloroethane	ug/L	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA
Volatile Organics-Field					
1,1-Dichloroethane	ug/L	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA
Semivolatile Organics (8270C)					
1,4-Dioxane	ug/L	2 U	NA	1 U	1.9 U
Volatile Organics (8260B) - SIM					
1,4-Dioxane	ug/L	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution					
1,4-Dioxane	ug/L	NA	NA	NA	NA
Volatile Organics (8260B)					
1,4-Dioxane	ug/L	NA	50 U	NA	NA
1,1-Dichloroethane	ug/L	0.5 U	0.2 U	NA	0.5 U
1,1-Dichloroethene	ug/L	0.5 U	0.2 U	NA	0.45 U
cis-1,2-Dichloroethene	ug/L	0.5 U	0.5 U	NA	0.65 U
Tetrachloroethene	ug/L	0.5 U	0.2 U	NA	0.34 U
Trichloroethene	ug/L	0.5 U	0.5 U	NA	0.28 U

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	7 7561/7571 15TH ST E AF Gravel 04/11/05	7 7561/7571 15TH ST E AF Gravel 05/11/05	7 7561/7571 15TH ST E AF Gravel 07/14/05	7 7561/7571 15TH ST E AF Gravel 03/08/06	7 7561/7571 15TH ST E AF Gravel 03/14/06	7 7561/7571 15TH ST E AF Gravel 03/16/06	7 7561/7571 15TH ST E AF Gravel 03/21/06	7 7561/7571 15TH ST E AF Gravel 12/20/06	7 7561/7571 15TH ST E AF Gravel 02/01/08
Volatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	1 U	1.9 U	1 U	1 U	1 U	1 U	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	1.1 J
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	0.76 U
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	50 U	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.2 U	NA	0.5 U	0.52 U	0.52 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.2 U	NA	0.5 U	0.45 U	0.45 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.5 U	NA	0.5 U	0.65 U	0.65 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.2 U	NA	0.5 U	0.34 U	0.34 U	0.34 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.5 U	NA	0.5 U	0.28 U	0.28 U	0.28 U	0.28 U	0.71 U

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	8 7602 16TH ST CT E AF Gravel 12/21/06	9 7604/7608 16TH ST E LSAS 05/20/04	9 7604/7608 16TH ST E LSAS 05/24/04	11 7616 16TH ST E AF Gravel 05/20/04	11 7616 16TH ST E AF Gravel 05/24/04	11 7616 16TH ST E AF Gravel 08/29/05	11 7616 16TH ST E AF Gravel 09/19/05	12 7620 16TH ST E AF Gravel 05/20/04	12 7620 16TH ST E AF Gravel 05/24/04	12 7620 16TH ST E AF Gravel 08/29/05
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	9.3 U	9.3 U	NA	NA	9.3 U
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	0.4 U	0.24 U	NA	NA	0.4 U
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	3.7	3	NA	NA	1.6
Tetrachloroethene	ug/L	NA	NA	NA	NA	0.17 U	0.26 U	NA	NA	0.17 U
Trichloroethene	ug/L	NA	NA	NA	NA	1.6	1.2	NA	NA	1.3
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	NA	0.23 U	0.38 I	0.23 U	0.36 I	NA	NA	0.23 U	0.29 I
1,1-Dichloroethene	ug/L	NA	0.95	1.1	0.7	0.98	NA	NA	0.23 U	0.19 U
cis-1,2-Dichloroethene	ug/L	NA	0.51	0.76	2.7	2.9	NA	NA	0.51	0.67
Tetrachloroethene	ug/L	NA	0.3 U	0.25 U	0.3 U	0.5 U	NA	NA	0.3 U	0.25 U
Trichloroethene	ug/L	NA	1.6	1.8	2.8	3.6	NA	NA	1.1	1.1
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	R	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	0.76 U	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.6 U	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	0.83 U	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	0.75 U	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	0.65 U	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	0.71 U	NA	NA	NA	NA	NA	NA	NA	NA

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	13 7624 16TH ST E LSAS 05/20/04	13 7624 16TH ST E LSAS 05/25/04	13 7624 16TH ST E LSAS 09/03/05	14,16 1507 TALLEVAST RD LSAS 09/09/05	17 7621 16TH ST E AF Gravel 05/20/04	17 7621 16TH ST E AF Gravel 05/24/04	17 7621 16TH ST E AF Gravel 08/31/05	20 1709 76TH AVE DR E AF Gravel 09/14/05	21 1710/1714 76TH AVE DR E LSAS 05/26/04	
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	9.3 U	9.3 U	NA	NA	9.3 U	9.3 U	NA
1,1-Dichloroethene	ug/L	NA	NA	0.22 U	0.22 U	NA	NA	0.37 I	11	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	2.4	2.3	NA	NA	0.41 I	22	NA
Tetrachloroethene	ug/L	NA	NA	0.09 U	0.09 U	NA	NA	0.09 U	0.09 U	NA
Trichloroethene	ug/L	NA	NA	1.1	0.84	NA	NA	1.2	57	NA
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	0.23 U	0.21 U	NA	NA	0.23 U	0.44 I	NA	NA	12
1,1-Dichloroethene	ug/L	0.23 U	0.19 U	NA	NA	0.23 U	0.19 U	NA	NA	98
cis-1,2-Dichloroethene	ug/L	3.2	3.6	NA	NA	0.25 U	0.28 I	NA	NA	5
Tetrachloroethene	ug/L	0.3 U	0.25 U	NA	NA	0.3 U	0.25 U	NA	NA	0.25 U
Trichloroethene	ug/L	1.4	1.7	NA	NA	0.22 U	0.21 U	NA	NA	15
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	22 7619 17TH ST E AF Gravel 05/25/04	22 7619 17TH ST E AF Gravel 09/14/05	24 1712 TALLEVAST RD USAS 09/09/05	25 7515 18TH ST E USAS 05/20/04	25 7515 18TH ST E USAS 05/25/04	25 7515 18TH ST E USAS 08/29/05	26 7519 18TH ST E AF Gravel 05/20/04	26 7519 18TH ST E AF Gravel 05/25/04	26 7519 18TH ST E AF Gravel 08/24/05	
Volatile Organics										
1,4-Dioxane	ug/L	NA	130	100	NA	NA	9.3 U	NA	NA	9.3 U
1,1-Dichloroethene	ug/L	NA	55	99	NA	NA	0.4 U	NA	NA	0.22 U
cis-1,2-Dichloroethene	ug/L	NA	14	50	NA	NA	0.17 U	NA	NA	0.11 U
Tetrachloroethene	ug/L	NA	0.09 U	1.8	NA	NA	0.17 U	NA	NA	0.09 U
Trichloroethene	ug/L	NA	41	730	NA	NA	0.27 U	NA	NA	0.14 U
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	0.21 U	NA	NA	0.23 U	0.21 U	NA	0.23 U	0.21 U	NA
1,1-Dichloroethene	ug/L	0.19 U	NA	NA	0.23 U	0.19 U	NA	0.23 U	0.19 U	NA
cis-1,2-Dichloroethene	ug/L	0.19 U	NA	NA	0.25 U	0.19 U	NA	0.25 U	0.19 U	NA
Tetrachloroethene	ug/L	0.25 U	NA	NA	0.3 U	0.25 U	NA	0.3 U	0.25 U	NA
Trichloroethene	ug/L	0.21 U	NA	NA	0.22 U	0.21 U	NA	0.22 U	0.21 U	NA
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	27 7609 18TH ST E AF Gravel 05/20/04	27 7609 18TH ST E AF Gravel 05/25/04	27 7609 18TH ST E AF Gravel 09/03/05	28 7611 18TH ST E AF Gravel 05/20/04	28 7611 18TH ST E AF Gravel 05/25/04	28 7611 18TH ST E AF Gravel 08/24/05	29 7615 18TH ST E LSAS 05/20/04	29 7615 18TH ST E LSAS 05/25/04	29 7615 18TH ST E LSAS 08/24/05	
Volatiles Organics										
1,4-Dioxane	ug/L	NA	NA	9.3 U	NA	NA	9.3 U	NA	NA	9.3 U
1,1-Dichloroethene	ug/L	NA	NA	0.22 U	NA	NA	0.22 U	NA	NA	0.22 U
cis-1,2-Dichloroethene	ug/L	NA	NA	0.11 U	NA	NA	1.8	NA	NA	17
Tetrachloroethene	ug/L	NA	NA	0.09 U	NA	NA	0.09 U	NA	NA	0.09 U
Trichloroethene	ug/L	NA	NA	0.2 I	NA	NA	4.8	NA	NA	120
Volatiles Organics (524)										
1,1-Dichloroethane	ug/L	0.23 U	0.21 U	NA	0.23 I	0.27 I	NA	0.7	0.21 U	NA
1,1-Dichloroethene	ug/L	0.23 U	0.19 U	NA	0.41 I	0.54	NA	3	3.2	NA
cis-1,2-Dichloroethene	ug/L	0.25 U	0.19 U	NA	1.1	1.5	NA	43	44	NA
Tetrachloroethene	ug/L	0.3 U	0.25 U	NA	0.3 U	0.25 U	NA	0.3 U	0.25 U	NA
Trichloroethene	ug/L	0.22 U	0.21 U	NA	2.3	3.3	NA	220 C	240	NA
Volatiles Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatiles Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatiles Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatiles Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatiles Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	30 7619 18TH ST E AF Gravel 05/20/04	30 7619 18TH ST E AF Gravel 05/25/04	30 7619 18TH ST E AF Gravel 08/29/05	30 7619 18TH ST E AF Gravel 09/19/05	31 1811 TALLEVAST RD LSAS 05/20/04	31 1811 TALLEVAST RD LSAS 05/25/04	31 1811 TALLEVAST RD LSAS 08/29/05	32 7624 19TH ST E AF Gravel 05/20/04	32 7624 19TH ST E AF Gravel 05/25/04	32 7624 19TH ST E AF Gravel 09/09/05
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	9.3 U	9.3 U	NA	NA	9.3 U	NA	9.3 U
1,1-Dichloroethene	ug/L	NA	NA	0.4 U	1.6	NA	NA	2.4	NA	4.8
cis-1,2-Dichloroethene	ug/L	NA	NA	7.1	24	NA	NA	30	NA	21
Tetrachloroethene	ug/L	NA	NA	0.17 U	0.26 U	NA	NA	0.17 U	NA	0.09 U
Trichloroethene	ug/L	NA	NA	44	150 L	NA	NA	160	NA	89
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	1	0.27 I	NA	NA	2.4	2.5	NA	2.7	2.9
1,1-Dichloroethene	ug/L	2.6	0.68	NA	NA	3.4	3.8	NA	7.5 C	7.6
cis-1,2-Dichloroethene	ug/L	22	9.2	NA	NA	26	32	NA	24	30
Tetrachloroethene	ug/L	0.33 I	0.25 U	NA	NA	0.3 U	0.25 U	NA	0.3 U	0.25 U
Trichloroethene	ug/L	180 C	110	NA	NA	180 C	240	NA	95 C	140
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	33 7600 19TH ST E LSAS 05/26/04	33 7600 19TH ST E LSAS 08/31/05	35 1911/1913 TALLEVAST RD AF Gravel 05/20/04	35 1911/1913 TALLEVAST RD AF Gravel 05/25/04	35 1911/1913 TALLEVAST RD AF Gravel 09/03/05	35 1911/1913 TALLEVAST RD AF Gravel 01/16/07
Volatile Organics						
1,4-Dioxane	ug/L	NA	120	NA	NA	9.3 U
1,1-Dichloroethene	ug/L	NA	1.8	NA	NA	0.22 U
cis-1,2-Dichloroethene	ug/L	NA	27	NA	NA	0.98
Tetrachloroethene	ug/L	NA	0.09 U	NA	NA	0.09 U
Trichloroethene	ug/L	NA	140	NA	NA	0.14 U
Volatile Organics (524)						
1,1-Dichloroethane	ug/L	1.6	NA	0.23 U	0.21 U	NA
1,1-Dichloroethene	ug/L	3.4	NA	0.23 U	0.19 U	NA
cis-1,2-Dichloroethene	ug/L	46	NA	1.2	1	NA
Tetrachloroethene	ug/L	0.25 U	NA	0.3 U	0.25 U	NA
Trichloroethene	ug/L	370	NA	0.42 I	0.23 I	NA
Volatile Organics-Field						
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution						
1,4-Dioxane	ug/L	NA	NA	NA	NA	5.1
Volatile Organics (8260B)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	0.6 U
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	0.83 U
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	0.75 UJ
Tetrachloroethene	ug/L	NA	NA	NA	NA	0.65 U
Trichloroethene	ug/L	NA	NA	NA	NA	0.71 UJ

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	36 1955 TALLEVAST RD AF Gravel 05/20/04	36 1955 TALLEVAST RD AF Gravel 05/25/04	36 1955 TALLEVAST RD AF Gravel 08/29/05	37 2003 TALLEVAST RD AF Gravel 05/20/04	37 2003 TALLEVAST RD AF Gravel 05/25/04	37 2003 TALLEVAST RD AF Gravel 08/29/05	37 2003 TALLEVAST RD AF Gravel 11/27/06	37 2003 TALLEVAST RD AF Gravel 12/20/06	
Volatile Organics									
1,4-Dioxane	ug/L	NA	NA	9.3 U	NA	NA	9.3 U	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	2	NA	NA	0.4 U	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	0.17 U	NA	NA	0.17 U	NA	NA
Tetrachloroethene	ug/L	NA	NA	0.17 U	NA	NA	0.17 U	NA	NA
Trichloroethene	ug/L	NA	NA	0.46 I	NA	NA	7.8	NA	NA
Volatile Organics (524)									
1,1-Dichloroethane	ug/L	2.5	0.54	NA	3.3	3.9	NA	NA	NA
1,1-Dichloroethene	ug/L	4	0.75	NA	5.5	5.9	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	0.56	0.19 U	NA	0.74	0.82	NA	NA	NA
Tetrachloroethene	ug/L	0.3 U	0.25 U	NA	0.3 U	0.25 U	NA	NA	NA
Trichloroethene	ug/L	0.57	0.26 I	NA	0.7	0.64	NA	NA	NA
Volatile Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	R	17 J
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	0.76 U	28
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	0.6 U	0.7 I
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	0.83 U	0.83 U
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	0.75 U	1.2 J
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	0.65 U	0.65 U
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	0.71 U	25 J

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HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	38 2105 TALLEVAST RD AF Gravel 08/24/05	38 2105 TALLEVAST RD AF Gravel 03/01/06	38 2105 TALLEVAST RD AF Gravel 03/27/06	38 2105 TALLEVAST RD AF Gravel 12/15/06	38 2105 TALLEVAST RD AF Gravel 02/01/08	40 1808 TALLEVAST RD AF Gravel 09/03/05	40 1808 TALLEVAST RD AF Gravel 11/27/06	41 1864 TALLEVAST RD AF Gravel 08/24/05	
Volatile Organics									
1,4-Dioxane	ug/L	66	NA	NA	NA	NA	320	NA	65
1,1-Dichloroethene	ug/L	10	NA	NA	NA	NA	130	NA	2.1
cis-1,2-Dichloroethene	ug/L	1.5	NA	NA	NA	NA	18	NA	0.35 I
Tetrachloroethene	ug/L	0.09 U	NA	NA	NA	NA	0.36 U	NA	0.09 U
Trichloroethene	ug/L	0.58	NA	NA	NA	NA	80	NA	0.58
Volatile Organics (524)									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	4.4 I	1.1 U	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	6.6 J	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	3.8	96	NA	0.76 U	NA
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	0.52 U	0.52 U	0.6 I	3.4	NA	0.6 U	NA
1,1-Dichloroethene	ug/L	NA	0.45 U	0.45 U	1.2	4.9	NA	0.83 U	NA
cis-1,2-Dichloroethene	ug/L	NA	0.65 U	0.65 U	0.75 U	0.93 I	NA	0.75 U	NA
Tetrachloroethene	ug/L	NA	0.34 U	0.34 U	0.65 U	0.5 U	NA	0.65 U	NA
Trichloroethene	ug/L	NA	0.28 U	0.28 U	0.71 U	0.5 U	NA	0.71 U	NA

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	42 1804 TALLEVAST RD AF Gravel 05/20/04	42 1804 TALLEVAST RD AF Gravel 05/25/04	42 1804 TALLEVAST RD AF Gravel 09/09/05	43 7715 17TH ST CT E LSAS 08/31/05	44 7716 17TH ST CT E LSAS 05/25/04	44 7716 17TH ST CT E LSAS 08/31/05	46 7819 17TH ST E LSAS 05/25/04	46 7819 17TH ST E LSAS 08/31/05	
Volatile Organics									
1,4-Dioxane	ug/L	NA	NA	330	390	NA	510	NA	480
1,1-Dichloroethene	ug/L	NA	NA	24	61	NA	61	NA	92
cis-1,2-Dichloroethene	ug/L	NA	NA	4.7	14	NA	20	NA	43
Tetrachloroethene	ug/L	NA	NA	0.48 I	5.1	NA	45	NA	0.09 U
Trichloroethene	ug/L	NA	NA	6.9	100	NA	270	NA	15
Volatile Organics (524)									
1,1-Dichloroethane	ug/L	18	12	NA	NA	100	NA	110	NA
1,1-Dichloroethene	ug/L	35 C	28	NA	NA	160	NA	94	NA
cis-1,2-Dichloroethene	ug/L	6.3	4.3	NA	NA	20	NA	50	NA
Tetrachloroethene	ug/L	1.9	2.7	NA	NA	30	NA	0.25 U	NA
Trichloroethene	ug/L	7.4 C	4.6	NA	NA	510	NA	11	NA
Volatile Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID:	47	47	47	47	47	47	47	47	47	47	47	50	52
Location ID:	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	7851 15TH ST E	1234 CLYDE JONES RD	7881 15TH ST E
Zone:	Floridan	Floridan	Floridan	Floridan	Floridan	Floridan	Floridan	Floridan	Floridan	Floridan	AF Gravel	USAS	AF Gravel
Date Collected:	06/18/04	11/01/04	05/26/05	03/09/06	03/14/06	03/16/06	03/21/06	12/20/06	01/30/08	01/30/08		12/27/06	03/20/06
Volatile Organics													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)													
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field													
1,1-Dichloroethane	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)													
1,4-Dioxane	ug/L	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	NA	1 U
Volatile Organics (8260B) - SIM													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	1.7 J	NA	NA	1.1 J
Volatile Organics (8260) - SIM Isotope Dilution													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	0.76 U	1.7	1.6	1.9 I
Volatile Organics (8260B)													
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.2 U	NA	NA	0.52 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.2 U	NA	NA	0.45 U	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	0.65 U	0.65 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.2 U	NA	NA	0.34 U	0.34 U	0.34 U	0.34 U	0.65 U	0.5 U	0.5 U	0.65 U	0.34 U
Trichloroethene	ug/L	0.5 U	NA	NA	1.6	1.6	2	1.6	1.2	1.9	2.2	0.71 U	0.28 U

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	53 8011 15TH ST E Lower AF - Floridan 07/06/05	53 8011 15TH ST E Lower AF - Floridan 04/14/06	53 8011 15TH ST E Lower AF - Floridan 12/20/06	53 8011 15TH ST E Lower AF - Floridan 01/31/07	55 7350 26TH CT E AF Gravel 12/27/06	56 7455 16TH ST E AF Gravel 04/05/06	56 7455 16TH ST E AF Gravel 12/21/06	
Volatile Organics								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)								
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field								
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)								
1,4-Dioxane	ug/L	2 U	1 U	NA	NA	NA	1 U	NA
Volatile Organics (8260B) - SIM								
1,4-Dioxane	ug/L	NA	NA	0.95 J	NA	R	NA	1.2 J [R]
Volatile Organics (8260) - SIM Isotope Dilution								
1,4-Dioxane	ug/L	NA	NA	0.76 U	0.76 U	0.76 U	NA	0.76 U [0.76 U]
Volatile Organics (8260B)								
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.5 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U	0.6 U [0.6 U]
1,1-Dichloroethene	ug/L	0.5 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U	0.83 U [0.83 U]
cis-1,2-Dichloroethene	ug/L	0.5 U	0.65 U	0.75 U	0.75 U	0.75 U	0.65 U	0.75 U [0.75 U]
Tetrachloroethene	ug/L	0.5 U	0.34 U	0.65 U	0.65 U	0.65 U	0.34 U	0.65 U [0.65 U]
Trichloroethene	ug/L	0.5 U	0.28 U	0.71 U	0.71 U	0.71 U	3.3	1.7 [1.8]

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	57 7500 26TH CT E AF Gravel 12/29/06	57 7500 26TH CT E AF Gravel 02/01/08	58 7501 15TH ST AF Gravel 04/04/06	58 7501 15TH ST AF Gravel 12/20/06	59 7524 COMMERCE PLACE AF Gravel 02/01/08	60 7575 COMMERCE CT AF Gravel 02/21/07	64 1107 TALLEVAST RD AF Gravel 12/27/06	64 1107 TALLEVAST RD AF Gravel 02/04/08	
Volatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	1 UJ	NA	NA	NA	NA	NA
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	R	NA	NA	R	NA	NA	R	NA
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	0.76 U	1 U	NA	0.76 U	1 U	1 I	0.76 U	1 U
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U	0.75 U	0.75 U	0.65 U
Tetrachloroethene	ug/L	0.65 U	0.5 U	0.34 U	0.65 U	0.5 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.71 U	0.5 U	0.28 U	0.71 U	0.5 U	0.71 U	0.71 U	0.5 U

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	65 1201 TALLEVAST RD AF Gravel 12/20/06	65 1201 TALLEVAST RD AF Gravel 02/04/08	70 1600 TALLEVAST RD AF Gravel 11/27/06	72 1611 TALLEVAST RD USAS 11/27/06	73 1619 TALLEVAST RD AF Gravel 08/31/05	73 1619 TALLEVAST RD AF Gravel 09/14/05	74 1701 BIOTECH WAY AF Gravel 04/03/06	74 1701 BIOTECH WAY AF Gravel 01/31/07	
Volatile Organics									
1,4-Dioxane	ug/L	NA	NA	NA	9.3 U	9.3 U	NA	NA	
1,1-Dichloroethene	ug/L	NA	NA	NA	16	20	NA	NA	
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	8.6	13	NA	NA	
Tetrachloroethene	ug/L	NA	NA	NA	0.09 U	0.09 U	NA	NA	
Trichloroethene	ug/L	NA	NA	NA	77	120	NA	NA	
Volatile Organics (524)									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Volatile Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	1 U	NA	
Volatile Organics (8260B) - SIM									
1,4-Dioxane	ug/L	R	NA	R	R	NA	NA	NA	
Volatile Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	0.76 U	1 U	0.76 U	0.76 U	NA	NA	0.76 U	
Volatile Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.6 U	0.52 U	0.6 U	0.6 U	NA	NA	0.52 U	
1,1-Dichloroethene	ug/L	0.83 U	0.45 U	0.83 U	0.83 U	NA	NA	0.45 U	
cis-1,2-Dichloroethene	ug/L	0.75 UJ	0.65 U	0.75 U	0.75 U	NA	NA	0.65 U	
Tetrachloroethene	ug/L	0.65 U	0.5 U	0.65 U	0.65 U	NA	NA	0.34 U	
Trichloroethene	ug/L	0.71 UJ	0.5 U	0.71 U	0.71 U	NA	NA	0.28 U	

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	75 1715 TALLEVAST RD AF Gravel 11/27/06	77 1807 TALLEVAST RD AF Gravel 11/27/06	78 1905 72ND DR E AF Gravel 04/14/06	79 1915 72ND DR E AF Gravel 04/10/06	79 1915 72ND DR E AF Gravel 04/21/06	79 1915 72ND DR E AF Gravel 12/20/06	80 2217 72ND AVE E AF Gravel 04/12/06	80 2217 72ND AVE E AF Gravel 04/19/06	80 2217 72ND AVE E AF Gravel 04/21/06	80 2217 72ND AVE E AF Gravel 12/20/06
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	NA	1 UJ	11	1 U	NA	1 U	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	79 J	2.8 J	NA	NA	NA	1.2 J	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	88	3.2	NA	NA	NA	0.76 U	NA	NA	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	6.4	0.6 U	0.52 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	5.8	0.83 U	0.45 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	1.9	1	0.65 U	0.65 U	0.65 U	0.75 UJ	0.65 U	0.65 U	0.75 UJ
Tetrachloroethene	ug/L	0.65 U	0.65 U	0.34 U	0.34 U	0.34 UJ	0.65 U	0.34 U	0.34 UJ	0.65 U
Trichloroethene	ug/L	3.5	0.71 UJ	0.28 U	1.2	0.87 I	0.71 UJ	0.28 U	0.28 U	1 J

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	81 2227 72ND AVE E AF Gravel 04/12/06	81 2227 72ND AVE E AF Gravel 12/20/06	82 2305 72ND AVE E AF Gravel 12/20/06	83 2337 72ND DR E AF Gravel 04/09/06	83 2337 72ND DR E AF Gravel 04/21/06	83 2337 72ND DR E AF Gravel 12/26/06	84 2400 TALLEVAST RD AF Gravel 08/29/05	84 2400 TALLEVAST RD AF Gravel 01/24/08	
Volatiles Organics									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	9.3 U	NA	
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	0.4 U	NA	
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	0.17 U	NA	
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	0.17 U	NA	
Trichloroethene	ug/L	NA	NA	NA	NA	NA	0.27 U	NA	
Volatiles Organics (524)									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Volatiles Organics-Field									
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	
Semivolatile Organics (8270C)									
1,4-Dioxane	ug/L	1 U	NA	NA	4 I	1 U	NA	NA	
Volatiles Organics (8260B) - SIM									
1,4-Dioxane	ug/L	NA	R	R	NA	NA	R	NA	
Volatiles Organics (8260) - SIM Isotope Dilution									
1,4-Dioxane	ug/L	NA	0.76 U	0.76 U	NA	NA	0.76 U	1 U [1 U]	
Volatiles Organics (8260B)									
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	
1,1-Dichloroethane	ug/L	0.52 U	0.6 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U [0.52 U]	
1,1-Dichloroethene	ug/L	0.45 U	0.83 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U [0.45 U]	
cis-1,2-Dichloroethene	ug/L	0.65 U	0.75 UJ	0.75 UJ	0.65 U	0.65 U	0.75 U	0.65 U [0.65 U]	
Tetrachloroethene	ug/L	0.34 U	0.65 U	0.65 U	0.34 U	0.34 UJ	0.65 U	0.5 U [0.5 U]	
Trichloroethene	ug/L	0.28 U	0.71 UJ	0.71 UJ	0.28 U	0.28 U	0.71 U	0.5 U [0.5 U]	

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	85 2411 TALLEVAST RD AF Gravel 03/23/06	85 2411 TALLEVAST RD AF Gravel 03/30/06	85 2411 TALLEVAST RD AF Gravel 12/12/06	85 2411 TALLEVAST RD AF Gravel 01/31/08	86 7061 15TH ST E AF Gravel 04/12/06	86 7061 15TH ST E AF Gravel 12/20/06	87 7116 24TH CT E AF Gravel 04/03/06	87 7116 24TH CT E AF Gravel 01/31/07	88 7126 24TH CT E AF Gravel 12/27/06	
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	19	1 U	NA	NA	1 U	NA	1 U	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	5.3 J	NA	NA	1.2 J	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	4.4	22	NA	0.76 U	NA	0.76 U	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.8 I	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.6 U	0.6 U
1,1-Dichloroethene	ug/L	0.82 I	1.1	0.83 U	0.55 I	0.45 U	0.83 U	0.45 U	0.83 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 UJ	0.65 U	12	5.7	0.65 U	0.75 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.34 U	0.65 U	0.5 U	0.34 U	0.65 U	0.34 U	0.65 U	0.65 U
Trichloroethene	ug/L	0.28 U	0.28 U	0.71 UJ	0.5 U	0.53 I	1.7	0.28 U	0.71 U	0.71 U

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	91 7205/7207 21ST ST E AF Gravel 04/09/06	91 7205/7207 21ST ST E AF Gravel 04/21/06	91 7205/7207 21ST ST E AF Gravel 12/27/06	92 7212 24TH CT E AF Gravel 04/09/06	92 7212 24TH CT E AF Gravel 04/21/06	92 7212 24TH CT E AF Gravel 12/27/06	93 7216 21ST ST E AF Gravel 04/09/06	93 7216 21ST ST E AF Gravel 04/21/06	93 7216 21ST ST E AF Gravel 12/27/06	94 7220 24TH CT E AF Gravel 04/09/06
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	15	1 U	NA	3.3 I	1 U	NA	4.6 I	1 U	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	NA	NA	R	NA	NA	R	NA	NA	R
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	NA	NA	0.76 U	NA	NA	0.76 U	NA	NA	0.76 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.34 U	0.34 UJ	0.65 U	0.34 U	0.34 UJ	0.65 U	0.34 U	0.34 UJ	0.65 U
Trichloroethene	ug/L	0.92 I	0.28 U	0.71 U	0.28 U	0.28 U	0.71 U	5.6	1.2	2.6

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID: Location ID: Zone: Date Collected:	95 7230 15TH ST E AF Gravel 12/20/06	96 7245 21ST ST E AF Gravel 04/10/06	96 7245 21ST ST E AF Gravel 04/21/06	96 7245 21ST ST E AF Gravel 12/27/06	97 7260 15TH ST E AF Gravel 04/05/06	97 7260 15TH ST E AF Gravel 12/20/06	99 7606 16TH ST CT E AF Gravel 11/27/06	100 7630 MATOAKA RD AF Gravel 12/20/06	102 8005 15TH ST E AF Gravel 02/01/08	103 8155 27TH ST E AF Gravel 12/20/06
Volatile Organics										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics (524)										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Volatile Organics-Field										
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)										
1,4-Dioxane	ug/L	NA	4.5 I	45	NA	1.1 U	NA	NA	NA	NA
Volatile Organics (8260B) - SIM										
1,4-Dioxane	ug/L	R	NA	NA	R	NA	0.96 J	2.7 J	1.2 J	NA
Volatile Organics (8260) - SIM Isotope Dilution										
1,4-Dioxane	ug/L	0.76 U	NA	NA	0.76 U	NA	0.76 U	1.9 I	0.76 U	1 U
Volatile Organics (8260B)										
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.6 U	0.52 U	0.52 U	0.6 U	0.52 U	0.6 U	0.6 U	0.6 U	0.52 U
1,1-Dichloroethene	ug/L	0.83 U	0.45 U	0.45 U	0.83 U	0.45 U	0.83 U	0.83 U	0.83 U	0.45 U
cis-1,2-Dichloroethene	ug/L	0.75 U	0.65 U	0.65 U	0.75 U	0.65 U	0.75 UJ	0.75 U	0.75 UJ	0.65 U
Tetrachloroethene	ug/L	0.65 U	0.34 U	0.34 UJ	0.65 U	0.34 U	0.65 U	0.65 U	0.65 U	0.5 U
Trichloroethene	ug/L	0.71 U	0.28 U	0.28 U	0.71 U	0.28 U	0.71 UJ	0.71 U	0.71 UJ	0.5 U

APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA

Map ID:	104	104	104	105	105	112
Location ID:	8161 15TH ST E	8161 15TH ST E	8161 15TH ST E	904 PONDEROSA PINE LANE	904 PONDEROSA PINE LANE	7609 16TH ST E Well #2
Zone:	USAS	USAS	USAS	AF Gravel	AF Gravel	LSAS
Date Collected:	08/18/05	04/05/06	12/21/06	04/04/06	12/27/06	08/31/05
Volatile Organics						
1,4-Dioxane	ug/L	NA	NA	NA	NA	9.3 U
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	6.8
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	5
Tetrachloroethene	ug/L	NA	NA	NA	NA	0.09 U
Trichloroethene	ug/L	NA	NA	NA	NA	40
Volatile Organics (524)						
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA
Volatile Organics-Field						
1,1-Dichloroethane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethene	ug/L	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	ug/L	NA	NA	NA	NA	NA
Tetrachloroethene	ug/L	NA	NA	NA	NA	NA
Trichloroethene	ug/L	NA	NA	NA	NA	NA
Semivolatile Organics (8270C)						
1,4-Dioxane	ug/L	1.9 U	1 U	NA	1 U	NA
Volatile Organics (8260B) - SIM						
1,4-Dioxane	ug/L	NA	NA	0.94 J	NA	R
Volatile Organics (8260) - SIM Isotope Dilution						
1,4-Dioxane	ug/L	NA	NA	0.76 U	NA	0.76 U
Volatile Organics (8260B)						
1,4-Dioxane	ug/L	NA	NA	NA	NA	NA
1,1-Dichloroethane	ug/L	0.5 U	0.52 U	0.6 U	0.52 U	0.6 U
1,1-Dichloroethene	ug/L	0.5 U	0.45 U	0.83 U	0.45 U	0.83 U
cis-1,2-Dichloroethene	ug/L	0.5 U	0.65 U	0.75 U	0.65 U	0.75 U
Tetrachloroethene	ug/L	0.5 U	0.34 U	0.65 U	0.34 U	0.65 U
Trichloroethene	ug/L	0.5 U	0.28 U	0.71 U	0.99 I	0.71 U

**APPENDIX D-4
HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL RESULTS**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA**

Notes:

AF Gravel = Arcadian Formation Gravels

LSAS = Lower Shallow Aquifer System

Lower AF = Lower Arcadia Formation

S&P Sand = Salt & Pepper sands

USAS = Upper Surficial Aquifer System

ug/L = micrograms per liter

C = Identification confirmed by gas chromatograph/mass spectrometer (GC/MS).

I = Detected but below reporting limit. Result is an estimated concentration.

J = Estimated value.

L = Estimated value, biased low.

R = Rejected.

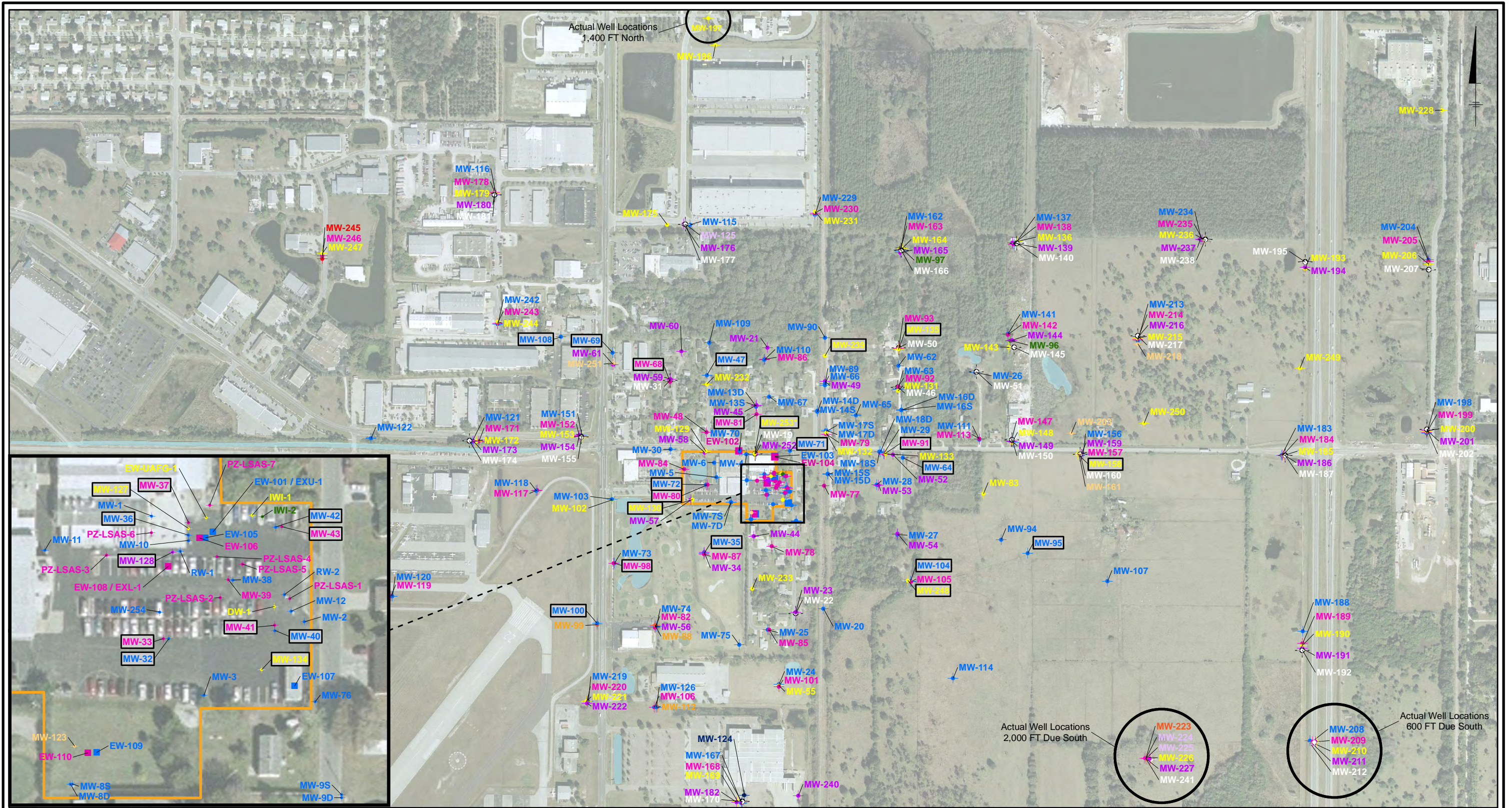
U = The analyte was analyzed for, but not detected.

UJ = The analyte was analyzed for, but not detected. The reporting limit is an estimated value.

[] = Duplicate sample result.

Appendix E

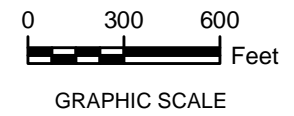
Concentration vs. Time Graphs for
Representative Wells



- Extraction Wells (USAS)
- Extraction Wells (LSAS)
- Extraction Wells (AF Gravel)
- ★ Monitoring Wells (USAS)
- ★ Monitoring Wells (Hardstreak Clay)
- ★ Monitoring Wells (LSAS)
- ★ Monitoring Wells (Venice Clay)
- ★ Monitoring Wells (Zone 1 Clay)
- ★ Monitoring Wells (AF Gravel)
- ★ Monitoring Wells (Zone 2 Clay)
- ★ Monitoring Wells (S&P Sand)
- ★ Monitoring Wells (Zone 3-4 Clay)
- ★ Monitoring Wells (Lower AF)
- ★ Monitoring Wells (Floridan)

- ▭ Former ABC Facility
- ▭ Ponds
- ▭ MW-100 Wells Selected for Concentration Versus Time Graphs

Note:
* Indicates well is planned for future graphing



FORMER AMERICAN BERYLLIUM COMPANY SITE
TALLEVAST, FLORIDA
2008 GROUNDWATER MONITORING REPORT

**REPRESENTATIVE WELLS SELECTED FOR
CONCENTRATION VERSUS TIME GRAPHS**

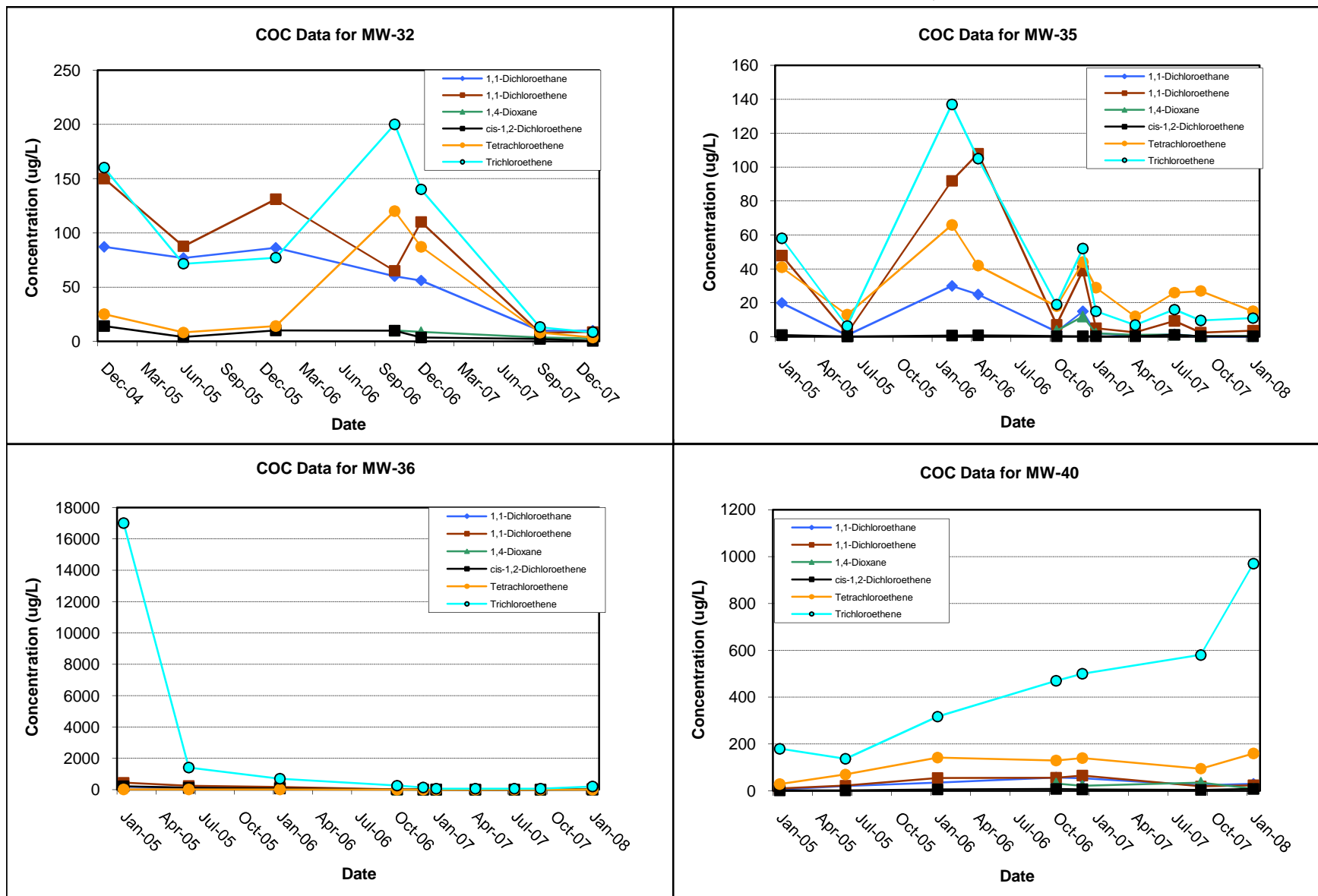
ARCADIS

FIGURE
E-1

DEN-080 MSS
Project #: B0038055.0000

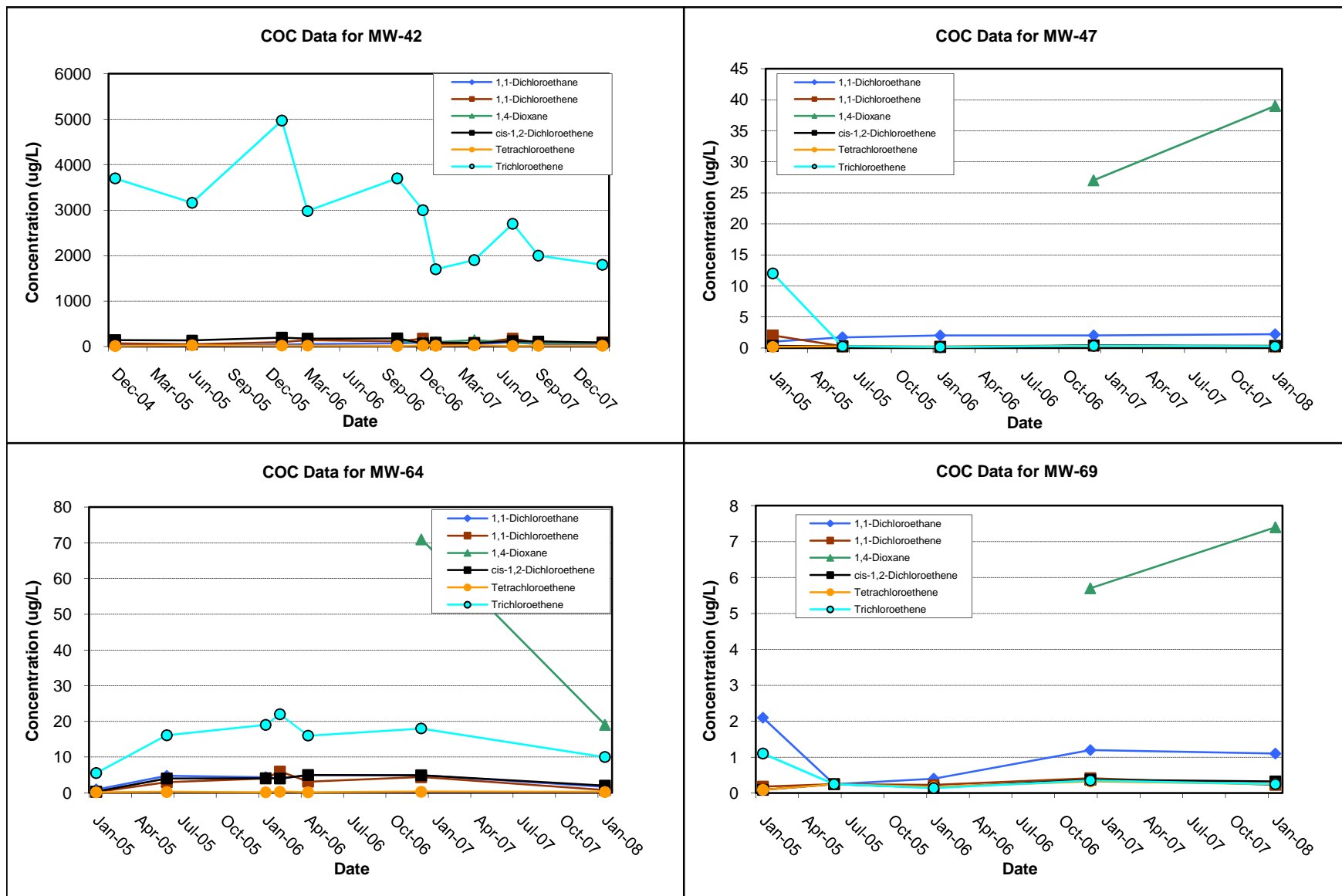
**APPENDIX E
CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
USAS ZONE**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



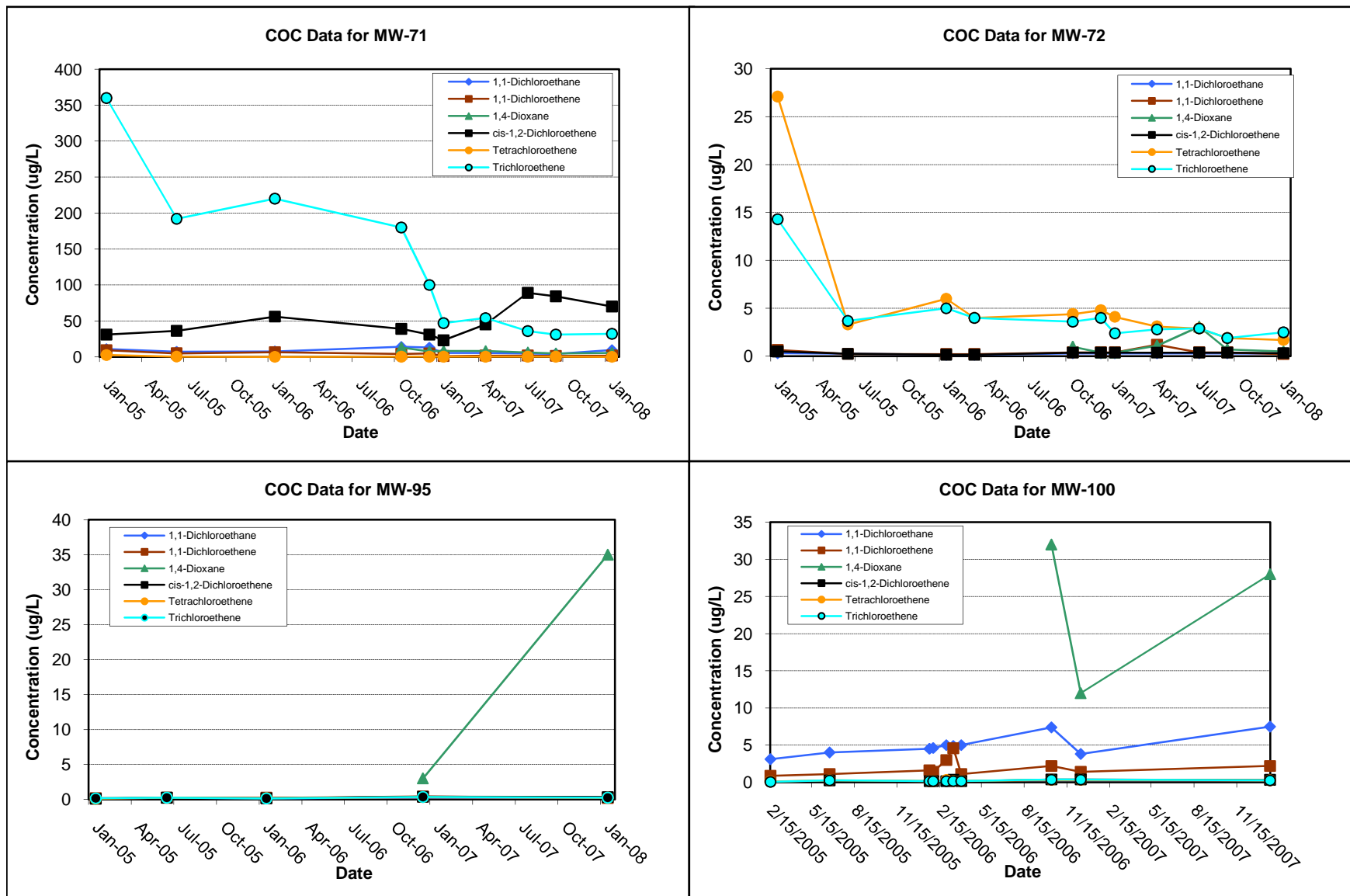
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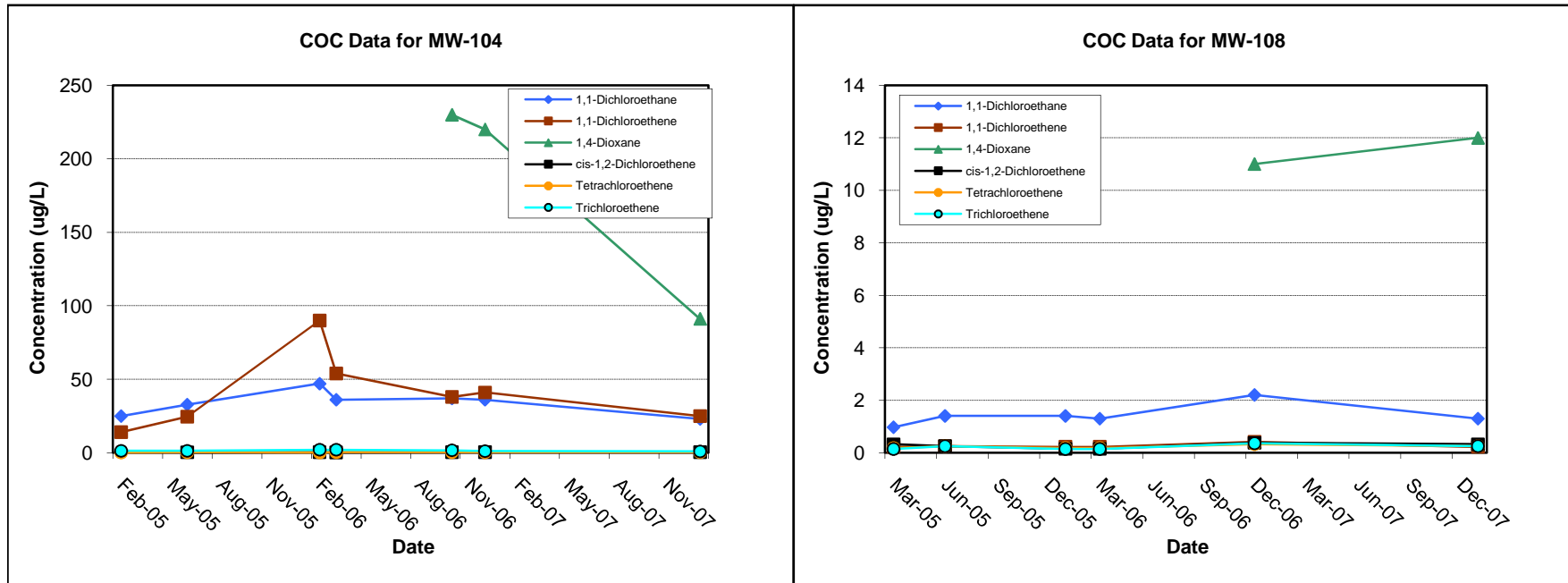
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USAS ZONE**

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FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



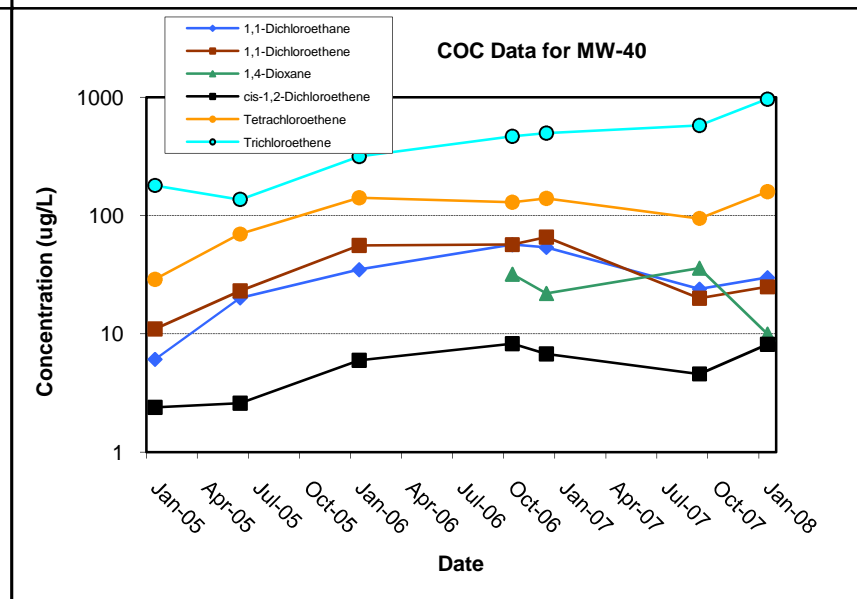
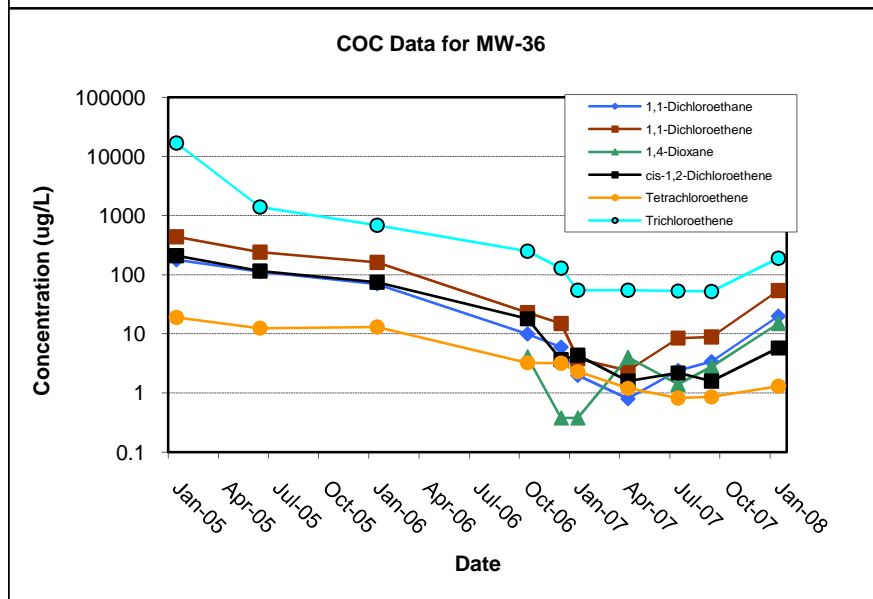
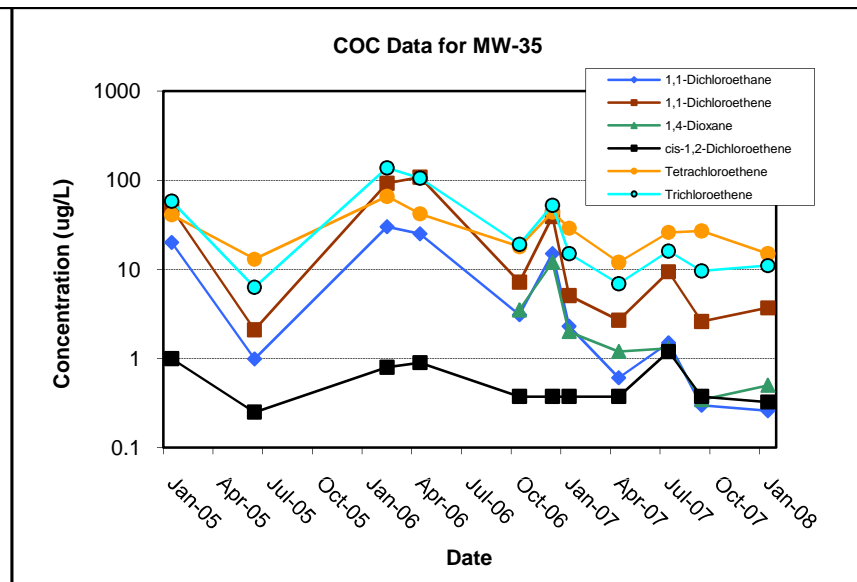
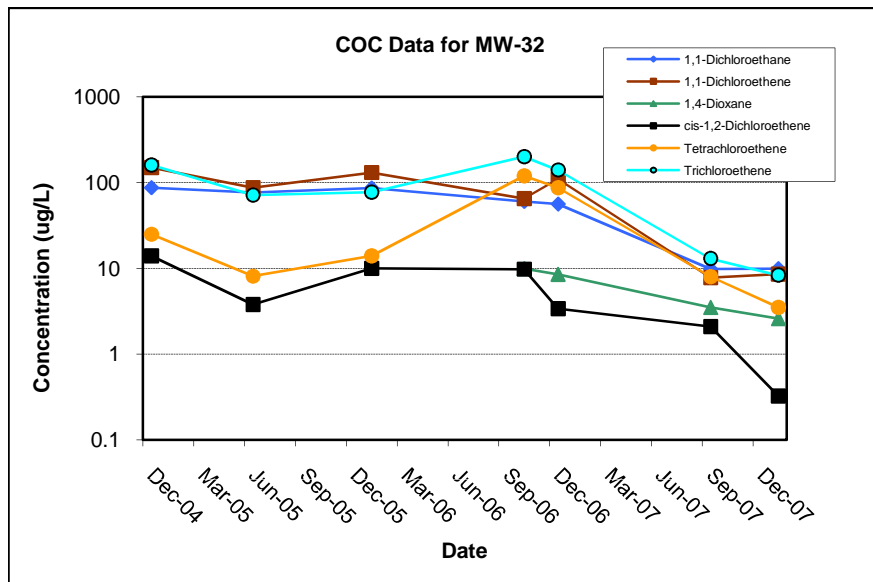
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CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
USAS ZONE**

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FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



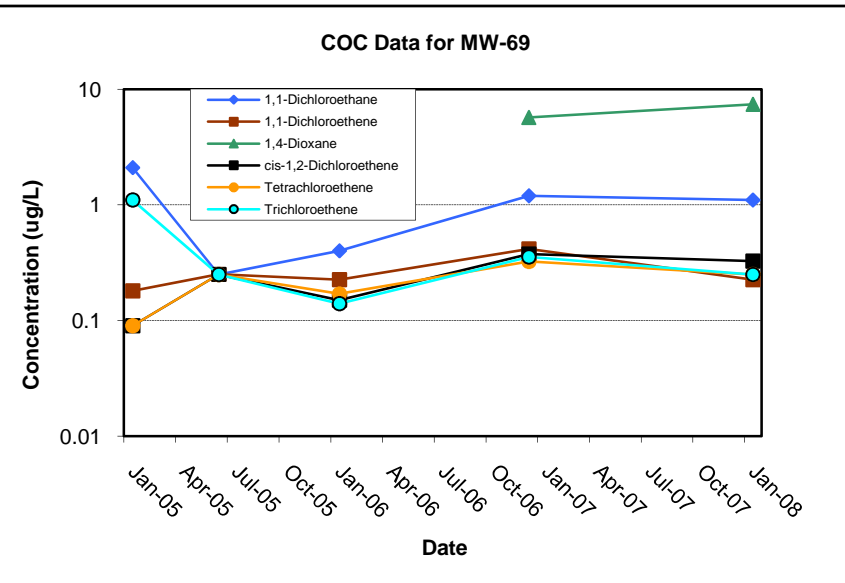
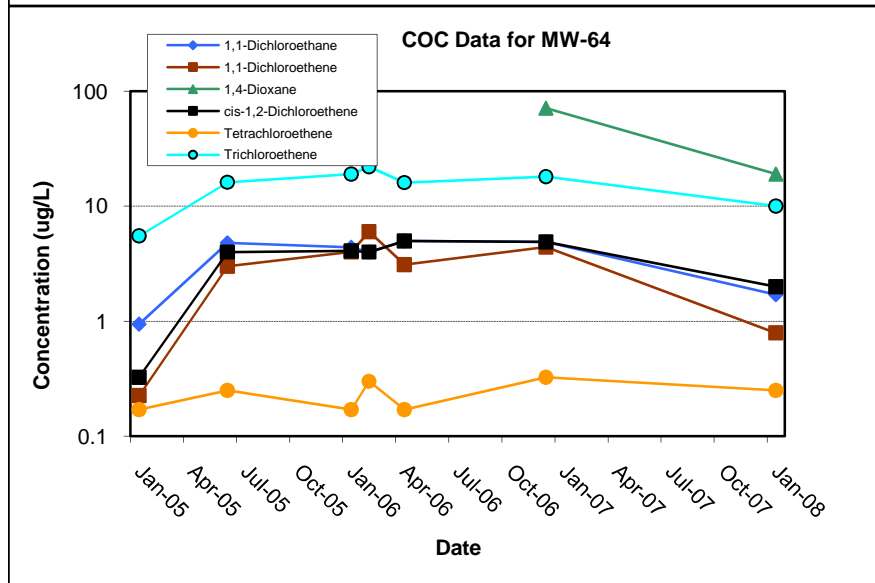
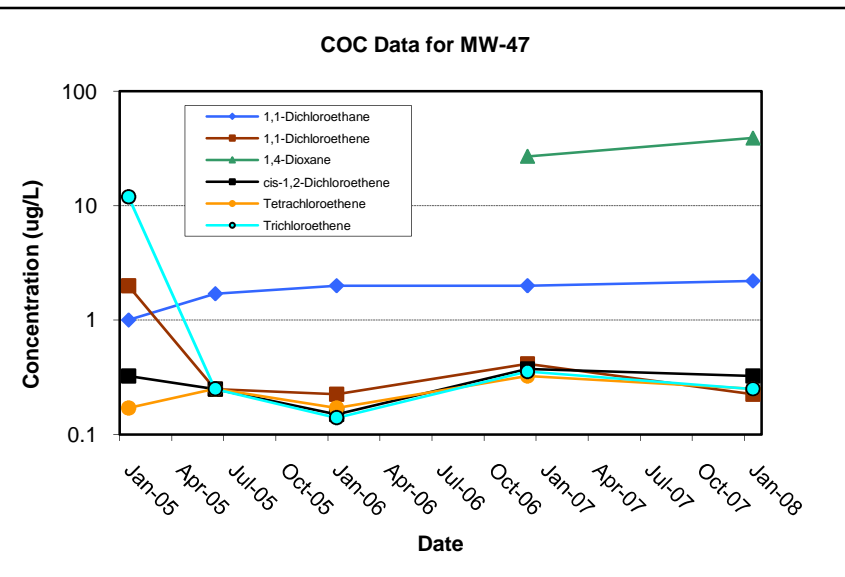
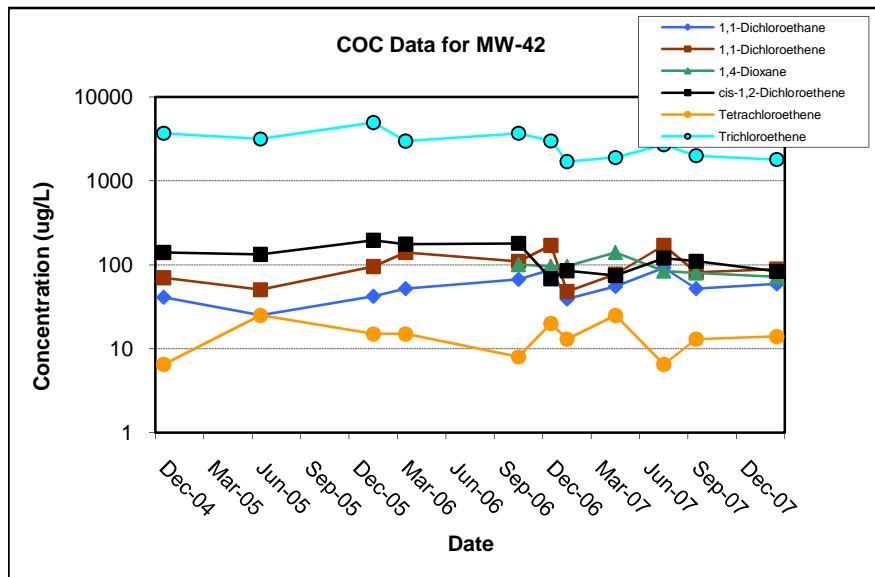
**APPENDIX E
LOG CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
USAS ZONE**

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FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



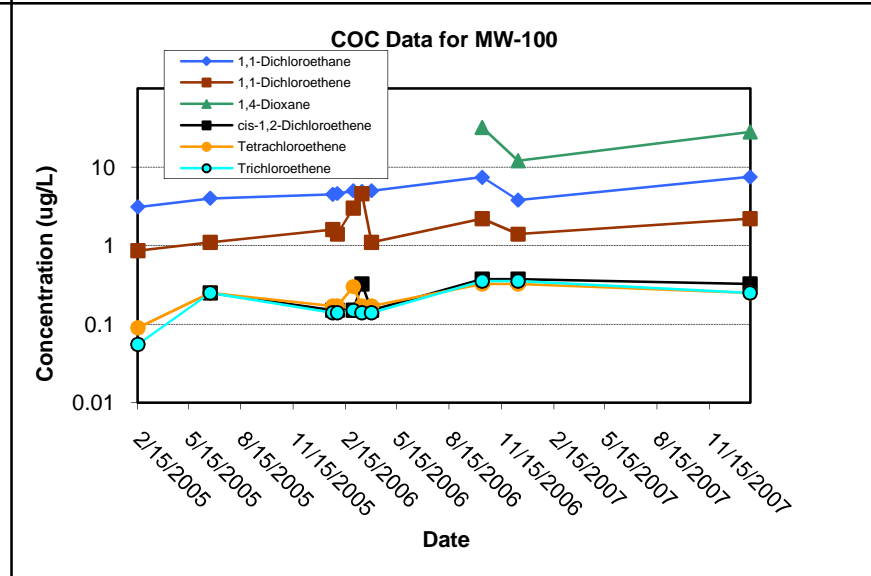
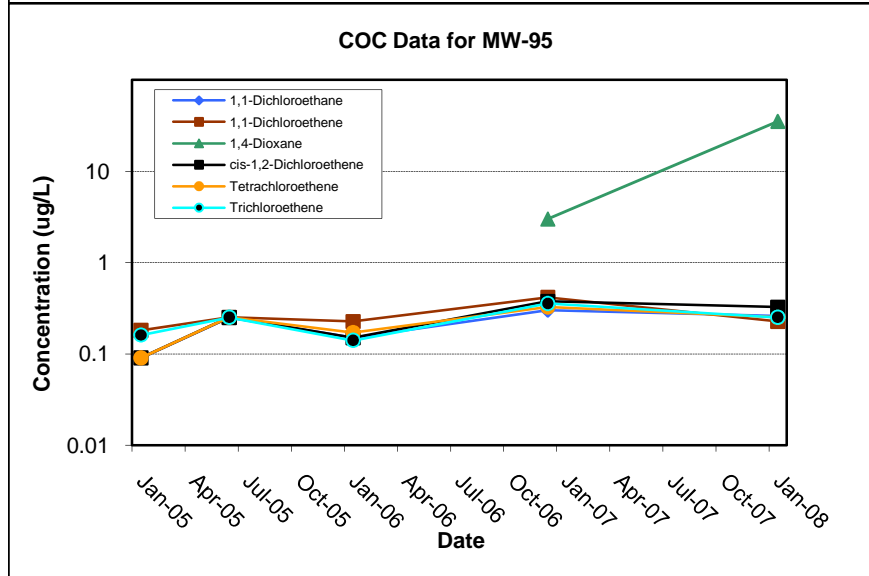
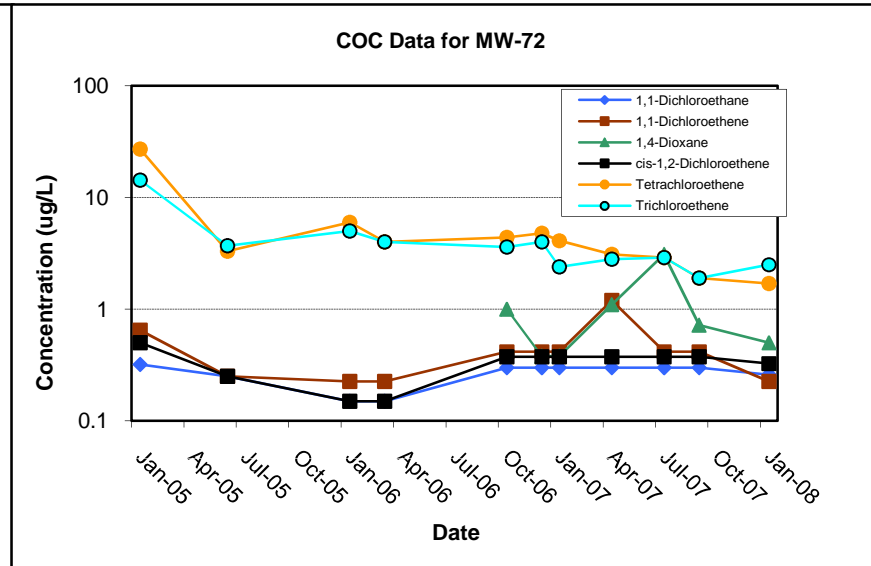
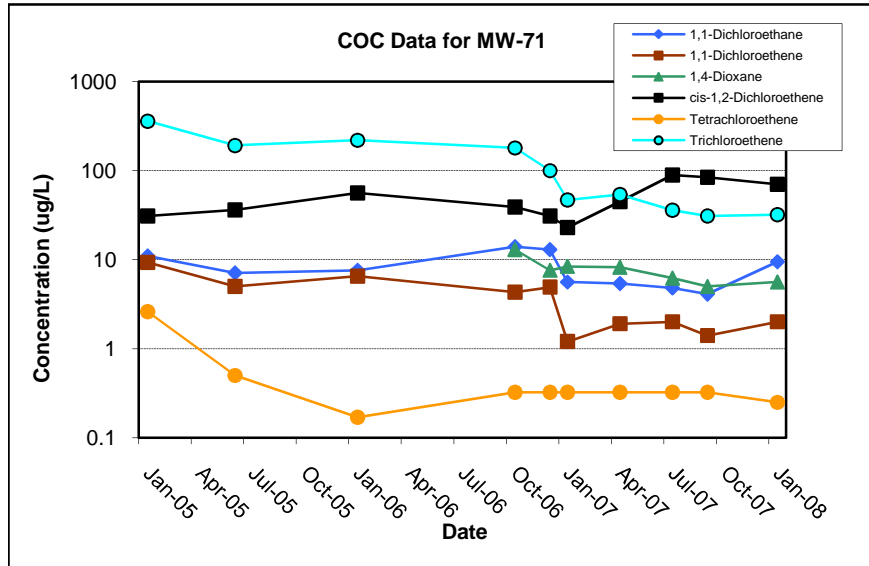
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USAS ZONE**

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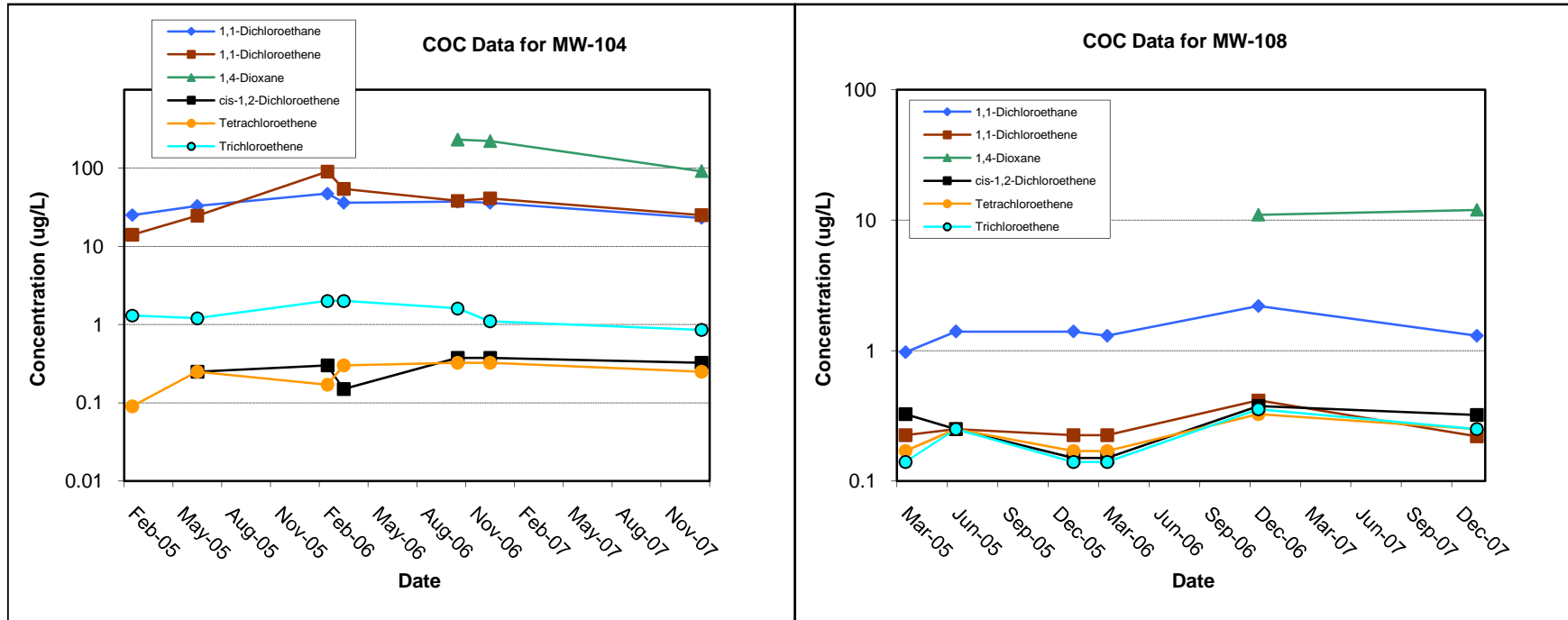
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USAS ZONE**

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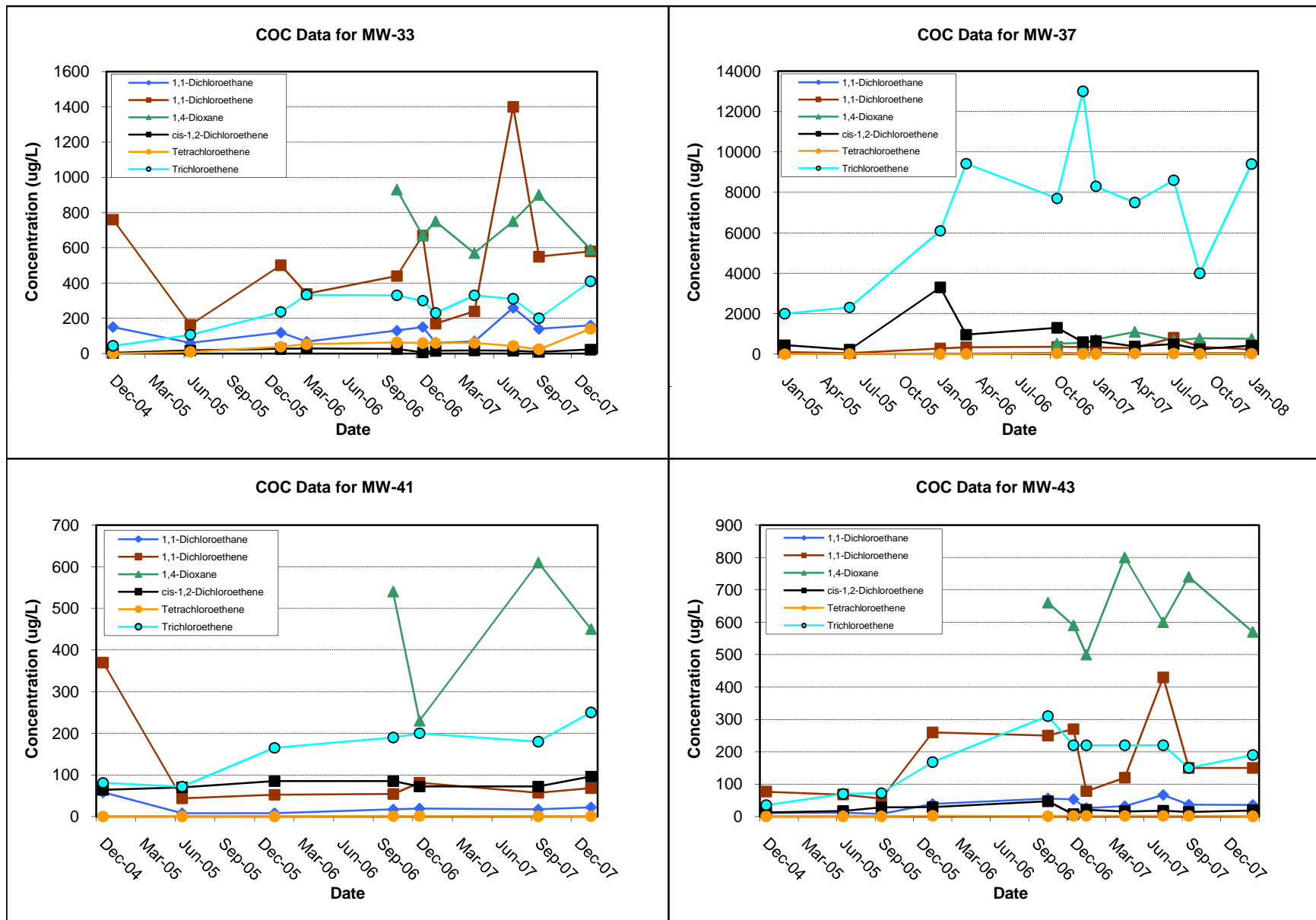
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USAS ZONE**

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FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



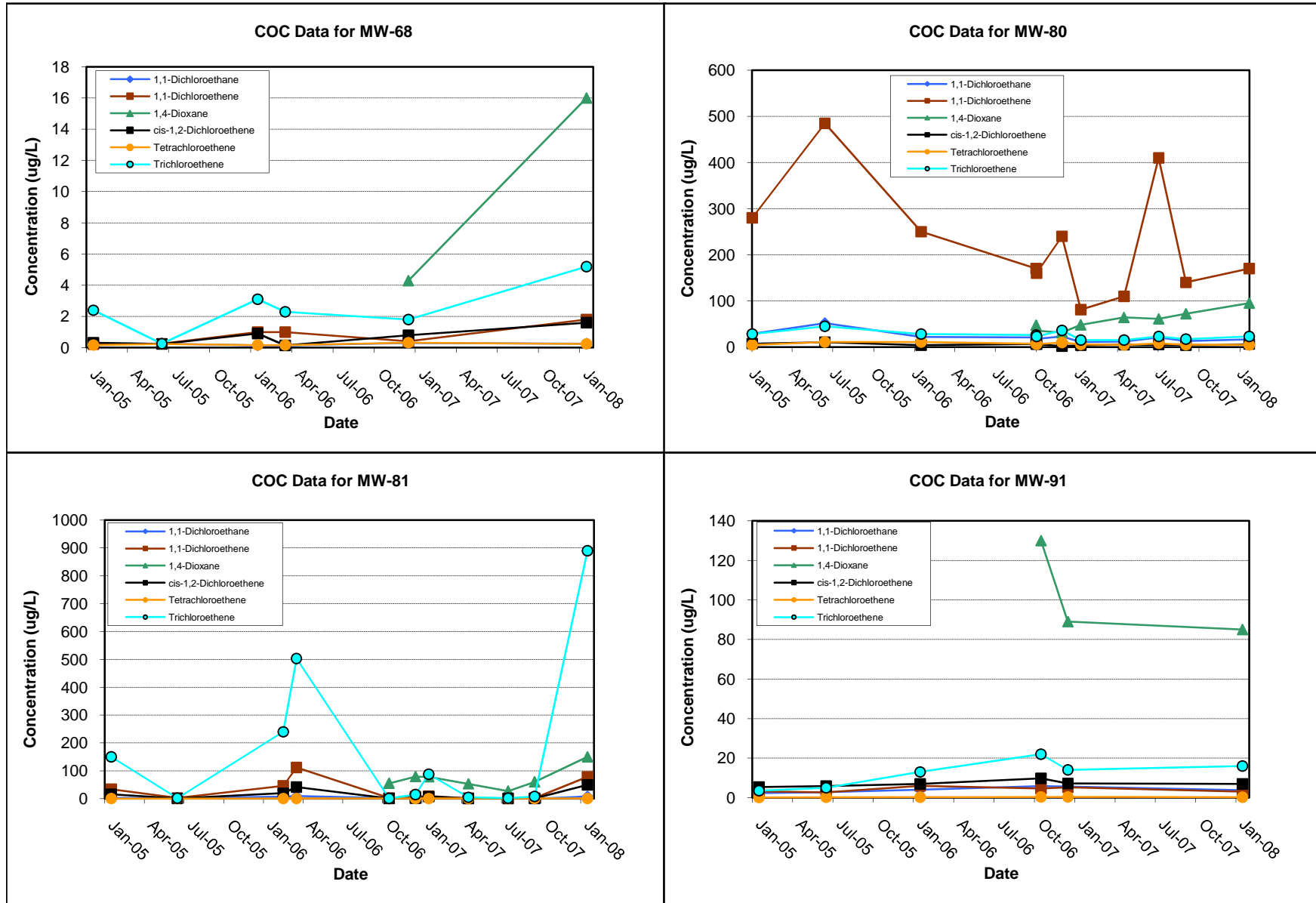
**APPENDIX E
CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
LSAS ZONE**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



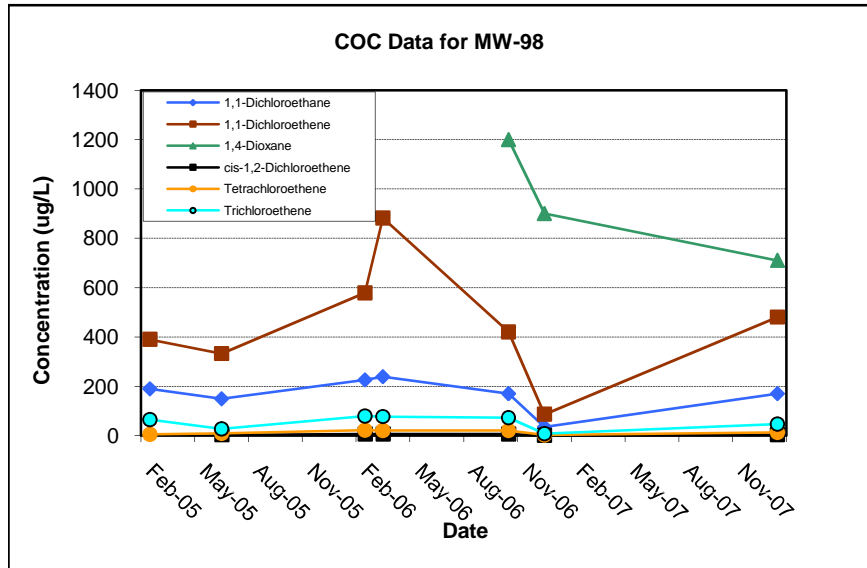
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LSAS ZONE**

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FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



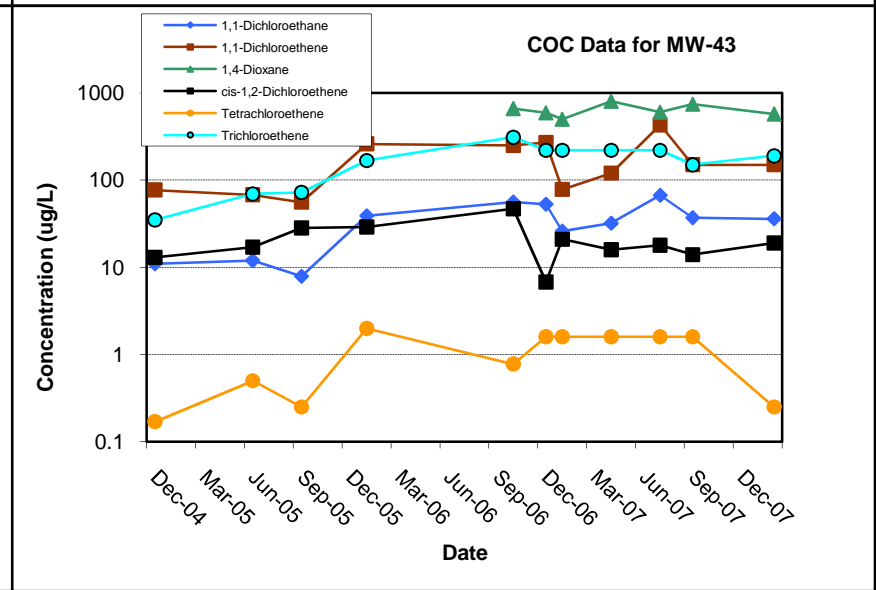
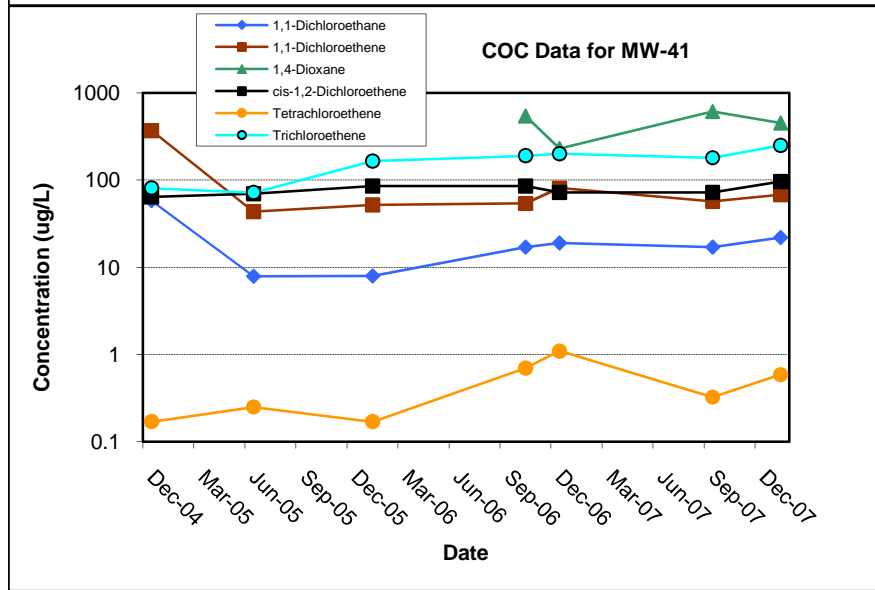
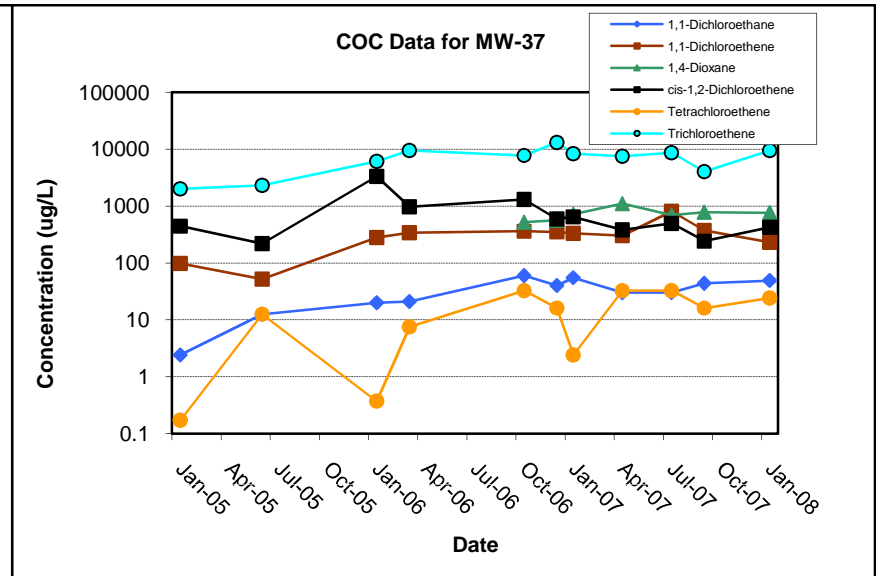
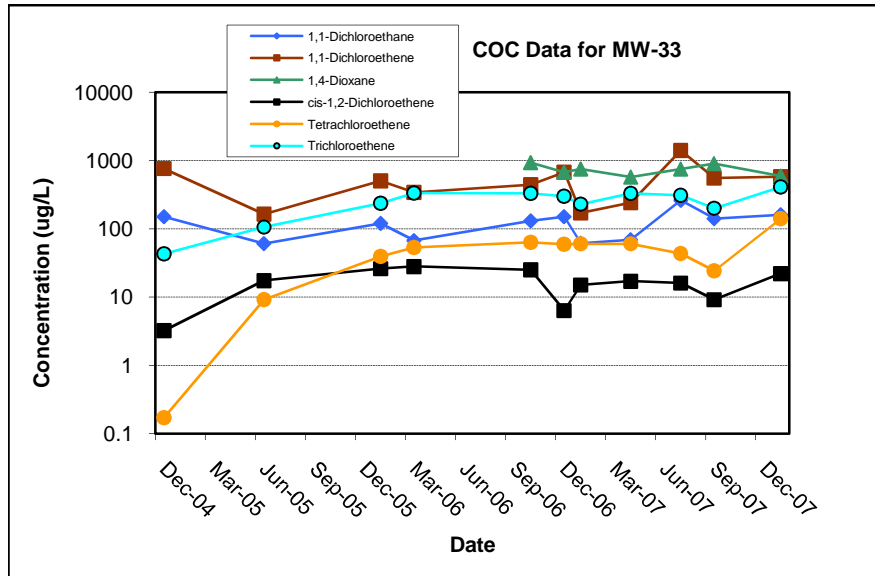
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CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
LSAS ZONE

2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA



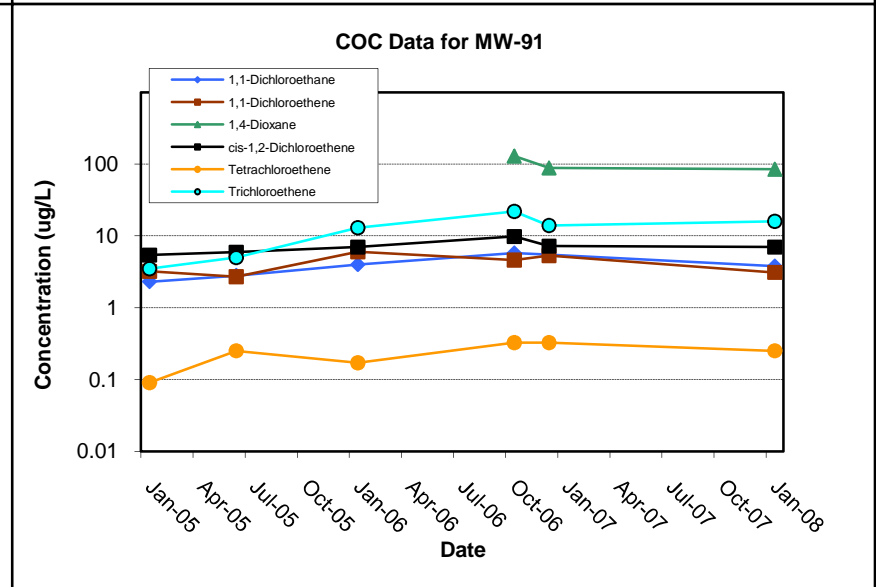
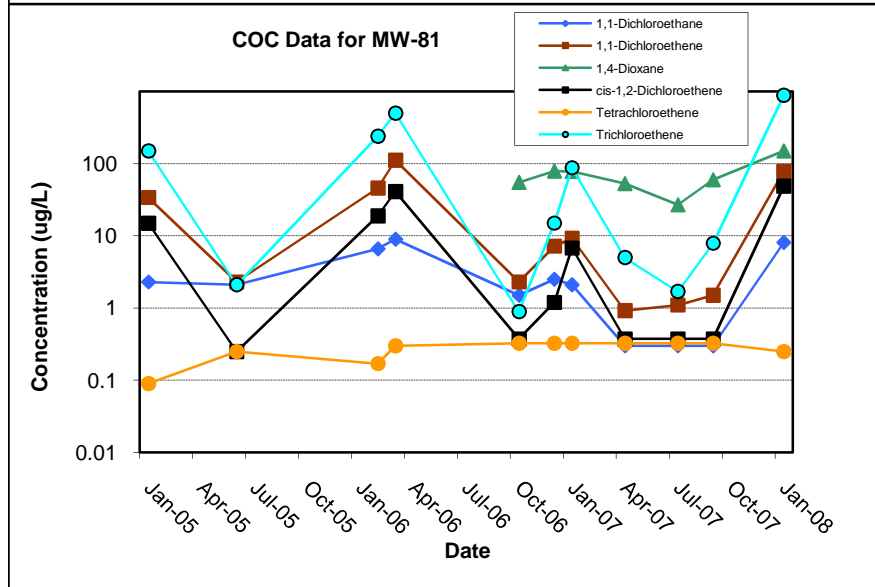
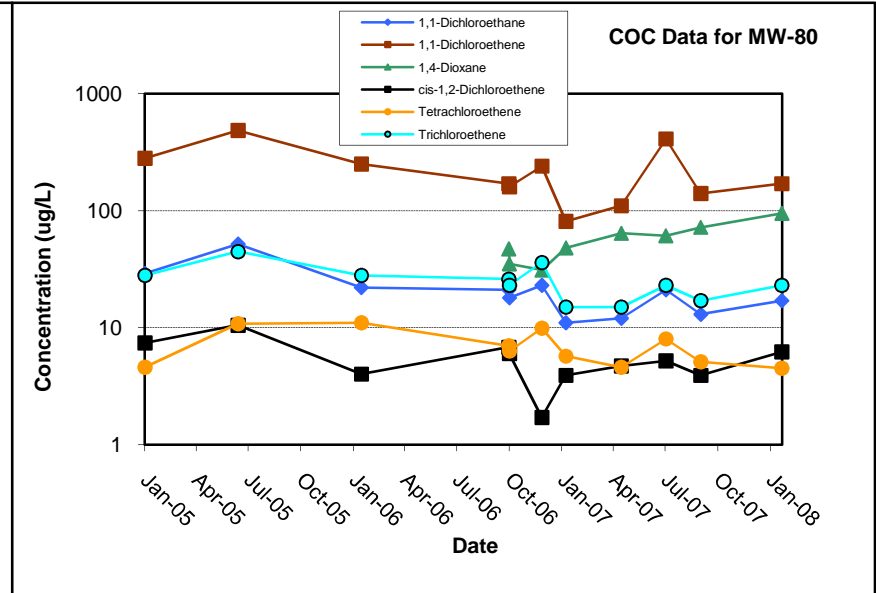
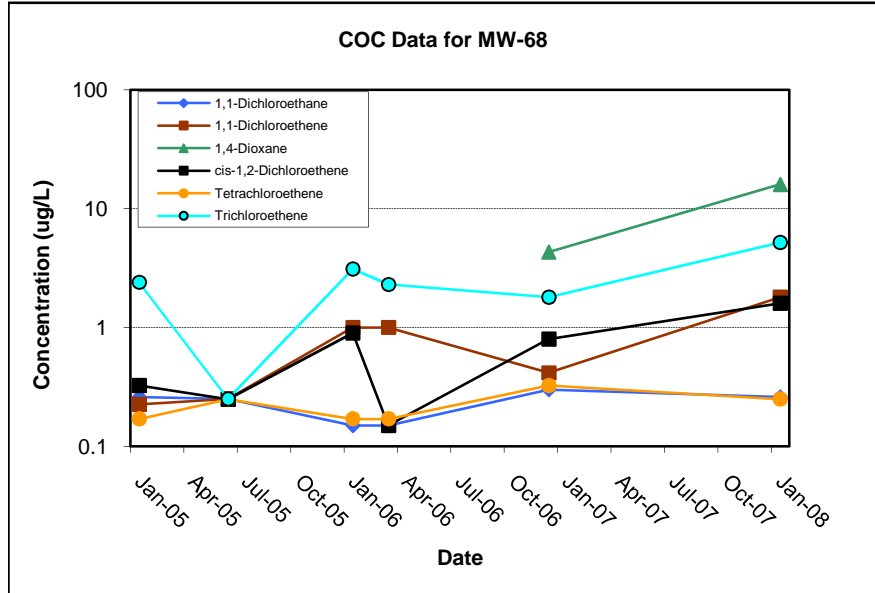
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LOG CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
LSAS ZONE**

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FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



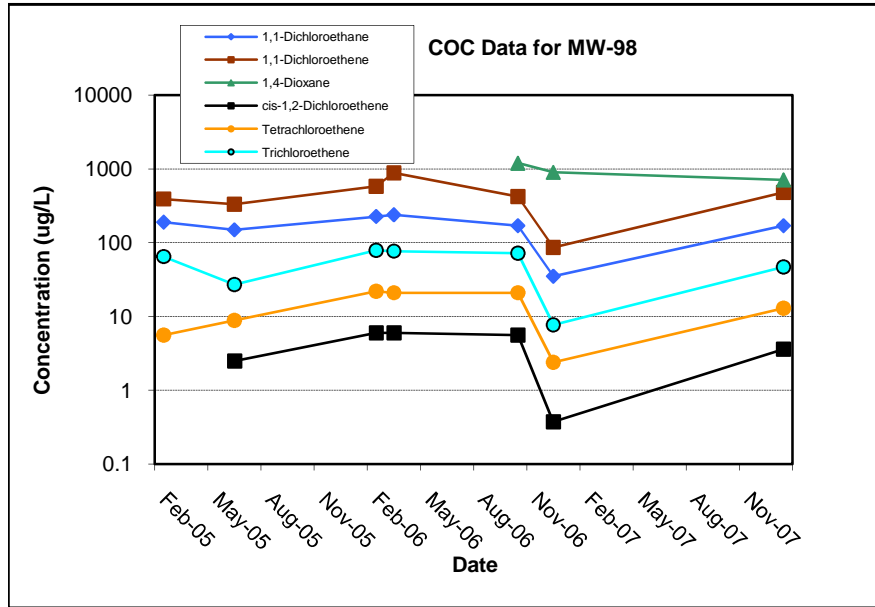
**APPENDIX E
LOG CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
LSAS ZONE**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



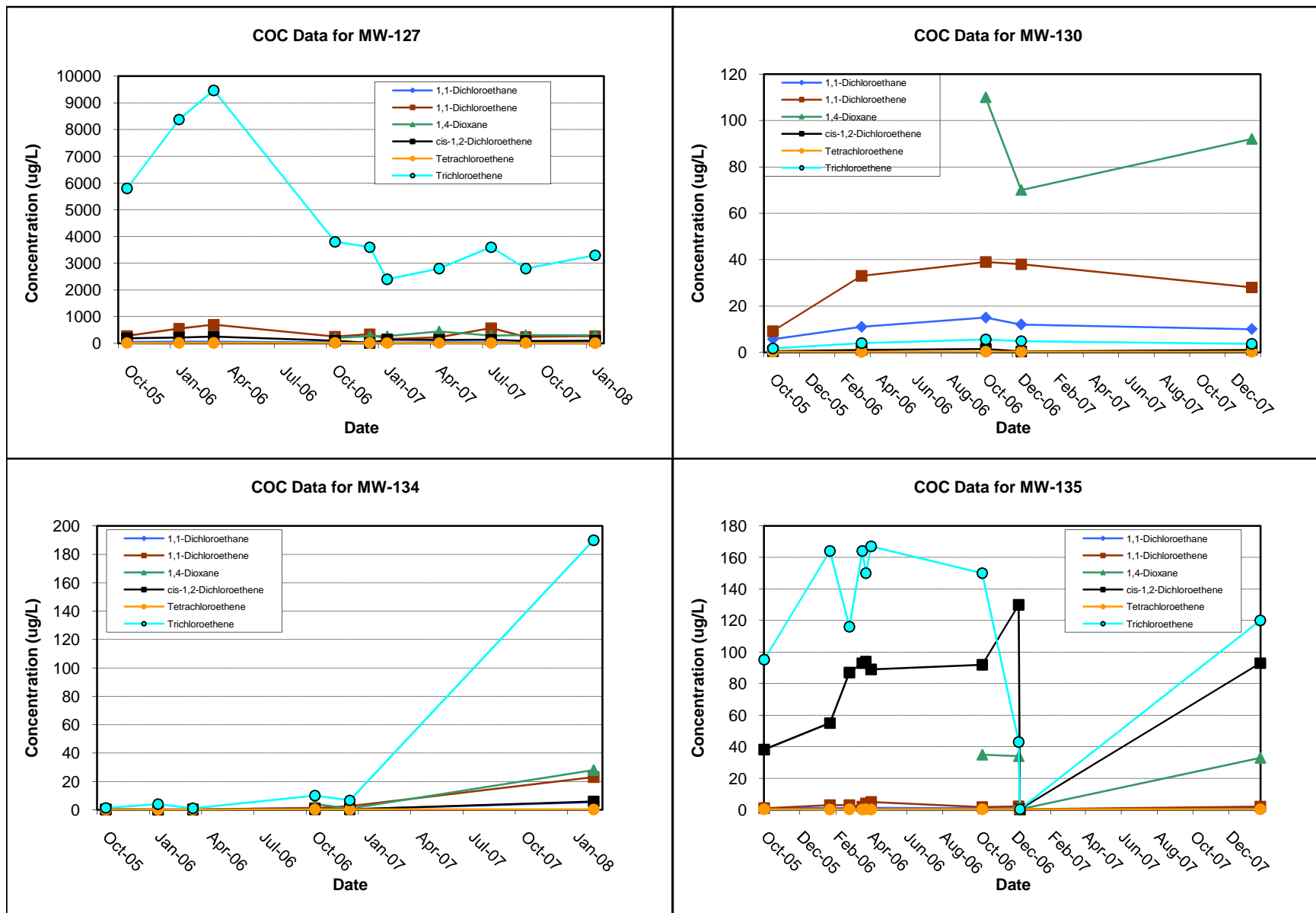
**APPENDIX E
LOG CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
LSAS ZONE**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



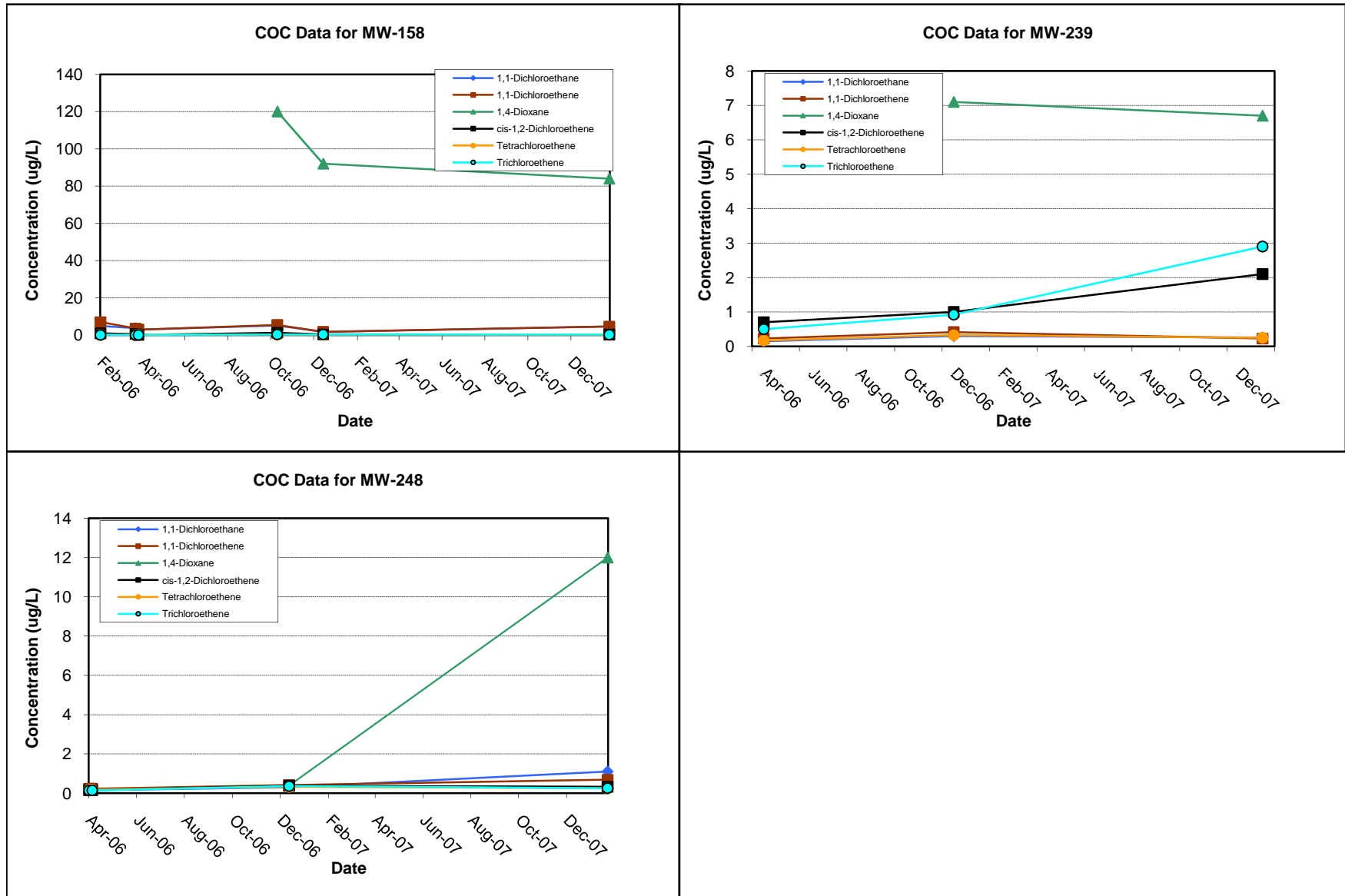
**APPENDIX E
CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
AF GRAVEL ZONE**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



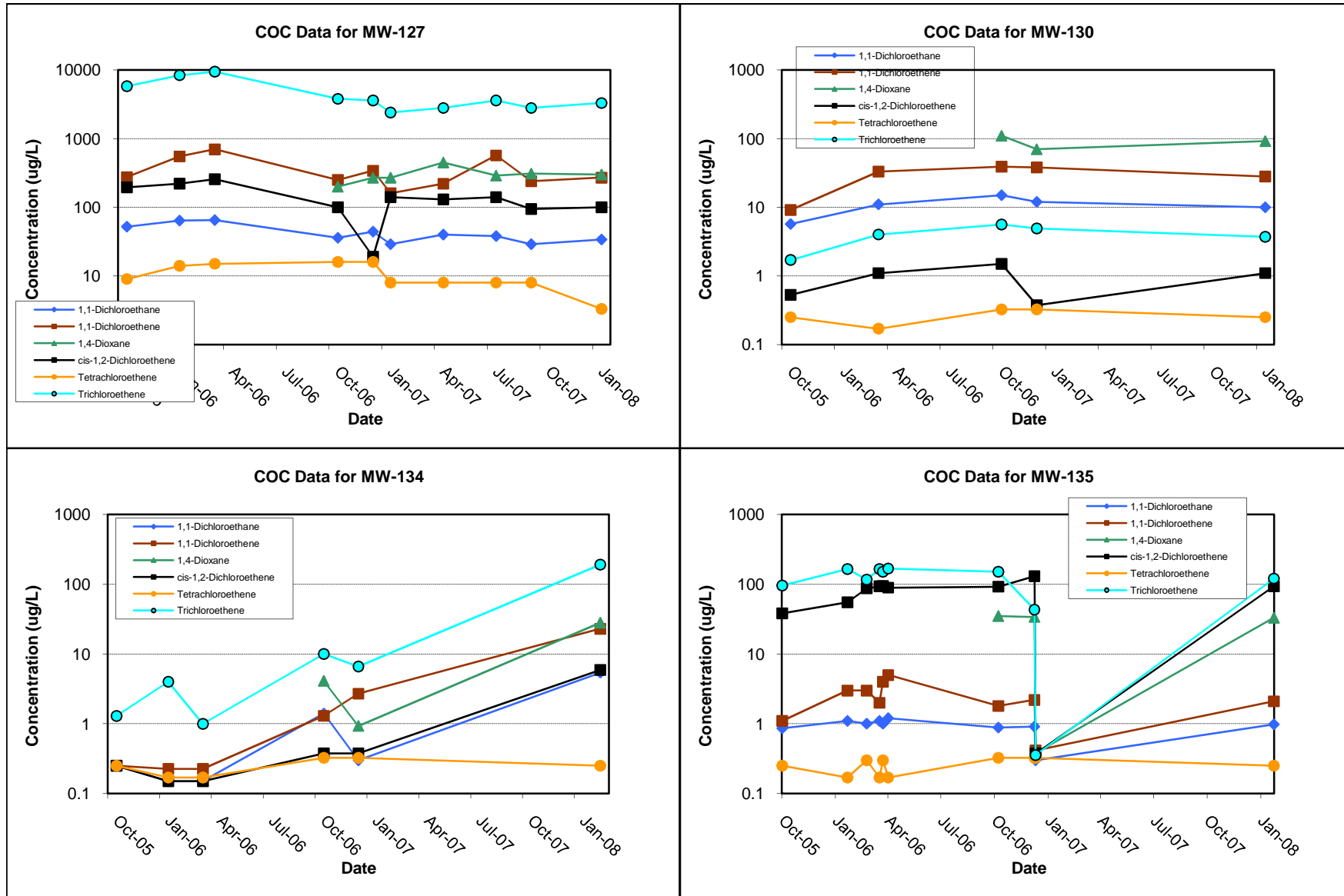
**APPENDIX E
CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
AF GRAVEL ZONE**

**2008 GROUNDWATER MONITORING REPORT
FORMER AMERICAN BERYLLIUM COMPANY SITE - TALLEVAST, FLORIDA**



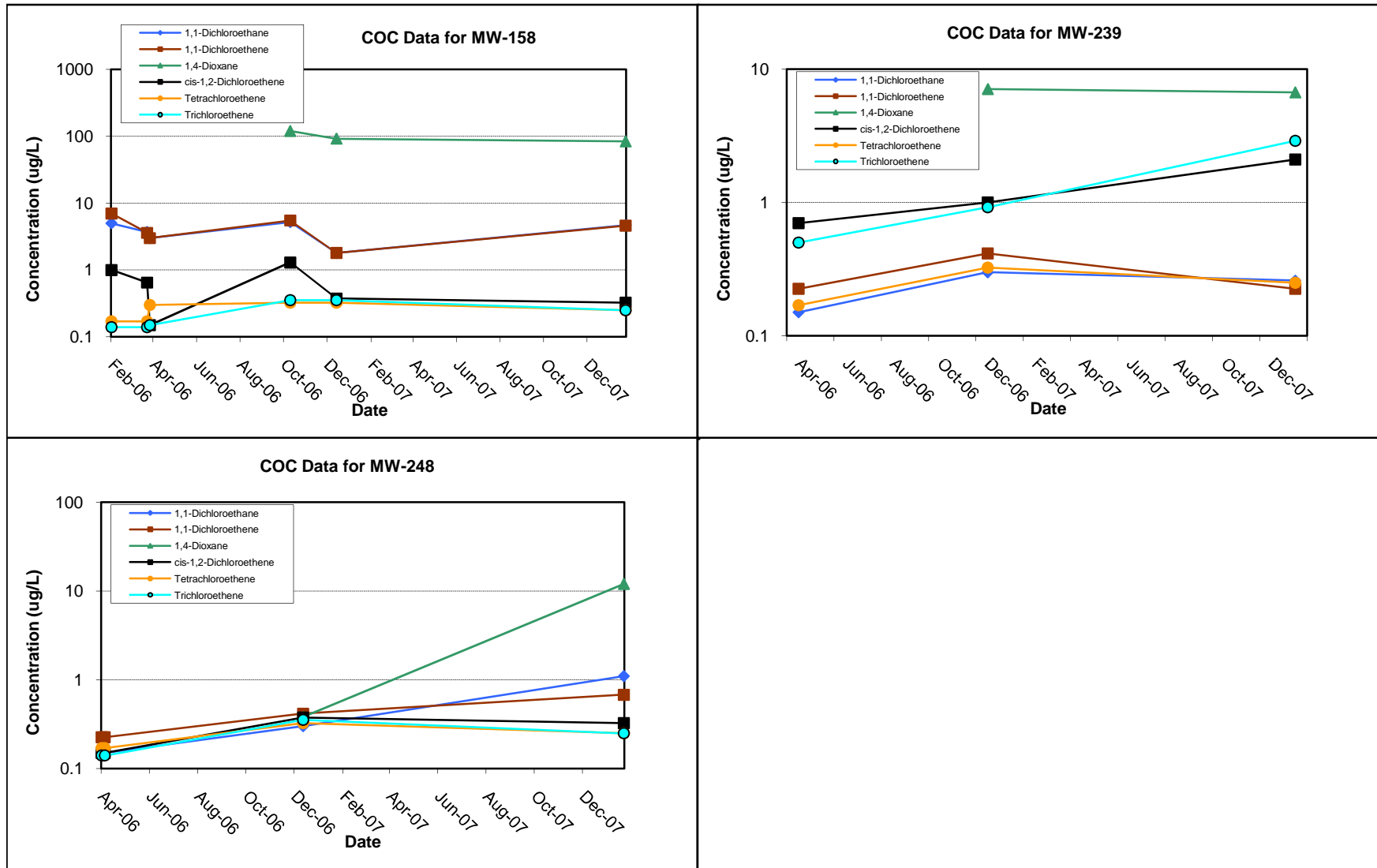
APPENDIX E
LOG CONCENTRATION VS. TIME GRAPHS FOR SELECTED MONITORING WELLS
AF GRAVEL ZONE

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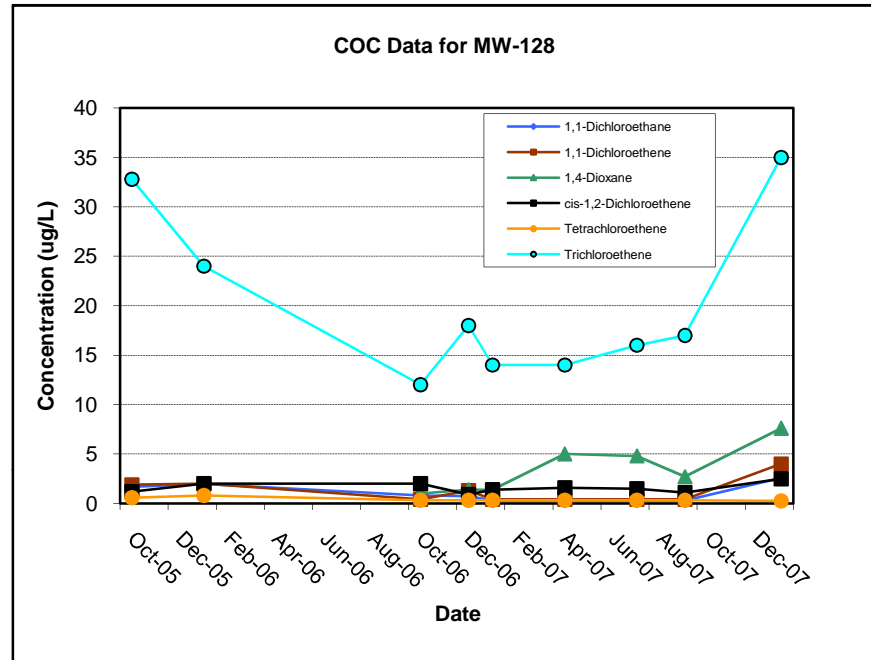
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