

TIDEWATER SEDIMENT DATA REPORT FORMER TIDEWATER FACILITY

Prepared for National Grid

Prepared by Anchor QEA, LLC

June 2009

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TABLE OF CONTENTS

1	INT	RODUCTION	.1
2	FIEI	D ACTIVITIES – SEDIMENT SAMPLING	.2
3	DAT	ГА	.5
	3.1	Physical Analysis	5
	3.2	Chemical Analysis	5
	3.3	Data Validation	6
4	REF	ERENCES	.7

List of Tables

Table 2-1	Tidewater As Builts
Table 3-1	Summary of Sediment Sample Analytical Results – Grain Size
Table 3-2	Summary of Sediment Sample Analytical Results – PAHs
Table 3-3	Summary of Sediment Sample Analytical Results – TPH
Table 3-4	Summary of Sediment Sample Analytical Results – VOCs
Table 3-5	Summary of Sediment Sample Analytical Results – Metals
Table 3-6	Summary of Sediment Sample Analytical Results – TOC, PAC, PSOL

List of Figures

Figure 1	Site Location
Figure 2	Station Locations
Figure 3	Station Locations Grain Size
Figure 4-A	Station Locations Total PAH Distribution
Figure 4-B	Station Locations Benzo(a)pyrene Distribution
Figure 4-C	Station Locations Napthalene Distribution
Figure 4-D	Station Locations Phenanthrene Distribution

List of Appendices

Appendix ACore LogsAppendix BData Validation Reports

LIST OF ACRONYMS AND ABBREVIATIONS

DEM	Department of Environmental Management's
MGP	Manufactured Gas Plant
PACN	Physiologically Available Cyanide
PAHs	Polycyclic Aromatic Hydrocarbons
SdIWP	Sediment Investigation Work Plan
TOC	Total Organic Carbon
TPHs	Total Petroleum Hydrocarbons
VOCs	Volatile Organic Compounds

1 INTRODUCTION

On behalf of The Narragansett Electric Company d/b/a National Grid (National Grid), Anchor QEA, LLC (Anchor QEA) and ARCADIS conducted a sediment sampling investigation in the Seekonk River near the former Tidewater Facility on Tidewater Street in Pawtucket, Rhode Island. The Tidewater Facility is the location of the former Tidewater Manufactured Gas Plant (MGP) and the Pawtucket No. 1 Power Station (Figure 1) and is hereinafter referred to as "the site". This sediment sampling investigation was intended to evaluate potential impacts to Seekonk River environmental conditions that may be associated with the site.

The objectives of this investigation were to collect information in the portions of the Seekonk River near the former Tidewater Facility in compliance with Rhode Island Department of Environmental Management's (DEM) February 2004 Rules and Regulations for the Investigation and Remediation of Hazardous Material Releases (Remediation Regulations) and the Sediment Investigation Work Plan (SdIWP) dated July 3, 2008.

The data collected during this site investigation will be used to preliminarily evaluate Seekonk River sediment for the presence of compounds that were detected in upland subsurface investigations performed previously at the site. This Tidewater Sediment Data Report (report) summarizes field activities and analytical data collected during the sediment sampling investigation. Details of the field and analytical methods are described in the SdIWP (ARCADIS 2008).

2 FIELD ACTIVITIES – SEDIMENT SAMPLING

To evaluate the potential impacts associated with the site and other potential contaminant sources to the Seekonk River sediments in the site vicinity, TG&B Marine Services, Inc. of Falmouth, Massachusetts collected 49 sediment cores from July 8, 2008 to July 16, 2008. A total of 48 sediment samples were collected from the sediment cores (which ranged in length from one to 14 feet) and submitted for physical and chemical analyses as described in Section 3. ARCADIS personnel observed the sampling, processed the sediment cores for sampling, and logged the cores.

Core locations were established prior to entering the field by placing transects approximately 200 feet apart along the shoreline of the site with two to three coring locations for each transect. An additional five locations were situated on the far side of the navigational channel from the site. Two locations were situated upstream of the site and additional sampling locations were situated downstream of the site. Several locations were adjusted in the field to compensate for difficulty in recovering sediment cores. Two additional cores were added to investigate potential MGP impacts. Figure 2 shows the sediment sampling locations. A summary of information associated with core collection and sampling is presented in Table 2-1.

The majority of cores had a five-foot target depth. Selected cores were chosen to be advanced to 20 feet or to refusal to provide a preliminary indication of deeper sediment stratigraphy and subsurface conditions.

The cores were collected by both push core and vibracore methods. Core locations were located from GPS coordinates. The workboat was anchored over the sample location. For the push cores, the acetate liner was cut to length and rinsed with river water. The piston was inserted into the bottom of the liner, and the line attached to the piston was drawn up through the top of the liner. The liner was then filled with river water. The top of the liner was attached to a pushrod. The assembly was lowered until the bottom of the acetate liner was approximately one inch above the sediment-water interface, and the piston line was tied off to keep the piston at the sediment-water interface. Pressure was exerted onto the pushrod to push the liner into the sediment. If the liner could not be pushed the full five feet into the sediment, a slide hammer was used to attempt full penetration. The liner was then pulled from the sediment. The bottom of the core was immediately capped and taped. The push rod was then detached. Standing water was drained from the top of the core, and the top was capped. The water depth, as-built coordinates, time and penetration depth were recorded. The cores were labeled and stored in an upright position.

Vibratory cores were collected by suspending a vibrating head from an A-frame from a work boat. The vibratory head has a 3-inch-diameter acetate-lined aluminum core barrel suspended below it. The head/core barrel apparatus was lowered on a cable until the core barrel bottom rim contacted the sediment. The core barrel penetrated the sediment by vibrating the core barrel and concurrently pushing the core barrel into the sediment. The process of sediment coring typically only minimally disturbs sediment structure, so sediment stratigraphy and other details were preserved and observed in the sediment cores.

After collection, the vibracores and push cores were brought to shore to be processed and stored in an upright position. Each core was cut lengthwise with power snips and split open on a table. The core was photographed and visually logged. Core logs are provided in Appendix A.

Sediment samples were collected from predetermined locations along the length of the core. Additional samples were collected from cores based on visual observations. Due to the number of analyses and the size of the core barrel, additional cores were required from each core location in order to collect enough material. The material from the same interval in each core was placed in a stainless steel bowl and homogenized. The homogenized sample was placed in laboratory-supplied glassware and labeled. The samples were then stored on ice until delivery to the laboratory under chain of custody.

For this preliminary sediment investigation, samples from the top six inches of the cores were collected for laboratory analyses. Samples from deeper intervals were collected and archived for later analysis, based on the results from the surface interval. Samples were collected from 6 to 12 inches and 12 to 18 inches for archival purposes. If a change in sediment stratigraphy was observed, a six-inch sample was collected at a deeper depth.

Excess sediment, decontamination water and additional waste material were stored in labeled drums and removed from the site by Clean Harbors, Inc. for off-site disposal at a permitted facility.

3 DATA

3.1 Physical Analysis

Forty-eight sediment samples were submitted to Geotesting Express of Boxborough, MA for analysis of grain size. Results are provided in Table 3-1. Grain-size distribution is depicted for each station on Figure 3. Grain size analyses were consistent with field observations in that the sediment varied from granular sand and gravel to lesser amounts of cohesive silt with sand. The percentage of sediment consisting of fine-grained particles is important because, in general, the fine-grained fraction combined with the total organic carbon (TOC) content of the sediment controls the extent that most sediment contaminants are adsorbed to the sediment matrix.

3.2 Chemical Analysis

Forty-eight sediment samples were submitted to Alpha Analytical of Mansfield, MA for analysis of polycyclic aromatic hydrocarbons (PAHs), total petroleum hydrocarbons (TPHs), 13 priority pollutant metals, volatile organic compounds (VOCs), physiologically available cyanide (PACN), and TOC. A subset of samples were also analyzed for an extended list of alkylated PAH compounds. Results are provided in Tables 3-2 through 3-6.

Distribution of Total PAHs and selected individual PAH compounds are depicted on Figures 4-A through 4-D. Total PAHs were detected in all samples and ranged from a minimum concentration of 6,100 μ g/kg at Station SC40 to a maximum concentration of 16,000,000 μ g/kg at Station SW2. The mean concentration was 438,744 μ g/kg.

A subset of samples was also analyzed for alkylated PAHs. The distribution of alkylated PAHs can aid in the determination of the type and source of the petroleum compounds present in a sample.

Out of a total of 65 VOCs that were analyzed, only 15 were detected, with most detected only in one sample (sample SW-2).

All 13 metals were detected in at least one sample, although many results were "J" qualified, indicating the result is an estimate. There were no detected concentrations of PACN.

In general, the field investigation findings indicate relatively elevated PAH concentrations are concentrated in localized areas. When compared in our experience to other New England properties that were formerly occupied by MGPs, the concentration and extent of organic compounds (PAHs and VOCs) in sediment was generally lower in magnitude (i.e., for PAH and VOC concentration) and occupied a smaller area in the river.

3.3 Data Validation

Data were validated according to the United States Environmental Protection Agency's (USEPA's) National Functional Guidelines of October 1999 (USEPA, 1999). The quality indicators of this limited data review included holding times, associated blanks, matrix spike/matrix spike duplicate analysis, field duplicates, laboratory control sample and surrogate recoveries. The data collected during this sediment investigation met overall system performance, and the overall data quality is within the guidelines specified in the method. Data validation reports are provided in Appendix B.

Data

4 REFERENCES

ARCADIS. 2008. *Sediment Investigation Work Plan, Former Tidewater Facility*. Prepared for National Grid (July 3, 2008).

United States Environmental Protection Agency. 1999. United States Environmental Protection Agency Contract Laboratory Program National Functional Guidelines for Organic Data Review. EPA540/R-99/008. October 1999.

TABLES

Table 2-1 Tidewater As Builts

Tidewater Data Report

Ctation	Data	Time	La	titude ¹	Lor	ngitude	DOW	Dan	Dee	Mudline	Sampled ^{3,4}			Additional Nation		
Station	Date		DD	MM.MMMM	DD	MM.MMMM	DOw	Pen	Rec	Elev. (ft)		Sa	mpled			Additional Notes
SC1	7/15/2008	1000	41 52	0.2189	71 22	0.9352	11.0	7.0	4.6	-9.35	S1	S2	S3	S6		MS/MSD, Alk PAH
SC2	7/15/2008	1130	41 52	0.1603	71 22	0.8832	12.6	8.0	9 ²	-11.69	S1	S2	S3	S12		Field Dup3
SC3	7/9/2008	830	41 52	0.1505	71 22	0.8560	11.5	6.0	3.1	-10.30	S1	S2	S3	S6		
SC4	7/9/2008	1030	41 52	0.1195	71 22	0.8408	9.6	5.0	3.8	-7.27	S 1	S2	S3			
SC5	7/9/2008	1350	41 52	0.0888	71 22	0.8233	12.0	5.0	2.9	-6.89	S1	S2	S3	S5		
SC6	7/9/2008	1430	41 52	0.0550	71 22	0.8228	7.4	6.0	6.1	-2.30	S1	S2	S3	S11		Alk PAH
SC7	7/14/2008	1130	41 51	0.8819	71 22	0.7842	1.0	4.0	3.7	0.04	S 1	S2	S3	S5		
SC8	7/14/2008	1100	41 51	0.8815	71 22	0.7693	3.2	5.0	3.8	-2.08	S 1	S2	S3	S4		Alk PAH
SC9	7/10/2008	950	41 52	0.0123	71 22	0.7997	5.5	5.0	4.8	-4.17	S1	S2	S3	S8		Field Dup1, Alk PAH
SC10	7/8/2008	1350	41 52	0.1861	71 22	0.9142	17.5	5.0	2.8	-12.53	S1	S2	S3			
SC11	7/15/2008	1310	41 52	0.1871	71 22	0.8996	9.4	10.0	5.4	-8.46	S1	S2	S3	S4		Alk PAH
SC12	7/8/2008	1430	41 52	0.1598	71 22	0.8774	14.2	4.5	2.4	-9.81	S 1	S2	S3			
SC13	7/15/2008	1420	41 52	0.1201	71 22	0.8328	12.0	10.0	5.4	-10.40	S1	S2	S3	S10		
SC13	7/15/2008	NA	41 52	0.1201	71 22	0.8328	12.0	16.0	6.4	-12.00						not sampled
SC13A	7/16/2008	1120	41 52	0.1187	71 22	0.8320	12.5	16.5	11.5	-11.28						not sampled
SC14	7/15/2008	1500	41 52	0.0875	71 22	0.8321	5.5	20.0	13.3	-3.43	S1	S2	S3	S7	S32	
SC15	7/16/2008	840	41 52	0.0523	71 22	0.8113	8.2	19.0	14.0	-4.52	S1	S2	S3	S4		
SC16	7/16/2008	1140	41 52	0.0205	71 22	0.8079	3.7	17.5	12.9	-2.69	S1	S2	S3	S11	S20	
SC17	7/10/2008	1050	41 52	0.0048	71 22	0.7893	6.9	5.0	5.0	-5.02	S1	S2	S3	S9		
SC18	7/14/2008	1030	41 51	0.8798	71 22	0.7474	10.0	5.5	5.5	-8.70	S1	S2	S3	S6		
SC19	7/10/2008	850	41 51	0.9955	71 22	0.7437	11.8	5.0	4.1	-10.83	S1	S2	S3	S8		Alk PAH
SC20	7/9/2008	1515	41 52	0.0541	71 22	0.7674	17.9	4.0	2.3	-13.22	S1	S2	S3			
SC21	7/8/2008	1150	41 52	0.2067	71 22	0.8799	5.5	5.5	5.4	-1.11	S1	S2	S3	S5		
SC22	7/9/2008	1140	41 52	0.2843	71 22	0.9971	17.3	2.0	4.3*	-13.83	S1	S2				S2 collected for GS
SC23	7/8/2008	1015	41 52	0.2578	71 22	0.9509	8.4	5.3	4.7	-5.88	S1	S2	S3	S9		
SC24	7/8/2008	1100	41 52	0.2636	71 22	0.9546	10.7	5.0	2.7	-7.23	S1	S2	S3			MS/MSD
SC25	7/8/2008	1500	41 52	0.1618	71 22	0.8656	12.8	5.0	1.7	-8.90	S1	S2	S3			
SC26	7/10/2008	815	41 52	0.1109	71 22	0.7987	13.8	6.0	5.6	-12.86	S1	S2	S3	S9		
SC27	7/14/2008	820	41 51	0.9391	71 22	0.7185	6.6	3.0	1.35	-3.71	S1	S2				
SC28	7/10/2008	1320	41 51	0.9796	71 22	0.7904	7.8	5.0	5.0	-3.73	S1	S2	S3	S9		
SC29	7/10/2008	1120	41 51	0.9802	71 22	0.7999	3.1	5.5	5.3	-0.88	S1	S2	S3			
SC30	7/10/2008	1350	41 51	0.9800	71 22	0.8107	3.4	3.0	2.2	1.20	S1	S2	S3	S4		
SC31	7/14/2008	1340	41 51	0.9513	71 22	0.8134	1.0	2.5	2.75	1.09	S1	S2	S3	S4		Alk PAH
SC32	7/11/2008	1100	41 51	0.9469	71 22	0.7992	1.9	4.5	4.0	-0.21	S1	S2	S3			
SC33	7/11/2008	1130	41 51	0.9479	71 22	0.7784	6.1	6.0	5.1	-4.12	S1	S2	S3	S10		Alk PAH
SC34	7/11/2008	1420	41 51	0.9305	71 22	0.7661	10.8	5.0	4.8	-6.74	S1	S2	S3	S5		
SC35	7/11/2008	1440	41 51	0.9303	71 22	0.7890	4.5	2.0	1.1	-0.24	S1	S2				
SC36	7/14/2008	1300	41 51	0.9327	71 22	0.7994	1.0	1.5	1.2	0.66	S1	S2				
SC37	7/14/2008	1320	41 51	0.9209	71 22	0.7824	5.1	4.5	3.25	-3.19	S1	S2	S3	S5		MS/MSD
SC38	7/11/2008	1340	41 51	0.9134	71 22	0.7776	5.6	5.5	5.4	-1.99	S1	S2	S3			
SC39	7/11/2008	1310	41 51	0.9129	71 22	0.7669	8.4	5.5	5.2	-5.25	S1	S2	S3			
SC40	7/9/2008	1000	41 52	0.1430	71 22	0.8334	13.7	5.0	4.8	-11.82	S1	S2	S3			

Table 2-1 **Tidewater As Builts**

Tidewater Data Report

Former Tidewater MGP Site, Pawtucket, Rhode Island

Station	Date	Time	La	titude ¹	Loi	ngitude	DOW	Pen	Rec	Mudline		60	Sampled ^{3,4}				
Station	Date		DD	MM.MMMM	DD	MM.MMMM	DOW	Fell	Neu	Elev. (ft)		5 a	mpled	L L		Additional Notes	
SC41	7/9/2008	920	41 52	0.0865	71 22	0.8016	11.2	6.0	5.8	-9.85	S1	S2	S3	S7			
SC42	7/10/2008	1020	41 52	0.0218	71 22	0.7782	8.5	5.0	4.7	-6.95	S 1	S2	S3				
SC43	7/10/2008	920	41 51	0.9817	71 22	0.7651	9.1	5.5	5.4	-7.94	S1	S2	S3				
SC44	7/14/2008	900	41 51	0.9151	71 22	0.7470	12.4	5.0	4.85	-10.10	S1	S2	S3	S8		Field Dup2	
SC45	7/14/2008	930	41 51	0.8819	71 22	0.7262	12.2	4.5	3.5	-10.25	S1	S2	S3	S5			
SC46	7/14/2008	1000	41 51	0.8555	71 22	0.7232	12.0	5.5	5.3	-10.43	S 1	S2	S3	S10			
SC47	7/11/2008	1030	41 52	0.1253	71 22	0.8449	10.5	5.0	3.7	-9.06	S1	S2	S3	S7			
SW-2	7/11/2008	848	41 52	0.1252	71 22	0.8511	NA	1.5	1.0	NA	S1	S2					

Notes:

NA = Not available

DOW = Depth of Water (ft)

Pen = Core Penetration (ft)

Rec = Core Recovery (ft)

¹ Datum WGS-84

² Additional material in core barrel due to suction from check valve
 ³ Samples in bold = analyzed; additional samples collected for archive purposes

⁴ Sampled Depths (ft):

	• • • •		
S1	0.0-0.5	S8	3.5-4.0
S2	0.5-1.0	S9	4.0-4.5
S3	1.0-1.5	S10	4.5-5.0
S4	1.5-2.0	S11	5.0-5.5
S5	2.0-2.5	S12	5.5-6.0
S6	2.5-3.0	S20	9.5-10.0
S7	3.0-3.5	S32	15.5-16.0

Table 3-1 Summary of Sediment Sample Analytical Results - Grain Size

Tidewater Data Report Former Tidewater MGP Site, Pawtucket, Rhode Island

				Geotech	
Location ID:	Date Collected:	Sample Name:	% Gravel	% Sand	% Silt/Clay
SC1	07/15/08	SC1-S1	3.3	92.9	3.8
SC2	07/15/08	SC2-S1	17.1	81.6	1.3
SC3	07/09/08	SC3-S1	8.5	89.4	2.1
SC4	07/09/08	SC4-S1	22.9	74.8	2.3
SC5	07/09/08	SC5-S1	4.5	94.5	1
SC6	07/09/08	SC6-S1	0	28.4	71.6
SC7	07/14/08	SC7-S1	3.7	63.9	32.4
SC7	07/14/08	SC7-S7	5.2	58.4	36.4
SC8	07/14/08	SC8-S1	0.8	30.8	68.4
SC9	07/10/08	SC9-S1	0.3	46.8	52.9
SC10	07/08/08	SC10-S1	5.7	89.8	4.5
SC11	07/15/08	SC11-S1	9	78.5	12.5
SC12	07/08/08	SC12-S2	16	83.2	0.8
SC13	07/15/08	SC13-S1	8.2	89.3	2.5
SC14	07/15/08	SC14-S1	0.4	41.9	57.7
SC15	07/16/08	SC15-S1	0.5	57.8	41.7
SC16	07/16/08	SC16-S1	0.2	41.1	58.7
SC17	07/10/08	SC17-S1	0.2	80.5	19.3
SC18	07/14/08	SC18-S1	0.2	21.9	77.9
SC19	07/10/08	SC19-S1	0.6	36.3	63.1
SC20	07/09/08	SC20-S1	0.5	98.8	0.7
SC21	07/08/08	SC21-S1	0.0	47.4	52.6
SC22	07/09/08	SC22-S2	1.2	97.8	1
SC23	07/08/08	SC23-S1	0	76.5	23.5
SC24	07/08/08	SC24-S1	5.8	92.1	2.1
SC25	07/08/08	SC25-S1	2.4	91	6.6
SC25	07/10/08	SC26-S1	0.2	97.5	2.3
SC20	07/14/08	SC27-S1	1.9	37.6	60.5
SC28	07/10/08	SC28-S1	0.7	95.2	4.1
SC28	07/10/08	SC29-S1	0.7	95.2 26.1	73.9
SC30	07/10/08	SC30-S1	0	42.3	57.7
			4.2		
SC31	07/14/08	SC31-S1		77.8	18
SC32 SC33	07/11/08 07/11/08	SC32-S1 SC33-S1	3.8 0	57.3 40.6	38.9 59.4
SC33 SC34		SC33-S1 SC34-S1			20.9
SC34 SC35	07/11/08		0.1	79	
	07/11/08	SC35-S1	0.3	46.3	53.4
SC36	07/14/08	SC36-S1	6.8	62.4	30.8
SC37	07/14/08	SC37-S1	0.9	46.9	52.2
SC38	07/11/08	SC38-S1	3.1	91.2	5.7
SC39	07/11/08	SC39-S1	0	15.8	84.2
SC40	07/09/08	SC40-S1	0.2	98.7	1.1
SC41	07/09/08	SC41-S1	0.1	97.5	2.4
SC42	07/10/08	SC42-S1	2.9	94.2	2.9
SC43	07/10/08	SC43-S1	1.7	86.4	11.9
SC44	07/14/08	SC44-S1	0.1	95.4	4.5
SC45	07/14/08	SC45-S1	0	97.5	2.5
SC46	07/14/08	SC46-S1	0.1	98.1	1.8
SC47	07/11/08	SC47-S1	25.6	72.4	2

Tidewater Data Report

Location ID:		SC1	SC2	SC3	SC4	SC5	SC6	SC7	SC8	SC9	SC10	SC11	SC12
Date Collected:	Units	07/15/08	07/15/08	07/09/08	07/09/08	07/09/08	07/09/08	07/14/08	07/14/08	07/10/08	07/08/08	07/15/08	07/08/08
PAHs													
Acenaphthene	µg/kg	150 J	300 J [50 J]	4,200	570	87	240	140	1,400	290 [310]	470	400	24
Acenaphthylene	µg/kg	350	320 [180]	6,100	920	250	960	610	2,000	1,000 [940]	220	480	170
Anthracene	µg/kg	1,000 J	3,200 J [280 J]	52,000	1,900	340	1,100	580	3,500	1,300 [1,400]	1,500	1,800	240
Benzo(a)anthracene	µg/kg	2,800	3,400 J [930 J]	55,000	3,800	730	4,000	1,700	8,300	4,200 [4,100]	3,900	3,900	700
Benzo(a)pyrene	µg/kg	2,000	2,600 J [800 J]	43,000	4,500	830	5,500	1,900	6,900	4,900 [4,800]	4,400	2,900	620
Benzo(b)fluoranthene	µg/kg	1,900	2,100 [790]	30,000	2,800	540	5,700	2,100	6,200	4,700 [4,500]	3,500	2,600	490
Benzo(g,h,i)perylene	µg/kg	1,200 J	1,300 [440]	18,000	1,900	610	4,200	1,400	4,500	3,500 [3,400]	2,300	1,700	310
Benzo(k)fluoranthene	µg/kg	2,200	2,100 J [690 J]	32,000	2,800	550	4,900	2,000	6,700	4,300 [4,300]	3,400	2,900	490
Chrysene	µg/kg	2,900	3,300 J [1,000 J]	45,000	3,600	720	5,400	2,000	8,800	4,900 [4,800]	3,800	4,000	630
Dibenzo(a,h)anthracene	µg/kg	360	410 J [130 J]	5,900	540	120	1,000	370	1,300	880 [860]	730	510	88
Fluoranthene	µg/kg	6,500	7,900 J [2,100 J]	140,000	9,200	1,300	11,000	3,300	18,000	9,500 [9,600]	7,100	8,700	1,400
Fluorene	µg/kg	260 J	520 J [69 J]	12,000	510	130	310	180	650	370 [380]	550	450	35
Indeno(1,2,3-cd)pyrene	µg/kg	1,400 J	1,400 J [480 J]	23,000	2,100	460	4,400	1,400	4,600	3,600 [3,600]	2,600	1,800	360
Naphthalene	µg/kg	300	480 J [94 J]	2,500	650	230	640	340	2,300	540 [620]	440	500	58
Phenanthrene	µg/kg	3,100	6,200 J [1,100 J]	120,000	5,400	710	3,200	1,300	6,300	3,500 [3,600]	4,400	4,800	480
Pyrene	µg/kg	5,700	6,600 [2,200]	110,000	8,300	1,400	9,000	3,600	18,000	8,500 [8,200]	6,100	7,400	1,200
Total PAHs	µg/kg	32,000	42,000 J [11,000 J]	700,000	50,000	9,000	61,000	23,000	99,000	56,000 [55,000]	45,000	45,000	7,200
Additional PAHs/Alkylated PAHs													
Perylene	µg/kg	540	NA	NA	NA	NA	1,400	NA	1,800	1,600 [1,500]	NA	740	NA
Benzo[e]pyrene	µg/kg	1,500	NA	NA	NA	NA	4,300	NA	5,100	3,700 [3,600]	NA	2,100	NA
C1-Chrysenes	µg/kg	1,200	NA	NA	NA	NA	2,600	NA	5,900	2,600 [2,400]	NA	1,900	NA
C1-Fluoranthenes/Pyrenes	µg/kg	2,700	NA	NA	NA	NA	4,300	NA	11,000	4,400 [4,200]	NA	3,600	NA
C1-Fluorenes	µg/kg	160	NA	NA	NA	NA	210	NA	980	240 [250]	NA	270	NA
C1-Naphthalenes	µg/kg	91	NA	NA	NA	NA	380	NA	1,100	370 [380]	NA	280	NA
C1-Phenanthrenes/Anthracenes	µg/kg	1,700	NA	NA	NA	NA	1,800	NA	7,000	2,100 [2,000]	NA	2,600	NA
C2-Chrysenes	µg/kg	520	NA	NA	NA	NA	1,600	NA	3,400	1,500 [1,200]	NA	1,000	NA
C2-Fluorenes	µg/kg	150	NA	NA	NA	NA	300	NA	1,500	360 [350]	NA	400	NA
C2-Naphthalenes	µg/kg	130	NA	NA	NA	NA	390	NA	1,500	450 [430]	NA	470	NA
C2-Phenanthrenes/Anthracenes	µg/kg	920	NA	NA	NA	NA	1,300	NA	5,900	1,400 [1,400]	NA	1,800	NA
C3-Chrysenes	µg/kg	380	NA	NA	NA	NA	1,300	NA	2,600	1,300 [1,100]	NA	810	NA
C3-Fluorenes	µg/kg	250	NA	NA	NA	NA	770	NA	1,800	780 [920]	NA	740	NA
C3-Naphthalenes	µg/kg	160	NA	NA	NA	NA	290	NA	2,300	360 [330]	NA	610	NA
C3-Phenanthrenes/Anthracenes	µg/kg	380	NA	NA	NA	NA	760	NA	3,400	890 [800]	NA	1,400	NA
C4-Chrysenes	µg/kg	150	NA	NA	NA	NA	790	NA	1,200	820 [720]	NA	410	NA
C4-Naphthalenes	µg/kg	95	NA	NA	NA	NA	220	NA	1,600	270 [260]	NA	560	NA
C4-Phenanthrenes/Anthracenes	µg/kg	110	NA	NA	NA	NA	390	NA	1,700	550 [470]	NA	1,100	NA

Table 3-2

Summary of Sediment Sample Analytical Results – PAHs

Tidewater Data Report

Location ID:		SC13	SC14	SC15	SC16	SC17	SC18	SC19	SC20	SC21	SC22	SC23	SC24
Date Collected:	Units	07/15/08	07/15/08	07/16/08	07/16/08	07/10/08	07/14/08	07/10/08	07/09/08	07/08/08	07/09/08	07/08/08	07/08/08
PAHs													
Acenaphthene	µg/kg	800	330	1,500	490	60	260	910	750	260	38	200	450
Acenaphthylene	µg/kg	370	1,300	1,100	1,100	250	780	1,200	190	620	200	270	280 J
Anthracene	µg/kg	2,100	1,300	3,900	1,700	310	910	2,800	2,200	1,000	330	640	2,100
Benzo(a)anthracene	µg/kg	5,800	4,700	7,900	4,700	1,100	2,900	4,300	3,100	3,600	790	2,000	2,700
Benzo(a)pyrene	µg/kg	5,900	5,500	7,600	4,800	1,200	3,300	4,500	2,800	4,100	740	2,100	2,700
Benzo(b)fluoranthene	µg/kg	5,100	5,700	6,500	4,900	1,400	3,500	3,300	2,100	4,000	580	2,000	1,800
Benzo(g,h,i)perylene	µg/kg	5,000	4,300	5,100	3,600	1,000	2,600	2,700	1,300	2,900	420	1,400	1,200
Benzo(k)fluoranthene	µg/kg	5,400	5,100	7,000	5,100	1,300	3,300	3,500	2,200	3,800	560	1,900	2,000
Chrysene	µg/kg	6,100	5,800	9,200	5,800	1,400	3,900	4,900	2,900	4,100	740	2,400	2,800
Dibenzo(a,h)anthracene	µg/kg	1,400	1,000	1,200	930	250	640	730	400	710	120	330	340
Fluoranthene	µg/kg	15,000	11,000	19,000	11,000	2,700	7,300	8,300	7,000	8,200	1,600	5,000	6,000
Fluorene	µg/kg	560	370	1,800	580	89	330	770	840	370	72	270	510
Indeno(1,2,3-cd)pyrene	µg/kg	4,900	4,000	4,900	3,500	1,100	2,500	2,700	1,500	3,000	460	1,400	1,300
Naphthalene	µg/kg	520	560	1,700	790	290	470	1,600	360	400	50	280	460
Phenanthrene	µg/kg	12,000	3,800	17,000	6,000	820	3,300	4,600	6,200	3,500	880	2,700	4,600
Pyrene	µg/kg	12,000	9,800	18,000	11,000	2,300	6,500	8,900	5,600	7,400	1,500	4,300	5,100
Total PAHs	µg/kg	83,000	64,000	110,000	66,000	16,000	42,000	56,000	39,000	48,000	9,100	27,000	34,000
Additional PAHs/Alkylated PAHs													
Perylene	µg/kg	NA	NA	NA	NA	NA	NA	990	NA	NA	NA	NA	NA
Benzo[e]pyrene	µg/kg	NA	NA	NA	NA	NA	NA	3,100	NA	NA	NA	NA	NA
C1-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	3,300	NA	NA	NA	NA	NA
C1-Fluoranthenes/Pyrenes	µg/kg	NA	NA	NA	NA	NA	NA	5,900	NA	NA	NA	NA	NA
C1-Fluorenes	µg/kg	NA	NA	NA	NA	NA	NA	880	NA	NA	NA	NA	NA
C1-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	790	NA	NA	NA	NA	NA
C1-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	5,400	NA	NA	NA	NA	NA
C2-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	2,200	NA	NA	NA	NA	NA
C2-Fluorenes	µg/kg	NA	NA	NA	NA	NA	NA	1,700	NA	NA	NA	NA	NA
C2-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	1,800	NA	NA	NA	NA	NA
C2-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	4,800	NA	NA	NA	NA	NA
C3-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	2,000	NA	NA	NA	NA	NA
C3-Fluorenes	µg/kg	NA	NA	NA	NA	NA	NA	2,200	NA	NA	NA	NA	NA
C3-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	2,600	NA	NA	NA	NA	NA
C3-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	3,900	NA	NA	NA	NA	NA
C4-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	1,200	NA	NA	NA	NA	NA
C4-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	2,600	NA	NA	NA	NA	NA
C4-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	3,800	NA	NA	NA	NA	NA

Summary of Seument Sample Analytical Result

Tidewater Data Report

Location ID:		SC25	SC26	SC27	SC28	SC29	SC30	SC31	SC32	SC33	SC34	SC35	SC36
Date Collected:	Units	07/08/08	07/10/08	07/14/08	07/10/08	07/10/08	07/10/08	07/14/08	07/11/08	07/11/08	07/11/08	07/11/08	07/14/08
PAHs													
Acenaphthene	µg/kg	2,500	45	11,000	160	220	760	990	2,500	220	51	650	3,200
Acenaphthylene	µg/kg	1,100	130	3,200	830	1,300	5,700	3,800	8,600	690	250	4,900	31,000
Anthracene	µg/kg	8,100	380	6,800	940	1,300	4,600	2,900	13,000	890	300	6,800	34,000
Benzo(a)anthracene	µg/kg	12,000	1,200	13,000	3,400	4,000	12,000	6,000	19,000	3,000	950	12,000	51,000
Benzo(a)pyrene	µg/kg	12,000	1,100	10,000	3,900	4,500	11,000	4,900	17,000	3,600	1,100	11,000	38,000
Benzo(b)fluoranthene	µg/kg	8,800	950	7,600	4,000	4,600	9,000	4,900	12,000	3,900	1,300	7,800	24,000
Benzo(g,h,i)perylene	µg/kg	5,900	620	6,000	3,100	3,500	7,500	3,400	10,000	2,900	910	6,900	20,000
Benzo(k)fluoranthene	µg/kg	8,400	910	8,200	4,000	4,600	9,600	4,800	13,000	3,600	1,100	8,800	28,000
Chrysene	µg/kg	11,000	1,200	13,000	4,200	4,900	11,000	6,200	20,000	4,000	1,200	13,000	53,000
Dibenzo(a,h)anthracene	µg/kg	1,700	180	1,600	800	930	2,200 J	1,000	3,200	720	240	2,100	7,200
Fluoranthene	µg/kg	24,000 D	2,700	23,000	6,800	7,900	20,000	12,000	30,000	7,500	2,400	19,000	73,000 D
Fluorene	µg/kg	3,100	75	3,900	240	340	1,300	2,600	5,800	280	79	1,800	23,000
Indeno(1,2,3-cd)pyrene	µg/kg	6,800	700	6,000	3,200	3,700	7,700	3,700	10,000	3,000	950	7,000	21,000
Naphthalene	µg/kg	3,800	53	8,000	530	1,000	4,300	9,700	17,000	530	130	4,300	21,000
Phenanthrene	µg/kg	25,000	1,200	16,000	2,400	3,300	9,500	5,500	33,000	2,700	840	18,000	110,000 D
Pyrene	µg/kg	23,000	2,100	27,000	6,800	8,600	22,000	11,000	35,000	6,500	2,100	23,000	94,000 D
Total PAHs	µg/kg	160,000	14,000	160,000	45,000	55,000	140,000	83,000	250,000	44,000	14,000	150,000	630,000
Additional PAHs/Alkylated PAHs													
Perylene	µg/kg	NA	NA	NA	NA	NA	NA	1,300	NA	1,000	NA	NA	NA
Benzo[e]pyrene	µg/kg	NA	NA	NA	NA	NA	NA	3,700	NA	3,100	NA	NA	NA
C1-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	4,500	NA	1,900	NA	NA	NA
C1-Fluoranthenes/Pyrenes	µg/kg	NA	NA	NA	NA	NA	NA	8,000	NA	3,100	NA	NA	NA
C1-Fluorenes	µg/kg	NA	NA	NA	NA	NA	NA	1,200	NA	170	NA	NA	NA
C1-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	2,500	NA	290	NA	NA	NA
C1-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	6,000	NA	1,400	NA	NA	NA
C2-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	2,500	NA	1,100	NA	NA	NA
C2-Fluorenes	µg/kg	NA	NA	NA	NA	NA	NA	1,000	NA	310	NA	NA	NA
C2-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	2,700	NA	340	NA	NA	NA
C2-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	4,000	NA	980	NA	NA	NA
C3-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	1,700	NA	1,100	NA	NA	NA
C3-Fluorenes	µg/kg	NA	NA	NA	NA	NA	NA	960	NA	630	NA	NA	NA
C3-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	1,800	NA	270	NA	NA	NA
C3-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	1,800	NA	620	NA	NA	NA
C4-Chrysenes	µg/kg	NA	NA	NA	NA	NA	NA	840	NA	680	NA	NA	NA
C4-Naphthalenes	µg/kg	NA	NA	NA	NA	NA	NA	1,000	NA	190	NA	NA	NA
C4-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA	NA	630	NA	320	NA	NA	NA

Tidewater Data Report

Former Tidewater MGP Site, Pawtucket, Rhode Island

Location ID:		SC37	SC38	SC39	SC40	SC41	SC42	SC43	SC44	SC45	SC46	SC47	SW2
Date Collected:	Units	07/14/08	07/11/08	07/11/08	07/09/08	07/09/08	07/10/08	07/10/08	07/14/08	07/14/08	07/14/08	07/11/08	07/11/08
PAHs													
Acenaphthene	µg/kg	1,200	310	180	21	27	150	210	37 [56]	540	69	52	260,000
Acenaphthylene	µg/kg	3,200	2,000	840	89	110	340	370	210 [130]	590	100	430	430,000
Anthracene	µg/kg	2,500	1,800	840	190	200	600	730	210 [190]	1,400	220	980	1,200,000
Benzo(a)anthracene	µg/kg	5,800	4,800	2,700	520	560	1,700	1,900	840 [740]	4,200	760	3,400	1,000,000
Benzo(a)pyrene	µg/kg	4,800	5,000	3,000	530	540	2,000	2,100	690 [610]	3,000	570	2,700	870,000
Benzo(b)fluoranthene	µg/kg	4,900	5,200	3,700	450	440	1,300	1,800	700 [640]	3,300	670	2,500	700,000
Benzo(g,h,i)perylene	µg/kg	4,000	4,100	2,700	310	300	1,100	1,300	510 [460]	2,500	470	1,800	450,000
Benzo(k)fluoranthene	µg/kg	5,200	5,100	3,200	460	430	1,400	1,900	750 [700]	3,100	690	2,400	670,000
Chrysene	µg/kg	6,000	5,400	3,400	540	520	1,800	2,000	960 [820]	4,400	920	3,000	840,000
Dibenzo(a,h)anthracene	µg/kg	1,200	1,100	680	93	90	270	340	140 [120]	630	130	530	110,000
Fluoranthene	µg/kg	11,000	8,900	6,200	1,100	1,100	3,600	4,500	1,800 [1,700]	11,000	2,200	6,600	3,500,000
Fluorene	µg/kg	1,200	480	240	40	48	190	250	66 [67]	690	75	96	840,000
Indeno(1,2,3-cd)pyrene	µg/kg	4,000	4,300	2,700	360	340	1,100	1,400	490 [480]	2,500	470	2,100	550,000
Naphthalene	µg/kg	2,500	1,400	510	42	43	440	310	58 [96]	610	52	220	100,000
Phenanthrene	µg/kg	5,400	4,000	1,900	530	540	1,900	2,500	610 [720]	7,100	1,000	1,500	2,400,000
Pyrene	µg/kg	12,000	9,400	5,700	870	870	3,600	3,800	1,600 [1,400]	9,300	1,800	6,000	2,400,000
Total PAHs	µg∕kg	75,000	63,000	39,000	6,100	6,100	21,000	25,000	9,600 [8,900]	54,000	10,000	34,000	16,000,000
Additional PAHs/Alkylated PAHs													
Perylene	µg/kg	NA	NA	NA	NA	NA							
Benzo[e]pyrene	µg/kg	NA	NA	NA	NA	NA							
C1-Chrysenes	µg/kg	NA	NA	NA	NA	NA							
C1-Fluoranthenes/Pyrenes	µg/kg	NA	NA	NA	NA	NA							
C1-Fluorenes	µg/kg	NA	NA	NA	NA	NA							
C1-Naphthalenes	µg/kg	NA	NA	NA	NA	NA							
C1-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA							
C2-Chrysenes	µg/kg	NA	NA	NA	NA	NA							
C2-Fluorenes	µg/kg	NA	NA	NA	NA	NA							
C2-Naphthalenes	µg/kg	NA	NA	NA	NA	NA							
C2-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA							
C3-Chrysenes	µg/kg	NA	NA	NA	NA	NA							
C3-Fluorenes	µg/kg	NA	NA	NA	NA	NA							
C3-Naphthalenes	µg/kg	NA	NA	NA	NA	NA							
C3-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA							
C4-Chrysenes	µg/kg	NA	NA	NA	NA	NA							
C4-Naphthalenes	µg/kg	NA	NA	NA	NA	NA							
C4-Phenanthrenes/Anthracenes	µg/kg	NA	NA	NA	NA	NA							
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Notes:

D – Compound quantitated using a secondary dilution.

J – Estimated value.

NA – Not analyzed.

Total PAH includes sum of 16 parent PAH compounds. Values in brackets are field duplicates.

Tidewater Data Report Former Tidewater MGP Site, Pawtucket, Rhode Island

		Total Petroleum
Location ID:	Date Collected:	Hydrocarbons
		mg/kg
SC1	07/15/08	1,290
SC2	07/15/08	520 [573]
SC3	07/09/08	1,790
SC4	07/09/08	980
SC5	07/09/08	611
SC6	07/09/08	7,280
SC7	07/14/08	2,750
SC8	07/14/08	6,890
SC9	07/10/08	5,820 [5,410]
SC10	07/08/08	1,980
SC11	07/15/08	5,390
SC12	07/08/08	406
SC12	07/15/08	1,730
SC14	07/15/08	9,320
SC14 SC15	07/16/08	9,630
SC16	07/16/08	5,010
SC17	07/10/08	4,330
SC18	07/14/08	12,200
SC18	07/10/08	13,600
SC20	07/09/08	924
SC20	07/08/08	-
		3,690
SC22	07/09/08	518
SC23 SC24	07/08/08	4,880
	07/08/08	2,260
SC25	07/08/08	1,620 594
SC26	07/10/08	
SC27	07/14/08	8,040
SC28	07/10/08	7,770
SC29	07/10/08	3,870
SC30	07/10/08	6,090
SC31	07/14/08	1,290
SC32	07/11/08	6,900
SC33	07/11/08	9,080
SC34	07/11/08	3,480
SC35	07/11/08	6,600
SC36	07/14/08	4,790
SC37	07/14/08	3,960
SC38	07/11/08	7,300
SC39	07/11/08	12,400
SC40	07/09/08	538
SC41	07/09/08	729
SC42	07/10/08	1,480
SC43	07/10/08	3,150
SC44	07/14/08	958 [1,260]
SC45	07/14/08	637
SC46	07/14/08	451
SC47	07/11/08	1,380
SW2	07/11/08	87,700

Tidewater Data Report

Location ID:		SC1	SC2	SC3	SC4	SC5	SC6	SC7	SC8	SC9	SC10	SC11
Date Collected:	Units	07/15/08	07/15/08	07/09/08	07/09/08	07/09/08	07/09/08	07/14/08	07/14/08	07/10/08	07/08/08	07/15/08
Volatile Organics												
1,1,1,2-Tetrachloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,1,1-Trichloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,1,2,2-Tetrachloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,1,2-Trichloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,1-Dichloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,1-Dichloroethene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,1-Dichloropropene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2,3-Trichlorobenzene	µg/kg	1.7 UJ	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2,3-Trichloropropane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2,4-Trichlorobenzene	µg/kg	1.7 UJ	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2,4-Trimethylbenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2-Dibromo-3-chloropropane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2-Dibromoethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2-Dichlorobenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.1 J
1,2-Dichloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,2-Dichloropropane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,3,5-Trimethylbenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,3-Dichlorobenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,3-Dichloropropane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
1,4-Dichlorobenzene	µg/kg	1.4 J	1.6 U [1.7 U]	2.0 U	2.9	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	7.6	3.6
2,2-Dichloropropane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
2-Butanone	µg/kg	2.3	1.6 U [1.7 U]	3.4	5.8	1.3 U	140	9.2	12	33 [76]	3.3	27
2-Chloroethylvinylether	µg/kg	4.3 U	4.1 U [4.2 U]	2.0 U	1.4 U	1.3 U	4.9 U	5.1 U	6.0 U	3.7 U [9.0 U]	1.6 U	4.0 U
2-Chlorotoluene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
2-Hexanone	µg/kg	4.3 U	4.1 U [4.2 U]	2.0 U	1.4 U	1.3 U	4.9 U	5.1 U	6.0 U	3.7 U [9.0 U]	1.6 U	4.0 U
4-Chlorotoluene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
4-Methyl-2-pentanone	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Acetone	µg/kg	24 U	12 U [12 U]	21 U	26 U	13 UJ	490 J	46 U	63 U	130 J [300]	21 U	130 J
Benzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.0 J
Bromobenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Bromodichloromethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Bromoform	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Bromomethane	µg/kg	1.7 UJ	1.6 UJ [1.7 UJ]	5.0 UJ	3.6 UJ	3.3 UJ	12 UJ	2.0 UJ	2.4 UJ	9.3 UJ [3.6 U]	3.9 UJ	1.6 UJ

Tidewater Data Report

Location ID:		SC1	SC2	SC3	SC4	SC5	SC6	SC7	SC8	SC9	SC10	SC11
Date Collected:	Units	07/15/08	07/15/08	07/09/08	07/09/08	07/09/08	07/09/08	07/14/08	07/14/08	07/10/08	07/08/08	07/15/08
Volatile Organics		•	•				•					
Carbon Disulfide	µg/kg	7.6	5.7 [4.6]	2.6	4.0	2.0	67	5.8	9.0	45 [67]	11	15
Carbon Tetrachloride	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Chlorobenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Chloroethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Chloroform	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Chloromethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
cis-1,2-Dichloroethene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
cis-1,3-Dichloropropene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Dibromochloromethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Dibromomethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Dichlorodifluoromethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Ethylbenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Hexachlorobutadiene	µg/kg	1.7 UJ	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Isopropylbenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	3.4	3.7 U [3.6 U]	1.6 U	4.8
Methyl tert-butyl ether	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Methylene Chloride	µg/kg	4.3 U	4.1 U [4.2 U]	5.0 U	3.6 U	3.3 U	12 U	5.1 U	6.0 U	9.3 U [9.0 U]	3.9 U	4.0 U
n-Butylbenzene	µg/kg	4.3 UJ	4.1 U [4.2 U]	2.0 U	1.4 U	1.3 U	4.9 U	5.1 U	6.0 U	3.7 U [9.0 U]	1.6 U	4.0 U
n-Propylbenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
o-Xylene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
p/m-Xylene	µg/kg	3.5 U	3.3 U [3.4 U]	4.0 U	2.9 U	2.6 U	9.9 U	4.1 U	4.8 U	7.5 U [7.2 U]	3.1 U	3.2 U
p-Isopropyltoluene	µg/kg	4.3 UJ	4.1 U [4.2 U]	2.0 U	1.4 U	1.3 U	4.9 U	5.1 U	6.0 U	3.7 U [9.0 U]	1.6 U	4.0 U
sec-Butylbenzene	µg/kg	1.7 UJ	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	2.5
Styrene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
tert-Butylbenzene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Tetrachloroethene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Toluene	µg/kg	0.87 J	1.6 U [1.2 J]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 J
trans-1,2-Dichloroethene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
trans-1,3-Dichloropropene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Trichloroethene	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Trichlorofluoromethane	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Vinyl Acetate	µg/kg	1.7 UJ	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U
Vinyl Chloride	µg/kg	1.7 U	1.6 U [1.7 U]	2.0 U	1.4 U	1.3 U	4.9 U	2.0 U	2.4 U	3.7 U [3.6 U]	1.6 U	1.6 U

Tidewater Data Report

Location ID:		SC12	SC13	SC14	SC15	SC16	SC17	SC18	SC19	SC20	SC21	SC22	SC23	SC24
Date Collected:	Units	07/08/08	07/15/08	07/15/08	07/16/08	07/16/08	07/10/08	07/14/08	07/10/08	07/09/08	07/08/08	07/09/08	07/08/08	07/08/08
Volatile Organics						•								
1,1,1,2-Tetrachloroethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,1,1-Trichloroethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,1,2,2-Tetrachloroethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,1,2-Trichloroethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,1-Dichloroethane	µg/kg	1.6 U	0.98 J	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,1-Dichloroethene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,1-Dichloropropene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2,3-Trichlorobenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2,3-Trichloropropane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2,4-Trichlorobenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2,4-Trimethylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2-Dibromo-3-chloropropane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2-Dibromoethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2-Dichlorobenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2-Dichloroethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,2-Dichloropropane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,3,5-Trimethylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,3-Dichlorobenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,3-Dichloropropane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
1,4-Dichlorobenzene	µg/kg	4.0	3.4	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.2 J
2,2-Dichloropropane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
2-Butanone	µg/kg	2.7	3.4	68	73	46	46 J	52	110	2.0	36	1.7	56	1.8 U
2-Chloroethylvinylether	µg/kg	1.6 U	3.8 U	10 U	12 U	10 U	6.5 U	14 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
2-Chlorotoluene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
2-Hexanone	µg/kg	1.6 U	3.8 U	10 U	12 U	10 U	6.5 U	14 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
4-Chlorotoluene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
4-Methyl-2-pentanone	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Acetone	µg/kg	20 U	17 U	240	270 J	190 J	210 J	250 J	380 J	17 UJ	130	9.6 UJ	200	14 U
Benzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Bromobenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Bromodichloromethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Bromoform	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Bromomethane	µg/kg	3.9 UJ	1.5 UJ	4.1 UJ	4.8 UJ	4.0 UJ	2.6 UJ	5.6 UJ	10 UJ	3.8 UJ	9.4 UJ	2.9 UJ	6.5 UJ	4.5 UJ

Tidewater Data Report

Location ID:		SC12	SC13	SC14	SC15	SC16	SC17	SC18	SC19	SC20	SC21	SC22	SC23	SC24
Date Collected:	Units	07/08/08	07/15/08	07/15/08	07/16/08	07/16/08	07/10/08	07/14/08	07/10/08	07/09/08	07/08/08	07/09/08	07/08/08	07/08/08
Volatile Organics		•	•											
Carbon Disulfide	µg/kg	7.4	4.1	27	40	26	21	32	45	11	21	0.64 J	18	3.8
Carbon Tetrachloride	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Chlorobenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Chloroethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Chloroform	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Chloromethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
cis-1,2-Dichloroethene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
cis-1,3-Dichloropropene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Dibromochloromethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Dibromomethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Dichlorodifluoromethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Ethylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Hexachlorobutadiene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Isopropylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	1.9 J	1.8 U
Methyl tert-butyl ether	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Methylene Chloride	µg/kg	3.9 U	3.8 U	10 U	12 U	10 U	6.5 U	14 U	10 U	3.8 U	9.4 U	2.9 U	6.5 U	4.5 U
n-Butylbenzene	µg/kg	1.6 U	3.8 U	10 U	12 U	10 U	6.5 U	14 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
n-Propylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
o-Xylene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
p/m-Xylene	µg/kg	3.1 U	3.0 U	8.2 U	9.5 U	8.0 U	5.2 U	11 U	8.3 U	3.0 U	7.5 U	2.3 U	5.2 U	3.6 U
p-lsopropyltoluene	µg/kg	1.6 U	3.8 U	10 U	12 U	10 U	6.5 U	14 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
sec-Butylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Styrene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
tert-Butylbenzene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Tetrachloroethene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Toluene	µg/kg	3.6	1.5 U	2.1 J	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.1 J	3.8 U	0.71 J	2.6 U	1.8 U
trans-1,2-Dichloroethene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
trans-1,3-Dichloropropene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Trichloroethene	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Trichlorofluoromethane	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Vinyl Acetate	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U
Vinyl Chloride	µg/kg	1.6 U	1.5 U	4.1 U	4.8 U	4.0 U	2.6 U	5.6 U	4.1 U	1.5 U	3.8 U	1.2 U	2.6 U	1.8 U

Tidewater Data Report

1,1,1-2:Tetrachloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1,1-Trichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1,2-Trichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1-Dichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1-Dichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,2,3-Trichloropropane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,2,4-Trichloropenpane μg/kg <	Location ID:		SC25	SC26	SC27	SC28	SC29	SC30	SC31	SC32	SC33	SC34	SC35	SC36
1,1,1-2:Tetrachloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1,1-Trichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1,2-Trichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1-Dichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1-Dichloroethane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,2,3-Trichloropropane μg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,2,4-Trichloropenpane μg/kg <	Date Collected:	Units	07/08/08	07/10/08	07/14/08	07/10/08	07/10/08	07/10/08	07/14/08	07/11/08	07/11/08	07/11/08	07/11/08	07/14/08
1,1-Trichloroethane µg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1,2-Trichloroethane µg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1-Dichloroethane µg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,1-Dichloroethane µg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 120 U 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,2,3-Trichloroporpane µg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 180 U 4,9 U 3,1 U 2,3 U 2,2 U 1,2,4-Trinethybenzene µg/kg 1,3 U 1,5 U 2,5 U 4,0 U 3,2 U 2,6 UJ 180 U 4,9 U 3,1 U 2,3 U 2,3 U 2,2 U 1,2-Dichorobenzene µg/kg 1,3 U <th>Volatile Organics</th> <th></th>	Volatile Organics													
1,1,2-2-Tetrachloroethane µg/kg 1.3,U 1.5,U 2.5,U 4.0,U 3.2,U 2.6,U,I 180,U 4.9,U 3.1,U 2.3,U 2.2,U 1,1,2-Trichloroethane µg/kg 1.3,U 1.5,U 2.5,U 4.0,U 3.2,U 2.6,U,I 180,U 4.9,U 3.1,U 2.3,U 2.2,U 1,1-Dichloroethane µg/kg 1.3,U 1.5,U 2.5,U 4.0,U 3.2,U 2.6,U,I 180,U 4.9,U 3.1,U 2.3,U 2.2,U 1,1-Dichloroptopene µg/kg 1.3,U 1.5,U 2.5,U 4.0,U 3.2,U 2.6,U,I 180,U 4.9,U 3.1,U 2.3,U 2.2,U 1,2,4-Trichloroptopane µg/kg 1.3,U 1.5,U 2.5,U 4.0,U 3.2,U 2.6,U,I 70,J 180,U 4.9,U 3.1,U 2.3,U 2.2,U 1,2,4-Trichlorobenzene µg/kg 1.3,U 1.5,U 2.5,U 4.0,U 3.2,U 2.6,U,I 180,U 4.9,U 3.1,U 2.3,U 2.2,U	1,1,1,2-Tetrachloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2-Trinchloroethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,1-Dichloroethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,1-Dichloropropene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U. 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Strichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U. 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Atrinethylbenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U. 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichorophopane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U. 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichorophopane µg/kg 1.3 U 1.5 U 2.5 U	1,1,1-Trichloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1, Dichloroethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1, Dichloropropene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1, Dichloropropene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,3-Trichloroporpane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichorobarzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 180 U 4.9 U 3.1 U 2.3 U <t< td=""><td>1,1,2,2-Tetrachloroethane</td><td>µg/kg</td><td>1.3 U</td><td>1.5 U</td><td>2.5 U</td><td>4.0 U</td><td>3.2 U</td><td>2.6 UJ</td><td>120 U</td><td>180 U</td><td>4.9 U</td><td>3.1 U</td><td>2.3 U</td><td>2.2 U</td></t<>	1,1,2,2-Tetrachloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1.1-Dichloroethene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,1-Dichloropropene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-3-Trichloroperaene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-4-Trinchlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Diorono-3-chloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichoropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U <td>1,1,2-Trichloroethane</td> <td>µg/kg</td> <td>1.3 U</td> <td>1.5 U</td> <td>2.5 U</td> <td>4.0 U</td> <td>3.2 U</td> <td>2.6 UJ</td> <td>120 U</td> <td>180 U</td> <td>4.9 U</td> <td>3.1 U</td> <td>2.3 U</td> <td>2.2 U</td>	1,1,2-Trichloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,1-Dichloropropene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,3-Trichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 70 J 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trimethylbenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 70 J 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dibromo-schloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorophane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorophopane µg/kg 1.3 U	1,1-Dichloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2,3-Trichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trichloropopane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trichloropopane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.0 J 1,2-Diorono-3-chloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Diorhoros-tenten µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorobenzene <td< td=""><td>1,1-Dichloroethene</td><td>µg/kg</td><td>1.3 U</td><td>1.5 U</td><td>2.5 U</td><td>4.0 U</td><td>3.2 U</td><td>2.6 UJ</td><td>120 U</td><td>180 U</td><td>4.9 U</td><td>3.1 U</td><td>2.3 U</td><td>2.2 U</td></td<>	1,1-Dichloroethene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2,3-Trichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trinethylbenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 70 J 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-bitromos-3-chloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-bitromos-3-chloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-bichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-bichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-bichlorobenzene	1,1-Dichloropropene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
μg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 70 J 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2,4-Trimethylbenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 230 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dibromo-schloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloropenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 U 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 180 U 4.9 U 3.1 U 2.3 U<	1,2,3-Trichlorobenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2,4-Trimethylbenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 230 180 U 4.9 U 3.1 U 2.3 U 2.0 J 1,2-Dibromo-3-chloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichloropropane µg/	1,2,3-Trichloropropane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2-Dibromo-3-chloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dibromoethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichloropropane µg/kg<	1,2,4-Trichlorobenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	70 J	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2-Dibromoethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloroperpane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloroperpane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichloroperpane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichloroporpane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,4-Dichloroporpane µg/kg	1,2,4-Trimethylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	230	180 U	4.9 U	3.1 U	2.3 U	2.0 J
1,2-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloroethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,2-Dichloropropane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 1,3-Dichlorobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 2.3 U 2.2 U 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 2.2 U 2.2 U 2.3 U 2.2 U 2.2 U	1,2-Dibromo-3-chloropropane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,2-Dichloroethaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,2-Dichloropropaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,3.5-Trimethylbenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ74 J180 U4.9 U3.1 U2.3 U2.2 U1,3.Dichlorobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,3.Dichlorobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,4.Dichlorobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneμg/kg1.3 U<	1,2-Dibromoethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1.2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 U120 U180 U4.9 U3.1 U2.3 U2.2 U1,3-5-Trimethylbenzeneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ74 J180 U4.9 U3.1 U2.3 U2.2 U1,3-Dichlorobenzeneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,3-Dichlorobenzeneµg/kg1.2 J1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,4-Dichlorobenzeneµg/kg1.2 J1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Butanoneµg/kg1.3 U1.	1,2-Dichlorobenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,3,5-Trimethylbenzeneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ74 J180 U4.9 U3.1 U2.3 U2.2 U1,3-Dichlorobenzeneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,3-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,4-Dichlorobenzeneµg/kg1.2 J1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1.3 U1.5 U <td>1,2-Dichloroethane</td> <td>µg/kg</td> <td>1.3 U</td> <td>1.5 U</td> <td>2.5 U</td> <td>4.0 U</td> <td>3.2 U</td> <td>2.6 UJ</td> <td>120 U</td> <td>180 U</td> <td>4.9 U</td> <td>3.1 U</td> <td>2.3 U</td> <td>2.2 U</td>	1,2-Dichloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,3-Dichlorobenzeneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,3-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U1,4-Dichlorobenzeneµg/kg1.2 J1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Butanoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1.3 U1.5 U2.5 U4	1,2-Dichloropropane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1,3-Dichloropropaneµg/kg1,3 U1,5 U2,5 U4,0 U3,2 U2,6 U120 U180 U4,9 U3,1 U2,3 U2,2 U1,4-Dichlorobenzeneµg/kg1,2 J1,5 U2,5 U4,0 U3,2 U2,6 UJ120 U180 U4,9 U3,1 U2,3 U2,2 U2,2-Dichloropropaneµg/kg1,3 U1,5 U2,5 U4,0 U3,2 U2,6 UJ120 U180 U4,9 U3,1 U2,3 U2,2 U2,2-Dichloropropaneµg/kg9,52,87,361 J4766 J120 U180 UJ85 J62 J6.3 J132-Chloroethylvinyletherµg/kg1,3 U1,5 U6.4 U9.9 U8.0 U6.4 UJ300 U450 U12 U7.7 U5.8 U5.5 U2-Chlorotolueneµg/kg1,3 U1,5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1,3 U1,5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1,3 U1,5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneµg/kg1,3 U1,5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneµg/kg4.6 U26 UJ32 U2	1,3,5-Trimethylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	74 J	180 U	4.9 U	3.1 U	2.3 U	2.2 U
1.4-Dichlorobenzeneµg/kg1.2 J1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2.2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Butanoneµg/kg9.52.87.361 J4766 J120 U180 UJ85 J62 J6.3 J132-Chloroteluylvinyletherµg/kg1.3 U1.5 U6.4 U9.9 U8.0 U6.4 UJ300 U450 U12 U7.7 U5.8 U5.5 U2-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UAcetoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UBenzeneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U <td< td=""><td>1,3-Dichlorobenzene</td><td>µg/kg</td><td>1.3 U</td><td>1.5 U</td><td>2.5 U</td><td>4.0 U</td><td>3.2 U</td><td>2.6 UJ</td><td>120 U</td><td>180 U</td><td>4.9 U</td><td>3.1 U</td><td>2.3 U</td><td>2.2 U</td></td<>	1,3-Dichlorobenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
2,2-Dichloropropaneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Butanoneµg/kg9.52.87.361 J4766 J120 U180 UJ85 J62 J6.3 J132-Chloroethylvinyletherµg/kg1.3 U1.5 U6.4 U9.9 U8.0 U6.4 UJ300 U450 U12 U7.7 U5.8 U5.5 U2-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Horotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Chlorotolueneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UAcetoneµg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3	1,3-Dichloropropane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
2-Butanone µg/kg 9.5 2.8 7.3 61 J 47 66 J 120 U 180 UJ 85 J 62 J 6.3 J 13 2-Chloroethylvinylether µg/kg 1.3 U 1.5 U 6.4 U 9.9 U 8.0 U 6.4 UJ 300 U 450 U 12 U 7.7 U 5.8 U 5.5 U 2-Chlorotoluene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 2-Hexanone µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 4-Chlorotoluene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 4-Methyl-2-pentanone µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U	1,4-Dichlorobenzene	µg/kg	1.2 J	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
2-Chloroethylvinyletherμg/kg1.3 U1.5 U6.4 U9.9 U8.0 U6.4 UJ300 U450 U12 U7.7 U5.8 U5.5 U2-Chlorotolueneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Chlorotolueneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UAcetoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UBenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140 U130 J4.9 U3.1 U2.3 U2.3 U1.3 JBromobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140 U130 J4.9 U3.1 U2.3 U2.3 U2.2 UBromodichloromethaneμg/kg1.3 U1.5 U </td <td>2,2-Dichloropropane</td> <td>µg/kg</td> <td>1.3 U</td> <td>1.5 U</td> <td>2.5 U</td> <td>4.0 U</td> <td>3.2 U</td> <td>2.6 UJ</td> <td>120 U</td> <td>180 U</td> <td>4.9 U</td> <td>3.1 U</td> <td>2.3 U</td> <td>2.2 U</td>	2,2-Dichloropropane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
2-Chlorotolueneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 U120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneμg/kg1.3 U1.5 U6.4 U9.9 U8.0 U6.4 U300 U450 U12 U7.7 U5.8 U5.5 U4-Chlorotolueneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UAcetoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UBenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140 U130 J4.9 U3.1 U2.3 U2.2 UBromobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140 U130 J4.9 U3.1 U2.3 U2.2 UBromobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140 U180 U4.9 U3.1 U2.3 U2.2 UBromodichloromethaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U<	2-Butanone	µg/kg	9.5	2.8	7.3	61 J	47	66 J	120 U	180 UJ	85 J	62 J	6.3 J	13
2-Chlorotolueneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U2-Hexanoneμg/kg1.3 U1.5 U6.4 U9.9 U8.0 U6.4 U300 U450 U12 U7.7 U5.8 U5.5 U4-Chlorotolueneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 U4-Methyl-2-pentanoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ120 U180 U4.9 U3.1 U2.3 U2.2 UAcetoneμg/kg46 U26 UJ32 U280 J180200 J300 U630 J370 J250 J4.4 UBenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140130 J4.9 U3.1 U2.3 U2.2 UAcetoneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140130 J4.9 U3.1 U2.3 U2.2 UBenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140130 J4.9 U3.1 U2.3 U2.3 U1.3 JBromobenzeneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ140 U180 U4.9 U3.1 U2.3 U2.2 UBromodichloromethaneμg/kg1.3 U1.5 U2.5 U4.0 U3.2 U2.6 UJ180 U	2-Chloroethylvinylether	µg/kg	1.3 U	1.5 U	6.4 U	9.9 U	8.0 U	6.4 UJ	300 U	450 U	12 U	7.7 U	5.8 U	5.5 U
4-Chlorotoluene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U 4-Methyl-2-pentanone µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Acetone µg/kg 46 U 26 UJ 32 U 280 J 180 200 J 300 U 630 J 370 J 250 J 44 UJ 63 J Benzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 2.3 U 2.2 U Benzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 2.	2-Chlorotoluene		1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
4-Methyl-2-pentanone µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Acetone µg/kg 46 U 26 UJ 32 U 280 J 180 200 J 300 U 630 J 370 J 250 J 44 UJ 63 J Benzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 63 J Benzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 1.3 U Bromobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3	2-Hexanone	µg/kg	1.3 U	1.5 U	6.4 U	9.9 U	8.0 U	6.4 U	300 U	450 U	12 U	7.7 U	5.8 U	5.5 U
Acetone µg/kg 46 U 26 UJ 32 U 280 J 180 200 J 300 U 630 J 370 J 250 J 44 UJ 63 J Benzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 1.3 J Bromobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromoform µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2	4-Chlorotoluene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Benzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 1.3 J Bromobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 140 130 J 4.9 U 3.1 U 2.3 U 1.3 J Bromobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromoform µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U	4-Methyl-2-pentanone	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Bromobenzene µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromoform µg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U	Acetone	µg/kg	46 U	26 UJ	32 U	280 J	180	200 J	300 U	630 J	370 J	250 J	44 UJ	63 J
Bromodichloromethane μg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromodichloromethane μg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromoform μg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U	Benzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	140	130 J	4.9 U	3.1 U	2.3 U	1.3 J
Bromodichloromethane μg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U Bromoform μg/kg 1.3 U 1.5 U 2.5 U 4.0 U 3.2 U 2.6 UJ 120 U 180 U 4.9 U 3.1 U 2.3 U 2.2 U	Bromobenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
	Bromodichloromethane		1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
	Bromoform	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
	Bromomethane		3.1 UJ	3.7 UJ	2.5 UJ	4.0 UJ	3.2 U	2.6 UJ	120 UJ	180 UJ	4.9 UJ	3.1 UJ	2.3 UJ	2.2 UJ

Tidewater Data Report

Location ID:		SC25	SC26	SC27	SC28	SC29	SC30	SC31	SC32	SC33	SC34	SC35	SC36
Date Collected:	Units	07/08/08	07/10/08	07/14/08	07/10/08	07/10/08	07/10/08	07/14/08	07/11/08	07/11/08	07/11/08	07/11/08	07/14/08
Volatile Organics													
Carbon Disulfide	µg/kg	6.9	1.7	3.0	29	58	30 J	120 U	180 U	43	46	3.3	3.4
Carbon Tetrachloride	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Chlorobenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Chloroethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Chloroform	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Chloromethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
cis-1,2-Dichloroethene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
cis-1,3-Dichloropropene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	R	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Dibromochloromethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Dibromomethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Dichlorodifluoromethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Ethylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Hexachlorobutadiene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Isopropylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Methyl tert-butyl ether	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Methylene Chloride	µg/kg	3.1 U	3.7 U	6.4 U	9.9 U	8.0 U	6.4 U	300 U	450 U	12 U	7.7 U	5.8 U	5.5 U
n-Butylbenzene	µg/kg	1.3 U	1.5 U	6.4 U	9.9 U	8.0 U	6.4 UJ	300 U	450 U	12 U	7.7 U	5.8 U	5.5 U
n-Propylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
o-Xylene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	97 J	180 U	4.9 U	3.1 U	2.3 U	2.2 U
p/m-Xylene	µg/kg	2.5 U	3.0 U	5.1 U	7.9 U	6.4 U	5.1 UJ	130 J	360 U	9.8 U	6.2 U	4.6 U	1.3 J
p-Isopropyltoluene	µg/kg	1.3 U	1.5 U	6.4 U	9.9 U	8.0 U	6.4 UJ	300 U	450 U	12 U	7.7 U	5.8 U	5.5 U
sec-Butylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Styrene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
tert-Butylbenzene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Tetrachloroethene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Toluene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	60 J	180 U	4.9 U	3.1 U	2.3 U	2.2 U
trans-1,2-Dichloroethene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
trans-1,3-Dichloropropene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	R	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Trichloroethene	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Trichlorofluoromethane	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Vinyl Acetate	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 UJ	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U
Vinyl Chloride	µg/kg	1.3 U	1.5 U	2.5 U	4.0 U	3.2 U	2.6 U	120 U	180 U	4.9 U	3.1 U	2.3 U	2.2 U

Tidewater Data Report

Location ID:		SC37	SC38	SC39	SC40	SC41	SC42	SC43	SC44	SC45	SC46	SC47	SW2
Date Collected:	Units	07/14/08	07/11/08	07/11/08	07/09/08	07/09/08	07/10/08	07/10/08	07/14/08	07/14/08	07/14/08	07/11/08	07/11/08
Volatile Organics													
1,1,1,2-Tetrachloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,1,1-Trichloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,1,2,2-Tetrachloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,1,2-Trichloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,1-Dichloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,1-Dichloroethene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,1-Dichloropropene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2,3-Trichlorobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2,3-Trichloropropane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2,4-Trichlorobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2,4-Trimethylbenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	7,800
1,2-Dibromo-3-chloropropane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2-Dibromoethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2-Dichlorobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2-Dichloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,2-Dichloropropane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,3,5-Trimethylbenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	3,800
1,3-Dichlorobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,3-Dichloropropane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
1,4-Dichlorobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	8.2	390 U
2,2-Dichloropropane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
2-Butanone	µg/kg	12 J	49 J	120 J	2.0	1.9	30	57 J	7.0 [15 J]	1.2 J	1.2 J	1.6 UJ	390 UJ
2-Chloroethylvinylether	µg/kg	6.0 UJ	9.6 U	13 U	1.2 U	1.3 U	1.4 U	2.1 UJ	3.9 U [4.6 UJ]	3.8 U	3.5 U	4.0 U	970 U
2-Chlorotoluene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
2-Hexanone	µg/kg	6.0 UJ	9.6 U	13 U	1.2 U	1.3 U	1.4 U	2.1 UJ	3.9 U [4.6 UJ]	3.8 U	3.5 U	4.0 U	970 U
4-Chlorotoluene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
4-Methyl-2-pentanone	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Acetone	µg/kg	53 UJ	210 J	470 J	14 U	17 U	110 J	190 J	44 UJ [97 J]	13 U	8.4 U	14 UJ	1,500 UJ
Benzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	230 J
Bromobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Bromodichloromethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Bromoform	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Bromomethane	µg/kg	2.4 UJ	3.8 UJ	5.0 UJ	2.9 UJ	3.2 UJ	3.6 UJ	5.3 UJ	1.6 UJ [1.8 UJ]	1.5 UJ	1.4 UJ	1.6 UJ	390 UJ

Tidewater Data Report

Former Tidewater MGP Site, Pawtucket, Rhode Island

Location ID:		SC37	SC38	SC39	SC40	SC41	SC42	SC43	SC44	SC45	SC46	SC47	SW2
Date Collected:	Units	07/14/08	07/11/08	07/11/08	07/09/08	07/09/08	07/10/08	07/10/08	07/14/08	07/14/08	07/14/08	07/11/08	07/11/08
Volatile Organics													
Carbon Disulfide	µg/kg	15 J	44	92	1.3	1.6	20	45 J	5.3 [8.9 J]	1.9	2.2	1.7	730
Carbon Tetrachloride	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Chlorobenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Chloroethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Chloroform	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Chloromethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
cis-1,2-Dichloroethene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
cis-1,3-Dichloropropene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Dibromochloromethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Dibromomethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Dichlorodifluoromethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Ethylbenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	330 J
Hexachlorobutadiene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Isopropylbenzene	µg/kg	3.2 J	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	280 J
Methyl tert-butyl ether	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Methylene Chloride	µg/kg	6.0 UJ	9.6 U	13 U	2.9 U	3.2 U	3.6 U	5.3 UJ	3.9 U [4.6 UJ]	3.8 U	3.5 U	4.0 U	970 U
n-Butylbenzene	µg/kg	6.0 UJ	9.6 U	13 U	1.2 U	1.3 U	1.4 U	2.1 UJ	3.9 U [4.6 UJ]	3.8 U	3.5 U	4.0 U	2,700
n-Propylbenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	410
o-Xylene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	620
p/m-Xylene	µg/kg	4.8 UJ	7.6 U	10 U	2.3 U	2.6 U	2.9 U	4.2 UJ	3.1 U [3.7 UJ]	3.1 U	2.8 U	3.2 U	1,000
p-Isopropyltoluene	µg/kg	6.0 UJ	9.6 U	13 U	1.2 U	1.3 U	1.4 U	2.1 UJ	3.9 U [4.6 UJ]	3.8 U	3.5 U	4.0 U	1,500
sec-Butylbenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	370 J
Styrene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
tert-Butylbenzene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Tetrachloroethene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Toluene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	6.8 B	270 J
trans-1,2-Dichloroethene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
trans-1,3-Dichloropropene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Trichloroethene	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Trichlorofluoromethane	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Vinyl Acetate	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U
Vinyl Chloride	µg/kg	2.4 UJ	3.8 U	5.0 U	1.2 U	1.3 U	1.4 U	2.1 UJ	1.6 U [1.8 UJ]	1.5 U	1.4 U	1.6 U	390 U

Notes:

B – Analyte was also detected in the associated method blank.

J – Estimated value.

R - Rejected.

U – The compound was analyzed for but not detected. The associated value is the compound quantitation limit.

Tidewater Data Report

Location ID:		SC1	SC2	SC3	SC4	SC5	SC6	SC7	SC8	SC9	SC10
Date Collected:	Units	07/15/08	07/15/08	07/09/08	07/09/08	07/09/08	07/09/08	07/14/08	07/14/08	07/10/08	07/08/08
Inorganics											
Antimony	mg/kg	0.283 J	0.169 J [0.273 J]	0.107 J	0.180 J	R	0.633 J	0.817 J	1.73 J	1.09 J [1.25 J]	0.345 J
Arsenic	mg/kg	4.71	5.18 [3.54]	4.03 J	5.32 J	2.58 J	16.8 J	7.50	16.3	15.7 J [15.9 J]	7.07 J
Beryllium	mg/kg	0.211 J	0.364 J [0.224 J]	0.200 J	0.333 J	0.385 J	1.11 J	0.529 J	0.711 J	0.982 J [1.04 J]	0.460 J
Cadmium	mg/kg	0.551	0.816 [0.716]	1.15 J	1.07 J	0.823 J	8.45 J	7.20	8.67	10.4 J [11.1 J]	1.37 J
Chromium	mg/kg	12.7 J	15.5 J [16.4 J]	16.2 J	15.7 J	10.4 J	140 J	103 J	326 J	154 J [158 J]	18.9 J
Copper	mg/kg	79.7	35.8 [29.0]	24.8	44.7	29.0	266	244	588	322 [314]	336
Lead	mg/kg	95.8 J	55.9 J [47.9 J]	71.0 J	79.2 J	56.0 J	242 J	177 J	376 J	308 J [302 J]	170 J
Mercury	mg/kg	0.0516	0.0311 [0.0300]	0.0591 J	0.0312 J	0.0650 J	0.718 J	0.404	1.76	0.903 J [0.942 J]	0.0529 J
Nickel	mg/kg	9.39 J	11.9 J [9.73 J]	10.1	13.1	10.1	35.9	34.9 J	37.0 J	41.0 [41.0]	15.7
Selenium	mg/kg	0.184 UJ	0.194 UJ [0.188 UJ]	0.230	0.448	0.199 U	2.39	0.667 J	1.48 J	1.99 [1.88]	0.241
Silver	mg/kg	0.142	0.105 [0.0997]	0.224	0.225	0.115	3.46	3.01	2.88	5.80 [6.10]	0.272
Thallium	mg/kg	0.0486	0.0645 [0.0377 U]	0.0531	0.0530	0.0398 U	0.224	0.104	0.222	0.180 [0.183]	0.0494
Zinc	mg/kg	83.2 J	75.6 J [76.1 J]	170	95.1	89.9	439	272 J	810 J	455 [446]	158

Tidewater Data Report

Location ID:		SC11	SC12	SC13	SC14	SC15	SC16	SC17	SC18	SC19	SC20	SC21	SC22	SC23	SC24
Date Collected:	Units	07/15/08	07/08/08	07/15/08	07/15/08	07/16/08	07/16/08	07/10/08	07/14/08	07/10/08	07/09/08	07/08/08	07/09/08	07/08/08	07/08/08
Inorganics															
Antimony	mg/kg	0.557 J	0.154 J	0.271 J	1.25 J	1.34 J	1.15 J	1.19 J	1.77 J	0.981 J	0.395 J	0.654 J	1.16 J	0.484 J	0.206 J
Arsenic	mg/kg	7.48	3.41 J	3.71	24.7	18.0	15.0	6.94 J	26.6	17.6 J	3.96 J	12.7 J	3.07 J	8.39 J	2.91 J
Beryllium	mg/kg	0.321 J	0.212 J	0.248 J	1.58 J	1.16 J	0.974 J	0.512 J	1.43 J	0.706 J	0.212 J	0.827 J	0.184 J	0.445 J	1.11 J
Cadmium	mg/kg	3.73	0.555 J	1.44	13.1	13.8	9.24	4.68 J	23.5	19.7 J	1.58 J	11.0 J	0.541 J	5.87 J	0.619 J
Chromium	mg/kg	66.4 J	10.1 J	17.8 J	233 J	203 J	136 J	67.9 J	345 J	280 J	18.3 J	130 J	11.6 J	85.0 J	10.4 J
Copper	mg/kg	202	25.8	47.2	430	401	293	158	668	773	54.4	249	68.3	173	45.6
Lead	mg/kg	193 J	40.1 J	214 J	355 J	352 J	277 J	158 J	428 J	383 J	128 J	253 J	43.6 J	178 J	60.4 J
Mercury	mg/kg	0.453	0.0141 UJ	0.0616	1.52	1.22	0.792	0.426 J	1.83	1.39 J	0.182 J	0.635 J	0.0492 J	0.374 J	0.0252 J
Nickel	mg/kg	19.4 J	11.7	12.4 J	51.6 J	50.0 J	37.0 J	20.4	76.8 J	49.6	9.71	38.4	10.4	24.9	13.4
Selenium	mg/kg	0.473 J	0.176 U	0.200 J	3.14 J	2.37 J	1.79 J	0.996	2.97 J	1.77	0.271	1.77	0.193 U	0.918	0.318
Silver	mg/kg	1.39	0.0951	0.211	6.02	6.94	8.23	2.76	10.6	5.25	0.436	4.00	0.129	1.94	0.120
Thallium	mg/kg	0.0778	0.0352 U	0.0400	0.333	0.228	0.219	0.0986	0.310	0.200	0.0398 U	0.247	0.0385 U	0.132	0.0388 U
Zinc	mg/kg	280 J	64.9	225 J	654 J	578 J	467 J	206	750 J	792	93.6	408	69.8	320	105

Tidewater Data Report

Location ID:		SC25	SC26	SC27	SC28	SC29	SC30	SC31	SC32	SC33	SC34	SC35	SC36	SC37	SC38	SC39	SC40
Date Collected:	Units	07/08/08	07/10/08	07/14/08	07/10/08	07/10/08	07/10/08	07/14/08	07/11/08	07/11/08	07/11/08	07/11/08	07/14/08	07/14/08	07/11/08	07/11/08	07/09/08
Inorganics																	
Antimony	mg/kg	5.13 J	0.372 J	3.56 J	0.928 J	0.848 J	1.40 J	0.350 J	1.66 J	1.38 J	0.482 J	1.19 J	1.14 J	0.815 J	1.06 J	1.52 J	0.208 J
Arsenic	mg/kg	5.39 J	1.79 J	25.2	14.7 J	13.2 J	13.5 J	7.19	11.7 J	17.1 J	7.07 J	20.1 J	19.1	12.6	17.5 J	16.8 J	2.39 J
Beryllium	mg/kg	0.283 J	0.143 J	0.671 J	1.09 J	1.04 J	1.35 J	0.376 J	1.53 J	1.18 J	0.526 J	0.610 J	0.483 J	1.50 J	1.58 J	1.34 J	0.165 J
Cadmium	mg/kg	2.57 J	0.551 J	2.40	11.0 J	10.8 J	6.46 J	0.620	6.93 J	10.0 J	4.50 J	10.4 J	4.77	11.1	18.2 J	16.1 J	0.673 J
Chromium	mg/kg	28.7 J	7.68 J	327 J	158 J	132 J	143 J	28.7 J	138 J	172 J	69.1 J	210 J	310 J	165 J	243 J	232 J	8.96 J
Copper	mg/kg	74.5	22.5	840	320	245	271	65.8	295	357	184	348	219	282	446	516	26.6
Lead	mg/kg	124 J	28.4 J	14,700 J	306 J	257 J	246 J	77.0 J	239 J	311 J	153 J	553 J	198 J	255 J	450 J	413 J	40.9 J
Mercury	mg/kg	0.0819 J	0.0300 J	4.91	0.875 J	0.678 J	0.882 J	0.240	1.23 J	0.922 J	0.617 J	0.914 J	0.673	0.723	1.58 J	1.35 J	0.0264 J
Nickel	mg/kg	20.6	4.91	20.2 J	39.8	38.6	61.7	21.6 J	33.7	39.1	21.8	50.3	23.8 J	152 J	71.7	55.0	7.41
Selenium	mg/kg	0.277	0.264	1.53 J	2.14	2.13	2.62	0.752 J	1.70	2.45	0.957	1.78	1.72 J	1.44 J	2.17	2.90	0.187 U
Silver	mg/kg	0.658	0.124	1.67	5.69	4.05	3.35	0.250	2.74	6.72	2.85	3.37	2.33	4.51	9.50	11.0	0.109
Thallium	mg/kg	0.0549	0.0374 U	0.358	0.214	0.192	0.179	0.0577	0.148	0.240	0.112	0.174	0.108	0.140	0.243	0.283	0.0375 U
Zinc	mg/kg	131	58.2	689 J	446	384	496	230 J	391	465	200	1,440	680 J	376 J	593	583	55.2

Tidewater Data Report

Former Tidewater MGP Site, Pawtucket, Rhode Island

Location ID:		SC41	SC42	SC43	SC44	SC45	SC46	SC47	SW2
Date Collected:	Units	07/09/08	07/10/08	07/10/08	07/14/08	07/14/08	07/14/08	07/11/08	07/11/08
Inorganics									
Antimony	mg/kg	0.102 J	0.139 J	0.334 J	0.172 J [0.155 J]	0.103 UJ	0.100 UJ	0.264 J	3.22 J
Arsenic	mg/kg	2.28 J	3.87 J	4.93 J	3.04 [2.79]	2.43	1.72	3.84 J	10.4 J
Beryllium	mg/kg	0.184 J	0.237 J	0.408 J	0.233 J [0.227 J]	0.181 J	0.155 J	0.254 J	0.144 J
Cadmium	mg/kg	0.752 J	2.76 J	2.95 J	1.26 [1.28]	1.05	0.814	0.847 J	1.88 J
Chromium	mg/kg	9.88 J	15.2 J	46.5 J	18.2 J [17.9 J]	12.6 J	9.05 J	12.4 J	108 J
Copper	mg/kg	80.1	44.4	117	39.5 [39.8]	26.6	21.7	30.3	283
Lead	mg/kg	28.3 J	83.1 J	112 J	48.9 J [48.3 J]	56.3 J	26.4 J	40.6 J	318 J
Mercury	mg/kg	0.0264 J	0.0542 J	0.292 J	0.104 [0.101]	0.0641	0.0577	0.0210 J	1.48 J
Nickel	mg/kg	7.99	10.8	16.3	8.87 J [8.57 J]	7.35 J	5.77 J	12.3	15.0
Selenium	mg/kg	0.192 U	0.346	0.707	0.340 J [0.333 J]	0.201 UJ	0.197 UJ	0.232	2.52
Silver	mg/kg	0.167	0.783	1.71	0.447 [0.444]	0.190	0.141	0.175	0.697
Thallium	mg/kg	0.0385	0.0441	0.115	0.0566 [0.0562]	0.0424	0.0394 U	0.0731	0.375
Zinc	mg/kg	64.4	992	176	89.3 J [84.4 J]	70.2 J	63.5 J	79.9	206

Notes:

J – Estimated value.

R – Rejected.

U – The compound was analyzed for but not detected. The associated value is the compound quantitation limit.

Table 3-6 Summary of Sediment Sample Analytical Results – TOC, PAC, PSOL

Tidewater Data Report

Location ID:		SC1	SC2	SC3	SC4	SC5	SC6	SC7	SC8	SC9	SC10	SC11
Date Collected:	Units	07/15/08	07/15/08	07/09/08	07/09/08	07/09/08	07/09/08	07/14/08	07/14/08	07/10/08	07/08/08	07/15/08
Miscellaneous												
Percent Solids	%	82.7	83 [84]	80.2	89.3	81.5	30.9	62.5	52.7	36.6 [36.2]	82.4	74.7
Physiologically Available Cyanide	mg/kg	0.27 U	0.3 U [0.29 U]	0.091 U	0.092 U	0.099 U	0.25 U	0.15 U	0.39 U	0.18 U [0.22 U]	0.075 U	0.32 U
Solids, Total (TPH sample)	%	79	87 [89]	83	89	85	31	59	49	35 [35]	72	70
Total Organic Carbon (Run 1)	%	NA	2.2 J [0.4 J]	1	NA	0.55	4.8	2.7	4.7	NA	NA	1.5 J
Total Organic Carbon (Run 2)	%	0.75	NA	NA	1.9	NA	NA	NA	NA	4.2	0.6	NA

Table 3-6 Summary of Sediment Sample Analytical Results – TOC, PAC, PSOL

Tidewater Data Report

Location ID:		SC12	SC13	SC14	SC15	SC16	SC17	SC18	SC19	SC20	SC21	SC22	SC23	SC24
Date Collected:	Units	07/08/08	07/15/08	07/15/08	07/16/08	07/16/08	07/10/08	07/14/08	07/10/08	07/09/08	07/08/08	07/09/08	07/08/08	07/08/08
Miscellaneous														
Percent Solids	%	86	80.9	30.5	30.6	38.3	52.9	24.3	35.7	77.6	42.5	80.2	59.3	82.6
Physiologically Available Cyanide	mg/kg	0.076 U	0.26 U	0.78 U	0.8 U	0.63 U	0.13 U	0.73 U	0.17 U	0.08 U	0.15 U	0.089 U	0.14 U	0.093 U
Solids, Total (TPH sample)	%	90	80	29	30	38	54	23	35	78	41	78	55	82
Total Organic Carbon (Run 1)	%	NA	NA	4.6	NA	4	NA	5.8	5.2	1.1	2.9	0.33	2.7	0.77 J
Total Organic Carbon (Run 2)	%	0.14	3.4 J	NA	3.8	NA	2.4	NA						

Table 3-6 Summary of Sediment Sample Analytical Results – TOC, PAC, PSOL

Tidewater Data Report

Location ID:		SC25	SC26	SC27	SC28	SC29	SC30	SC31	SC32	SC33	SC34	SC35	SC36	SC37
Date Collected:	Units	07/08/08	07/10/08	07/14/08	07/10/08	07/10/08	07/10/08	07/14/08	07/11/08	07/11/08	07/11/08	07/11/08	07/14/08	07/14/08
Miscellaneous														
Percent Solids	%	76.7	83.5	50.8	34.6	42.1	53.8	75.4	60.1	31.1	53.9	55.1	63.3	52
Physiologically Available Cyanide	mg/kg	0.089 U	0.092 U	0.47 U	0.23 U	0.17 U	0.15 U	1.5	0.13 U	0.25 U	0.13 U	0.12 U	3.8	0.36 U
Solids, Total (TPH sample)	%	72	81	53	36	43	54	73	57	29	49	52	61	49
Total Organic Carbon (Run 1)	%	NA	NA	7.3	4.8	3.8	NA	NA	NA	5.1	1.7	NA	NA	4.5
Total Organic Carbon (Run 2)	%	1.6	0.39	NA	NA	NA	10 J	5.3	10	NA	NA	5.4	6.4	NA

Table 3-6 Summary of Sediment Sample Analytical Results – TOC, PAC, PSOL

Tidewater Data Report

Former Tidewater MGP Site, Pawtucket, Rhode Island

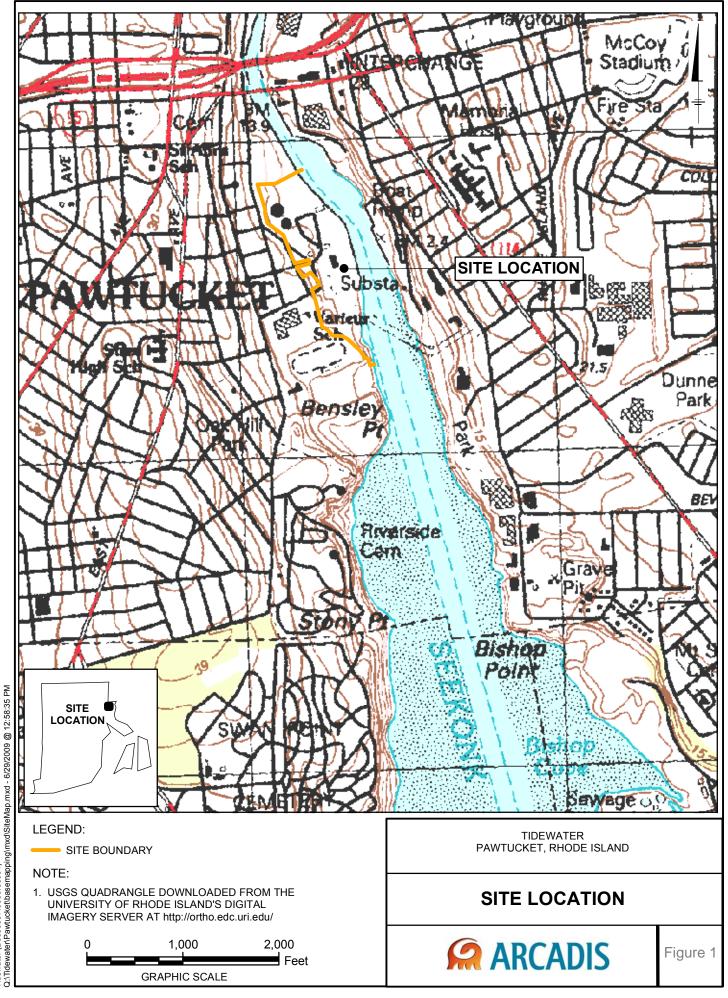
Location ID:		SC38	SC39	SC40	SC41	SC42	SC43	SC44	SC45	SC46	SC47	SW2
Date Collected:	Units	07/11/08	07/11/08	07/09/08	07/09/08	07/10/08	07/10/08	07/14/08	07/14/08	07/14/08	07/11/08	07/11/08
Miscellaneous												
Percent Solids	%	34.5	23.7	84.4	81.3	77.3	62	76 [73.4]	77.8	78.4	83.2	39.6
Physiologically Available Cyanide	mg/kg	0.21 U	0.32 U	0.089 U	0.076 U	0.1 U	0.11 U	0.23 U [0.3 U]	0.19 U	0.27 U	0.091 U	0.19 U
Solids, Total (TPH sample)	%	36	23	82	78	78	62	73 [72]	74	78	83	53
Total Organic Carbon (Run 1)	%	NA	6.5	NA	NA	0.92	NA	1.2	0.76	0.34	NA	34
Total Organic Carbon (Run 2)	%	4.7	NA	1	1.3	NA	1.8	NA	NA	NA	1.4	NA

Notes:

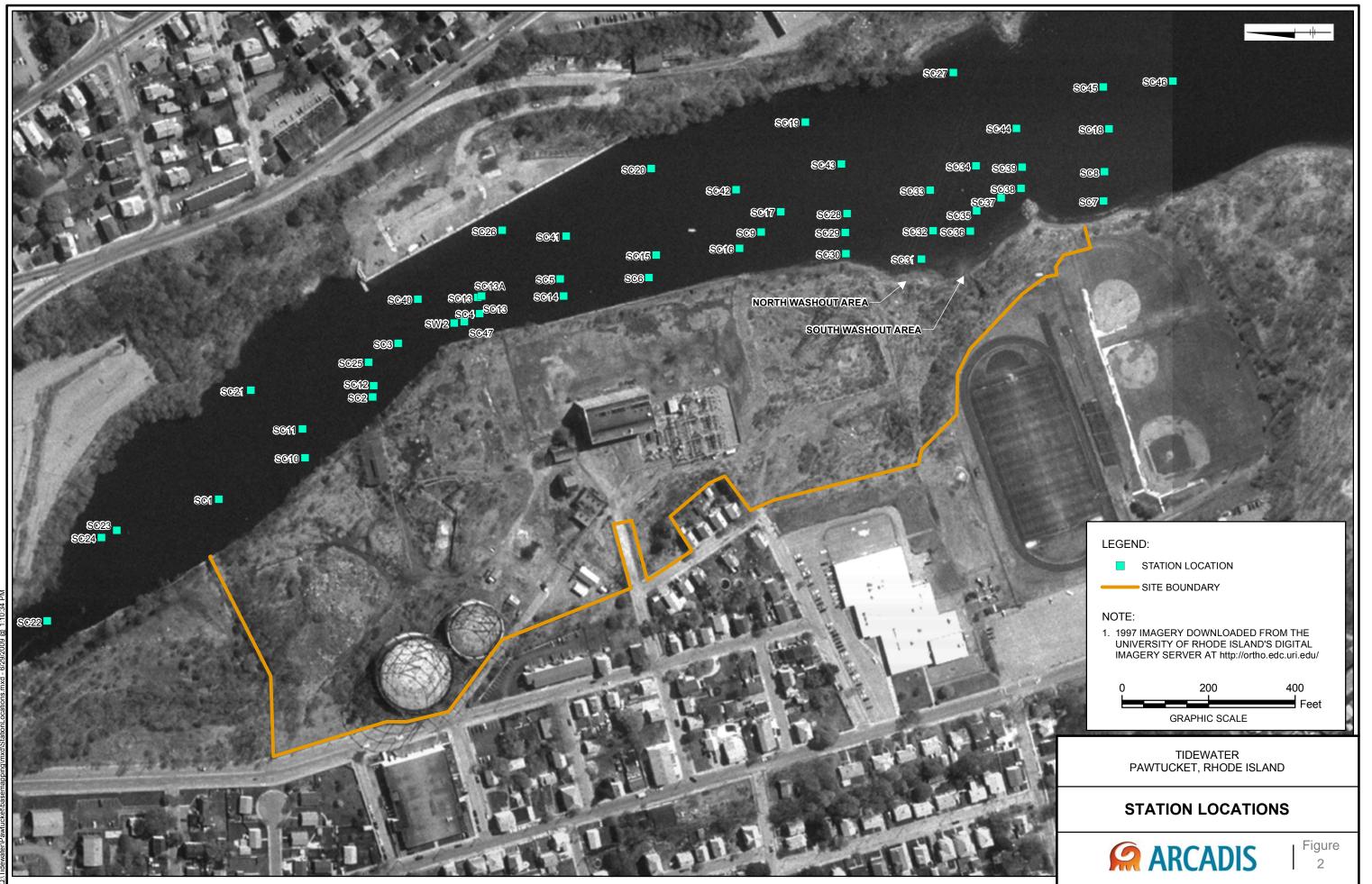
J – Indicates an estimated value.

U – The compound was analyzed for but not detected.

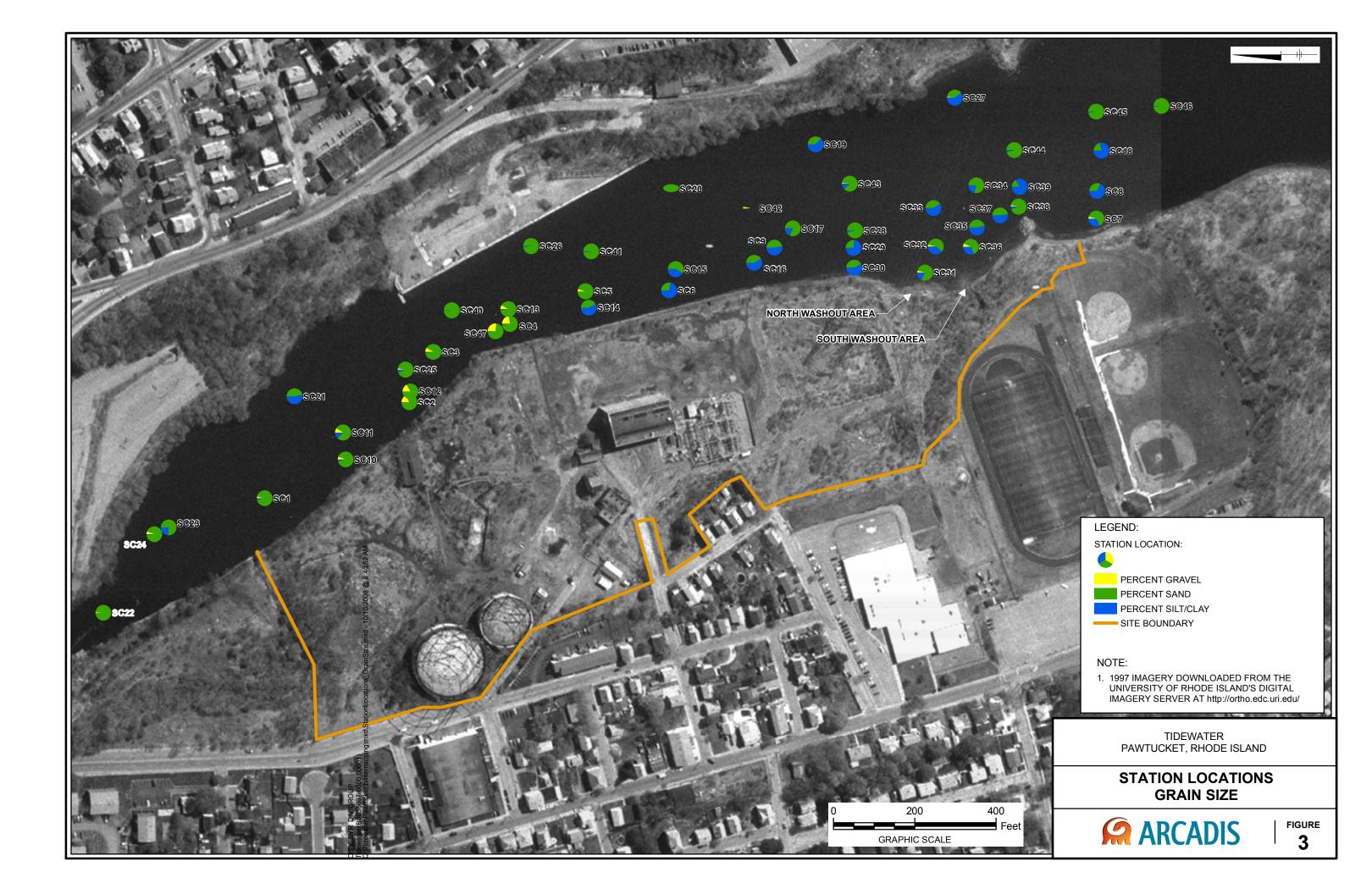
FIGURES

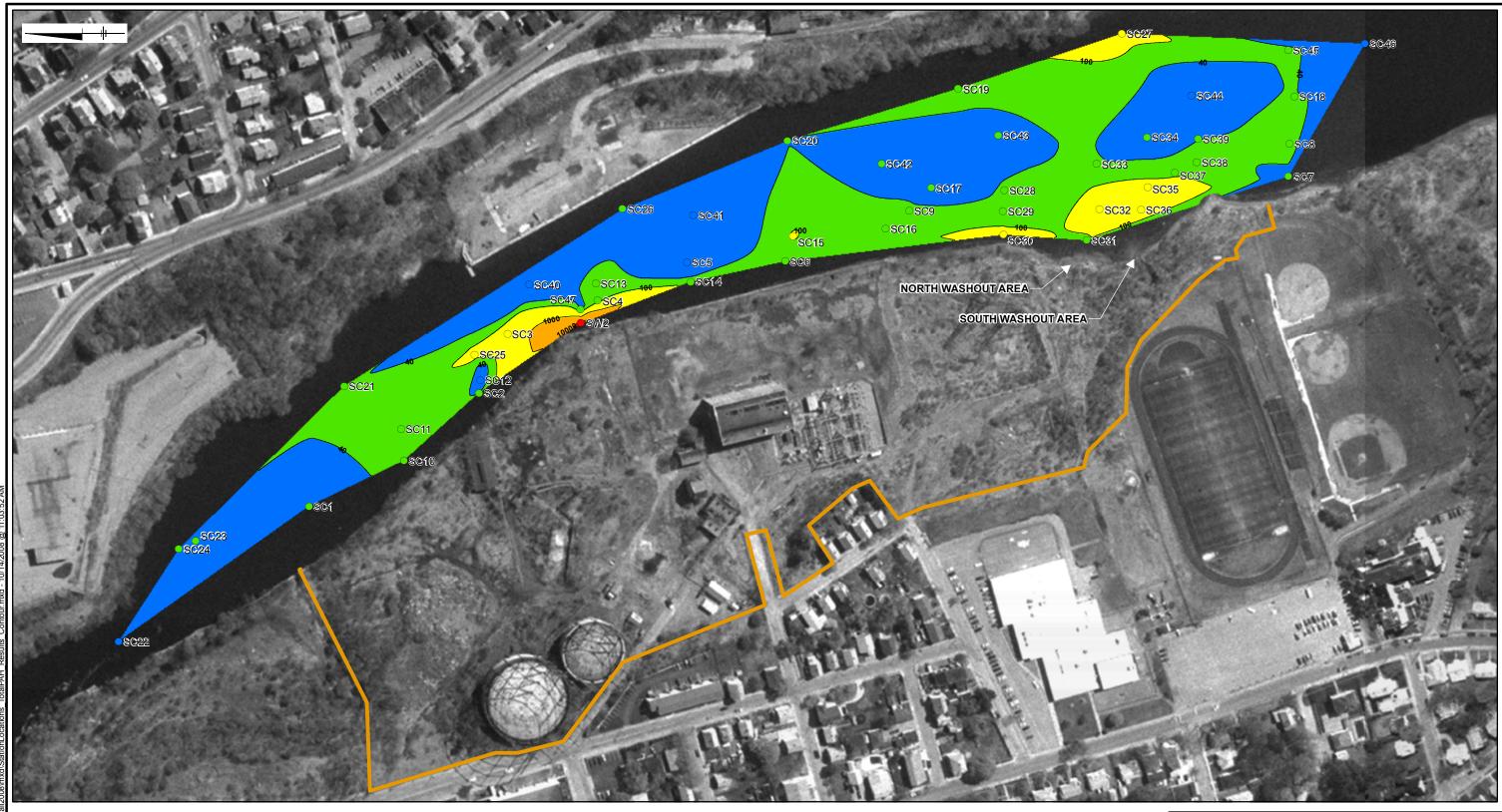


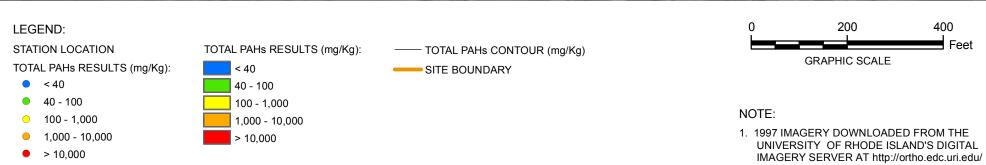
CITY: SYR DIV/GROUP: AIT DB: KEW JCR Tidewater (B0036697.0000.00001) Q:/Tidewaten/Pawtucket\basemapping/mxd/SiteMap.mxd - 6/29/2009 @ 12:58:35 PM



TY: SYR DIV/GROUP: AIT DB: KEW JCR PM: HD Jewater (B0036697.0000.00001)







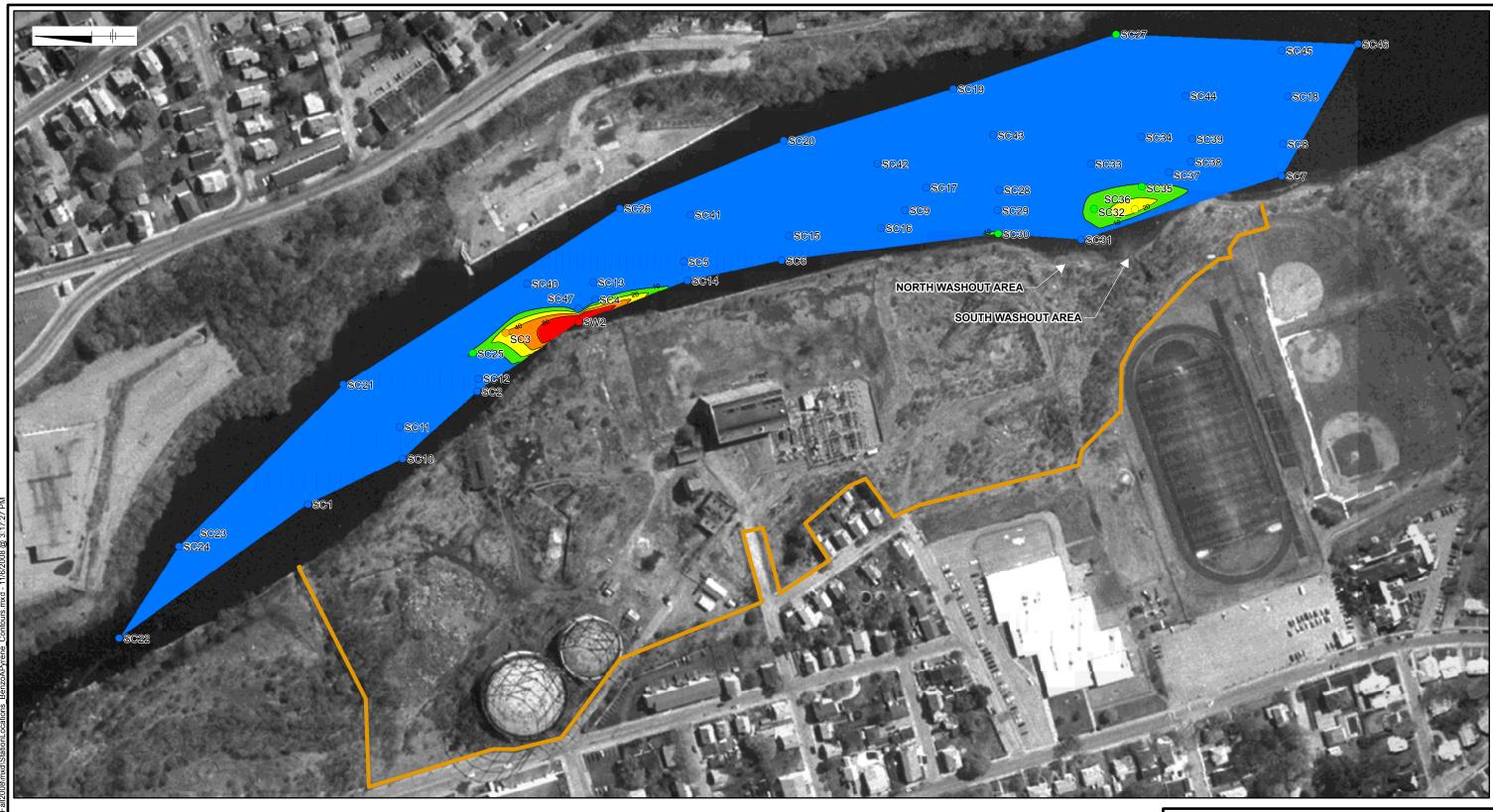
GR

∑ TO E



TIDEWATER PAWTUCKET, RHODE ISLAND

STATION LOCATIONS





LEGEND: 200 BENZO(A)PYRENE RESULTS (mg/Kg): STATION LOCATION ------ BENZO(A)PYRENE CONTOUR (mg/Kg) BENZO(A)PYRENE RESULTS (mg/Kg): < 10 SITE BOUNDARY GRAPHIC SCALE • < 10 10 - 20 10 - 20 20 - 40 20 - 40 0 40 - 80 NOTE: 9 40 - 80 > 80 • > 80

1. 1997 IMAGERY DOWNLOADED FROM THE UNIVERSITY OF RHODE ISLAND'S DIGITAL IMAGERY SERVER AT http://ortho.edc.uri.edu/

400

Feet

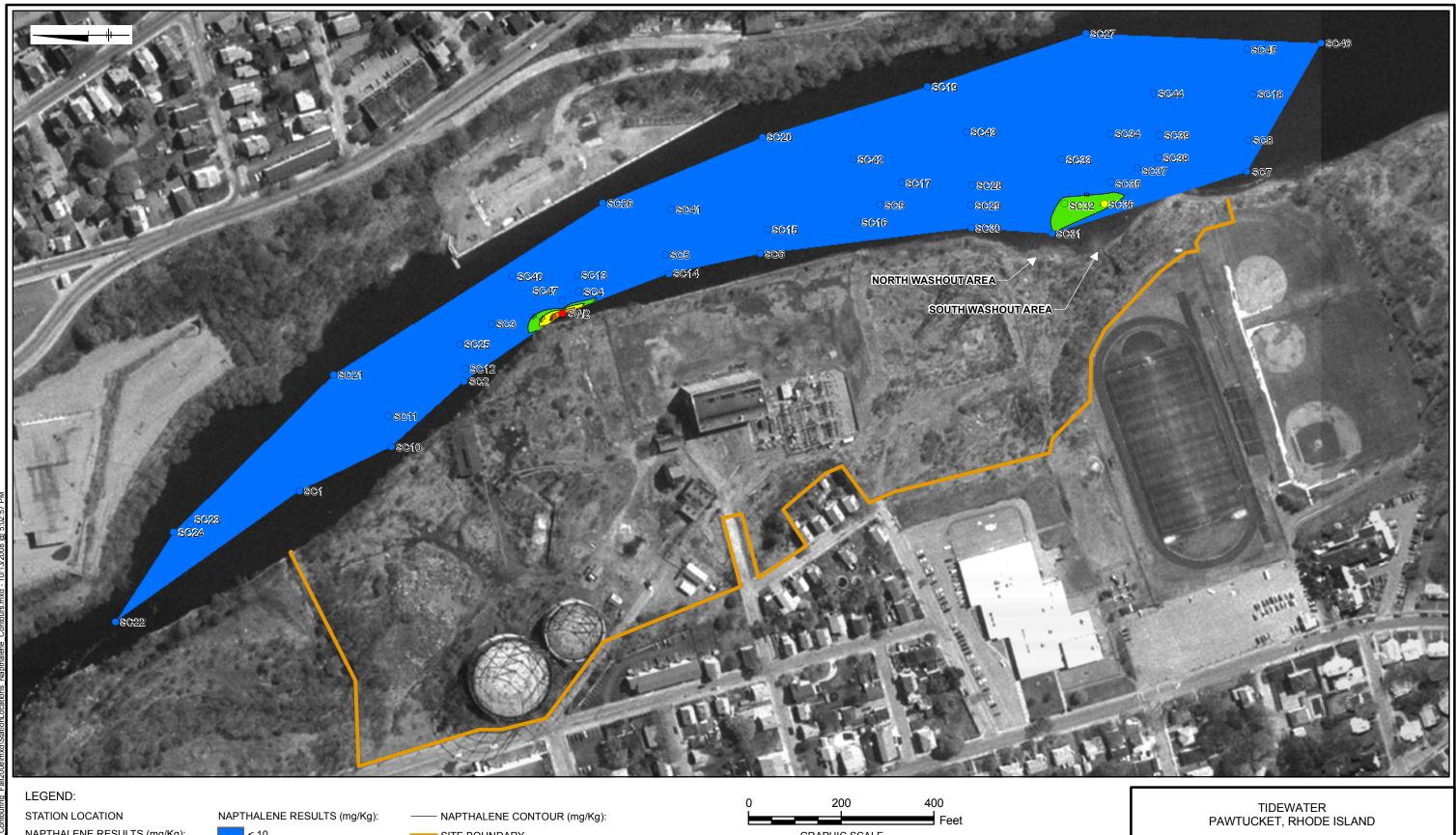
Idewater/Pawtucket/Res



FIGURE

STATION LOCATIONS BENZO(A)PYRENE DISTRIBUTION

TIDEWATER PAWTUCKET, RHODE ISLAND





<u>C</u>R

В

∑ TO E

NAPTHALENE RESULTS (mg/Kg): < 10</p> • 10 - 20 20 - 40 40 - 80 0 > 80

9 40 - 80 • > 80

< 10 10 - 20 20 - 40



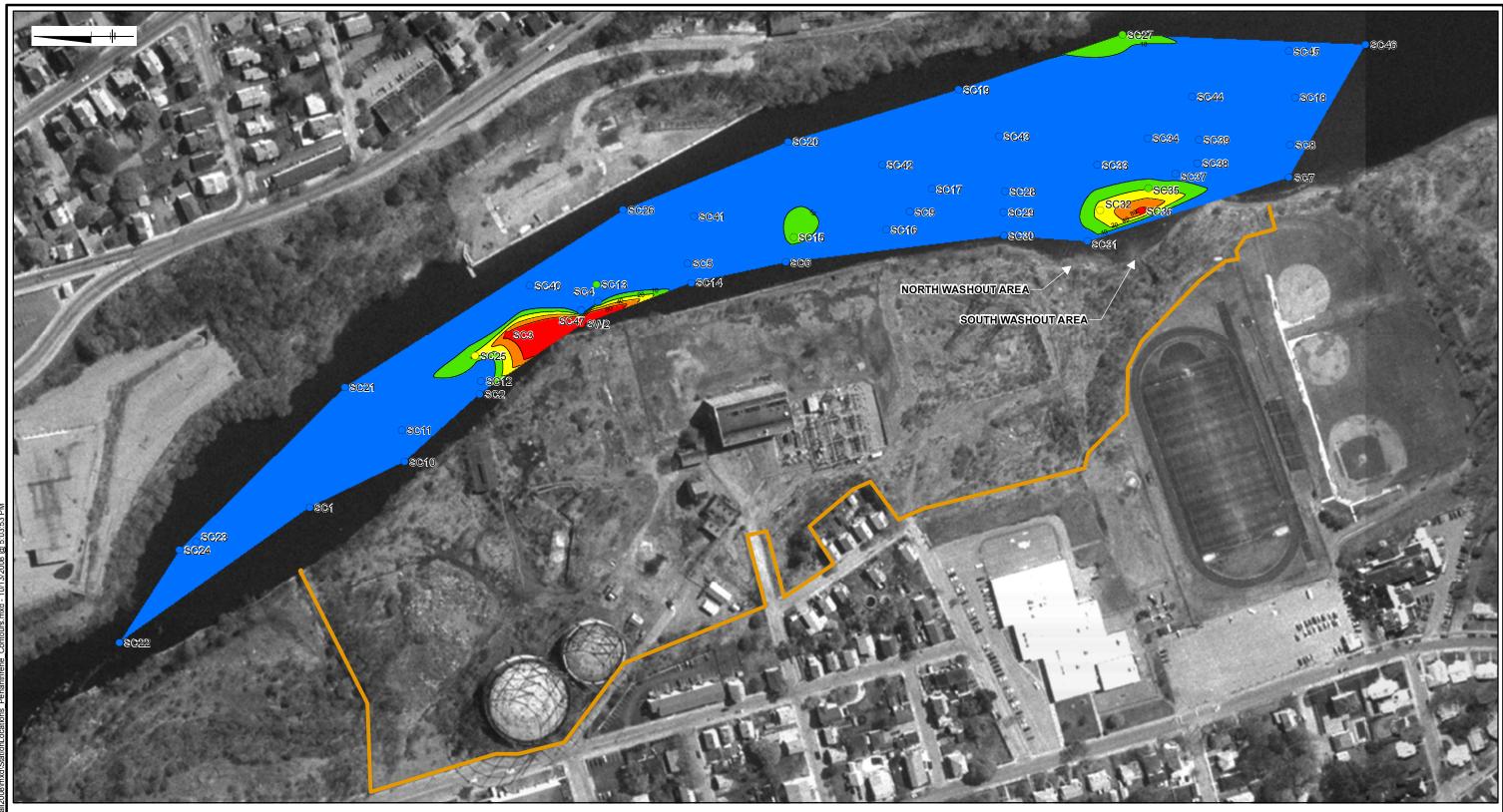
NOTE:

1. 1997 IMAGERY DOWNLOADED FROM THE UNIVERSITY OF RHODE ISLAND'S DIGITAL IMAGERY SERVER AT http://ortho.edc.uri.edu/

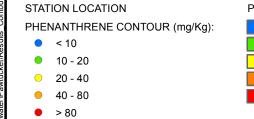
SITE BOUNDARY

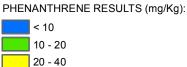
STATION LOCATIONS NAPTHALENE DISTRIBUTION











40 - 80

> 80

PHENANTHRENE CONTOUR (mg/Kg):
 SITE BOUNDARY



NOTE:

1. 1997 IMAGERY DOWNLOADED FROM THE UNIVERSITY OF RHODE ISLAND'S DIGITAL IMAGERY SERVER AT http://ortho.edc.uri.edu/



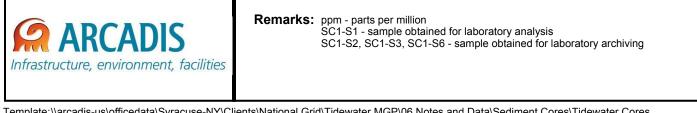
FIGURE

STATION LOCATIONS PHENANTHRENE DISTRIBUTION

TIDEWATER PAWTUCKET, RHODE ISLAND

APPENDIX A CORE LOGS

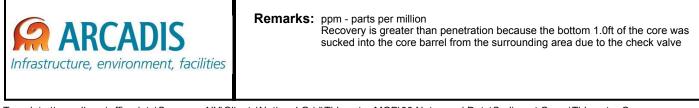
D D D	rilling riller's rilling	tart/Finish: Company s Name: Method: ation:	: T M V	I. Avak ibraco	larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.2189 N 71 22.9352 W 11.0 ft -9.35 ft 4.6 ft B. Thibault	Well/Boring ID: SC1 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
	_								
	- 10 -	SC1-S1 SC1-S2 SC1-S3 SC1-S6	15.2		coarse gravel, and b 3.0ft grain size coars be coarse rounded of	rick fragments; thin layer sens to 20-30% coarse ro	of dark gray silt at 0. bunded gravel; at base ravel and medium sar	and fine gravel; <10% silt, Oft; 1/2" coal at 0.4ft; starting at e of core sediment appears to nd; piece of concrete at 4.5ft; sweet odor.	
- 	- -15 -								
-	_								
- 10 - -	-20 -								- 1 _ _



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 Data File:SC1.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

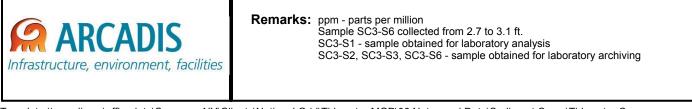
Di Di Di	rilling riller'៖ rilling	tart/Finish: Company s Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1603 N 71 22.8832 W 12.6 ft -11.69 ft 9.0 ft B. Thibault	Well/Boring ID: SC2 Client: National (Location: Former Tidewa Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraph	ic Description		Well/Boring Construction
5		SC2-S1 SC2-S2 SC2-S3 SC2-S12	8.0		glass fragments; 1" i	ND (SP); 20-30% coarse ece of coal at 3.0ft; piece ist; dark gray; no appare	e of black and white	gravel; <10% coarse gravel and pottery at 4.8ft; 2" piece of wood ning; musty odor.	
- 10 -	_								



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 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

D D D)rilling)riller's	art/Finish: Company s Name: Method: ation:	: T M V		Marine Services, Inc. tian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1505 N 71 22.8560 W 11.5 ft -10.30 ft 3.1 ft B. Thibault	Well/Boring ID: SC3 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- - - -	- 10 -	SC3-S1 SC3-S2 SC3-S3	1.0		bricks and glass frag fine material increase	ments; amount of coarse	e material decreases mately 1.0ft compose	rse sand; <10% fine gravel, with depth and percentage of ed of medium sand, <10% fine c grav to black	
5	- 15 -	SC3-S6			/	e and medium sand and	-	; black; medium plasticity; bluish	
- - - 10	-20 -								-



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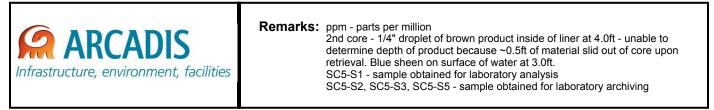
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 Date:August 11, 2008
 Project: 36697
 Page: 1 of 1

D D D)rilling)riller':)rilling	tart/Finish: յ Company s Name: յ Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1195 N 71 22.8408 W 9.6 ft -7.27 ft 3.8 ft B. Thibault	Well/Boring ID: SC4 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraph	ic Description		Well/Boring Construction
-	-5 -								-
	-10 -	SC4-S1 SC4-S2 SC4-S3	11.8		layer of dark grayish	brown silt at 0.0ft; gradu ine gravel at 0.8 to 1.3ft;	ally transitions to me	ravel; <10% medium sand; thin dium angular sand, 10-20% uish sheen on water in top of	0
- 5	-								-5
- 10	-15 - -								- 1
	-								



Remarks: ppm - parts per million SC4-S1 - sample obtained for laboratory analysis SC4-S2, SC4-S3 - sample obtained for laboratory archiving

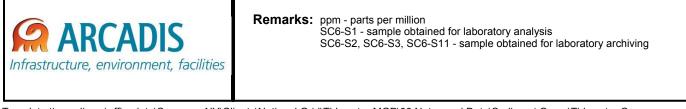
D D D	rilling riller'៖ rilling	art/Finish: Company s Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0888 N 71 22.8233 W 12.0 ft -6.89 ft 2.9 ft B. Thibault	Well/Boring ID: SC5 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	-5 -	SC5-S1	3.5	••••	Poorly graded mediu	ım angular SAND (SP); 1	10-20% coarse sand	; 4.5% gravel; moist; gray; no	
	-10 -	SC5-S2 SC5-S3 SC5-S5			Apparent oil/tar odor SILT (ML); 30% vege sand from 2.6 to 2.7 2.0ft, and 2.2ft; organ	etation; 10-20% fine grav 5ft; soft; moist; black; me	vel; <10% fine and m dium plasticity; spot	nedium sand; silt and medium of iridescent sheen at 1.9ft,	
- 5	_								-
 - 10	- 15 - - -								-



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 Date:August 11, 2008
 Project: 36697
 Page: 1 of 1

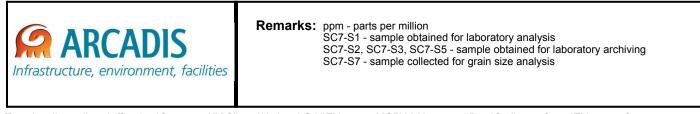
D D D	rilling riller'៖ rilling	tart/Finish: Company s Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0550 N 71 22.8228 W 7.4 ft -2.30 ft 6.1 ft B. Thibault	Well/Boring ID: SC6 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	0-								
	-5 -	SC6-S1 SC6-S2 SC6-S3	0.3		becomes more firm (soft) with depth; fine san ; layer of leaves at 6.0ft;	d layer at 3.9ft, 4.1ft,	s; very soft at surface and 4.4 to 4.5ft, 4.9 to 5.1ft, 5.4 to plasticity; spotty small sheen	
5 - - -		SC6-S11							-5
- 10	-								- 1



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 Data File:SC6.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller's rilling	tart/Finish: ı Company s Name: ı Method: ation:	: T J. V	G&B N	, 2008 Marine Services, Inc. Ion and J. Balmer re	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.8819 N 71 22.7842 W 1.0 ft 0.04 ft 3.7 ft B. Thibault	Well/Boring ID: SC7 Client: National G Location: Former Tidewate Pawtucket, RI	-
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- 0		SC7-S1 SC7-S2 SC7-S3 SC7-S5 SC7-S5 SC7-S7	0.2		soft; black to dark gr SILTY SAND (SM); 2 to gray with depth; h	ay; nonplastic to low plas 20-30% silt; medium san eavy iridescent sheen fro	ticity; <10% bluish d layer from 2.7 to 2 m 2.3 to 2.65ft; no	naterial; moist to wet; very soft to spotty sheens; organic odor. 2.9ft; moist; dark gray transitioning apparent oil/tar odor. and light gray; no apparent oil/tar	
5 - - - - - - - - - 10-	-5 - - - - - - - - - - - - - - - - - - -								5



10

Di Di Di	rilling riller's	art/Finish: Company Name: Method: ation:	: T J. V		Narine Services, Inc. Ion and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.8815 N 71 22.7693 W 3.2 ft -2.08 ft 3.8 ft B. Thibault	Well/Boring ID: SC8 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
		SC8-S1 SC8-S2 SC8-S3 SC8-S4	1.5		depth and is medium nonplastic to low plas throughout. SILT (ML); 10-20% fi 20% spotty bluish sh Medium SAND (SP);	n sand with 30-40% fine s sticity; heavy iridescent s ine to medium sand; 1" c ieen; organic odor. ; 20-30% coarse sand an 2.93ft and 3.0 to 3.1ft; m	and by the base of heen from 0.8 to 0.8 linker at 1.2ft; moist d fine gravel; <10%	ount of sand increases with the unit; wet; soft; black; 85ft; 10-20% spotty bluish sheen ;; soft; black; low plasticity; 10- fine sand and coarse gravel; silt ange staining at 3.5ft; no	



Remarks: ppm - parts per million SC8-S1 - sample obtained for laboratory analysis SC8-S2, SC8-S3, SC8-S4 - sample obtained for laboratory archiving 10

Di Di Di	rilling riller'៖ rilling	tart/Finish: Company s Name: Method: ation:	: T M V		Narine Services, Inc.	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0123 N 71 22.7997 W 5.5 ft -4.17 ft 4.8 ft B. Thibault	Well/Boring ID: SC9 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- - - -	-5-	SC9-S1 SC9-S2 SC9-S3	2.5		sand layer at 1.1ft, 1. 2.5ft, 3.8ft, and 4.0ft; 0.9 to 4.8ft; soft; blac	3 to 1.55ft, 2.25 to 2.35f 30-40% organic materia k; nonplastic to low plas at 0.80 to 0.90ft and 2.8	t, 3.05 to 3.3ft, and 4 Il from 3.0 to 4.3ft; we ticity; pin-head sized	naterial; 0.3% gravel; medium 5 to 4.65ft; leaf layer at 1.6ft, et from 0.0 to 0.9ft; moist from droplets of brown product on from 1.9 to 2.3ft; organic odor,	
- 5 		SC9-S8							Ę
- 10 	- - - -								-



Remarks: ppm - parts per million SC9-S1 - sample obtained for laboratory analysis SC9-S2, SC9-S3, SC9-S8 - sample obtained for laboratory archiving

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 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D)rilling)riller':)rilling	tart/Finish: յ Company s Name: յ Method: ation:	: T N V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1861 N 71 22.9142 W 17.5 ft 12.53 ft 2.8 ft B. Thibault	Well/Boring ID: SC10 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	-10 -								
	- - -15 -	SC21-S1 SC21-S2 SC21-S3	0.6		Poorly graded mediur gravel; <10% coal, gla apparent oil/tar odor o	ass, and brick fragments	0-20% coarse sand; s; thin layer of dark gr	4.5% silt; 5.7% fine and coarse ay silt at 0.0ft; moist; gray; no	-
	-								
- - -	- 20 - -								-
- 10	-								-

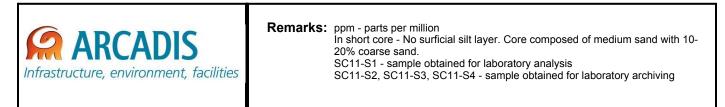


Remarks: ppm - parts per million SC10-S1 - sample obtained for laboratory analysis SC10-S2, SC10-S3 - sample obtained for laboratory archiving

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 Date:August 11, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller's	art/Finish: Company s Name: Method: ation:	: T M V	I. Avak ibraco	Aarine Services, Inc.	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1871 N 71 22.8996 W 9.4 ft -8.46 ft 5.4 ft B. Thibault	Well/Boring ID: SC11 Client: National G Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	_								-
	-10	SC11-S1 SC11-S2 SC11-S3 SC11-S4	23.3		water above core; of Medium angular SA glass fragments; sec	rganic odor. ND (SP); 20-30% coarse diment coarsens with dep 30% fine gravel and medi	sand; 10-20% fine g	10% spotty iridescent sheen in ravel; <10% coarse gravel and e core is a layer of coarse parse sand; moist; gray; no	
-	-15 - -								-
- 10	-20 -								

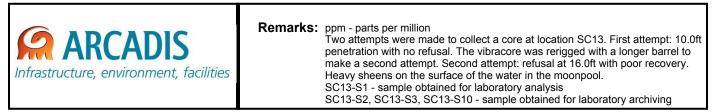


D D D	rilling riller'៖ rilling	tart/Finish: Company s Name: Method: ation:	: Т М V		larine Services, Inc. tian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1598 N 71 22.8774 W 14.2 ft -9.81 ft 2.4 ft B. Thibault	Well/Boring ID: SC12 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- -	-10 -	SC21-S1	0.9						C
-		SC21-S1 SC21-S2 SC21-S3	0.9		coarse gravel, glass a	and brick fragments; 0.8	% silt; piece of hard p	16% gravel; <10% fine sand, blastic at 2.3 ft; clinker at 2.1 ft; t oil/tar odor or staining.	
- - 5	-								-5
-	- 15 -								-
-	_								
- 10_ -	-20 -								



Remarks: ppm - parts per million SC12-S1 - sample obtained for laboratory analysis SC12-S2, SC12-S3 - sample obtained for laboratory archiving

Di Di Di	rilling riller'៖ rilling	tart/Finish: Company s Name: Method: ation:	: T N V		larine Services, Inc. tian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1201 N 71 22.8328 W 12.0 ft -10.40 ft 5.4 ft B. Thibault	Well/Boring ID: SC13 Client: National G Location: Former Tidewat Pawtucket, RI	-
DЕРТН	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
_	-10 -								-
0		SC13-S1 SC13-S2 SC13-S3	48.6		gravel at 0.0ft; 3" piec		; 2" piece of black p	r of silt, vegetation,and fine lastic at 3.9ft; moist; black; r.	-
- 5	-15 -	SC13-S10			visible product; faint p Catcher (9.9-10.0ft):	petroleum-like odor.	nd and organic mate	k; low plasticity; no sheens or erial; moist; soft; black; low	-
- 10	- 20 -				sand, glass, and coal 5.1-5.9ft: Black SILT 5.9-6.4ft: Medium any fine sand; moist; blac odor. Catcher (15.9-16.0ft) moist; soft; silver-blac	ngular SAND; 20-30% cd ; piece of 1/2" rubber tul gular SAND; 20-30% coa k; brown product coating : Fine SILTY SAND; 20-	bing at 4.5ft; moist; b arse sand and fine g g sediment from 6.3 30% fine rounded gr covered with silver s	gravel; <10% coarse gravel, fine black; faint petroleum-like odor. ravel; <10% coarse gravel and to 6.4ft; strong petroleum-like avel; 10-20% medium sand; sheen; brown product smears on	-



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 Data File:SC13.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

Drillin Driller Drillin	Start/Finish: g Company 's Name: g Method: ration:	: TO M. Vil		larine Services, Inc. ian and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1187 N 71 22.8320 W 12.5 ft -11.28 ft 11.5 ft B. Thibault	Well/Boring ID: SC13A Client: National Gr Location: Former Tidewater Pawtucket, RI	
DEPTH ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-10 - 					metal end of light bulb a		nd and fine gravel; thin layer of 2" spot of iridescent sheen at	-
- <i>20</i> - - 10 -				or sheen; faint musty SILT (ML); <10% fine moderate petroleum- Medium SAND (SP); 2" pieces of gravel at from 8.6 to 8.8ft; pin- SILT (ML); 20-30% fil of coal at 9.0ft; moist; Medium grained SILT visible oil/tar on sedir Catcher: Medium gra	petroleum-like odor. to medium sand; moist like odor. 20-30% coarse sand; < 8.1 to 8.3ft; moist; black head size smears of bro ne to medium sand and soft; black; low plasticit Y SAND (SM); moist; si nent; strong petroleum-l	soft; black; low plas 10% fine gravel and ; <10% spotty irides wn product on glove organic material; <1 y; no apparent stain ver-black; sediment ke odor. moist; silver-black;	sticity; no apparent oil/tar staining sticity; no apparent sheen; coal and glass fragments; 1" and scent sheen; heavy silver sheen s; moderate petroleum-like odor. 0% rounded coarse sand; piece ing or sheen; organic odor. t covered with silver sheen; no sediment covered with silver tar-like odor.	-

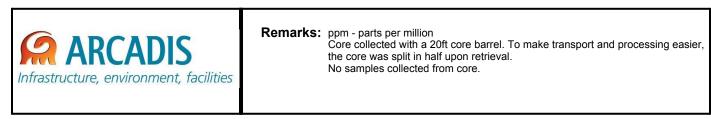


Remarks: ppm - parts per million Core collected with a 20ft core barrel. To make transport and processing easier, the core was split in half upon retrieval. No samples collected from core.

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 Data File:SC13A.dat
 Date:August 21, 2008
 Project: 36697
 Page: 1 of 2

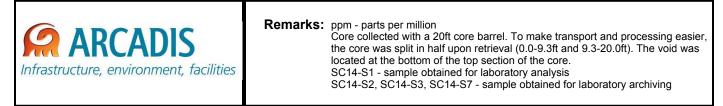
Drill Drill Drill	ling C ler's l	rt/Finish: Company Name: Aethod: ion:	: T N V	G&B I	, 2008 Marine Services, Inc. kian and J. Balmer ore	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1187 N 71 22.8320 W 12.5 ft -11.28 ft 11.5 ft B. Thibault	Well/Boring ID: SC13A Client: National (Location: Former Tidewa Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- 25 - 15 -	5								- 15
- 30 - 20 - -	-								- 20



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 Data File:SC13A.dat
 Date:August 21, 2008
 Project: 36697
 Page: 2 of 2

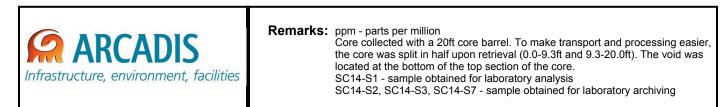
	Drilling Driller' Drilling	tart/Finish: y Company s Name: y Method: ation:	: T M V		Narine Services, Inc.	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0875 N 71 22.8321 W 5.5 ft -3.43 ft 13.3 ft B. Thibault	Well/Boring ID: SC14 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	-								-
- 0 - - - - 5	-5	SC14-S1 SC14-S2 SC14-S3 SC14-S3	85.8		4.7ft; wet from 0.0 to	1.2ft, moist from 1.2 to 6	6.4ft; very soft from 0	terial; fine sand layer from 4.5 to .0 to 1.2ft, soft from 1.2 to 6.4ft; /8" iridescent sheen at 2.9ft and	-5
-	-10 -				Void SILT (ML); <10% find staining.	e to medium sand; moist	soft; black; low plas	ticity; no apparent oil/tar odor or	-
- 10	-15 -					ND (SP); <10% coarse sa	and and fine gravel; r	noist; gray; no apparent oil/tar	- 1



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 Data File:SC14.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 2

D D D	rilling riller' rilling	tart/Finish: g Company s Name: g Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0875 N 71 22.8321 W 5.5 ft -3.43 ft 13.3 ft B. Thibault	Well/Boring ID: SC14 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- - 15	-	SC14-S32			Medium angular SAN odor or staining.	ID (SP); <10% coarse sa	and and fine gravel; r	noist; gray; no apparent oil/tar	- 15
-	-20 -								-
- 20	- -25 - -								■ - 20 - - -



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 Data File:SC14.dat
 Date:August 20, 2008
 Project: 36697
 Page: 2 of 2

D D D	rilling riller's	art/Finish: Company s Name: Method: ation:	: T M V		Marine Services, Inc. kian and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0523 N 71 22.8113 W 8.2 ft -4.52 ft 14.0 ft B. Thibault	Well/Boring ID: SC15 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
0	-5	SC15-S1 SC15-S2	3.1		and 2.7 to 3.1ft, remain	ainder is moist; very soft	from 0.0 to 0.5ft, soft	l; wet (soupy); from 0.0 to 0.5ft t from 0.5 to 3.0ft; black;	- - -
5	-	SC15-S3 SC15-S4			Medium SAND (SP); SILT (ML); 10-20% f plasticity; no apparei	ent oil/tar odor or staining ; 30-40% fine sand; mois ine to medium sand; 20-3 nt oil/tar staining; chemic	t; gray; no apparent o		-
	- 10 - - -				plasticity; no apparen SILT (ML); 10-20% f 7.0 to 9.0ft, moist fro	ine to medium sand; 20-; nt oil/tar staining; chemic ine to medium sand; 20-;	al-like odor. 30% organic material (soupy) from 7.0 to 9	l; <10% coarse gravel; wet from 9.0ft, soft from 9.0 to 9.85ft;	-
10	-15 - -					<10% spotty iridescent sl		ivel; grain size coarsens with smears of brown product on	-



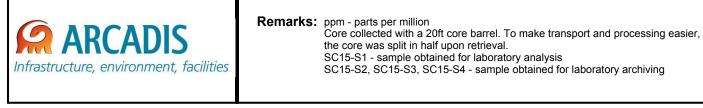
Remarks: ppm - parts per million Core collected with a 20ft core barrel. To make transport and processing easier, the core was split in half upon retrieval.

SC15-S1 - sample obtained for laboratory analysis

SC15-S2, SC15-S3, SC15-S4 - sample obtained for laboratory archiving

Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores
Page: 1 of 2
Page: Date:August 20, 2008 Project: 36697

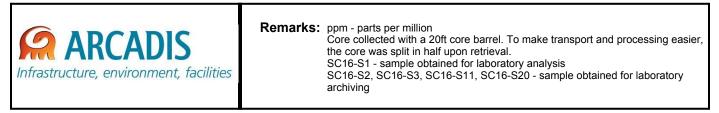
)rilling)riller's	tart/Finish: J Company s Name: J Method: ation:	: T M V		larine Services, Inc. ian and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0523 N 71 22.8113 W 8.2 ft -4.52 ft 14.0 ft B. Thibault	Well/Boring ID: SC15 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- - - 20	-20 -				and coarse gravel; 3"	by 2" brick fragment at	14.0ft; moist; no appa	arent oil/tar odor or staining.	- 15



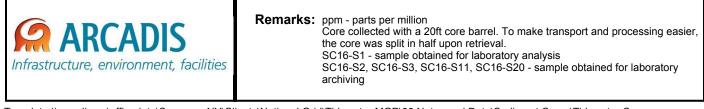
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 Data File:SC15.dat
 Date:August 20, 2008
 Project: 36697
 Page: 2 of 2

Dr Dr Dr	rilling riller's rilling	tart/Finish: յ Company s Name: յ Method: ation:	: T M V	I. Avak ibracoi	larine Services, Inc. ian and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0205 N 71 22.8079 W 3.7 ft -2.69 ft 12.9 ft B. Thibault	Well/Boring ID: SC16 Client: National G Location: Former Tidewat Pawtucket, RI	-
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraph	ic Description		Well/Boring Construction
-	0 -								-
- 0	-5 -	SC16-S1 SC16-S2 SC16-S3	0.5		sand layer at 3.1ft; 4 soft from 0.0 to 1.0ft,	.2 to 4.5ft, and 4.97ft; we	et from 0.0 to 3.75ft, n ack; nonplastic to low	terial; leaf layer at 5.0ft; medium noist from 3.75 to 5.6ft; very plasticity; heavy silver sheen sty odor at 5.6ft.	
- 5	-	SC16-S11			6.8ft; soft; moist; blac	ck; low plasticity; 10-20%	5 1/4" iridescent shee	all piece of rounded gravel at n; pin-head size smears of nt; moderate musty petroleum-	5
- - 10	- 10 -	SC16-S20			Medium SAND (SP);	oble from 10.6 to 10.9ft;	pravel and coarse sar gray transitoning to y	nd; 30-40% fine sand from 11.25 ellow-brown at 12.0ft; moist; no	- 1
	- 15 -					-			



Dri Dri Dri	illing iller's illing	art/Finish: Company Name: Method: tion:	: T N V	G&B I 1. Aval ïbracc	, 2008 Marine Services, Inc. kian and J. Balmer ore (refusal)	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0205 N 71 22.8079 W 3.7 ft -2.69 ft 12.9 ft B. Thibault	Well/Boring ID: SC16 Client: National (Location: Former Tidewa Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- 15 	- - - 20 -								- 15
- 20	-								- 20
	25 -								



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 Data File:SC16.dat
 Date:August 21, 2008
 Project: 36697
 Page: 2 of 2

Dril Dril Dril	lling Iler's Iling	art/Finish: Company Name: Method: ntion:	: T M V		Marine Services, Inc. kian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0048 N 71 22.7893 W 6.9 ft -5.02 ft 5.0 ft B. Thibault	Well/Boring ID: SC17 Client: National G Location: Former Tidewat Pawtucket, RI	
DЕРТН	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- - - - -		SC17-S1 SC17-S2 SC17-S3 SC17-S9	4.2		0.2ft, 1.1 to 1.2ft, 1.4	to 1.55ft, 2.1 to 2.2ft, 3.4	45 to 3.65ft, 4.25 to 4	erial; medium sand layer at 0.0 to 4.4ft, and 4.8 to 4.9ft; moist; ty sheens throughout; organic	
-5 -1 - - - - - - 10-1	-								

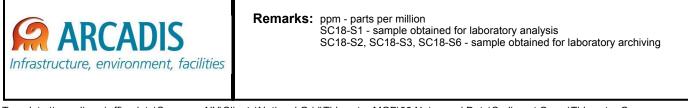


Remarks: ppm - parts per million SC17-S1 - sample obtained for laboratory analysis SC17-S2, SC17-S3, SC17-S9 - sample obtained for laboratory archiving

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 Data File:SC17.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller': rilling	tart/Finish: Company s Name: Method: ation:	: T J. V		Narine Services, Inc. Ion and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.8798 N 71 22.7474 W 10.0 ft -8.70 ft 5.5 ft B. Thibault	Well/Boring ID: SC18 Client: National G Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	_								
	- 10 -	SC18-S1 SC18-S2 SC18-S3 SC18-S6	71.4		layer at 0.35ft, 1.7ft, 2 0.6 to 5.5ft; very soft	.2% gravel; 21.9% sand 2.15 to 2.3ft, and 4.75ft; to approximately 2.0ft an at oil/tar odor or staining.	leaf layer at 0.8ft; we nd then firms to soft;	terial; medium grained sand et from 0.0 to 0.6ft, moist from black; nonplastic to low	
	_								-5
- - -	- 15 - - -								
10 	-20 -								-



D D D	Date Start/Finish:July 10, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:M. Avakian and J. ScanlonDrilling Method:VibracorePenetration:5.0 ft					Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9955 N 71 22.7437 W 11.8 ft -10.83 ft 4.1 ft B. Thibault	Well/Boring ID: SC19 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
	- - - 10 -								
- 0 - -		SC19-S1 SC19-S2 SC19-S3	1.9		gravel at 4.1ft; moist).6% gravel; 36.3% sand ; soft; black; low plasticity duct at 0.7ft on water in c	; small spotty bluish	aterial; layer of fine rounded n sheen from 0.0 to 4.1ft; 1/8"	
 5	- 15 -	SC19-S8		<u>·</u> - <u>·</u> - · -					-5
	- -20 -								
- 10 -	_								-1



Remarks: ppm - parts per million SC19-S1 - sample obtained for laboratory analysis SC19-S2, SC19-S3, SC19-S8 - sample obtained for laboratory archiving

 Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores

 Data File:SC19.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

Date Start/Finish:July 9, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:M. Avakian and J. ScanlonDrilling Method:VibracorePenetration:4.0 ft (refusal)					larine Services, Inc. ian and J. Scanlon re	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0541 N 71 22.7674 W 17.9 ft -13.22 ft 2.3 ft B. Thibault	Well/Boring ID: SC20 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	-								-
-0	- 15 -	SC20-S1 SC20-S2 SC20-S3	1.8		with no distinct layers	; at 2.3 ft the sediment h	has graded to fine gra	coarser gradation with depth avel, 10-20% coarse gravel and pparent oil/tar odor or staining.	-
-	_								-
-5 	- - 20 -								-
	-								-
- 10	- 25 -								



Remarks: ppm - parts per million SC20-S1 - sample obtained for laboratory analysis SC20-S2, SC20-S3 - sample obtained for laboratory archiving

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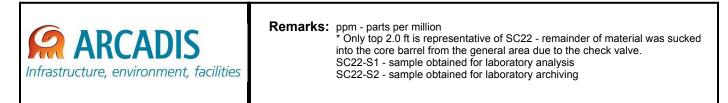
 Data File:SC20.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	Date Start/Finish:July 8, 2008Drilling Company:TG&B Marine Services,Driller's Name:M. Avakian and J. ScarDrilling Method:VibracorePenetration:5.5 ft					Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.2067 N 71 22.8799 W 5.5 ft -1.11 ft 5.4 ft B. Thibault	Well/Boring ID: SC21 Client: National C Location: Former Tidewat Pawtucket, RI			
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction		
- - - - - -	0- - - -5-	SC21-S1 SC21-S2 SC21-S3 SC21-S5	NA		iridescent sheen; org SILT (ML); soft; mois	NDY SILT (ML); 47.4% sand; very soft; wet; very dark grayish brown; nonplastic; spotty lescent sheen; organic odor. .T (ML); soft; moist; has a blocky texture when split; black to dark gray; medium plasticity; otty bluish sheen from 1.7 to 4.4ft (~ dozen); leather-like odor.					
5 - - 	- 10								-5		



Remarks: ppm - parts per million SC21-S1 - sample obtained for laboratory analysis SC21-S2, SC21-S3, SC21-S5 - sample obtained for laboratory archiving 10

D D D	rilling riller'៖ rilling	art/Finish: Company Name: Method: ation:	: T M V		Narine Services, Inc.	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.2843 N 71 22.9971 W 17.3 ft -13.83 ft 4.3 ft* B. Thibault	Well/Boring ID: SC22 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-									
-0	- 15 -	SC22-S1 SC22-S2	2.5		Poorly graded mediu layer of gray silt at 0. apparent oil/tar odor	Oft; coarser gradation wi	.2% gravel; 10-20% th depth - no distinct	coarse sand; 1.0% silt; thin layers; moist; brownish gray; no	
- 5	- 20 -				Fine to coarse GRAV	/EL (GW); rounded to su	bangular; gray; no ap	oparent oil/tar odor or staining.	
-	-								-
- 10 	-25 -								



D D D	rilling riller'៖ rilling	tart/Finish: Company s Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.2578 N 71 22.9509 W 8.4 ft -5.88 ft 4.7 ft B. Thibault	Well/Boring ID: SC23 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	-5-								
- 0	-	SC23-S1 SC23-S2 SC23-S3	0.9		from 3.4 to 4.0ft; laye 4.7ft; black to very da 0.0 to 2.6ft, layer of in	r of vegetation (organics ark gray; medium plastici	s) at 4.7ft; wet from 0 ty (not in sand layer) t 3.4ft; small spots of	leaf litter; layer of fine sand .0 to 0.6ft, moist from 0.6 to ; spotty iridescent sheen from f brown product on gloves from	-
- 5	- 10 -	SC23-S9							
- 10	- - - 15 -								-
- 1U	_								



Remarks: ppm - parts per million SC23-S1 - sample obtained for laboratory analysis SC23-S2, SC23-S3, SC23-S9 - sample obtained for laboratory archiving

	Drilling Driller's	tart/Finish: J Company s Name: J Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.2636 N 71 22.9546 W 10.7 ft -7.23 ft 2.7 ft B. Thibault	Well/Boring ID: SC24 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	-5 -								
- 0 - - -	-10 -	SC24-S1 SC24-S2 SC24-S3	1.1		Poorly graded mediu gravel and fine sand; apparent oil/tar odor o	2.1% silt; thin layer of d	SP); 5.8% gravel; 10- ark gray silt at 0.0 ft; ı	20% coarse sand; <10% fine moist; brownish gray; no	
- 5	_								-5
-	- 15 - -								-
- 10	-								



Remarks: ppm - parts per million SC24-S1 - sample obtained for laboratory analysis SC24-S2, SC24-S3 - sample obtained for laboratory archiving

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 Data File:SC24.dat
 Date:August 11, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller': rilling	tart/Finish: J Company s Name: J Method: ation:	: T N V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1618 N 71 22.8656 W 12.8 ft -8.90 ft 1.7 ft B. Thibault	Well/Boring ID: SC25 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	-								
	-10 -	SC25-S1 SC25-S2 SC25-S3	15.3		Poorly graded mediu sand; 2.4% gravel; 6. oil/tar odor or staining	.6% fines; thin layer of ve	i-10% coarse sand ar ery dark brown silt at (nd brick fragments; 10-20% fine 0.0 ft; moist; gray; no apparent	
- - 5	-	_							- - - 5
-	-15 -								-
- 10	_								
	-20 -								



Remarks: ppm - parts per million SC25-S1 - sample obtained for laboratory analysis SC25-S2, SC25-S3 - sample obtained for laboratory archiving

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 Data File:SC25.dat
 Date:August 11, 2008
 Project: 36697
 Page: 1 of 1

D D D)rilling)riller':)rilling	tart/Finish: յ Company s Name: յ Method: ation:	: T N V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1109 N 71 22.7987 W 13.8 ft -12.86 ft 5.6 ft B. Thibault	Well/Boring ID: SC26 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	-10 -								
- 0 - - - -	-15 -	SC26-S1 SC26-S2 SC26-S3	4.8		SILT (ML); <10% fine 3.7ft; medium grained from 2.5 to 5.6ft; very	0.0ft; moist; gray; no app e and medium sand; 10-2 d sand layers at 2.3ft, 3.3	arent oil/tar odor or st 20% organic material; 2 to 3.35ft, and 4.55ft	0.2% gravel; 2.3% fines; thin aining. leaf layer at 3.0ft, 3.35ft, and ; wet from 0.65 to 2.5ft; moist ck; nonplastic to low plasticity;	
- 5 -	-	SC26-S9							
	-20 - - -								-
- 10	-								



Remarks: ppm - parts per million SC26-S1 - sample obtained for laboratory analysis SC26-S2, SC26-S3, SC26-S9 - sample obtained for laboratory archiving

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 Data File:SC26.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller's	cart/Finish: Company s Name: Method: ation:	: T J. V	. Scan ibraco	Narine Services, Inc. Ion and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9391 N 71 22.7185 W 6.6 ft -3.71 ft 1.35 ft B. Thibault	Well/Boring ID: SC27 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	-	SC27-S1	0.4		SANDY SILT (ML); 1	.9% gravel; 37.6% fine t	o medium sand; <109	% organic material; layer with	(
-	-5 -	SC27-S2		<u> </u>	medium sand, 20-30	% coarse sand, 10-20% gray banding from 1.25 t	fine gravel and fine s	0.15ft layer of brown-gray and at 1.35ft; moist; soft; black faint spotty bluish sheen	
- 5 	-10 -								Ę
- 10	- - -15 -								-



Remarks: ppm - parts per million SC27-S1 - sample obtained for laboratory analysis SC27-S2 - sample obtained for laboratory archiving

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 Data File:SC27.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller': rilling	tart/Finish: Company s Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9796 N 71 22.7904 W 7.8 ft -3.73 ft 5.0 ft B. Thibault	Well/Boring ID: SC28 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	-								-
 - -	-5-	SC28-S1 SC28-S2 SC28-S3	1.9		Poorly graded SAND nonplastic to low plas	(SP); 0.7% gravel; 4.1% ticity; small spotty bluish	5 fines; 10-20% organ n sheen throughout.	nic material; moist; soft; black;	-
- 5	- 10 -	SC28-S9							-
-	-								-
- 10 	- 15 -								-



Remarks: ppm - parts per million SC28-S1 - sample obtained for laboratory analysis SC28-S2, SC28-S3, SC28-S9 - sample obtained for laboratory archiving

D D D	rilling riller's rilling	art/Finish: Company s Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.9802 N 71 22.7999 W 3.1 ft -0.88 5.3 ft B. Thibault	Well/Boring ID: SC29 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- - - - - -	-5 -	SC29-S1 SC29-S2 SC29-S3	0.7		4.25ft, and 5.2ft; ver 1.0ft, black from 1.0 core; <10% pin-head	y soft from 0.0 to 1.0ft; so to 5.3ft; nonplastic to low	oft from 1.0 to 5.3ft; n plasticity; bluish sm t on gloves from sec	n sand layer at 2.4 to 2.6ft, noist; grayish brown from 0.0 to all spotty sheens throughout liment thoughout core; no visible 5.3ft.	0
	- 10 -								- 1



Remarks: ppm - parts per million SC29-S1 - sample obtained for laboratory analysis SC29-S2, SC29-S3 - sample obtained for laboratory archiving

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 Data File:SC29.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

Di Di Di	rilling riller'៖ rilling	art/Finish: Company s Name: Method: ation:	: T N V	I. Avak ibraco	Marine Services, Inc.	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9800 N 71 22.8107 W 3.4 ft 1.20 ft 2.2 ft B. Thibault	Well/Boring ID: SC30 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
	-								_
0	0-	SC30-S1 SC30-S2 SC30-S3 SC30-S4	0.9		bluish spotty sheen the Fine SILTY SAND (S	hroughout; slight organic	: odor.	ack; low plasticity; small (<1/8") 1.55 to 1.7ft); moist; black; light organic odor.	(
- 5	_								
5	-5 - -								{
10	_								-
	-10 -								_



Remarks: ppm - parts per million SC30-S1 - sample obtained for laboratory analysis SC30-S2, SC30-S3, SC30-S4 - sample obtained for laboratory archiving

Di Di Di	rilling riller's	art/Finish: Company S Name: Method: ation:	: T J. V	. Scan ibraco	<i>l</i> larine Services, Inc. Ion and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9513 N 71 22.8134 W 1.0 ft 1.09 ft 2.75 ft B. Thibault	Well/Boring ID: SC31 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-									-
-0- -		SC31-S1 SC31-S2 SC31-S3 SC31-S4	153		small gravel from 0.9 0.4ft, and black from petroleum-like odor. SILT (ML); 10-20% fil	to 1.0ft and 1.25 to 1.4f 0.4 to 1.4ft; small spotty ne to medium sand; 1/2"	t; moist; brown from (bluish sheen at 0.7ft coal at 2.1ft; 20-30%	6 sand from 2.0 to 2.1ft; moist;	-
	-			••••	<u> </u>	city; bluish spotty sheen; ; 30-40% fine sand; mois		/	
-5	-5 -								-5
	-								-
- 10	_								-
· _	-10 -								

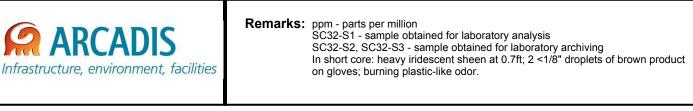


Remarks: ppm - parts per million SC31-S1 - sample obtained for laboratory analysis SC31-S2, SC31-S3, SC31-S4 - sample obtained for laboratory archiving

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 Data File:SC31.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

Di Di Di	rilling riller's	art/Finish: Company Name: Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon re	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9469 N 71 22.7992 W 1.9 ft -0.21 ft 4.0 ft B. Thibault	Well/Boring ID: SC32 Client: National G Location: Former Tidewat Pawtucket, RI	
DЕРТН	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
0	0-	SC32-S1 SC32-S2 SC32-S3	3.0		nonplastic; silver blu	ish spotty sheens throug D (SM); 20-30% fine san	hout. d and organic materi	wet; very soft; grayish brown;	
- 5	-5 -			••••	plastīcitý; spotty blui: 0.87ft. Medium angular SAI gravel; moist; gray; r SILT (ML); 20-30% f	ND (SP); 20-30% coarse no apparent oil/tar odor o	I/8" spot of brown pro sand; 10-20% fine g r staining.	a gray; nonplastic to low oduct in very fine sand layer at ravel; <10% coarse rounded moist; soft; grayish brown; low	
- 10	- - -10 -				plasticity; no appared Void. SILT (ML); 20-30% f with water); soft; gra Medium angular SAI	nt oil/tar odor or sheen. ine to medium sand; <10 yish brown; nonplastic to	% organic material; r low plasticity; no ap	t; soft; grayish brown; low moist to wet (due to void filled parent oil/tar odor or sheen.	-
- 10	_								



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 Data File:SC32.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller' rilling	tart/Finish: y Company s Name: y Method: ation:	: T M V		larine Services, Inc. tian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9479 N 71 22.7784 W 6.1 ft -4.12 ft 5.1 ft B. Thibault	Well/Boring ID: SC33 Client: National G Location: Former Tidewate Pawtucket, RI	
DЕРТН	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	_								-
-	-5 -	SC33-S1 SC33-S2 SC33-S3	0.0		from 0.0 to 1.7ft, mois low plasticity; iridesce	st from 1.7 to 5.1ft; very	soft grading to soft w t; bluish sheen at 2.21	nt has a blocky texture; wet ith depth; black; nonplastic to ft, 3.1 to 3.65ft, and 4.2ft; <1/16" 3ft; slight organic odor.	
	-10 -	SC33-S10		· · · · · ·					5
-	-								-
10 	- - 15 - -								

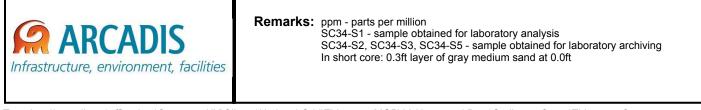


Remarks: ppm - parts per million SC33-S1 - sample obtained for laboratory analysis SC33-S2, SC33-S3, SC33-S10 - sample obtained for laboratory archiving

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 Data File:SC33.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

Di Di Di	rilling riller': rilling	tart/Finish: J Company s Name: J Method: ation:	: T M V		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9305 N 71 22.7661 W 10.8 ft -6.74 ft 4.8 ft B. Thibault	Well/Boring ID: SC34 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- - - - - -	-55	SC34-S1 SC34-S2 SC34-S3 SC34-S5	2.2		texture; wet to moist;	avel; 20.9% fines; 10-20 very soft grading to soft sh sheens throughout.	% organic material; with depth; black; r	leaf layer at 0.85ft; blocky nonplastic to low plasticity;	
- 5									- 1



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 Data File:SC34.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	rilling riller's	art/Finish: Company Name: Method: ation:	: T M V	I. Aval ibraco	Marine Services, Inc. kian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9303 N 71 22.7890 W 4.5 ft -0.24 ft 1.1 ft B. Thibault	Well/Boring ID: SC35 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
		SC35-S1 SC35-S2	0.1		0.1 to 0.15ft; inclusio moist from 0.3 to 0.5 scattered spots of blu Medium angular SAN	n of yellow/brown mediu ft; very soft grading to so uish sheen; organic odor	m grained sand at 0.2 oft with depth; gray; no sand; very fine sand	terial; inclusion of fine sand at 2 to 0.25ft; wet from 0.0 to 0.3ft, onplastic to low plasticity; <1/4" inclusion from 0.6 to 0.8ft;	



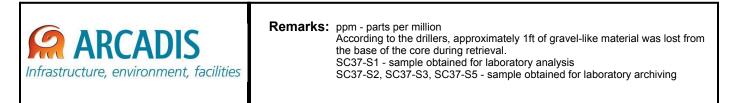
Remarks: ppm - parts per million SC35-S1 - sample obtained for laboratory analysis SC35-S2 - sample obtained for laboratory archiving

D D D	rilling riller'៖ rilling	art/Finish: Company Name: Method: ation:	: T J. V	. Scanl ibraco	Narine Services, Inc. Ion and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9327 N 71 22.7994 W 1.0 ft 0.66 ft 1.2 ft B. Thibault	Well/Boring ID: SC36 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-	-	SC36-S1	2.4	•••		8 90/ grouple 20 90/ cite 4		<10% organic material; moist;	C
-	0 -	SC36-S2			orange-brown at 0.0f odor. SILT (ML); 10-20% fi odor or staining; orga Medium SAND (SP);	t transitioning to gray with ne to medium sand and anic odor. 10-20% coarse sand; <	th depth; no apparent organic material; soft 10% fine gravel and c	; oil/tar odor or staining; organic	 - -
- 5	_ _ _5 _				gravel at 1.2π, moist; iridescent sheen from	orange-brown at 0.0ft tr n 0.9 to 1.0ft; organic od	ansitioning to gray wi	itn deptn; <10% spotty	
	_								
- 10 - -	- 10 -								



Remarks: ppm - parts per million SC36-S1 - sample obtained for laboratory analysis SC36-S2 - sample obtained for laboratory archiving

D D D	rilling riller'៖ rilling	tart/Finish: J Company s Name: J Method: ation:	: T J V		larine Services, Inc. on and J. Balmer re	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9209 N 71 22.7824 W 5.1 ft -3.19 ft 3.25 ft B. Thibault	Well/Boring ID: SC37 Client: National G Location: Former Tidewate Pawtucket, RI	-
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
-									-
-0	-5 -	SC37-S1 SC37-S2 SC37-S3	0.3		material from 1.5 to 1	.8ft; 2" clinker at 1.5ft; m	noist; soft; brownish g	ganic material; 30-40% organic rray; nonplastic to low plasticity; 37ft and 1.8 to 1.85ft; slight	-
	-	SC37-S5			coarse gravel; 20-30%	10-20% fine to coarse s % fine gravel at 2.15ft; a ent oil/tar odor or stainir	mount of coarse mate	oarse sand with depth; 10-20% erial decreases with depth;	
- 5	_								- I
-	-10 -								-
	_								
-10	_								-
	-15 -								-



	Drilling Driller's	tart/Finish: J Company s Name: J Method: ation:	: T M P		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9134 N 71 22.7776 W 5.6 ft -1.99 ft 5.4 ft B. Thibault	Well/Boring ID: SC38 Client: National G Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	0 -								
	-	SC38-S1 SC38-S2 SC38-S3	0.4		SAND (SP); 3.1% gra with depth; black; non from 2.55 to 2.8ft; org	plastic to low plasticity;	organic material; m scattered spotty blui	oist; very soft grading to soft ish sheens throughout, heaviest	(
	-5 - - -				4.5ft the amount of co coarse gravel; fine sa	arse material increases nd layer from 4.6 to 5.0 m; gray; spotty iridesce	to 20-30% coarse s t; silt layer with 10-2	nd and fine gravel; from 4.0 to and and fine gravel, 10-20% .0% fine to medium sand from 5.4ft; organic odor, possible	
-	- 10 -							/	-
- 10 -	-								



Remarks: ppm - parts per million SC38-S1 - sample obtained for laboratory analysis SC38-S2, SC38-S3 - sample obtained for laboratory archiving

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 Data File:SC38.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D)rilling)riller's	tart/Finish: Company s Name: Method: ation:	: T M P		larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9129 N 71 22.7669 W 8.4 ft -5.25 ft 5.2 ft B. Thibault	Well/Boring ID: SC39 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- - -	-5 -	SC39-S1 SC39-S2 SC39-S3	0.5		moist from 1.2 to 5.2ft	t; very soft from 0.0 to 0 astic to low plasticity; 1/	.9ft, soft from 0.9 to 5	/ texture; wet from 0.0 to 1.2ft, 5.2ft; fine gray sand layer at 0.9 it sheen at 0.9ft; spotty bluish	C
- - 5	-10 -								5
- 10	-15 -								-



Remarks: ppm - parts per million SC39-S1 - sample obtained for laboratory analysis SC39-S2, SC39-S3 - sample obtained for laboratory archiving

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 Data File:SC39.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	Date Start/Finish:July 9, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:M. Avakian and J. ScanlonDrilling Method:VibracorePenetration:5.0 ft				larine Services, Inc. ian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1430 N 71 22.8334 W 13.7 ft -11.82 ft 4.8 ft B. Thibault	Well/Boring ID: SC40 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- - -	-10 -	SC40-S1	1.7		Poorly graded mediu	m angular SAND (SD). 1	10-20% fine to coord	e sand; 0.2% gravel; 1.1% silt;	
	-15 -	SC40-S2 SC40-S3			two thin layers of black gray; no apparent oil/ SILT (ML); <10% fine 4.0ft, 4.3ft, 4.6ft, and	ck (color change - no col (tar odor or staining. e and medium sand; 10-2 4.8ft; medium grained s	mposition change) b 20% organic materia and layer from 3.0 to	/ ////////////////////////////////////	
- 5	-			· · · · · ·					
- - -10	-20 - - - -								-



Remarks: ppm - parts per million SC40-S1 - sample obtained for laboratory analysis SC40-S2, SC40-S3 - sample obtained for laboratory archiving

Date Start/Finish: Drilling Company: Driller's Name: Drilling Method: Penetration:	July 9, 2008 TG&B Marine Services, Inc. M. Avakian and J. Scanlon Vibracore 6.0 ft	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.0865 N 71 22.8016 W 11.2 ft -9.85 ft 5.8 ft B. Thibault	Well/Boring Client: Location:	g ID: SC41 National Grid Former Tidewater MGP Site Pawtucket, RI
--	--	--	--	-------------------------------------	--

						- -
	SC41-S1 SC41-S2	1.5		Poorly graded medium angular SAND (SP); 0.1% gravel; 10-20% coarse sand; 2.4% fines; moist; gray; no apparent oil/tar odor or staining.		
	SC41-S3			SILT (ML); <10% fine and medium sand; 1" diameter piece of wood at 1.1ft; layer of leaves at 1.3ft; very soft; moist; black; nonplastic; no apparent oil/tar odor or staining; organic odor.		-
	SC41-S7		· · · · · · · · · · · · · · · · · · ·	Poorly graded medium SAND (SP); <10% fine and coarse grained sand; moist; gray; no apparent oil/tar odor or staining.		-
	3041-37			SILT (ML); 20-30% organic material (leaves and twigs); <10% medium and fine sand; soft; moist; black; low plasticity; small copper-colored sheen at 3.1ft; no apparent oil/tar odor or staining; organic odor.		-
-5 -5 -			·	Poorly graded medium SAND (SP); 30-40% fine sand; <10% coarse sand and fine gravel, brick fragments, and organics; moist; black; no apparent oil/tar odor or staining.	-	-5
				SILT (ML); 20-30% organic material (leaves and twigs); 10-20% fine and medium sand; soft; moist; black; low plasticity; no apparent oil/tar odor or staining.		
				Poorly graded medium angular SAND (SP); <10% fine and coarse grained sand; moist gray; no apparent oil/tar odor or staining.		
						-
- 10- <i>10</i> -					_	10



Remarks: ppm - parts per million SC41-S1 - sample obtained for laboratory analysis SC41-S2, SC41-S3, SC41-S7 - sample obtained for laboratory archiving

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 Data File:SC41.dat
 Date:August 11, 2008
 Project: 36697
 Page: 1 of 1

Date Start/Finish:	July 10, 2008	Latitude:	41 52.0218 N	Well/Boring	ID: SC42
Drilling Company:	TG&B Marine Services, Inc.	Longitude:	71 22.7782 W	Client:	National Grid
Driller's Name:	M. Avakian and J. Scanlon	Water Depth:	8.5 ft	Location:	Former Tidewater MGP Site
Drilling Method:	Vibracore	Mudline Elevation:	-6.95 ft		Pawtucket, RI
Penetration:	5.0 ft	Recovery:	4.7 ft		
		Geologist:	B. Thibault		

DEPTH ELEVATION	Sample ID Headspace (ppm)	ogic Column	Stratigraphic Description	Well/Boring Construction
--------------------	------------------------------	-------------	---------------------------	-----------------------------

5 -						
	SC42-S1 SC42-S2	5.1		Poorly graded medium angular SAND (SP); 2.9% gravel; 10-20% coarse sand; 2.9% fines; moist; gray; no apparent oil/tar odor or staining.		
	SC42-S3			SILT (ML); 10-20% fine to medium sand and organic material; wet; soft; moist; black; nonplastic; no apparent oil/tar odor or staining.		
10 -				Poorly graded medium angular SAND (SP); <10% coarse sand; moist; gray; no apparent oil/tar odor or staining.		-
				SILT (ML); 10-20% fine to medium sand and organic material; moist; soft; black; low plasticity; no apparent oil/tar odor or staining.		-
-5 -			<u> </u>	Poorly graded medium angular SAND (SP); <10% coarse sand; moist; gray; no apparent oil/tar odor or staining.		5
				SILT (ML); 10-20% fine to medium sand and organic material; layer of medium sand at 4.7ft; soft; moist; black; low plasticity; faint small (<1/8") bluish spotty sheens; organic odor.		-
						1
15 -						1
_ 10 _					_	10



Remarks: ppm - parts per million SC42-S1 - sample obtained for laboratory analysis SC42-S2, SC42-S3 - sample obtained for laboratory archiving

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 Data File:SC42.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

D D D	Date Start/Finish:July 10, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:M. Avakian and J. ScanlonDrilling Method:VibracorePenetration:5.5 ft				larine Services, Inc. tian and J. Scanlon	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9817 N 71 22.7651 W 9.1 ft -7.94 ft 5.4 ft B. Thibault	Well/Boring ID: SC43 Client: National Grid Location: Former Tidewater M Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraph	ic Description		Well/Boring Construction
-0	_5 _ _ _ _ _ _	SC43-S1 SC43-S2 SC43-S3	0.6		gray; no apparent oil	tar odor or staining.	_	coarse sand; 11.9% fines; moist;	-
- 5	-10 -				Poorly graded mediu odor or staining. SILT (ML); <10% fine staining.	e to medium sand; soft; r m angular SAND (SP); <	<10% coarse sand; r noist; black; low plas	or staining. moist; gray; no apparent oil/tar sticity; no apparent oil/tar odor or moist; gray, black band at 2.15ft;	-
-	- -15 - - -				from 3.5 to 3.9ft; low Poorly graded mediu odor or staining. SILT (ML); 20-30% fi	plasticity; no apparent o m angular SAND (SP); <	<pre>il/tar odor or staining <10% coarse sand; r 20% organic materia</pre>	moist; black, dark brownish gray g. moist; gray; no apparent oil/tar al; moist; soft; black; low	-
- 10	_								_



Remarks: ppm - parts per million SC43-S1 - sample obtained for laboratory analysis SC43-S2, SC43-S3 - sample obtained for laboratory archiving

 Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores

 Data File:SC43.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

Dri Dri Dri	Date Start/Finish:July 14, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:J. Scanlon and J. BalmerDrilling Method:VibracorePenetration:5.0 ft				larine Services, Inc. on and J. Balmer	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.9151 N 71 22.7470 W 12.4 ft -10.10 ft 4.85 ft B. Thibault	Well/Boring ID: SC44 Client: National C Location: Former Tidewat Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- - - - - - - - - - - - - - - - - - -	-	SC44-S1 SC44-S2 SC44-S3 SC44-S8	1.9		No apparent oil/tar of SILT (ML); 10-20% f 0.9ft; leaf layer at 1.2 organic odor. Fine SAND (SP); <10 sheen; no apparent of SILT (ML); 10-20% f	dor or staining. ine to medium sand and 25ft; wet to moist; soft; bl 0% medium sand and or pil/tar odor or staining. ine to medium sand and ayer at 3.5ft, 4.35ft, and 4	organic material; fir ack; nonplastic to lo ganic material; mois organic material; 20	material; 4.5% fines; moist; gray; he gray sand layer from 0.7 to w plasticity; spotty bluish sheen; st; gray; <10% spotty iridescent -30% organic material from 4.0 ck; low plasticity; 10-20% spotty	



Remarks: ppm - parts per million SC44-S1 - sample obtained for laboratory analysis SC44-S2, SC44-S3, SC44-S9 - sample obtained for laboratory archiving

 Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores

 Data File:SC44.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

Dril Dril Dril	Date Start/Finish:July 14, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:J. Scanlon and J. BalmerDrilling Method:VibracorePenetration:4.5 ft				Latitude:41 51.8819 NLongitude:71 22.7262 WWater Depth:12.2 ftMudline Elevation:-10.25 ftRecovery:3.5 ftGeologist:B. Thibault			
DEPTH	ELEVATION Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
- - - -	- - - - - - - - - - - - - - - - - - -	150		with 10-20% coarse odor or staining.	sand at 0.5ft; thin layer o	f black silt at 0.0ft; n	organic material; 2.5% silt; layer noist; gray; no apparent oil/tar	
-	SC45-S5			Apparent oil/tar odoro Medium SAND (SW) 30% coarse sand at SILT (ML); 10-20% fi	or staining. ; 30-40% fine sand; <10 ⁰ 1.85ft, 2.2ft, and 2.45ft; ç	% coarse sand and o gray; no apparent oil organic material; blo	bist; soft; black; low plasticity; no organic material; layer with 20- /tar odor or staining.	
	-			1				Ę



Remarks: ppm - parts per million SC45-S1 - sample obtained for laboratory analysis SC45-S2, SC45-S3, SC45-S5 - sample obtained for laboratory archiving

 Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores

 Data File:SC45.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

Date Start/Finish: Drilling Company: Driller's Name: Drilling Method: Penetration:	July 14, 2008 TG&B Marine Services, Inc. J. Scanlon and J. Balmer Vibracore 5.5 ft	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 51.8555 N 71 22.7232 W 12.0 ft -10.43 ft 5.3 ft B. Thibault	Well/Boring Client: Location:	g ID: SC46 National Grid Former Tidewater MGP Sit Pawtucket, RI

DEPTH ELEVATION Sample ID Headspace (ppn	Stratigraphic Description	Well/Boring Construction
---	---------------------------	-----------------------------

-10 -	-)
-	SC46-S1 SC46-S2	16.2	· · · · · ·	Medium SAND (SW); 0.1% gravel; 30-40% fine sand; <10% coarse sand; 1.8% fines; moist; gray; no apparent oil/tar odor or staining.		
-	SC46-S3			SILT (ML); 10-20% fine to coarse sand and organic material; moist; soft; black; low plasticity; no apparent oil/tar odor or staining.		
-			•••• • •	Medium SAND (SW); 30-40% fine sand; <10% coarse sand; moist; gray; no apparent oil/tar odor or staining.		
-	-			Medium SAND (SP); 20-30% fine and coarse sand and organic material; <10% fine gravel; possible plastic footing; moist; black; no apparent oil/tar odor or staining.	-	
-15 -	SC46-S10		· ·	SILT (ML); 10-20% fine to medium sand and organic material; medium grained sand layer from 2.85 to 2.9ft and 3.1 to 3.45ft; moist; soft; black; low plasticity; no apparent oil/tar odor or staining.	5	5
-				Medium SAND (SW); 30-40% fine sand; <10% coarse sand; moist; gray, 0.1ft layer of black at 4.4ft; no apparent oil/tar odor or staining.	-	
-				SILT (ML); 10-20% fine to medium sand and organic material; moist; soft; black; low plasticity; no apparent oil/tar odor or staining.	-	
-					-	
-20 -						10
-	-					10



Remarks: ppm - parts per million SC46-S1 - sample obtained for laboratory analysis SC46-S2, SC46-S3, SC46-S10 - sample obtained for laboratory archiving

 Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores

 Data File:SC46.dat
 Date:August 20, 2008
 Project: 36697
 Page: 1 of 1

D D D	Date Start/Finish:July 11, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:M. Avakian and J. ScanlonDrilling Method:VibracorePenetration:5.0 ft				Marine Services, Inc.	Latitude: Longitude: Water Depth: Mudline Elevation: Recovery: Geologist:	41 52.1253 N 71 22.8449 W 10.5 ft -9.06 ft 3.7 ft B. Thibault	Well/Boring ID: SC47 Client: National G Location: Former Tidewate Pawtucket, RI	
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	c Description		Well/Boring Construction
-	_								
 	- 10 -	SC47-S1 SC47-S2 SC47-S3	3.5		material; 1/2" diamet gray; no apparent oil	er slag at 0.15ft; 2" shell /tar odor or staining. 	at 0.05ft; layer of bla	e sand; 2.0% silt; <10% organic ack silt at 0.0ft; moist; brownish and brick fragments; moist; gray;	
-5	- 15 -	SC47-S7							
	-								-
- 10	-20 -								_



Remarks: ppm - parts per million SC47-S1 - sample obtained for laboratory analysis SC47-S2, SC47-S3, SC47-S7 - sample obtained for laboratory archiving

 Template:\\arcadis-us\officedata\Syracuse-NY\Clients\National Grid\Tidewater MGP\06 Notes and Data\Sediment Cores\Tidewater Cores

 Data File:SC47.dat
 Date:August 19, 2008
 Project: 36697
 Page: 1 of 1

Di Di Di	Date Start/Finish:July 11, 2008Drilling Company:TG&B Marine Services, Inc.Driller's Name:M. Avakian and J. ScanlonDrilling Method:Push CorePenetration:1.5 ft			Marine Services, Inc.	-				
DEPTH	ELEVATION	Sample ID	Headspace (ppm)	Geologic Column		Stratigraphi	ic Description		Well/Boring Construction
- - - - - - - - - - - - - - - - - - -		SW-2-S1 SW-2-S2	71.2		moist; soft; black; nor visible product on sec SILT (ML); 30-40% or plasticity; sediment co	plastic; sediment cover liment; heavy possible p ganic material; 10-20%	ed with bluish sheen petroleum-like odor. fine to medium sand sheen; brown produ	; 10-20% silt; 10% coarse sand; ; brown product on gloves; no d; moist; soft; black; low ict on gloves; no product visible	



Remarks: ppm - parts per million SW-2-S1 - sample obtained for laboratory analysis SW-2-S2 - sample obtained for laboratory archiving

APPENDIX B DATA VALIDATION REPORTS

DATA REVIEW FOR

TIDEWATER MGP

PAWTUCKET, RHODE ISLAND

SDG #0807059

VOLATILE, SEMIVOLATILE, METALS, AND MISCELLANEOUS ANALYSES

Analyses performed by:

Alpha Woods Hole Labs Mansfield, Massachusetts

Review performed by:



Syracuse, New York Report #8839

Summary

The following is an assessment of the data package for sample delivery group (SDG) #0807059 for sampling from the Tidewater MGP Site. Included with this assessment are the corrected sample results and chain of custody. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample			Analy	ysis		
			Date	voc	svoc	ТРН	MET	PCBs	MISC
SC23-S1	0807059-01	Sediment	7/8/2008	Х	Х		Х		Х
SC24-S1	0807059-02	Sediment	7/8/2008	Х	Х		Х		Х
SC21-S1	0807059-03	Sediment	7/8/2008	Х	Х		Х		Х
SC10-S1	0807059-04	Sediment	7/8/2008	Х	Х		Х		Х
SC12-S1	0807059-05	Sediment	7/8/2008	Х	Х		Х		Х
SC25-S1	0807059-06	Sediment	7/8/2008	Х	Х		Х		Х
SC3-S1	0807059-07	Sediment	7/9/2008	Х	Х		Х		Х
SC41-S1	0807059-08	Sediment	7/9/2008	Х	Х		Х		Х
SC40-S1	0807059-09	Sediment	7/9/2008	Х	Х		Х		Х
SC4-S1	0807059-10	Sediment	7/9/2008	Х	Х		Х		Х
SC22-S1	0807059-11	Sediment	7/9/2008	Х	Х		Х		Х
SC5-S1	0807059-12	Sediment	7/9/2008	Х	Х		Х		Х
SC6-S1	0807059-13	Sediment	7/9/2008	Х	Х		Х		Х
SC20-S1	0807059-14	Sediment	7/9/2008	Х	Х		Х		Х
SC26-S1	0807059-15	Sediment	7/10/2008	Х	Х		Х		Х
SC19-S1	0807059-16	Sediment	7/10/2008	Х	Х		Х		Х
SC43-S1	0807059-17	Sediment	7/10/2008	Х	Х		Х		Х
SC42-S1	0807059-18	Sediment	7/10/2008	Х	Х		Х		Х
SC9-S1	0807059-19	Sediment	7/10/2008	Х	Х		Х		Х
DUP-1	0807059-20	Sediment	7/10/2008	Х	Х		Х		Х
SC29-S1	0807059-21	Sediment	7/10/2008	Х	Х		Х		Х
SC17-S1	0807059-22	Sediment	7/10/2008	Х	Х		Х		Х
SC28-S1	0807059-23	Sediment	7/10/2008	Х	Х		Х		Х
SC30-S1	0807059-24	Sediment	7/10/2008	Х	Х		Х		Х
SC47-S1	0807059-25	Sediment	7/11/2008	Х	Х		Х		Х
SC32-S1	0807059-26	Sediment	7/11/2008	Х	Х		Х		Х
SC33-S1	0807059-27	Sediment	7/11/2008	Х	Х		Х		Х
SW2-S1	0807059-28	Sediment	7/11/2008	Х	Х		Х		Х
SC39-S1	0807059-29	Sediment	7/11/2008	Х	Х		Х		Х
SC38-S1	0807059-30	Sediment	7/11/2008	Х	Х		Х		Х
SC34-S1	0807059-31	Sediment	7/11/2008	Х	Х		Х		Х
SC35-S1	0807059-32	Sediment	7/11/2008	Х	Х		Х		Х
Trip Blank	0807059-33	Water	7/2/2008	Х					

Notes:

- Miscellaneous parameters include total organic carbon, physiologically available cyanide, and oil and 1. grease. Sample location DUP-1 is the field duplicate of parent sample location SC9-S1.
- 2.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United Stated Environmental Protection Agency) USEPA SW-846 Method 8260. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999. The quality indicators of this limited data review included holding times, associated blanks, matrix spike/matrix spike duplicate (MS/MSD) analysis, field duplicates, laboratory control sample and surrogate recoveries.

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 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
	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
SW-846 8260	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	
Trip Blank	Analysis Completed	22 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.

	Qualification	
Criteria	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
SC32-S1 SW2-S1	Bromomethane	Detected sample results <rl <bal<="" and="" td=""><td>"U" at the PQL</td></rl>	"U" at the PQL
SC24-S1 SC10-S1 SC12-S1 SC25-S1 SC3-S1 SC41-S1 SC40-S1 SC4-S1 SC22-S1 SC20-S1 SC20-S1 SC26-S1 SC26-S1 SC47-S1 SW2-S1 SC35-S1	Acetone	Detected sample results >RL and <bal< td=""><td>"U" at detected sample concentration</td></bal<>	"U" at detected sample concentration

RL = reporting limit

3. 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
SC43-S1	1,2-Dichloroethane-d4	< LL but > 10%
	4-Bromofluorobenzene	AC
	Dibromofluoromethane	AC
	Toluene-d8	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	Non-detect		
calibration curve due to the high concentration of a target compounds	Detect	No Action	

4. 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 were the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

All compounds associated with sample location SC30-S1 except dichlorodifluoromethane, chloromethane, vinyl chloride, methylene chloride, methyl tert-butyl ether, 1,1-dichloroethane, chloroform, 1,2-dichlorpropane, methyl isobutyl ketone, and 2-hexanone had one or both MS/MSD recoveries below the lower control limit; therefore the associated sample result was qualified as estimated (J).

Sample locations associated with the MS/MSD exhibiting recoveries below 10% are presented in the following table.

Sample Locations	Compound	MS Recovery	MSD Recovery
SC30-S1	Acetone	<10%	<10%
	cis-1,3-Dichloropropene	<10%	<ll but=""> 10%</ll>
	trans-1,3-Dichloropropene	<10%	<ll but=""> 10%</ll>

AC = Acceptable

The criteria used to evaluate the MS/MSD recoveries are presented in the following table. In the case of an MS/MSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper control limit (UL)	Non-detect	No Action
	Detect	J
< the lower control limit (LL) but > 10%	Non-detect	UJ
	Detect	J

Control Limit	Sample Result	Qualification
< 10%	Non-detect	R
< 10%	Detect	J
Parent sample concentration > four times the MS/MSD	Detect	No Action
spiking solution concentration (D).	Non-detect	

Sample locations associated with MS/MSD recoveries exhibiting an RPD greater than of the control limit presented in the following table.

Sample Locations	Compound
	Bromomethane
	Vinyl acetate
SC30-S1	2-chloroethylvinyl ether
3030-31	cis-1,3-Dichloropropene
	trans-1,3-Dichloropropene
	Styrene

The criteria used to evaluate the RPD between the MS/MSD recoveries are presented in the following table. In the case of an RPD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
S10	Non-detect	J
> UL	Detect	J

5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/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.

Sample locations associated with LCS/LCSD analysis exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Compound	LCS Recovery	LCSD Recovery
SC23-S1 SC24-S1 SC21-S1 SC10-S1 SC12-S1 SC25-S1 SC3-S1 SC41-S1	Bromomethane	<ll but=""> 10%</ll>	<ll but=""> 10%</ll>

Sample Locations	Compound	Compound LCS LCSD Recovery Recover	
SC40-S1 SC4-S1 SC22-S1 SC5-S1 SC6-S1 SC20-S1	Bromomethane	<ll but=""> 10%</ll>	<ll but=""> 10%</ll>
SC20-S1 SC26-S1 SC19-S1 SC43-S1 SC42-S1 SC9-S1	Vinyl acetate		AC
SC30-S1	Bromomethane	<ll but=""> 10%</ll>	AC
	Vinyl acetate	>UL	AC
SC17-S1 SC28-S1 SC47-S1 SC32-S1 SC33-S1 SW2-S1	Bromomethane	<ll but=""> 10%</ll>	AC
SW2-S1 SC39-S1 SC38-S1 SC34-S1 SC35-S1 Trip Blank		>UL	AC
	Dichlorodifluoromethane	>UL	>UL
DUP-1 SC29-S1	Vinyl chloride	>UL	>UL
0020-01	Carbon disulfide	>UL	>UL

The criteria used to evaluate the LCS/LCSD recoveries are presented in the following table. In the case of an LCS deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper central limit (LLL)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (1.1.) but > 10%	Non-detect	J
< the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
< 10 <i>/</i> 0	Detect	J

Sample locations associated with LCS/LCSD recoveries exhibiting an RPD greater than of the control limit presented in the following table.

Sample Locations	Compound
SC40-S1 SC4-S1 SC22-S1 SC5-S1	Acetone
SC6-S1	

Sample Locations	Compound
SC20-S1	
SC26-S1	
SC19-S1	
SC43-S1	
SC42-S1	
SC9-S1	
SC30-S1	Acetone
3030-31	2-Butanone
SC17-S1	
SC28-S1	Acetone
SC47-S1	
SC32-S1	
SC33-S1	
SW2-S1	2-Butanone
SC39-S1	
SC38-S1	
SC34-S1	Bromomothene
SC35-S1	Bromomethane
Trip Blank	

The criteria used to evaluate the RPD between the MS/MSD recoveries are presented in the following table. In the case of an RPD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	J
	Detect	J

6. 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 and 100% for soil 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
	2-Butanone (MEK)	33.4	75.5	77.3 %
SC9-S1/DUP-1	Acetone	129	298	79.1 %
	Carbon disulfide	45.3	67.4	39.2 %

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two 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.

7. 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
SW2-S1	Naphthalene	39700 E	34200 D	34200 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

8. 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.

SEMI-VOLATILE ORGANIC COMPOUND (SVOC) ANALYSES

Introduction

Analyses were performed according to (United Stated Environmental Protection Agency) USEPA SW-846 Method 8270 selective ion monitoring (SIM). Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999. The quality indicators of this limited data review included holding times, associated blanks, matrix spike/matrix spike duplicate (MS/MSD) analysis, field duplicates, laboratory control sample and surrogate recoveries.

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 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 8270 SIM	Water	7 days from collection to extraction and 40 days from extraction to analysis	Cooled @ 4 °C
3W-040 0270 3IW	Soil	14 days from collection to extraction and 40 days from extraction to analysis	Cooled @ 4 °C

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance blanks (i.e., method 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. 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. 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. SVOC analysis requires that two of the three SVOC surrogate compounds within each fraction exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

4. Matrix Spike/Matrix Spike Duplicate (MS/MSD)/Laboratory Duplicate 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 were the compounds concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

Sample locations associated with the MS/MSD exhibiting recoveries outside of the control limits are presented in the following table.

Sample Locations	Compound MS Recov		MSD Recovery
SC24-S1	Acenaphthylene	<10%	> UL
SC30-S1	Dibenz(a,h)anthracene	> UL	> UL

AC = Acceptable

The criteria used to evaluate the MS/MSD recoveries are presented in the following table. In the case of an MS/MSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper central limit (LLL)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (1.1.) but > 100/	Non-detect	J
< the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
~ 1070	Detect	J

Sample locations associated with MS/MSD recoveries exhibiting an RPD greater than of the control limit presented in the following table.

Sample Locations	Compound
SC24-S1	Acenaphthylene

The criteria used to evaluate the RPD between the MS/MSD recoveries are presented in the following table. In the case of an RPD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	J
2 0L	Detect	J

5. 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.

6. 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 and 100% for soil matrices is applied to the RPD between the parent sample and the field duplicate.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
	Acenaphthene	286	305	6.4 %
	Acenaphthylene	1020	943	7.8 %
	Anthracene	1310	1350	3.0 %
	Benz[a]anthracene	4170	4090	1.9 %
	Benzo[a]pyrene	4900	4780	2.4 %
	Benzo[b]fluoranthene	4720	4530	4.1 %
	Benzo[e]pyrene	3730	3620	2.9 %
	Benzo[g,h,i]perylene	3460	3380	2.3 %
	Benzo[k]fluoranthene	4280	4260	0.4 %
	C1-Chrysenes	2640	2370	10.7 %
	C1-Fluoranthenes/Pyrenes	4370	4210	3.7 %
SC9-S1/DUP-1	C1-Fluorenes	235	246	4.5 %
	C1-Naphthalenes	366	382	4.2 %
	C1-Phenanthrenes/Anthracenes	2090	2020	3.4 %
	C2-Chrysenes	1490	1240	18.3 %
	C2-Fluorenes	362	348	3.9 %
	C2-Naphthalenes	449	425	5.4 %
	C2-Phenanthrenes/Anthracenes	1420	1350	5.0 %
	C3-Chrysenes	1260	1110	12.6 %
	C3-Fluorenes	781	916	15.9 %
	C3-Naphthalenes		334	6.9 %
	C3-Phenanthrenes/Anthracenes	888	804	9.9 %
	C4-Chrysenes	815	724	11.8 %

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
	C4-Naphthalenes	270	263	2.6 %
	C4-Phenanthrenes/Anthracenes	549	469	15.7 %
	Chrysene	4850	4790	1.2 %
	Dibenz[a,h]anthracene	881	861	2.2 %
	Fluoranthene	9520	9600	0.8 %
	Fluorene	370	382	3.1 %
	Indeno[1,2,3-cd]pyrene	3560	3550	0.2 %
	Naphthalene	541	615	12.8 %
	Perylene	1610	1530	5.0 %
	Phenanthrene	3450	3580	3.6 %
	Pyrene	8450	8200	3.0 %

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two 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.

7. 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
SC-25-S1	Fluoranthene	28800 E	24400 D	24400 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

Reported Sample Results	Qualification
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

8. 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.

METALS ANALYSES

Introduction

Analyses were performed according to United States Environmental Protection Agency (USEPA) SW-846 Method 6000/7000. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1994. The quality indicators of this limited data review included holding times, associated blanks, matrix spike (MS) and duplicate analysis, serial dilution analysis, field duplicate, and laboratory control sample recoveries.

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 that it was already 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 the USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The analyte was analyzed for but not detected. The associated value is the analyte instrument detection limit.
 - B The reported value was obtained from a reading less than the contract-required detection limit (CRDL), but greater than or equal to the instrument detection limit (IDL).
- Quantitation (Q) Qualifiers
 - E The reported value is estimated due to the presence of interference.
 - N Spiked sample recovery is not within control limits.
 - * Duplicate analysis is not within control limits.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - 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 6020	Water	180 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	180 days from collection to analysis	Cooled @ 4 °C.
SW-846 7474	Soil	28 days from collection to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method or rinse blanks), are prepared to identify any contamination that may have been introduced into the samples during sample preparation or field activity. Method blanks (including initial and continuing calibration blanks, and preparation blanks) measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected analyte in an associated blank is calculated for QA blanks containing concentrations greater than the IDL. The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No analytes were detected above the reporting limit in the associated blanks.

3. Matrix Spike (MS)/Laboratory Duplicate Analysis

MS and laboratory duplicate data are used to assess the precision and accuracy of the analytical method.

3.1 MS Analysis

All metal analytes must exhibit a percent recovery within the established acceptance limits of 75% to 125%. The MS recovery control limits do not apply for MS performed on sample locations where the analyte's concentration detected in the parent sample exceeds the MS concentration by a factor of four or greater. In instance where this is true, the data will not be qualified even if the percent recovery does not meet the control limits and the laboratory qualifier "N" will be removed.

All analytes associated with MS recoveries were within control limits with the exception of the

following analytes present in the table below.

Sample Location	Analytes	MS Recovery
SC24-S1	Antimony	61%
SC20 S1	Antimony	27%
SC30-S1	Mercury	128%

The criteria used to evaluate MS recoveries are presented in the following table. In the case of an MS deviation, the sample results are qualified. The qualifications are applied to all sample results associated with this SDG.

Control limit	Sample Result	Qualification
MS percent recovery 30% to	Non-detect	UJ
74%	Detect	J
MS percent recovery <30%	Non-detect	R
	Detect	J
MS percent	Non-detect	No Action
recovery >125%	Detect	J

3.2 Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the CRDL. A control limit of 20% for water matrices and 35% for soil matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the CRDL, a control limit of one times the CRDL is applied for water matrices and two times the CRDL for soil matrices.

All analytes associated with laboratory duplicate RPD were within the control limit, with the exception of the analytes presented in the following table.

Sample Location	Analytes	Laboratory RPD
	Arsenic	21%
	Beryllium	136%
SC24-S1	Cadmium	63%
	Chromium	39%
	Lead	142%
SC30-S1	Antimony	25%

The criteria used to evaluate laboratory duplicate RPD are presented in the following table. In the case of a laboratory duplicate RPD deviation, the sample results are qualified. The qualifications are applied to the all sample results associated with this SDG.

Sample Concentration	Control Limit	Sample Result	Qualification
Parent sample and laboratory sample concentration >5 times	Soil 35%	Non-detect	UJ
CRDL	3011 35 %	Detect	J
Parent sample and/or laboratory duplicate sample result ≤ five	Soil two times RL	Non-detect	UJ
times the RL and difference between samples >RL		Detect	J

4. 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 and 100% for soil 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
	Antimony	1.09	1.25	13.6 %
	Arsenic	15.7	15.9	1.2 %
	Beryllium	0.982	1.04	5.7 %
	Cadmium	10.4	11.1	6.5 %
	Chromium	154	158	2.5 %
	Copper	322	314	2.5 %
SC9-S1/DUP-1	Lead	308	302	1.9 %
	Nickel	41	41	0 %
	Selenium	1.99	1.88	5.6 %
	Silver	5.8	6.1	5.0 %
	Thallium	0.18	0.183	1.6 %
	Zinc	455	446	1.9 %
	Mercury	0.903	0.942	4.2 %

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two 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.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences.

The LCS analysis exhibited recoveries within the control limits.

6. Serial Dilution

The serial dilution analysis is used to assess if a significant physical or chemical interference exists due to sample matrix. Analytes exhibiting concentrations greater than 50 times the MDL in the undiluted sample are evaluated to determine if matrix interference exists. These analytes are required to have less than a 10% difference (%D) between sample results from the undiluted (parent) sample and results associated with the same sample analyzed with a five-fold dilution.

All serial dilutions were within control limits, with the exception of the analytes presented in the following table. The sample locations associated with the deviant %D are also presented in the following table.

Sample Locations	Analytes	%D
SC30-S1	Mercury	16%

The criteria used to evaluate the serial dilution are presented in the following table. In the case of a serial dilution deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	UJ
	Detect	J

7. 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.

MISCELLANEOUS ANALYSES

Introduction

Analyses were performed according to United States Environmental Protection Agency (USEPA) SW-846 methods 9010, 9060, and 9071b. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1994. The quality indicators of this limited data review included holding times, associated blanks, matrix spike (MS) and laboratory duplicate analysis, field duplicate, and laboratory control sample recoveries.

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 that it was already 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 the USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The analyte was analyzed for but not detected. The associated value is the analyte instrument detection limit.
 - B The reported value was obtained from a reading less than the contract-required detection limit (CRDL), but greater than or equal to the instrument detection limit (IDL).
- Quantitation (Q) Qualifiers
 - N Spiked sample recovery is not within control limits.
 - * Duplicate analysis is not within control limits.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - 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 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
Physiologically Available Cyanide by SW-846 9010	Soil	14 days from collection to analysis	Cooled @ 4 °C
Total Organic Carbon by 9060	Soil	28 days from collection to analysis	Cooled @ 4 °C
Oil and Grease by 9071b	Soil	28 days from collection to analysis	Cooled @ 4 °C

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method or rinse blanks), are prepared to identify any contamination that may have been introduced into the samples during sample preparation or field activity. Method blanks (including initial and continuing calibration blanks, and preparation blanks) measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected analyte in an associated blank is calculated for QA blanks containing concentrations greater than the MDL. The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No analytes were detected above the reporting limit in the associated blanks.

3. Matrix Spike (MS)/Laboratory Duplicate Analysis

MS and laboratory duplicate data are used to assess the precision and accuracy of the analytical method.

3.1 MS Analysis

All analyses must exhibit a percent recovery within the established acceptance limits of 75% to 125%. The MS recovery control limits do not apply for MS performed on sample locations where the analyte's concentration detected in the parent sample exceeds the MS concentration by a factor of four or greater

All analytes associated with MS recoveries were within control limits with the exception of the following analytes present in the table below.

Sample Location	Analytes	MS Recovery
SC24-S1	TOC	128%
SC30-S1	TOC	272%

The criteria used to evaluate MS recoveries are presented in the following table. In the case of an MS deviation, the sample results are qualified. The qualifications are applied to all sample results associated with this SDG.

Control limit	Sample Result	Qualification
MS percent recovery 30% to	Non-detect	UJ
74%	Detect	J
MS percent recovery <30%	Non-detect	R
	Detect	J
MS percent	Non-detect	No Action
recovery >125%	Detect	J

3.2 Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the CRDL. A control limit of 20% for water matrices and 35% for soil matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the CRDL, a control limit of one times the CRDL is applied for water matrices and two times the CRDL for soil matrices.

All analytes associated with laboratory duplicate RPD were within the control limit, with the exception of the analytes presented in the following table.

Sample Location	Analytes	Laboratory RPD
SC30-S1	TOC	35%

The criteria used to evaluate laboratory duplicate RPD are presented in the following table. In the case of a laboratory duplicate RPD deviation, the sample results are qualified. The qualifications are applied to the all sample results associated with this SDG.

Sample Concentration	Control Limit	Sample Result	Qualification
Parent sample and laboratory sample concentration >5 times	Soil 35%	Non-detect	UJ
CRDL	3011 33 76	Detect	J
Parent sample and/or laboratory duplicate sample result ≤ five	Soil two times RL	Non-detect	UJ
times the RL and difference between samples >RL		Detect	J

4. 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 and 100% for soil 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
	Total Organic Carbon (Run 1)	4.1	3.9	4.9 %
	Total Organic Carbon (Run 2)	4.2	3.6	15.3 %
SC9-S1/DUP-1	TPH, HEM-SGT	5820	5410	7.3 %
	Physiologically Available Cyanide	ND	ND	AC

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two 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.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences.

The LCS analysis exhibited recoveries within the control limits.

6. System Performance and Overall Assessment

Upon validation of the TOC analysis performed within this SDG it was discovered that all samples were analyzed in duplicate. Technical compliance of method SW-846 9060 requires analysis to be performed in quadruplicate. The purpose of the quadruplicate analysis is to establish data precision. Since the laboratory performed the sample analysis in duplicate for each sample location, for data qualification and usability purposes the sample analysis precision of the two reported runs for each sample location were evaluated against laboratory duplicate criterion:

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the reporting limit (RL). A control limit of 20% for water matrices and 35% for soil matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of one times the RL is applied for water matrices and two times the RL for soil matrices.

The criteria used to evaluate laboratory duplicate RPD (in this case the two sample runs) are presented in the following table. In the case of a laboratory duplicate RPD deviation, the sample results are qualified.

Sample Concentration	Control Limit	Sample Result	Qualification
Parent sample and laboratory	Water 20%	Non-detect	UJ
sample concentration >5 times RL	Soil 35%	Detect	J
Parent sample and/or laboratory		Non-detect	UJ
duplicate sample result ≤ five times the RL and difference between samples >RL	Water one times RL Soil two times RL	Detect	J

The duplicate sample results exhibited RPD within the control limit for all sample locations within this SDG; therefore no results were qualified due to the method deviation.

CORRECTED SAMPLE ANALYSIS DATA SHEETS

N/A



ARCADIS Tidewater MGP

SC23-S1

N/A SDG: Sediment Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-01 Associated Blank: VS071808B16 Concentration Units: _µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	<u>, e</u> /st
07/08/08	07/11/08	07/19/08	59.3	6.53	5	1	MLI	
Parameter		F	Result	Parameter	<u> </u>		Result	
Dichlorodifluoro	methane		2.58 U	2-Hexanone			2.58	U
Chloromethane			2.58 U	Tetrachloroeth	ene		2.58	U
Vinyl chloride			2.58 U	1,3-Dichlorop	ropane		2.58	Ų
Bromomethane			6.46 U T	Dibromochlor	omethane		2.58	U
Chloroethane			2,58 U	1,2-Dibromoe	thane		2.58	U
Trichlorofluoror	nethane		2.58 U	Chlorobenzen	3		2.58	U
Acetone			201 B	1,1,1,2-Tetrac	hloroethane		2.58	U
1,1-Dichloroethe	ene		2.58 U	Ethylbenzene			2.58	U
Carbon disulfide	;		18.1	p/m-Xylene			5.17	U
Methylene chlor	ide		6.46 U	o-Xylene			2,58	U
Methyl tert-buty	1 ether (MTBE)		2.58 U	Styrene			2,58	U
trans-1,2-Dichlo			2.58 U	Bromoform			2,58	U
1,1-Dichloroetha	ane		2.58 U	Isopropylbenz	ene		1.86	J
Vinyl acetate			2.58 U	1,1,2,2-Tetrac	hloroethane		2.58	U
2-Butanone (ME	EK)		56.4	Bromobenzene		2.58	U	
cis-1.2-Dichloro			2.58 U	1,2,3-Trichlor	opropane		2.58	U
2,2-Dichloropro	pane	and a second state of the second state of	2.58 U	n-Propylbenze	ne		2.58	U
Chloroform			2.58 U	2-Chlorotolue	ne		2.58	
1,1,1-Trichloroe	thane		2.58 U	1,3,5-Trimethy	ylbenzene		2.58	υ
1,1-Dichloropro			2.58 U	4-Chlorotolue	ne		2.58	U
Carbon tetrachlo	oride		2.58 U	tert-Butylbenz	2.58	U		
Benzene			2.58 U	1,2,4-Trimethy			2.58	U
1,2-Dichloroetha	ane		2.58 U	sec-Butylbenz	ene		2.58	U
Trichloroethene	and her and a		2.58 U	1,3-Dichlorob	enzene		2.58	U
1,2-Dichloropro	pane		2.58 U	p-Isopropyltol			2.58	
Dibromomethan			2.58 U	1,4-Dichlorob			2.58	
Bromodichloron			2.58 U	n-Butylbenzer			2.58	U
2-Chloroethylvin			2.58 U	1,2-Dichlorob			2.58	
	ketone (MIBK)		2.58 U		3-chloropropane		2.58	
cis-1,3-Dichloro		ar a 11	2.58 U	1,2,4-Trichlor			2.58	
Toluene	A ,		2.58 U	Hexachlorobu			2.58	U
trans-1,3-Dichlo	ropropene		2.58 U	Naphthalene			2,58	Ū
1,1,2-Trichloroe			2.58 U	1,2,3-Trichlor	obenzene		2.58	Ū

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	104	70-130
1,2-Dichloroethane-d4	108	70-130
Toluene-d8	93	70-130
4-Bromofluorobenzene	95	70-130

- 6

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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07/31/08 16:11

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Page 7 / 277

N/A



ARCADIS Tidewater MGP

SDG:

: SC-24-S1 N/A

Matrix: Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-02 Associated Blank: VS071808B16 Concentration Units: µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	/st
07/08/08	07/11/08	07/19/08	82.6	6.77	5	1	MLF	
Parameter		F	Result	Parameter			Result	
Dichlorodifluoro	methane		1.79 U	2-Hexanone			1.79	U
Chloromethane			1.79 U	Tetrachloroeth	iene		1.79	U
Vinyl chloride			1.79 U	1,3-Dichlorop	ropane		1.79	U
Bromomethane			4.47 U J	Dibromochlor	omethane		1.79	U
Chloroethane			1.79 U	1,2-Dibromoe	thane		1.79	U
Trichlorofluoror	nethane		1.79 U	Chlorobenzene	0		1.79	U
Acetone			13.7 BU	1,1,1,2-Tetrac	hloroethane		1.79	U
1,1-Dichloroethe	ene		1.79 U	Ethylbenzene			1.79	U
Carbon disulfide			3.81	p/m-Xylene				U
Methylene chlor	ide		4.47 U	o-Xylene			1.79	U
Methyl tert-buty	l ether (MTBE)		1.79 U	Styrene			1,79	υ
trans-1,2-Dichlo			1.79 U	Bromoform			1.79	U
1,1-Dichloroetha	ane		1.79 U	Isopropylbenzene			1.79	U
Vinyl acetate			1.79 U	1,1,2,2-Tetrac	hloroethane		1.79	U
2-Butanone (ME	EK)		1.79 U	Bromobenzen	e		1.79	U
cis-1,2-Dichloro	ethene		1.79 U	1,2,3-Trichloropropane			1.79	U
2,2-Dichloropro	pane		1.79 U	n-Propylbenzene			1,79	U
Chloroform			1.79 U	2-Chlorotoluene			1.79	U
1,1,1-Trichloroe	thane		1.79 U	1,3,5-Trimethy	ylbenzene		1.79	U
1,1-Dichloropro	pene		1.79 U	4-Chlorotolue	ne		1.79	U
Carbon tetrachlo	oride		1,79 U	tert-Butylbenz		1.79	U	
Benzene			1.79 U	1,2,4-Trimethy	ylbenzene		1.79	U
1,2-Dichloroetha	ane		1.79 U	sec-Butylbenz			1.79	U
Trichloroethene			1.79 U	1,3-Dichlorob	enzene		1.79	U
1,2-Dichloropro	· · · · · · · · · · · · · · · · · · ·		1.79 U	p-Isopropyltoluene			1.79	U
Dibromomethan	e		1.79 U	1,4-Dichlorobenzene			1.22	J
Bromodichloron	nethane		1.79 U	n-Butylbenzer	ne		1.79	U
2-Chloroethylvii	nyl ether		1.79 U	1,2-Dichlorob	enzene		1.79	U
Methyl isobutyl	ketone (MIBK)		1.79 U	1,2-Dibromo-	3-chloropropane		1.79	U
cis-1,3-Dichloro			1.79 U	1,2,4-Trichlorobenzene			1.79	U
Toluene			1.79 U	Hexachlorobutadiene			1.79	U
trans-1,3-Dichlo	ropropene		1.79 U	Naphthalene			1.11	J
1,1,2-Trichloroe			1.79 U	1,2,3-Trichlor	obenzene		1.79	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	100	70-130
1,2-Dichloroethane-d4	94	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	85	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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Page 8 / 277

N/A



ARCADIS

Tidewater MGP

SC-21-S1 SDG;

Matrix: Sediment

- r

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-03 Associated Blank: VS071808B16 Concentration Units: µg/Kg

				Sample	Final	Dilution		
Date Collected			Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	yst
07/08/08	07/11/08	07/19/08	42.5	6.24	5	1	ML	R
Parameter		1	Result	Parameter			Result	
Dichlorodifluoro	methane		3.77 U	2-Hexanone			3.77	U
Chloromethane			<u>3.77 U</u>	Tetrachloroeth	ene		3.77	U
Vinyl chloride			<u>3.77 U</u>	1,3-Dichloropi	ropane		3.77	U
Bromomethane			9.42 U J	Dibromochloro	omethane		3.77	U
Chloroethane			3.77 U	1,2-Dibromoet	hane		3.77	U
Trichlorofluoron	nethane		3.77 U	Chlorobenzene	>		3.77	U
Acetone			134 B	1,1,1,2-Tetracl	nloroethane		3.77	U
1,1-Dichloroethe	ne		3.77 U	Ethylbenzene			3.77	U
Carbon disulfide			20.9	p/m-Xylene			7.53	Ŭ
Methylene chlori		-	9.42 U	o-Xylene			3.77	Ū
Methyl tert-butyl			3.77 U	Styrene			3.77	Ū
trans-1,2-Dichlor	oethene		3.77 U	Bromoform	······································		3.77	Ŭ
1,1-Dichloroetha	ne		3,77 U	Isopropylbenze	ene		3.77	U
Vinyl acetate	~		3.77 U	1,1,2,2-Tetracl	loroethane		3.77	U
2-Butanone (ME	K)		35.6	Bromobenzene			3.77	U
cis-1,2-Dichlorod	thene		3.77 U	1,2,3-Trichloropropane			3.77	U
2,2-Dichloroprop	ane		3.77 U	n-Propylbenze	ne		3.77	Ū
Chloroform			3.77 U	2-Chlorotoluer	ıe		3.77	U
1,1,1-Trichloroet	hane		3.77 U	1,3,5-Trimethy	lbenzene	•	3.77	U
1,1-Dichloroprop	ene		3.77 U	4-Chlorotoluer	1e		3.77	υ
Carbon tetrachlor	ride		3.77 U	tert-Butylbenze	ene		3.77	U
Benzene			3.77 U	1,2,4-Trimethy	lbenzene		3.77	U
1,2-Dichloroetha	ne		3.77 U	sec-Butylbenze	ene		3.77	U
Trichloroethene			3.77 U	1,3-Dichlorobe	nzene		3.77	Ū
1,2-Dichloroprop	ane		3.77 U	p-Isopropyltoh	iene		3.77	U
Dibromomethane	;		3.77 U	1,4-Dichlorobe	nzene		3.77	U
Bromodichlorom	ethane		3.77 U	n-Butylbenzen	e		3.77	Ū
2-Chloroethylvin	yl ether		3.77 U	1,2-Dichlorobe		an Ber andere and an	3.77	U
Methyl isobutyl l	etone (MIBK)		3.77 U	monthermout and the state of th	-chloropropane		3.77	U
cis-1,3-Dichlorop						3.77	Ŭ	
Тоluепе			3.77 U	Hexachlorobut			3.77	U
trans-1,3-Dichlor	opropene		3.77 U	Naphthalene			3.77	Ŭ
1,1,2-Trichloroet	hane		3.77 U	1,2,3-Trichloro	benzene		3.77	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	98	70-130
1,2-Dichloroethane-d4	93	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	83	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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ARCADIS **Tidewater MGP**

SC-10-S1 N/A

SDG: N/A

Matrix: Sediment Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-04 Associated Blank: VS071808B16 **Concentration Units:** µg/Kg

	· · · · · · · · · · · · · · · · · · ·			Sample	Final	Dilution		g/ng
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	/st
07/08/08	07/11/08	07/19/08	82.4	7.80	5	1	MLI	2
Parameter		F	Result	Parameter	L	l,,,,,	Result	
Dichlorodifluoro	methane	- 1 · · · · · · · · · · · · · · · · ·	1.56 U	2-Hexanone			1.56	U
Chloromethane			<u>1.56 U</u>	Tetrachloroeth			1.56	U
Vinyl chloride			1.56 U	1,3-Dichlorop	ropane		1.56	U
Bromomethane			<u>3.89 U T</u>	Dibromochlor	omethane		1.56	U
Chloroethane			1.56 U	1,2-Dibromoet	hane		1.56	U
Trichlorofluoron	nethane		1.56 U	Chlorobenzene	e		1.56	U
Acetone			20.9 B 🗸	1,1,1,2-Tetrac	hloroethane		1.56	U
1,1-Dichloroethe	ene		1.56 U	Ethylbenzene			1.56	U
Carbon disulfide			11.1	p/m-Xylene			3.11	U
Methylene chlor	ide		3.89 U	o-Xylene			1.56	U
Methyl tert-buty	l ether (MTBE)		1.56 U	Styrene			1.56	U
trans-1,2-Dichlo			1.56 U	Bromoform		1,56	U	
1,1-Dichloroetha	ine		1.56 U	Isopropylbenzene			1.56	U
Vinyl acetate			1.56 U	1,1,2,2-Tetrac	1,1,2,2-Tetrachloroethane			U
2-Butanone (ME	K)		3.27	Bromobenzene			1.56	U
cis-1,2-Dichloro	ethene		1.56 U	1,2,3-Trichloropropane			1.56	U
2,2-Dichloroprop	oane		1.56 U	n-Propylbenzene			1.56	U
Chloroform			1.56 U	2-Chlorotoluer	1,56	U		
1,1,1-Trichloroe	thane		1.56 U	1,3,5-Trimethylbenzene			1.56	U
1,1-Dichloroproj	oene		1,56 U	4-Chlorotoluene			1,56	U
Carbon tetrachlo	ride		1.56 U	tert-Butylbenz	ene		1.56	U
Benzene			1.56 U	1,2,4-Trimethy			1.56	U
1,2-Dichloroetha	ine		1.56 U	sec-Butylbenz	ene		1.56	U
Trichloroethene			1.56 U	1,3-Dichlorob	enzene		1.56	U
1,2-Dichloroprop	oane		1.56 U	p-Isopropyltol	uene		1.56	U
Dibromomethan	e		1.56 U	1,4-Dichlorob	enzene		7,60	
Bromodichloron	nethane		1.56 U	n-Butylbenzen	e		1.56	U
2-Chloroethylvir	nyl ether		1.56 U	1,2-Dichlorob			1.56	U
Methyl isobutyl	ketone (MIBK)		1.56 U	1,2-Dibromo-2	3-chloropropane		1.56	U
cis-1,3-Dichloro	propene		1.56 U	1,2,4-Trichlor	obenzene		1.56	U
Toluene			1.56 U	Hexachlorobu	tadiene		1.56	U
trans-1,3-Dichlo	ropropene		1.56 U	Naphthalene			2.15	
1,1,2-Trichloroe			1.56 U	1,2,3-Trichlor	obenzene		1.56	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	101	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	83	70-130

- 14

N/A - Not Applicable

B - Found in associated blank as well as sample.

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U - The analyte was analyzed for but not detected at the sample specific level reported.

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N/A



ARCADIS **Tidewater MGP**

SC-12-S1 N/A

SDG: Sediment

Matrix: <u>т</u>

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-05 Associated Blank: VS071808B16 Concentration Units: µg/Kg

		T		<u> </u>		Concentration	$\mu g/\kappa g$
Data Collected	Date Received	Date Analyzed	Percent Solid	Sample	Final Values (m1)	Dilution	
and as a last treatment of the second s		T		Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/19/08	86.0	7.46	5	1	MLR
Parameter		F	Result	Parameter			Result
Dichlorodifluor	methane		<u>1.56 U</u>	2-Hexanone			1.56 U
Chloromethane			1.56 U	Tetrachloroeth	ene		1.56 U
Vinyl chloride			<u>1.56 U</u>	1,3-Dichlorop	ropane		1.56 U
Bromomethane			3.90 U J	Dibromochlor			1.56 U
Chloroethane			1.56 U	1,2-Dibromoet	hane		1.56 U
Trichlorofluoror	nethane		<u>1.56 U</u>	Chlorobenzene	3		1.56 U
Acetone			<u>20.0 B U</u>	1,1,1,2-Tetracl	hloroethane		1.56 U
1,1-Dichloroethe	······································	· · · · · · · · · · · · · · · · · · ·	1.56 U	Ethylbenzene			1.56 U
Carbon disulfide			7.39	p/m-Xylene			3.12 U
Methylene chlor			3.90 U	o-Xylene			1.56 U
Methyl tert-buty	l ether (MTBE)		1.56 U	Styrene			1.56 U
trans-1,2-Dichlo	roethene		1.56 U	Bromoform			1.56 U
1,1-Dichloroetha	ane		1.56 U	Isopropylbenzene			1.56 U
Vinyl acetate			1.56 U	1,1,2,2-Tetrachloroethane			1.56 U
2-Butanone (ME	EK)		2.69	Bromobenzene			1.56 U
cis-1,2-Dichloro	ethene		1.56 U	1,2,3-Trichloropropane			1.56 U
2,2-Dichloropro	pane		1.56 U	n-Propylbenzene			1.56 U
Chloroform			1.56 U	2-Chlorotoluer	2-Chlorotoluene		
1,1,1-Trichloroe			1.56 U	1,3,5-Trimethy	/lbenzene		<u>1.56 U</u> 1.56 U
1,1-Dichloropro	pene		1.56 U	4-Chlorotoluene			1.56 U
Carbon tetrachlo	ride		1.56 U	tert-Butylbenzene			1.56 U
Benzene			1.56 U	1,2,4-Trimethy	/Ibenzene		1.56 U
1,2-Dichloroetha	ine		1.56 U	sec-Butylbenzo	ene		1.56 U
Trichloroethene			1.56 U	1,3-Dichlorobe	enzene		1.56 U
1,2-Dichloropro	pane		1.56 U	p-Isopropyltol	uene		1.56 U
Dibromomethan	e		1.56 U	1,4-Dichlorobe	enzene		3.97
Bromodichloron	nethane		1.56 U	n-Butylbenzen	e		1.56 U
2-Chloroethylvin	nyl ether		1.56 U	1,2-Dichlorobe	enzene		1.56 U
Methyl isobutyl	ketone (MIBK)		1,56 U	1,2-Dibromo-3	-chloropropane		1.56 U
cis-1,3-Dichloro			1.56 U	1,2,4-Trichloro			1.56 U
Toluene			3.63	Hexachlorobut			1.56 U
trans-1,3-Dichlo			1.56 U	Naphthalene			1.45 J
1,1,2-Trichloroe			1.56 U	1,2,3-Trichloro	benzene		1.56 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	104	70-130
1,2-Dichloroethane-d4	103	70-130
Toluene-d8	91	70-130
4-Bromofluorobenzene	84	70-130

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N/A - Not Applicable

B - Found in associated blank as well as sample.

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J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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ARCADIS **Tidewater MGP**

SC-25-S1

N/A

SDG: N/A

Matrix: Sediment Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-06 Associated Blank: VS071808B16 **Concentration Units:** µg/Kg

Data Callastad	Dete Bresterd			Sample	Final	Dilution	
		Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/19/08	76.7	10.43	5	1	MLR
Parameter		I	Result	Parameter			Result
Dichlorodifluoro	methane		<u>1.25 U</u>	2-Hexanone			1.25 U
Chloromethane			1.25 U	Tetrachloroeth			1.25 U
Vinyl chloride			<u>1.25 U</u>	1,3-Dichlorop	ropane		1.25 U
Bromomethane			3.13 U J	Dibromochlor	omethane		1.25 U
Chloroethane	······································	·····	1.25 U	1,2-Dibromoe	thane		1.25 U
Trichlorofluoron	nethane		1.25 U	Chlorobenzene	Э		1.25 U
Acetone			45.6 B V	1,1,1,2-Tetracl	hloroethane		1.25 U
1,1-Dichloroethe	me		1.25 U	Ethylbenzene			1.25 U
Carbon disulfide			6.88	p/m-Xylene			2.50 U
Methylene chlor			<u>3.13 U</u>	o-Xylene			1.25 U
Methyl tert-buty	l ether (MTBE)		1.25 U	Styrene			1.25 U
trans-1,2-Dichlo			<u>1.25 U</u>	Bromoform			1.25 U
1,1-Dichloroetha	ine		1.25 U	Isopropylbenz			1.25 U
Vinyl acetate			1.25 U	1,1,2,2-Tetracl	nloroethane		1.25 U
2-Butanone (ME			9.51	Bromobenzene	3		1.25 U
cis-1,2-Dichloro			1.25 U	1,2,3-Trichlord	opropane		1.25 U
2,2-Dichloroproj	oane		<u>1.25 U</u>	n-Propylbenze	ne		1.25 U
Chloroform			1.25 U	2-Chlorotoluer	10		1.25 U
1,1,1-Trichloroe			1.25 U	1,3,5-Trimethy	/lbenzene		1.25 U
1,1-Dichloroprop			<u>1.25 U</u>	4-Chlorotoluer		1.25 U	
Carbon tetrachlo	ride		<u>1.25 U</u>	tert-Butylbenz		1.25 U	
Benzene			1.25 U	1,2,4-Trimethy	/lbenzene		1.25 U
1,2-Dichloroetha	ne		1.25 U	sec-Butylbenze			1.25 U
Trichloroethene		· · · · · · · · · · · · · · · · · · ·	1.25 U	1,3-Dichlorobe	enzene		1.25 U
1,2-Dichloroprop			1.25 U	p-Isopropyltol	uene		1.25 U
Dibromomethan			<u>1.25 U</u>	1,4-Dichlorobenzene			1.21 J
Bromodichlorom			1.25 U	n-Butylbenzen			1.25 U
2-Chloroethylvir			1.25 U	1,2-Dichlorobe			1.25 U
Methyl isobutyl		····	1.25 U		-chloropropane		1.25 U
cis-1,3-Dichloro	propene		1.25 U	1,2,4-Trichlord	obenzene		1.25 U
Toluene			1.25 U	Hexachlorobut	adiene		1.25 U
trans-1,3-Dichlor			1.25 U	Naphthalene			1.25 U
1,1,2-Trichloroet	thane		1.25 U	1,2,3-Trichlord	benzene		1.25 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	105	70-130
1,2-Dichloroethane-d4	107	70-130
Toluene-d8	95	70-130
4-Bromofluorobenzene	89	70-130

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N/A - Not Applicable

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B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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N/A



ARCADIS Tidewater MGP

SDG:

D: SC3-S1 N/A

Matrix: Sediment

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Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-07 Associated Blank: VS071808B16 Concentration Units: µg/Kg

				Sample	Final	Dilution	
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/19/08	80.2	6.29	5	1	MLR
Parameter		1	Result	Parameter			Result
Dichlorodifluoro	methane		1.98 U	2-Hexanone			1.98 U
Chloromethane			1.98 U	Tetrachloroeth			<u>1.98 U</u>
Vinyl chloride	······································		<u>1.98 U</u>	1,3-Dichlorop	opane		<u>1.98</u> U
Bromomethane			<u>4.96 U J</u>	Dibromochlore	omethane		1.98 U
Chloroethane			1.98 U	1,2-Dibromoet			1.98 U
Trichlorofluoror	nethane		<u>1.98 U</u>	Chlorobenzene			1.98 U
Acetone			21.3 BU	1,1,1,2-Tetracl	loroethane		1.98 U
1,1-Dichloroethe	ene		<u>1.98 U</u>	Ethylbenzene			1.98 U
Carbon disulfide)		2.64	p/m-Xylene			3.97 U
Methylene chlor	ide		4.96 U	o-Xylene			1.98 U
Methyl tert-buty	l ether (MTBE)		<u>1.98 U</u>	Styrene			1.98 U
trans-1,2-Dichlo	roethene		1.98 U	Bromoform			1.98 U
1,1-Dichloroetha	ane		1.98 U	Isopropylbenz	ene		1.98 U
Vinyl acetate			1.98 U	1,1,2,2-Tetracl	loroethane		1.98 U
2-Butanone (ME	EK)		3.41	Bromobenzene	;		1.98 U
cis-1,2-Dichloro	ethene		1.98 U	1,2,3-Trichloro	propane		1.98 U
2,2-Dichloropro	pane		1.98 U	n-Propylbenze	ne		1.98 U
Chloroform			1,98 U	2-Chlorotoluer	ne		1.98 U
1,1,1-Trichloroe	thane		1.98 U	1,3,5-Trimethy	/lbenzene		1.98 U
1,1-Dichloropro	pene		1,98 U	4-Chlorotoluer	10		1.98 U
Carbon tetrachlo	ride		1.98 U	tert-Butylbenz	ene		1.98 U
Benzene			1.98 U	1,2,4-Trimethy	/lbenzene		1.98 U
1,2-Dichloroetha	ane		1.98 U	sec-Butylbenze	ene		1.98 U
Trichloroethene			1.98 U	1,3-Dichlorob	enzene		1.98 U
1,2-Dichloropro	pane		1.98 U	p-Isopropyltol	uene		1.98 U
Dibromomethan	e		1.98 U	1,4-Dichlorob	enzene		1.98 U
Bromodichloron	nethane		1.98 U	n-Butylbenzen	e		1.98 U
2-Chloroethylvii	nyl ether		1.98 U	1,2-Dichlorobe	enzene		1.98 U
Methyl isobutyl	ketone (MIBK)		1.98 U	1,2-Dibromo-3	-chloropropane		1.98 U
cis-1,3-Dichloro	propene		1.98 U	1,2,4-Trichlor	obenzene		1.98 U
Toluene			1.98 U	Hexachlorobut	adiene		1,98 U
trans-1,3-Dichlo	ropropene		1.98 U	Naphthalene			1.98 U
1,1,2-Trichloroe			1.98 U	1,2,3-Trichlord	obenzene		1.98 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	102	70-130
Toluene-d8	94	70-130
4-Bromofluorobenzene	85	70-130

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N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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N/A



ARCADIS Tidewater MGP

SC41-S1 N/A SDG:

Matrix: Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-08 Associated Blank: VS071808B16 Concentration Units: µg/Kg

				Sample	Final	Dilution	
Date Collected	Date Received		Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/19/08	81.3	9.60	5	1	MLR
Parameter		I	Result	Parameter			Result
Dichlorodifluoro	omethane		1.28 U	2-Hexanone			1.28 U
Chloromethane			1.28 U	Tetrachloroeth			<u>1.28 U</u>
Vinyl chloride			1.28 U	1,3-Dichloropi	ropane		1.28 U
Bromomethane			<u>3.20 U J</u>	Dibromochloro		· · · · · · · · · · · · · · · · · · ·	<u>1.28 U</u>
Chloroethane			1.28 U	1,2-Dibromoet	hane		1.28 U
Trichlorofluoror	nethane		<u>1.28 U</u>	Chlorobenzene			1.28 U
Acetone			16.8 BU	1,1,1,2-Tetracl	nloroethane		1.28 U
1,1-Dichloroethe	ene		<u>1,28 U</u>	Ethylbenzene			1.28 U
Carbon disulfide			1.63	p/m-Xylene			2.56 U
Methylene chlor	ide		<u>3.20 U</u>	o-Xylene			1.28 U
Methyl tert-buty	l ether (MTBE)		1.28 U	Styrene			1.28 U
trans-1,2-Dichlo	roethene		1.28 U	Bromoform	West of the Marcel Sec		1.28 U
1,1-Dichloroetha	ine		1.28 U	Isopropylbenz	ene		1.28 U
Vinyl acetate			1.28 U	1,1,2,2-Tetracl	nloroethane		1.28 U
2-Butanone (MF	K)		1.90	Bromobenzene	<u>;</u>		1.28 U
cis-1,2-Dichloro		4 Mar and a 19 191 1-191	1.28 U	1,2,3-Trichloro	opropane		1.28 U
2,2-Dichloropro	pane		1.28 U	n-Propylbenze	ne		1.28 U
Chloroform			<u>1.28 U</u>	2-Chlorotoluer	ne		1.28 U
1,1,1-Trichloroe	thane		1.28 U	1,3,5-Trimethy	lbenzene		1,28 U
1,1-Dichloropro	pene		1.28 U	4-Chlorotoluer	ne		1.28 U
Carbon tetrachlo	ride		1.28 U	tert-Butylbenz	ene		1.28 U
Benzene			1.28 U	1,2,4-Trimethy			1.28 U
1,2-Dichloroetha	ine		1.28 U	sec-Butylbenze	ene		1.28 U
Trichloroethene			1.28 U	1,3-Dichlorobe	enzene		1.28 U
1,2-Dichloropro	pane		1.28 U	p-Isopropyltol	uene		1.28 U
Dibromomethan	e		1.28 U	1,4-Dichlorobe	enzene		1.28 U
Bromodichloron	nethane		1.28 U	n-Butylbenzen	e		1.28 U
2-Chloroethylvin	nyl ether		1.28 U	1,2-Dichlorobe	enzene		1.28 U
Methyl isobutyl	ketone (MIBK)		1.28 U	1,2-Dibromo-3	-chloropropane		1.28 U
cis-1,3-Dichloro			1.28 U	1,2,4-Trichlord	obenzene		1.28 U
Toluene			1.28 U	Hexachlorobut	tadiene		1.28 U
trans-1,3-Dichlo	ropropene		1.28 U	Naphthalene			1.28 U
1,1,2-Trichloroe			1.28 U	1,2,3-Trichlord	obenzene		1.28 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	105	70-130
1,2-Dichloroethane-d4	104	70-130
Toluene-d8	91	70-130
4-Bromofluorobenzene	86	70-130

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. هور N/A - Not Applicable

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B - Found in associated blank as well as sample,

U - The analyte was analyzed for but not detected at the sample specific level reported.

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N/A



Matrix:

ARCADIS Tidewater MGP

SC40-S1 N/A S

N/A SDG: Sediment Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-09 Associated Blank: VS072108B21 Concentration Units: µg/Kg

				Sample	Final	Dilution	
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/21/08	84.4	10.17	5	1	MLR
Parameter		I	Result	Parameter			Result
Dichlorodifluoro	methane		<u>1.17 U</u>	2-Hexanone			1.17 U
Chloromethane			1.17 U	Tetrachloroeth	iene		<u>1.17 U</u>
Vinyl chloride			<u>1.17 U</u>	1,3-Dichlorop	ropane		1.17 U
Bromomethane			2.91 U J	Dibromochlor	omethane		<u>1.17 U</u>
Chloroethane			<u>1.17 U</u>	1,2-Dibromoe			1.17 U
Trichlorofluoron	nethane		<u>1.17 U</u>	Chlorobenzene	Contraction of the American State of the Ame		<u>1.17 U</u>
Acetone			<u>13.7 メリ</u> ろ	1,1,1,2-Tetracl	hloroethane		1.17 U
1,1-Dichloroethe	ene		1.17 U	Ethylbenzene			1.17 U
Carbon disulfide)		1.29	p/m-Xylene			2.33 U
Methylene chlor	ide		2.91 U	o-Xylene			1.17 U
Methyl tert-buty	l ether (MTBE)		<u>1.17 U</u>	Styrene			<u>1.17 U</u>
trans-1,2-Dichlo	roethene		<u>1.17 U</u>	Bromoform			<u>1.17 U</u>
1,1-Dichloroetha	ine		1.17 U	Isopropylbenz	ene		1.17 U
Vinyl acetate			<u>1.17 U</u>	1,1,2,2-Tetrac	hloroethane		1.17 U
2-Butanone (ME	EK)		2.03	Bromobenzen			<u>1.17 U</u>
cis-1,2-Dichloro	ethene		<u>1.17 U</u>	1,2,3-Trichlor			1.17 U
2,2-Dichloropro	pane		<u>1.17 U</u>	n-Propylbenze			<u>1.17 U</u>
Chloroform			<u>1.17 U</u>	2-Chlorotolue	*****		1.17 U
1,1,1-Trichloroe	thane		1.17 U	1,3,5-Trimethy			<u>1.17 U</u>
1,1-Dichloropro			<u>1.17 U</u>	4-Chlorotolue		·	1.17 U
Carbon tetrachle	oride		<u>1.17 U</u>	tert-Butylbenz			1.17 U
Benzene			1.17 U	1,2,4-Trimethy			<u>1.17 U</u>
1,2-Dichloroetha	ane		<u>1.17 U</u>	sec-Butylbenz			<u>1.17 U</u>
Trichloroethene			1.17 U	1,3-Dichlorob			<u>1.17 U</u>
1,2-Dichloropro	pane		<u>1.17 U</u>	p-Isopropyltol			<u>1.17 U</u>
Dibromomethan	e		1.17 U	1,4-Dichlorob	enzene		<u>1.17 U</u>
Bromodichloron	nethane		1.17 U	n-Butylbenzer		***	1.17 U
2-Chloroethylvii			1.17 U	1,2-Dichlorob			1.17 U
Methyl isobutyl			1.17 U		3-chloropropane		<u>1.17 U</u>
cis-1,3-Dichloro	propene		<u>1.17 U</u>	1,2,4-Trichlor			1.17 U
Toluene			1.17 U	Hexachlorobu	tadiene		1.17 U
trans-1,3-Dichlo	ropropene		1.17 U	Naphthalene		·····	1.17 U
1,1,2-Trichloroe	thane		1.17 U	1,2,3-Trichlor	obenzene		1.17 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	111	70-130
1,2-Dichloroethane-d4	112	70-130
Toluene-d8	93	70-130
4-Bromofluorobenzene	86	70-130

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N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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N/A



Matrix:

ARCADIS **Tidewater MGP**

SC4-S1 N/A

SDG: Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-10 Associated Blank: VS072108B21 Concentration Units: µg/Kg

Date Collected	Data Passiyad	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	vot
07/09/08	07/11/08	07/21/08	89.3	7.76	5	1	ML	
Parameter	07/11/08		Result	Parameter			Result	<u> </u>
Dichlorodifluoro	methane	-	1.44 U	2-Hexanone			1.44	TI
Chloromethane	moulane		1.44 U	Tetrachloroeth	lene		1.44	U
Vinyl chloride			1.44 U	1,3-Dichlorop			1.44	
Bromomethane			3.61 U J	Dibromochlor			1.44	Ū
Chloroethane		Ar as a diameter	1.44 U	1,2-Dibromoe			1,44	Ū
Trichlorofluoror	nethane		1.44 U	Chlorobenzen			1.44	Ŭ
Acetone			25.5 BUJ				1,44	U
1,1-Dichloroethe	ene	Line	1.44 U	Ethylbenzene			1.44	U
Carbon disulfide			4.01	p/m-Xylene			2.89	U
Methylene chlor	ide		3.61 U	o-Xylene			1.44	U
Methyl tert-buty	l ether (MTBE)		1.44 U	Styrene			1.44	U
rans-1,2-Dichlo			1.44 U	Bromoform			1.44	U
1,1-Dichloroetha	ane		1.44 U	Isopropylbenz	ene		1.44	U
Vinyl acetate			1.44 U	1,1,2,2-Tetrac	hloroethane		1.44	U
2-Butanone (ME	EK)		5.81	Bromobenzen	e		1.44	U
cis-1,2-Dichloro	ethene		1.44 U	1,2,3-Trichlor	opropane		1.44	U
2,2-Dichloropro	pane		<u>1.44 U</u>	n-Propylbenze	ene		1.44	U
Chloroform			<u>1,44 U</u>	2-Chlorotolue	ne		1.44	U
1,1,1-Trichloroe	thane		<u>1.44 U</u>	1,3,5-Trimethy	ylbenzene		1.44	U
1,1-Dichloropro	pene		<u>1.44 U</u>	4-Chlorotolue	ne		1.44	<u>U</u>
Carbon tetrachlo	oride		<u>1.44 U</u>	tert-Butylbenz			1.44	
Benzene			<u>1.44 U</u>	1,2,4-Trimeth	ylbenzene		1.44	U
1,2-Dichloroetha	ane		<u>1.44 U</u>	sec-Butylbenz	ene		1.44	U
Trichloroethene			<u>1.44 U</u>	1,3-Dichlorob	enzene		1.44	U
1,2-Dichloropro	pane		<u>1.44 U</u>	p-Isopropyltol			1.44	U
Dibromomethan	1 / / · · · · · · · · · · · · · · · · ·		<u>1.44 U</u>	1,4-Dichlorob			2.94	
Bromodichloron	and the second sec		1.44 U	n-Butylbenzer			1.44	
2-Chloroethylvii	nyl ether		1.44 U	1,2-Dichlorob			1.44	
Methyl isobutyl			1.44 U		3-chloropropane		1.44	
cis-1,3-Dichloro	propene		1,44 U	1,2,4-Trichlor			1.44	
Toluene		ur e) — v v v v v v v v v v v v v v v	1.44 U	Hexachlorobu	tadiene		1.44	U
trans-1,3-Dichlo	ropropene		1.44 U	Naphthalene			2.84	
1,1,2-Trichloroe	thane		1.44 U	1,2,3-Trichlor	obenzene		1,44	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	109	70-130
1,2-Dichloroethane-d4	108	70-130
Toluene-d8	93	70-130
4-Bromofluorobenzene	93	70-130

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N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

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ARCADIS Tidewater MGP

SC22-S1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-11 Associated Blank: VS072108B21 Concentration Units: µg/Kg

07/09/0807/11/0807/21/08ParameterResultDichlorodifluoromethane1.15Chloromethane1.15Vinyl chloride1.15Bromomethane2.88	U U U U U J	Amount (g) 10.84 Parameter 2-Hexanone Tetrachloroeth 1,3-Dichloroph Dibromochloro		Factor 1	Analyst MLR Result 1.15 U 1.15 U
ParameterResultDichlorodifluoromethane1.15Chloromethane1.15Vinyl chloride1.15Bromomethane2.88	U U U U U	Parameter 2-Hexanone Tetrachloroeth 1,3-Dichloropi	ene	1	Result 1.15 U
Dichlorodifluoromethane1.15Chloromethane1.15Vinyl chloride1.15Bromomethane2.88	U U U U U	2-Hexanone Tetrachloroeth 1,3-Dichloropi			1.15 U
Chloromethane1.15Vinyl chloride1.15Bromomethane2.88	U U U J	Tetrachloroeth 1,3-Dichloropi			
Vinyl chloride1.15Bromomethane2.88	U UJ	1,3-Dichloropi			1 15 TT
Bromomethane 2.88	υΣ		onana		1.15 0
where a second sec		Dibromochlor	opane		1.15 U
Chloroethane 115	U	Dioronitoritori	omethane		1.15 U
Chloroethane 1.15		1,2-Dibromoet	hane		1.15 U
Trichlorofluoromethane 1.15	U	Chlorobenzene	e		1.15 U
Acetone 9.63	B O 2	1,1,1,2-Tetracl	nloroethane		1.15 U
1,1-Dichloroethene 1.15	U	Ethylbenzene			1.15 U
Carbon disulfide 0.64	J	p/m-Xylene			2.30 U
Methylene chloride 2.88	Ų	o-Xylene			1.15 U
Methyl tert-butyl ether (MTBE) 1.15	U	Styrene			1.15 U
trans-1,2-Dichloroethene 1.15	U	Bromoform			1.15 U
1,1-Dichloroethane 1.15	U	Isopropylbenz	ene		1.15 U
Vinyl acetate 1.15	U	1,1,2,2-Tetracl			1.15 U
2-Butanone (MEK) 1.72		Bromobenzene)		1.15 U
cis-1,2-Dichloroethene 1,15	U	1,2,3-Trichlord	opropane		1.15 U
2,2-Dichloropropane 1.15	U	n-Propylbenze	ne		1.15 U
Chloroform 1.15	U	2-Chlorotoluer	ne		1.15 U
1,1,1-Trichloroethane 1.15	U	1,3,5-Trimethy	/lbenzene		1.15 U
1,1-Dichloropropene 1.15	U	4-Chlorotoluer	10		1.15 U
Carbon tetrachloride 1.15		tert-Butylbenz	ene		1.15 U
Benzene 1.15	U	1,2,4-Trimethy	/lbenzene		1.15 U
1,2-Dichloroethane 1.15	U	sec-Butylbenz	ene		1.15 U
Trichloroethene 1.15	U	1,3-Dichlorobe	enzene		1.15 U
1,2-Dichloropropane 1.15	U	p-Isopropyltol	uene		1.15 U
Dibromomethane 1.15	U	1,4-Dichlorobe			1.15 U
Bromodichloromethane 1.15	U	n-Butylbenzen	C		1.15 U
2-Chloroethylvinyl ether 1.15	U	1,2-Dichlorobe	enzene		1.15 U
Methyl isobutyl ketone (MIBK) 1.15	U	1,2-Dibromo-3	-chloropropane		1.15 U
cis-1,3-Dichloropropene 1.15	U	1,2,4-Trichlord			1.15 U
Toluene 0.71	J	Hexachlorobut	tadiene	10 00 0 00 0 0 0 0 0 0 0 0 0 0 0 0 0 0	1.15 U
trans-1,3-Dichloropropene 1.15		Naphthalene		a West all	1.15 U
1,1,2-Trichloroethane 1.15		1,2,3-Trichloro	obenzene		1.15 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	101	70-130
1,2-Dichloroethane-d4	98	70-130
Toluene-d8	94	70-130
4-Bromofluorobenzene	86	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 16:13



ARCADIS Tidewater MGP

SC5-S1

SDG: N/A

Sediment

N/A

Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-12 Associated Blank: VS072108B21 Concentration Units: µg/Kg

Date CollectedDate ReceivedDate AnalyzedPercent SolidAmount (g)Volume (ml)Factor07/09/0807/11/0807/21/08 $\$1.5$ 9.32511ParameterResultParameterParameterFDichlorodifluoromethane1.32U2-Hexanone1Chloromethane1.32UTetrachloroethene1Vinyl chloride1.32U1,3-Dichloropropane1Bromomethane3.29UDibromochloromethane1Chloroethane1.32U1,2-Dibromochane1Trichlorofluoromethane1.32UChlorobenzeneAcetone13.4 \cancel{P} (U)1,1,1,2-Tetrachloroethane1,1-Dichloroethene1.32UStyreneI,1-Dichloroethene1.32UStyreneMethyl tert-butyl ether (MTBE)1.32U1,1-Dichloroethane1.32U1,1-Dichloroethane1.32U1,1-Dichloroethane1.32U1,1-Dichloroethane1.32U1,1-Dichloroethane1.32U1,1-Dichloroethane1.32U1,1-Dichloroethane1.32U1,2-Dichloroethane1.32U1,1-Dichloroethane1.32U1,2-Dichloroethane1.32U2,2-Dichloropropane1.32U2,2-Dichloropropane1.32U2,2-Dichloropropane1.32U2,2-Dichloropropane1.32U<	Analyst MLR Result 1.32 U 1.32 U
ParameterResultParameterFDichlorodifluoromethane1.32U2-HexanoneChloromethane1.32UTetrachloroetheneVinyl chloride1.32U1,3-DichloropropaneBromomethane3.29UDibromochloromethaneChloroethane1.32U1,2-DibromochlaromethaneChloroethane1.32U1,2-DibromochlaromethaneChloroethane1.32UChlorobenzeneAcetone13.4ØJ1,1-Dichloroethene1.32UCarbon disulfide2.03p/m-XyleneMethylene chloride3.29UO-Xylene1.32UStyrene1.32UStyrene1.32UCarbon disulfide3.29UO-Xylene1.32UStyrene1.32UStyrene1.32UStyrene1.32UStyrene1.1-Dichloroethene1,1-Dichloroethene1.32UStyrene1.1.2,2-Tetrachloroethane1,1-Dichloroethene1.32UStyrene1.1.2,2-Tetrachloroethane2-Butanone (MEK)1.32U1,32U1,2,3-Trichloroethane2,2-Dichloroethene1.32U2,2-Dichloroethene1.32U2,2-Dichloroethene1.32U2,2-Dichloroethene1.32U2,2-Dichloroethene1.32U2,2-Dichloroethene1.32U	Result 1.32 U
Dichlorodifluoromethane1.32U2-HexanoneChloromethane1.32UTetrachloroethene1Vinyl chloride1.32U1,3-DichloropropaneBromomethane3.29UJDibromochloromethaneChloroethane1.32U1,2-DibromoethaneTrichlorofluoromethane1.32UChlorobenzeneAcetone13.4B'UJ1,1,1,2-Tetrachloroethane1,1-Dichloroethene1.32UEthylbenzene2,1-Dichloroethene1.32UOMethylene chloride3.29Uo-XyleneMethyl tert-butyl ether (MTBE)1.32UStyrene1,1-Dichloroethane1.32UStyrenetrans-1,2-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32UIsopropylbenzene2.Butanone (MEK)1.32UBromobenzene2.2-Dichloroethene1.32Un-Propylbenzene2.3-Dichloroethene1.32UIsopropylbenzene1.32U1,2,2-Tetrachloroethane1.322.Butanone (MEK)1.32Un-Propylbenzene2.2-Dichloroethene1.32Un-Propylbenzene2.2-Dichloroethene1.32Un-Propylbenzene2.2-Dichloroethene1.32Un-Propylbenzene2.2-Dichloroethene1.32Un-Propylbenzene2.2-Dichloroethene1.32Un-Propylbenzene2.2-Dichloroethene1.32U	1.32 U
Chloromethane 1.32 UTetrachloroetheneVinyl chloride 1.32 U $1,3$ -DichloropropaneBromomethane 3.29 UDibromochloromethaneChloroethane 1.32 U $1,2$ -DibromocthaneChlorofluoromethane 1.32 U $1,2$ -DibromocthaneTrichlorofluoromethane 1.32 U $1,2$ -DibromocthaneAcetone 1.34 B' U $1,1,1,2$ -Tetrachloroethane $1,1$ -Dichloroethene 1.32 UEthylbenzene $1,1$ -Dichloroethene 1.32 UEthylbenzeneMethylene chloride 3.29 U o -XyleneMethyl tert-butyl ether (MTBE) 1.32 UStyrene $1,1$ -Dichloroethane 1.32 UBromoform $1,1$ -Dichloroethane 1.32 UBromoform $1,1$ -Dichloroethane 1.32 UIsopropylbenzene 2 -Butanone (MEK) 1.32 UBromobenzene $2,2$ -Dichloroethene 1.32 Un-Propylbenzene $2,2$ -Dichloroethene 1.32 Un-Propylbenzene $2,2$ -Dichloroethene 1.32 U $1,2,3$ -Trichloropropane $2,2$ -Dichloroethene 1.32 U 2 -Chlorotoluene	1.32 U
Vinyl chloride1.32U1,3-DichloropropaneBromomethane 3.29 UDibromochloromethaneChloroethane 1.32 U1,2-DibromochloromethaneTrichlorofluoromethane 1.32 UChlorobenzeneAcetone 13.4 $B' \cup J$ $1,1,1,2$ -Tetrachloroethane1,1-Dichloroethene 1.32 UEthylbenzene1,1-Dichloroethene 1.32 UEthylbenzeneCarbon disulfide 2.03 $p/m-Xylene$ Methylene chloride 3.29 U $o-Xylene$ Methyl tetr-butyl ether (MTBE) 1.32 UI,1-Dichloroethene 1.32 UJopropylbenzene 1.32 UJopropylbenzene 1.32 UStyrene 1.32 UStyrene 1.32 UStyrene 1.32 UJopropylbenzene 1.32 UJopropylbe	1.32 U
Bromomethane3.29UDibromochloromethaneChloroethane1.32U1,2-DibromocthaneTrichlorofluoromethane1.32UChlorobenzeneAcetone13.4BU1,1,1,2-Tetrachloroethane1,1-Dichloroethene1.32UEthylbenzene1,1-Dichloroethene1.32UEthylbenzeneCarbon disulfide2.03p/m-XyleneMethylene chloride3.29Uo-XyleneMethyl tert-butyl ether (MTBE)1.32UStyrene1,1-Dichloroethene1.32UBromoform1,1-Dichloroethene1.32UBromoform1,1-Dichloroethene1.32UIsopropylbenzene2-Butanone (MEK)1.32UBromobenzene2,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-Propylbenzene1.32U1,2,2-Tetrachloropropane1.32U1,2,3-Trichloropropane	1.32 U
Chloroethane1.32U1,2-DibromoethaneTrichlorofluoromethane1.32UChlorobenzeneAcetone13.4 $\not{B}' \cup J$ 1,1,1,2-Tetrachloroethane1,1-Dichloroethene1.32UEthylbenzeneQarbon disulfide2.03 p/m -XyleneMethylene chloride3.29UMethyl tert-butyl ether (MTBE)1.32U1,1-Dichloroethene1.32UMethyl tert-butyl ether (MTBE)1.32U1,1-Dichloroethene1.32U1,1-Dichloroethene1.32U1,2-Dichloroethene1.32U1,1-Dichloroethane1.32U1,2-Dichloroethane1.32U1,1-Dichloroethane1.32U1,2-Dichloroethane1.32U2-Butanone (MEK)1.32U2,2-Dichloroethene1.32U1,32U1,2,3-Trichloroethane2,2-Dichloropropane1.32U1,32U2-Chlorotoluene	1.32 U 1.32 U 1.32 U 1.32 U 2.63 U 1.32 U
Trichlorofluoromethane1.32UChlorobenzeneAcetone13.4 $\mathbb{B}' \cup \mathbb{J}$ 1,1,1,2-Tetrachloroethane1,1-Dichloroethene1.32UEthylbenzeneCarbon disulfide2.03 p/m -XyleneMethylene chloride3.29UOrthyl tert-butyl ether (MTBE)1.32Utrans-1,2-Dichloroethene1.32U1,1-Dichloroethene1.32Utrans-1,2-Dichloroethene1.32U1,1-Dichloroethene1.32U1,2-Dichloroethene1.32U1,2-Dichloroethene1.32U1,32UIsopropylbenzeneVinyl acetate1.32U2-Butanone (MEK)1.32U2,2-Dichloroethene1.32U1,32UPromobenzene2,2-Dichloropropane1.32U1,32U2-Chloropropane1,32U2-Chlorotoluene	1.32 U 1.32 U 1.32 U 2.63 U 1.32 U
Acetone13.4B'UI,1,1,2-Tetrachloroethane1,1-Dichloroethene1.32UEthylbenzeneCarbon disulfide2.03p/m-XyleneMethylene chloride3.29Uo-XyleneMethyl tert-butyl ether (MTBE)1.32UStyrenetrans-1,2-Dichloroethene1.32UBromoform1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32UIsopropylbenzene2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32UI.2,3-Trichloropropane1.32U1.2,2-Tetrachloropropane1.32U1.2,2-Tetrachloropropane1.32U2.3-Trichloropropane1.32U2.2-Chlorotoluene	1.32 U 1.32 U 2.63 U 1.32 U
1,1-Dichloroethene1.32UEthylbenzeneCarbon disulfide2.03p/m-XyleneMethylene chloride3.29Uo-XyleneMethyl tert-butyl ether (MTBE)1.32UStyrenetrans-1,2-Dichloroethene1.32UBromoform1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32UIsopropylbenzene2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32Un-Propylbenzene2-Butanone (MEK)1.32UI.2,3-Trichloropropane1.32U1.2,2-Cethoropropane1.2,2-Cethorotoluene	1.32 U 2.63 U 1.32 U 1.32 U 1.32 U 1.32 U 1.32 U 1.32 U
Carbon disulfide2.03p/m-XyleneMethylene chloride3.29Uo-XyleneMethyl tert-butyl ether (MTBE)1.32UStyrenetrans-1,2-Dichloroethene1.32UBromoform1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32UIsopropylbenzene2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32Un-Propylbenzene2,2-Dichloropropane1.32U2.3-Trichloropropane1.32U2.2-Chlorotoluene1.32U	2.63 U 1.32 U 1.32 U 1.32 U 1.32 U 1.32 U 1.32 U
Methylene chloride3.29Uo-XyleneMethyl tert-butyl ether (MTBE)1.32UStyrenetrans-1,2-Dichloroethene1.32UBromoform1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32U1,1,2,2-Tetrachloroethane2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	1.32 U 1.32 U 1.32 U 1.32 U 1.32 U
Methyl tert-butyl ether (MTBE)1.32UStyrenetrans-1,2-Dichloroethene1.32UBromoform1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32U1,1,2,2-Tetrachloroethane2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	1.32 U 1.32 U 1.32 U
trans-1,2-Dichloroethene1.32UBromoform1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32U1,1,2,2-Tetrachloroethane2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	1.32 U 1.32 U
1,1-Dichloroethane1.32UIsopropylbenzeneVinyl acetate1.32U1,1,2,2-Tetrachloroethane2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	1.32 U
Vinyl acetate1.32U1,1,2,2-Tetrachloroethane2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	
2-Butanone (MEK)1.32UBromobenzenecis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	1.32 U
cis-1,2-Dichloroethene1.32U1,2,3-Trichloropropane2,2-Dichloropropane1.32Un-PropylbenzeneChloroform1.32U2-Chlorotoluene	
2,2-Dichloropropane 1.32 U n-Propylbenzene Chloroform 1.32 U 2-Chlorotoluene	1.32 U
Chloroform 1.32 U 2-Chlorotoluene	1.32 U
	1.32 U
1,1,1-Trichloroethane 1.32 U 1,3,5-Trimethylbenzene	1.32 U
	1.32 U
1,1-Dichloropropene 1.32 U 4-Chlorotoluene	1.32 U
Carbon tetrachloride 1.32 U tert-Butylbenzene	1.32 U
Benzene 1.32 U 1,2,4-Trimethylbenzene	1.32 U
1,2-Dichloroethane 1.32 U sec-Butylbenzene	1.32 U
Trichloroethene 1.32 U 1,3-Dichlorobenzene	1.32 U
1,2-Dichloropropane 1.32 U p-Isopropyltoluene	1.32 U
Dibromomethane 1.32 U 1,4-Dichlorobenzene	1.32 U
Bromodichloromethane 1.32 U n-Butylbenzene	1.32 U
2-Chloroethylvinyl ether 1.32 U 1,2-Dichlorobenzene	1.32 U
Methyl isobutyl ketone (MIBK) 1.32 U 1,2-Dibromo-3-chloropropane	1.32 U
cis-1,3-Dichloropropene 1.32 U 1,2,4-Trichlorobenzene	1.32 U
Toluene 1.32 U Hexachlorobutadiene	1.32 U
trans-1,3-Dichloropropene 1.32 U Naphthalene	
1,1,2-Trichloroethane 1.32 U 1,2,3-Trichlorobenzene	1.32 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	101	70-130
Toluene-d8	93	70-130
4-Bromofluorobenzene	86	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

07/31/08 16:13



ARCADIS Tidewater MGP

SC6-S1

Sediment

N/A

Matrix:

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-13 Associated Blank: VS072108B21 Concentration Units: µg/Kg

				······	·····	Concentration	Units: μ	g/Kg
				Sample	Final	Dilution		
Date Collected	Date Received		Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/09/08	07/11/08	07/21/08	30.9	6.56	5	1	MLI	3
Parameter		F	Result	Parameter			Result	
Dichlorodifluoro	omethane		4.94 U	2-Hexanone			4.94	U
Chloromethane			4.94 U	Tetrachloroeth	ene		4.94	U
Vinyl chloride			4.94 U	1,3-Dichlorop	ropane		4.94	U
Bromomethane			<u>12.3 U J</u>	Dibromochlor	omethane		4.94	U
Chloroethane			4.94 U	1,2-Dibromoe	thane		4,94	U
Trichlorofluoror	nethane		4.94 U	Chlorobenzen	3	1	4.94	U
Acetone			490 B J	1,1,1,2-Tetrac	hloroethane		4.94	U
1,1-Dichloroethe			<u>4.94 U</u>	Ethylbenzene			4.94	U
Carbon disulfide			67.2	p/m-Xylene			9.87	U
Methylene chlor	and the second sec		12.3 U	o-Xylene			4.94	U
Methyl tert-buty			4,94 U	Styrene	· · · · · · · · · · · · · · · · · · ·		4.94	U
	,2-Dichloroethene 4.94 U Bromoform			4.94	U			
	-Dichloroethane 4.94 U		Isopropylbenz			4.94	U	
Vinyl acetate			<u>4.94 U</u>	1,1,2,2-Tetrac			4.94	U
2-Butanone (ME	the second secon	.)-	138	Bromobenzen			4.94	
cis-1,2-Dichloro			4.94 U	1,2,3-Trichlor			4.94	U
2,2-Dichloropro	pane	·····	<u>4.94 U</u>	n-Propylbenze			4,94	U
Chloroform			<u>4.94 U</u>	2-Chlorotoluene			4,94	U
1,1,1-Trichloroe			<u>4.94 U</u>	1,3,5-Trimethylbenzene			4.94	Ų
1,1-Dichloropro			4,94 U	4-Chlorotoluene			4.94	U
Carbon tetrachlo	oride		4.94 U	tert-Butylbenzene			4.94	U
Benzene			<u>4.94 U</u>	1,2,4-Trimethylbenzene			4.94	U
1,2-Dichloroetha	ane	······································	4.94 U	sec-Butylbenzene			4.94	U
Trichloroethene			<u>4.94 U</u>	1,3-Dichlorobenzene			4.94	U
1,2-Dichloropro	pane		<u>4.94 U</u>	p-Isopropyltoluene			4.94	U
Dibromomethan	e		4.94 U	1,4-Dichlorobenzene			4,94	U
Bromodichloron	nethane		<u>4.94 U</u>	n-Butylbenzene			4.94	U
	-Chloroethylvinyl ether 4.94 U		1,2-Dichlorobenzene			4.94		
Methyl isobutyl			4.94 U	1,2-Dibromo-3-chloropropane			4.94	
cis-1,3-Dichloro	propene		<u>4.94 U</u>	1,2,4-Trichlorobenzene			4,94	U
Toluene			4.94 U	Hexachlorobutadiene			4.94	U
trans-1,3-Dichlo			4.94 U	Naphthalene			4.94	U
1,1,2-Trichloroe	thane		4.94 U	1,2,3-Trichlor	obenzene		4.94	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	100	70-130
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	82	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 16:13



ARCADIS **Tidewater MGP**

SC20-S1

Sediment

N/A

Matrix:

SDG: N/A Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-14 Associated Blank: VS072108B21 Concentration Units: $\mu g/Kg$

						Concontration	
			1	Sample	Final	Dilution	
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/21/08	77.6	8.49	5	1	MLR
Parameter	<u> </u>	F	Result	Parameter			Result
Dichlorodifluor	methane		1.52 U	2-Hexanone	 _ _ _ _ _ _ _ _		1.52 U
Chloromethane			<u>1.52 U</u>	Tetrachloroeth			1.52 U
Vinyl chloride			1.52 U	1,3-Dichlorop			1.52 U
Bromomethane			3.79 U J	Dibromochlor		L.	1.52 U
Chloroethane			1.52 U	1,2-Dibromoe			1.52 U
Trichlorofluoron	nethane		1.52 U	Chlorobenzen			1.52 U
Acetone			16.8 BUJ	1,1,1,2-Tetrac	hloroethane		1.52 U
1.1-Dichloroeth	ene		1.52 U	Ethylbenzene	· · · ·		1.52 U
Carbon disulfide			10.6	p/m-Xylene			3.03 U
Methylene chlor			3.79 U	o-Xylene			1.52 U
Methyl tert-buty			1.52 U	Styrene			1.52 U
trans-1,2-Dichlo			1.52 U	Bromoform			1.52 U
1,1-Dichloroeth				Isopropylbenz			<u>1.52 U</u> 1.52 U
Vinyl acetate			1,1,2,2-Tetrac	1,1,2,2-Tetrachloroethane			
2-Butanone (MI	EK)		2.02	Bromobenzen	e		1.52 U
cis-1,2-Dichloro			1.52 U	1,2,3-Trichlor	opropane		1,52 U
2,2-Dichloropro			1.52 U	n-Propylbenzene			1.52 U
Chloroform			1.52 U	2-Chlorotolue	ne		1.52 U
1,1,1-Trichloroe	thane		1.52 U	1,3,5-Trimeth	ylbenzene		<u>1.52 U</u>
1,1-Dichloropro			1.52 U	4-Chlorotoluene			<u>1.52 U</u>
Carbon tetrachle			1.52 U	tert-Butylbenzene			1.52 U
Benzene			1.52 U	1,2,4-Trimeth	ylbenzene		1.52 U
1.2-Dichloroeth	ane		1.52 U	sec-Butylbenz			<u>1.52 U</u> 1.52 U
Trichloroethene	the second se		1.52 U	1,3-Dichlorob	1,3-Dichlorobenzene		
1,2-Dichloropro			1.52 U	p-Isopropylto	luene		1.52 U
Dibromometha			1.52 U	1,4-Dichlorobenzene			<u>1.52 U</u>
Bromodichloro			1.52 U	n-Butylbenzene			<u>1.52 U</u>
2-Chloroethylvi			1.52 U	1,2-Dichlorob			1.52 U
	ketone (MIBK)		1.52 U	1,2-Dibromo-	3-chloropropane		1.52 U
cis-1,3-Dichlor			1.52 U		robenzene		1.52 U
Toluene			1.07 J	Hexachlorobu	itadiene		1.52 U
······································	oropropene		1.52 U	Naphthalene			1.52 U
1,1,2-Trichloro			1.52 U	1,2,3-Trichlor			1.52 U
1,1,2-111011010							

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	107	70-130
1.2-Dichloroethane-d4	101	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	91	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 16:13



ARCADIS Tidewater MGP

SC26-S1

N/A

Sediment

Matrix:

SDG: N/A

Lab Code: **MA00030** ETR: **0807059** Lab ID: **0807059-15** Associated Blank: **VS072108B21** Concentration Units: μg/Kg

[Sample	Final	Dilution	Amalyast
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/21/08	83.5	8.07	5	1	MLR
Parameter		F	Result	Parameter			Result
Dichlorodifluoro	methane		1,48 U	2-Hexanone			1.48 U
Chloromethane			1,48 U	Tetrachloroeth			<u>1.48 U</u>
Vinyl chloride			<u>1,48 U</u>	1,3-Dichlorop			<u>1.48 U</u>
Bromomethane			3.71 U J	Dibromochlor			1.48 U
Chloroethane			1.48 U	1,2-Dibromoe	thane		1.48 U
Trichlorofluoron	nethane		1.48 U	Chlorobenzen			1.48 U
Acetone			25.8 B U J	1,1,1,2-Tetrac	hloroethane		1.48 U
1,1-Dichloroethe	ene		1.48 U	Ethylbenzene			1,48 U
Carbon disulfide			1.65	p/m-Xylene			2.97 U
Methylene chlor			3.71 U	o-Xylene			1,48 U
Methyl tert-buty			1.48 U	Styrene			1.48 U
trans-1.2-Dichlo			1.48 U	Bromoform			1.48 U
1.1-Dichloroetha			1.48 U	Isopropylbenzene			1.48 U
Vinyl acetate			1.48 U	1,1,2,2-Tetrachloroethane			<u>1.48 U</u>
2-Butanone (MI	 EK)		2.79	Bromobenzene			1.48 U
cis-1,2-Dichloro			1.48 U	1,2,3-Trichlor	opropane		1.48 U
2,2-Dichloropro			1.48 U	n-Propylbenzo			1.48 U
Chloroform			1.48 U	2-Chlorotolue	ene		1.48 U
1,1,1-Trichloroe	ethane		1.48 U	<u>1,3,5-Trimeth</u>			1.48 U
1,1-Dichloropro			1.48 U	4-Chlorotolue			1.48 U
Carbon tetrachle			1.48 U	tert-Butylben:	zene	<u></u>	1.48 U
Benzene			1.48 U	1,2,4-Trimeth	ylbenzene		1.48 U
1.2-Dichloroeth	ane		1.48 U	sec-Butylben:			1.48 U
Trichloroethene			1.48 U	1,3-Dichlorot			1.48 U
1,2-Dichloropro			1.48 U	p-Isopropylto			1.48 U
Dibromomethar			1.48 U	1,4-Dichlorot			1.48 U
Bromodichloron			1.48 U	n-Butylbenze	ne		1.48 U
2-Chloroethylvi			1,48 U	1,2-Dichlorol			1.48 U
	ketone (MIBK)		1.48 U	1,2-Dibromo-3-chloropropane			1.48 U
cis-1,3-Dichlor			1.48 U	1,2,4-Trichlorobenzene			<u>1,48 U</u>
Toluene			1.48 U	Hexachloroby	utadiene	· · · · · · · · · · · · · · · · · · ·	1.48 U
trans-1.3-Dichle	oropropene		1,48 U	Naphthalene		······	1.48 U
1,1,2-Trichloro			1.48 U	1,2,3-Trichlo	robenzene		1.48 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	104	70-130
1.2-Dichloroethane-d4	103	70-130
Toluene-d8	93	70-130
4-Bromofluorobenzene	87	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 16:13



ARCADIS **Tidewater MGP**

SC19-S1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-16 Associated Blank: VS072108B21 Concentration Units: µg/Kg

					·	Concentration	unta. µg/ng
				Sample	Final	Dilution	
Date Collected	Date Received		Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/21/08	35.7	6.79	5	1	MLR
Parameter		I	Result	Parameter			Result
Dichlorodifluor	omethane		<u>4.13 U</u>	2-Hexanone			4.13 U
Chloromethane			4.13 U	Tetrachloroeth			4.13 U
Vinyl chloride			4.13 U	1,3-Dichlorop			4.13 U
Bromomethane			10.3 UJ	Dibromochlor			4.13 U
Chloroethane			4.13 U	1,2-Dibromoe			4,13 U
Trichlorofluoro	methane		4.13 U	Chlorobenzen			4.13 U
Acetone			379 B J	1,1,1,2-Tetrac	hloroethane		4.13 U
1,1-Dichloroeth	ene		4.13 U	Ethylbenzene			4.13 U
Carbon disulfide	e		44.6	p/m-Xylene			8.26 U
Methylene chlor	ride		10.3 U	o-Xylene			4.13 U
Methyl tert-buty	l ether (MTBE)		4.13 U	Styrene		L 11	4.13 U
trans-1,2-Dichle	roethene		4.13 U	Bromoform			4.13 U
1,1-Dichloroeth	,1-Dichloroethane 4.13 U		Isopropylbenzene			4.13 U	
Vinyl acetate			4.13 U	1,1,2,2-Tetrachloroethane			<u>4.13 U</u>
2-Butanone (MI	EK)		110	Bromobenzene			4.13 U
cis-1,2-Dichloro	ethene		4.13 U	1,2,3-Trichloropropane			4.13 U
2,2-Dichloropro	pane		4.13 U	n-Propylbenzene			4.13 U
Chloroform			<u>4.13 U</u>	2-Chlorotoluene			4.13 U
1,1,1-Trichloroe	ethane		4.13 U	1,3,5-Trimethylbenzene			4.13 U
1,1-Dichloropro	pene		4.13 U	4-Chlorotoluene			4.13 U
Carbon tetrachle			4.13 U	tert-Butylbenzene			4.13 U
Benzene			4.13 U	1,2,4-Trimethylbenzene			4.13 U
1,2-Dichloroeth	ane		4.13 U	sec-Butylbenz			4.13 U
Trichloroethene			4.13 U	1,3-Dichlorobenzene			4.13 U
1,2-Dichloropro	pane		4.13 U	p-Isopropyltoluene			4.13 U
Dibromomethan			4.13 U	1,4-Dichlorobenzene			4.13 U
Bromodichloro	nethane		4.13 U	n-Butylbenzene			4.13 U
2-Chloroethylvi	Chloroethylvinyl ether 4.13 U		1,2-Dichlorobenzene			4.13 U	
Methyl isobutyl	ethyl isobutyl ketone (MIBK) 4.13 U		1,2-Dibromo-3-chloropropane			4.13 U	
cis-1,3-Dichloro			4.13 U	1,2,4-Trichlor	obenzene		4.13 U
Toluene			4.13 U	Hexachlorobu	tadiene		4.13 U
trans-1,3-Dichlo			4.13 U	Naphthalene			4.13 U
1,1,2-Trichloroe			4.13 U	1,2,3-Trichlor	obenzene		4.13 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	100	70-130
1,2-Dichloroethane-d4	101	70-130
Toluene-d8	91	70-130
4-Bromofluorobenzene	99	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SC43-S1

Sediment

N/A

Matrix:

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-17 Associated Blank: VS072108B21 Concentration Units: µg/Kg

				Sample	Final	Dilution	
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/21/08	62.0	7.62	5	1	MLR
Parameter		I	Result	Parameter			Result
Dichlorodifluoro	methane		2.12 UJ	2-Hexanone			<u>2.12 U J</u>
Chloromethane			2.12 U	Tetrachloroeth	iene		2.12 U
Vinyl chloride			2.12 U	1,3-Dichlorop			2.12 U
Bromomethane			5.29 U	Dibromochlor	omethane	a a cara ta a cara a	<u>2.12 U</u>
Chloroethane			<u>2.12 U</u>	1,2-Dibromoe	thane		<u>2.12 U</u>
Trichlorofluoron	nethane		2.12 U	Chlorobenzen			2.12 U
Acetone			190 B	1,1,1,2-Tetrac	hloroethane		2.12 U
1,1-Dichloroethe	ene		2.12 U	Ethylbenzene			2.12 U
Carbon disulfide			45.0	p/m-Xylene			4.23 U
Methylene chlor	ide		5.29 U	o-Xylene			2.12 U
Methyl tert-buty	l ether (MTBE)		2,12 U	Styrene			2.12 U
trans-1,2-Dichlo	roethene		2.12 U	Bromoform			2.12 U
1,1-Dichloroetha	oethane 2.12 U Isopropylbenzene				2.12 U		
Vinyl acetate			2.12 U	1.1.2.2-Tetrachloroethane			2.12 U
2-Butanone (ME	EK)		57.3	Bromobenzene		2.12 U	
cis-1,2-Dichloro	ethene		2.12 U	1,2,3-Trichlor	opropane		2.12 U
2,2-Dichloropro	pane		2.12 U	n-Propylbenze	ene		2.12 U
Chloroform			2,12 U	2-Chlorotolue	ne		2.12 U
1,1,1-Trichloroe	thane		2.12 U	1,3,5-Trimethy	ylbenzene		2.12 U
1,1-Dichloropro	pene		2.12 U	4-Chlorotolue	ne		2,12 U
Carbon tetrachlo	ride		2.12 U	tert-Butylbenz	ene		2.12 U
Benzene			2.12 U	1,2,4-Trimethy	ylbenzene		2.12 U
1,2-Dichloroetha	ine		2.12 U	sec-Butylbenz	ene		2.12 U
Trichloroethene			2.12 U	1,3-Dichlorob	enzene		2.12 U
1,2-Dichloropro	pane		2.12 U	p-Isopropyltol	uene		2.12 U
Dibromomethan	e		2.12 U	1,4-Dichlorobenzene			2.12 U
Bromodichloron	nethane		2.12 U	n-Butylbenzene			2.12 U
2-Chloroethylvin	nyl ether		2.12 U	1,2-Dichlorobenzene		2.12 U	
Methyl isobutyl			2.12 U	1,2-Dibromo-:	3-chloropropane		2.12 U
cis-1,3-Dichloro			2.12 U	1,2,4-Trichlor			2.12 U
Toluene			2.12 U	Hexachlorobu	tadiene		2.12 U
trans-1,3-Dichlo	ropropene		2.12 U	Naphthalene			2.12 U
1,1,2-Trichloroe			2.12 U 🖢	1,2,3-Trichlor	obenzene		2.12 U

Surrogate	% Recovery		Acceptance Range (%)
Dibromofluoromethane	94		70-130
1,2-Dichloroethane-d4	82		70-130
Toluene-d8	84		70-130
4-Bromofluorobenzene	64	§	70-130

N/A - Not Applicable

 \S - Surrogate value outside of acceptable range.

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 16:13



ARCADIS Tidewater MGP

SC42-S1

Sediment

N/A

Matrix:

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-18 Associated Blank: VS072108B21 Concentration Units: µg/Kg

				Sample	Final	Dilution		
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst	
07/10/08	07/11/08	07/21/08	77.3	8.98	5	1	MLR	
Parameter		I	Result	Parameter			Result	
Dichlorodifluoro	methane		1.44 U	2-Hexanone			1.44 U	I
Chloromethane			<u>1.44 U</u>	Tetrachloroeth	ene		<u>1.44 U</u>	J
Vinyl chloride		1015	<u>1.44 U</u>	1,3-Dichlorop	opane		<u> </u>	1
Bromomethane			<u>3.60 U J</u>	Dibromochlor	omethane		<u>1.44 U</u>	J
Chloroethane			1.44 U	1,2-Dibromoe	hane		<u>1.44 U</u>	J
Trichlorofluoror	nethane		1.44 U	Chlorobenzene	;		<u>1.44 U</u>	J
Acetone			111 B J	1,1,1,2-Tetracl	nloroethane		1.44 U	J
1,1-Dichloroethe	ene		1.44 U	Ethylbenzene			1.44 U	J
Carbon disulfide	;		19.9	p/m-Xylene			2.88 U	I
Methylene chlor	ide		3.60 U	o-Xylene			1.44 U	ľ
Methyl tert-buty	l ether (MTBE)		1.44 U	Styrene			1.44 U	J
trans-1,2-Dichlo	roethene		1.44 U	Bromoform			1.44 U	J
1,1-Dichloroetha						1.44 U	J	
Vinyl acetate			1.44 U	1,1,2,2-Tetrachloroethane			1.44 U	I
2-Butanone (ME	EK)		30.4	Bromobenzene			1.44 Ŭ	J
cis-1,2-Dichloro	ethene		1.44 U	1,2,3-Trichlor	opropane		<u>1.44 U</u>	J
2,2-Dichloropro	pane		1.44 U	n-Propylbenze	ne		1.44 U	J
Chloroform			1.44 Ŭ	2-Chlorotolue	<u>1e</u>		1.44 U	J
1,1,1-Trichloroe	thane		1.44 U	1,3,5-Trimethy			<u>1.44</u> U	J
1,1-Dichloropro	pene		1.44 U	4-Chlorotolue	ne		1.44 U	J
Carbon tetrachlo	oride		<u>1.44 U</u>	tert-Butylbenz	ene		<u>1.44 U</u>	J
Benzene			1.44 U	1,2,4-Trimethy	lbenzene		<u>1.44 U</u>	J
1,2-Dichloroetha	ane		1.44 U	sec-Butylbenz	ene		<u>1.44 U</u>	<u>)</u>
Trichloroethene			<u>1.44 U</u>	1,3-Dichlorob	enzene	1 * * * * * * * * * * * * * * * * * * *	<u>1.44 U</u>	J
1,2-Dichloropro	pane		<u>1.44 U</u>	p-Isopropyltol	uene		1.44 U	J
Dibromomethan	e		1.44 U 1,4-Dich		enzene		<u>1.44 U</u>	J
Bromodichloron	nethane		<u>1.44 U</u>	n-Butylbenzene			<u> </u>	
2-Chloroethylvii	nyl ether		1.44 U	1,2-Dichlorob			<u>1.44 U</u>	
Methyl isobutyl	ketone (MIBK)		1.44 U	1,2-Dibromo-3-chloropropane			<u>1.44 U</u>	
cis-1,3-Dichloro	propene		1.44 U	1,2,4-Trichlorobenzene			<u> </u>	
Toluene			1.44 U	Hexachlorobu	tadiene		1.44 U	J
trans-1,3-Dichlo	ropropene		1.44 U	Naphthalene			1.44 U	J
1,1,2-Trichloroe	thane		1.44 U	1,2,3-Trichlor	obenzene		<u>1.44 U</u>	J

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	90	70-130
1,2-Dichloroethane-d4	82	70-130
Toluene-d8	91	70-130
4-Bromofluorobenzene	79	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS **Tidewater MGP**

SC9-S1

Sediment

N/A

Matrix:

SDG:

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-19 Associated Blank: VS072108B21 **Concentration Units:** µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	vet
07/10/08	07/11/08	07/22/08	36.6	7,31	5	1	MLI	
Parameter		l I	Result	Parameter			Result	
Dichlorodifluor	omethane		3.74 U	2-Hexanone			3.74	U
Chloromethane			3.74 U	Tetrachloroeth	nene		3.74	U
Vinyl chloride			3.74 U	1,3-Dichlorop	ropane		3.74	U
Bromomethane			9.34 U J	Dibromochlor	omethane		3.74	
Chloroethane			3.74 U	1,2-Dibromoe	thane		3.74	U
Trichlorofluoror	nethane		<u>3.74 U</u>	Chlorobenzen	e		3.74	U
Acetone			129 B J	1,1,1,2-Tetrac	hloroethane		3.74	U
1,1-Dichloroethe	ene		3.74 U	Ethylbenzene			3.74	U
Carbon disulfide	;		45.3	p/m-Xylene			7.47	U
Methylene chlor	ide		9.34 U	o-Xylene			3.74	U
Methyl tert-buty	l ether (MTBE)		3.74 U	Styrene			3.74	U
trans-1,2-Dichlo	roethene		3.74 U	Bromoform			3.74	U
1,1-Dichloroetha	ane		3.74 U	Isopropylbenz	ene		3.74	U
Vinyl acetate			3.74 U	1,1,2,2-Tetrac	hloroethane		3.74	υ
2-Butanone (ME	EK)		33.4	Bromobenzen	e		3.74	U
cis-1,2-Dichloro	ethene		3.74 U	1,2,3-Trichlor	opropane		3.74	U
2,2-Dichloropro	pane		3.74 U	n-Propylbenze			3.74	U
Chloroform			3.74 U	2-Chlorotolue	ne		3.74	U
1,1,1-Trichloroe	thane		3.74 U	1,3,5-Trimeth	ylbenzene		3,74	U
1,1-Dichloropro	репе		3.74 U	4-Chlorotolue	ne		3.74	U
Carbon tetrachlo	oride		3.74 U	tert-Butylbenz	ene		3.74	U
Benzene			3.74 U	1,2,4-Trimeth	ylbenzene		3.74	U
1,2-Dichloroetha	ane		3.74 U	sec-Butylbenz	ene		3.74	U
Trichloroethene			3.74 U	1,3-Dichlorob	enzene		3.74	U
1,2-Dichloropro	pane		3.74 U	p-Isopropyltol	uene		3.74	U
Dibromomethan	e		3.74 U	1,4-Dichlorob	enzene		3.74	U
Bromodichloron	nethane		3.74 U	n-Butylbenzer	1e	-	3.74	U
2-Chloroethylvi	nyl ether		3.74 U	1,2-Dichlorob	enzene		3.74	U
Methyl isobutyl	ketone (MIBK)		3.74 U	1,2-Dibromo-	3-chloropropane		3.74	U
cis-1,3-Dichloro			3.74 U	1,2,4-Trichlor			3.74	U
Toluene			3.74 U	Hexachlorobu	tadiene		3.74	U
trans-1,3-Dichlo	ropropene		3.74 U	Naphthalene			3.74	U
1,1,2-Trichloroe			3.74 U	1,2,3-Trichlor	obenzene		3.74	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	97	70-130
1,2-Dichloroethane-d4	87	70-130
Toluene-d8	84	70-130
4-Bromofluorobenzene	70	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS **Tidewater MGP**

DUP-1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-20 Associated Blank: VS072308B09 Concentration Units: μ<u>g</u>/Kg

	······································			Sample	Final	Dilution		g/itg
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	/st
07/10/08	07/11/08	07/23/08	36.2	7.68	5	1	MLI	3.
Parameter	· · · · · · · · · · · · · · · · · · ·	F	Result	Parameter			Result	
Dichlorodifluoro	methane		3.60 U	2-Hexanone			9.00	
Chloromethane			<u>3.60 U</u>	Tetrachloroeth	iene		3,60	U
Vinyl chloride			<u>3.60 U</u>	1,3-Dichlorop			3.60	
Bromomethane			3.60 U	Dibromochlor	omethane		3.60	U
Chloroethane			3.60 U	1,2-Dibromoe	thane	·····	3.60	
Trichlorofluoror	nethane		3.60 U	Chlorobenzen			3.60	
Acetone			298	1,1,1,2-Tetrac	hloroethane		3.60	
1,1-Dichloroethe	ene		3.60 U	Ethylbenzene			3.60	
Carbon disulfide	;		67.4	p/m-Xylene			7,20	U
Methylene chlor	ide		9.00 U	o-Xylene			3.60	U
Methyl tert-buty	l ether (MTBE)		3.60 U	Styrene			3.60	U
trans-1,2-Dichlo	roethene		3.60 U	Bromoform			3.60	U
1,1-Dichloroetha	ane		3.60 U	Isopropylbenz	ene		3.60	U
Vinyl acetate			3.60 U	1,1,2,2-Tetrac	hloroethane		3.60	U
2-Butanone (ME			75.5	Bromobenzen	e		3.60	U
cis-1,2-Dichloro	ethene		3.60 U	1,2,3-Trichloropropane			3.60	U
2,2-Dichloropro	pane		3.60 U	n-Propylbenzene			3.60	U
Chloroform			3.60 U	2-Chlorotoluene			3.60	U
1,1,1-Trichloroe	thane		3.60 U	1,3,5-Trimethylbenzene			3.60	U
1,1-Dichloropro	pene		3.60 U	4-Chlorotoluene			3.60	U
Carbon tetrachlo			3.60 U	tert-Butylbenzene			3.60	U
Benzene			3.60 U	1,2,4-Trimethy	ylbenzene		3.60	U
1,2-Dichloroetha			3.60 U	sec-Butylbenz			3.60	U
Trichloroethene			3.60 U	1,3-Dichlorob	enzene		3.60	U
1,2-Dichloropro			3.60 U	p-Isopropyltol	uene		9.00	
Dibromomethan			3.60 U	1,4-Dichlorob	enzene		3.60	U
Bromodichloron			3.60 U	n-Butylbenzer	-		0.00	U
2-Chloroethylvin			9.00 U	1,2-Dichlorob			2 (0	
Methyl isobutyl			3.60 U	1,2-Dibromo-	3-chloropropane		3.60	
cis-1,3-Dichloro			3.60 U	1,2,4-Trichlor	obenzene		3.60	U
Toluene			3.60 U	Hexachlorobu	tadiene		3.60	U
trans-1,3-Dichlo	ropropene		3.60 U	Naphthalene			9.00	
1,1,2-Trichloroe			3.60 U	1,2,3-Trichlor	obenzene		3.60	

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	90	70-130
1,2-Dichloroethane-d4	83	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	78	70-130

N/A - Not Applicable

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:16



ARCADIS Tidewater MGP

SC29-S1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-21 Associated Blank: VS072308B09 Concentration Units: µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
07/10/08	07/11/08	07/23/08	42.1	7.45	5	1	MLR
Parameter		J	Result	Parameter			Result
Dichlorodifluoro	methane		3.19 U	2-Hexanone			7.97 U
Chloromethane			3.19 U	Tetrachloroeth	ene		<u>3,19 U</u>
Vinyl chloride			3.19 U	1,3-Dichlorop			3.19 U
Bromomethane			3.19 U	Dibromochlor			<u>3.19 U</u>
Chloroethane			3.19 U	1,2-Dibromoe			3.19 U
Trichlorofluoron	nethane		<u>3.19 U</u>	Chlorobenzen			<u>3.19 U</u>
Acetone			182	1,1,1,2-Tetrac	hloroethane		3.19 U
1,1-Dichloroethe	ne		3.19 U	Ethylbenzene		and the first of groups and the	3.19 U
Carbon disulfide			57.8	p/m-Xylene			6.37 U
Methylene chlor	ide		7.97 U	o-Xylene			3.19 U
Methyl tert-buty	ether (MTBE)		3.19 U	Styrene			3.19 U
trans-1,2-Dichlor	roethene		3.19 U	Bromoform			3.19 U
1,1-Dichloroetha	ine		3.19 U	Isopropylbenz	ene		3.19 U
Vinyl acetate			3.19 U	1,1,2,2-Tetrac	hloroethane		3.19 U
2-Butanone (ME	X)		47.1	Bromobenzen	3		3.19 U
cis-1,2-Dichloro		-	3.19 U	1,2,3-Trichlor	opropane		3.19 U
2,2-Dichloroprop			3.19 U	n-Propylbenze	ne		3.19 U
Chloroform			3.19 U	2-Chlorotolue	ne		3.19 U
1,1,1-Trichloroe	thane		3.19 U	1,3,5-Trimethy	lbenzene		3.19 U
1,1-Dichloropro	oene		3.19 U	4-Chlorotolue	ne		3.19 U
Carbon tetrachlo	ride		3.19 U	tert-Butylbenz	ene		3.19 U
Benzene			3.19 U	1,2,4-Trimethy	vlbenzene		3.19 U
1,2-Dichloroetha	me		3.19 U	sec-Butylbenz	ene		3.19 U
Trichloroethene			3.19 U	1,3-Dichlorob	enzene		3.19 U
1,2-Dichloroproj			3.19 U	p-Isopropyltol	uene		7.97 U
Dibromomethan			3.19 U	1,4-Dichlorob			3.19 U
Bromodichlorom	nethane		3.19 U	n-Butylbenzer	e		7.97 U
2-Chloroethylvin	ivl ether		7.97 Ŭ	1,2-Dichlorob			3.19 U
Methyl isobutyl			3.19 U	1,2-Dibromo-	3-chloropropane		3.19 U
cis-1,3-Dichloro		· · · · · · · · · · · · · · · · · · ·	3.19 U	1,2,4-Trichlor			3.19 U
Toluene	F	- 1	3.19 U	Hexachlorobu	tadiene		3.19 U
trans-1,3-Dichlo	ropropene	4.,/14 (a) (a) (a)	3.19 U	Naphthalene			7.97 U
1,1,2-Trichloroe		v	3,19 U	1,2,3-Trichlor	obenzene		3.19 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	87	70-130
1,2-Dichloroethane-d4	78	70-130
Toluene-d8	89	70-130
4-Bromofluorobenzene	73	70-130

N/A - Not Applicable

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:17



ARCADIS Tidewater MGP

SC17-S1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-22 Associated Blank: VS072408B15

µg/Kg

Concentration Units:

Data Callestad	Deta Passiwad	Data Analyzad	Domo	nt Calid	Sample	Final	Dilution		<u>6/115</u>
Date Collected		Date Analyzed		nt Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/10/08	07/11/08	07/24/08		2.9	7.28	5	1	BAS	5
Parameter		F	Result		Parameter			Result	
Dichlorodifluoro	methane		2.60	U	2-Hexanone			6,49	Ų
Chloromethane			2.60	<u>U</u>	Tetrachloroeth	ene		2.60	U
Vinyl chloride			2.60	U	1,3-Dichloropi			2.60	U
Bromomethane			2.60	<u> </u>	Dibromochlor	omethane		· 2.60	U
Chloroethane			2.60	U	1,2-Dibromoet	hane		2.60	U
Trichlorofluoron	nethane		2.60	U	Chlorobenzene	<u> </u>		2.60	U
Acetone			205	<u>P</u>	1,1,1,2-Tetracl	nloroethane		2.60	U
1,1-Dichloroethe	ene		2.60	<u>´Ů</u>	Ethylbenzene			2.60	U
Carbon disulfide			20.7		p/m-Xylene			5.19	U
Methylene chlor	ide '		6.49	U	o-Xylene			2.60	U
Methyl tert-buty	l ether (MTBE)		2.60	U	Styrene			2.60	U
trans-1,2-Dichlo	roethene		2.60	U	Bromoform			2.60	U
1,1-Dichloroetha	ine		2.60	U	Isopropylbenze	ene		2.60	υ
Vinyl acetate			2.60	U	1,1,2,2-Tetracl	nloroethane		2.60	U
2-Butanone (ME	EK)		46.1	J	Bromobenzene	3		2.60	U
cis-1,2-Dichloro	ethene		2.60	U	1,2,3-Trichloro	opropane		2.60	U
2,2-Dichloropro	pane		2,60	U	n-Propylbenze	ne		2.60	U
Chloroform			2.60	U	2-Chlorotoluer	ie		2.60	U
1,1,1-Trichloroe	thane		2.60	U	1,3,5-Trimethy	lbenzene		2.60	U
1,1-Dichloropro	pene		2.60	U	4-Chlorotoluer	ne		2.60	U
Carbon tetrachlo	ride		2.60	U	tert-Butylbenz	ene		2.60	U
Benzene			2.60	U	1,2,4-Trimethy	lbenzene		2.60	U
1,2-Dichloroetha	ine		2,60	U	sec-Butylbenze	ene		2.60	U
Trichloroethene			2.60	U	1,3-Dichlorobe	enzene		2.60	U
1,2-Dichloropro	pane		2.60	U	p-Isopropyltoh	uene		6.49	U
Dibromomethan	e		2.60	U	1,4-Dichlorobe	enzene		2.60	U
Bromodichloron	nethane		2.60	U	n-Butylbenzen	e		6.49	U
2-Chloroethylvir	yl ether		6.49	U	1,2-Dichlorobe	enzene		2.60	U
Methyl isobutyl	ketone (MIBK)		2.60	U	1,2-Dibromo-3	-chloropropane		2.60	U
cis-1,3-Dichloro	propene		2.60	U	1,2,4-Trichloro			2.60	U
Toluene			2.60	U	Hexachlorobut	adiene		2.60	υ
trans-1,3-Dichlo	ropropene		2.60	U	Naphthalene			6.49	U
1,1,2-Trichloroe			2.60	Ų	1,2,3-Trichlord	benzene		2.60	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	105	70-130
1,2-Dichloroethane-d4	103	70-130
Toluene-d8	87	70-130
4-Bromofluorobenzene	77	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:25



ARCADIS **Tidewater MGP**

SC28-S1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-23 Associated Blank: VS072408B15 Concentration Units: µg/Kg

Date Collected	Date Received	Data Analyzad	Daraca	nt Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	<u>, 156</u>
		Date Analyzed 07/24/08			7.31	5			
07/10/08	07/11/08	07/24/08	،د	4.6	/.31	3	1	BAS	<u> </u>
Parameter		F	Result		Parameter			Result	
Dichlorodifluoro	methane		· · · · · · · · · · · · · · · · · · ·	U	2-Hexanone			9.88	U
Chloromethane				U	Tetrachloroeth			3.95	
Vinyl chloride				U	1,3-Dichlorop			3,95	
Bromomethane				UJ	Dibromochlor			3.95	
Chloroethane			3.95	U	1,2-Dibromoe	thane		3.95	<u>U</u>
Trichlorofluoron	nethane		3,95	U	Chlorobenzen	8		3.95	U
Apatona			280	<u>B</u> T	1,1,1,2-Tetrac	hloroethane		3.95	U
1,1-Dichloroethe	ene		<u>3.95</u>	<u>U</u>	Ethylbenzene			3.95	U
Carbon disulfide			28.7		p/m-Xylene			7.90	U
Methylene chlor	ide		9.88	U	o-Xylene			3.95	U
Methyl tert-buty	l ether (MTBE)		3.95	U	Styrene			3.95	U
trans-1,2-Dichlo	roethene		3.95	U	Bromoform			3.95	U
1,1-Dichloroetha	me		3.95	U	Isopropylbenz	ene		3.95	U
Vinyl acetate			3.95	U ·	1,1,2,2-Tetrachloroethane			3.95	U
2-Butanone (ME	EK)		60.9	J	Bromobenzen	e		3,95	U
cis-1,2-Dichloro	ethene		3.95	U	1,2,3-Trichlor	opropane		3.95	U
2,2-Dichloropro	pane		3.95	U	n-Propylbenze	ene		3.95	U
Chloroform			3,95	U	2-Chlorotolue	ne		3.95	U
1,1,1-Trichloroe	thane		3.95	U	1,3,5-Trimethy	ylbenzene		3,95	U
1,1-Dichloropro	pene		3.95	U	4-Chlorotolue	ne		3.95	U
Carbon tetrachlo	ride		3.95	U	tert-Butylbenz	ene		3.95	U
Benzene			3.95	U	1,2,4-Trimethy	ylbenzene		3.95	U
1,2-Dichloroetha	ine		3.95	U	sec-Butylbenz	ene		3.95	ືບ
Trichloroethene			3.95	U	1,3-Dichlorob	enzene		3.95	U
1,2-Dichloropro	pane		3.95	U	p-Isopropyltol	uene		9.88	U
Dibromomethan	e		3.95	U	1,4-Dichlorob	enzene		3.95	
Bromodichloron	nethane		3.95	U	n-Butylbenzer	ne		9.88	U
2-Chloroethylvii	nyl ether		9.88	U	1,2-Dichlorobenzene			3.95	U
Methyl isobutyl			3.95	U	1,2-Dibromo-	3-chloropropane		3.95	U
cis-1,3-Dichloro			3.95	U	1,2,4-Trichlor			3.95	U
Toluene			3.95	U	Hexachlorobu	tadiene		3.95	U
trans-1,3-Dichlo	ropropene		3.95	U	Naphthalene			9.88	U
1,1,2-Trichloroe			3.95	U	1,2,3-Trichlor	obenzene		3.95	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	103	70-130
1,2-Dichloroethane-d4	105	70-130
Toluene-d8	88	70-130
4-Bromofluorobenzene	80	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SC30-S1

N/A

Matrix:

SDG: N/A

Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-24 Associated Blank: VS072308B14 Concentration Units: µg/Kg

m. 14 . .

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
07/10/08	07/11/08	07/24/08	53.8	7.29	5	1	BAS
Parameter	l	1	Result	Parameter			Result
Dichlorodifluor	omethane		2.55 U	· 2-Hexanone			6.37 U
Chloromethane			2.55 U	Tetrachloroeth	iene		2.55 U J
Vinyl chloride			2.55 U	1,3-Dichlorop	ropane		2.55 U
Bromomethane			2.55 U J	Dibromochlor	omethane		<u>2.55 U</u>
Chloroethane			2.55 U	1,2-Dibromoe	thane		2.55 U
Trichlorofluoror	nethane		<u>2.55 U</u>	Chlorobenzen			<u>2.55 U</u>
Acetone			197	1,1,1,2-Tetrac	hloroethane	· · · · .	2.55 U
1,1-Dichloroethe	ene		<u>2.55 U 🐓</u>	Ethylbenzene			2.55 U
Carbon disulfide			<u>29.8</u> T	p/m-Xylene	annan an a		5.10 U
Methylene chlor	ide		6.37 U	o-Xylene		¢	2.55 U
Methyl tert-buty	l ether (MTBE)		2.55 U	Styrene			2.55 U
trans-1,2-Dichlo			<u>2.55 U J</u>	Bromoform			2.55 U
1,1-Dichloroetha	ane		2.55 U	Isopropylbenz	ene		2.55 U
Vinyl acetate			2.55 U J	1,1,2,2-Tetrac	hloroethane		2.55 U
2-Butanone (ME	EK)		<u>66.3</u> T	Bromobenzen	e		2.55 U
cis-1,2-Dichloro			2.55 U J	1,2,3-Trichlor	opropane		2,55 U
2,2-Dichloropro	pane		2.55 U J	n-Propylbenze	ene		2.55 U
Chloroform			<u>2.55 U</u>	2-Chlorotolue	ne		2.55 U
1,1,1-Trichloroe	thane		2.55 U J	1,3,5-Trimethy	ylbenzene		2.55 U
1,1-Dichloropro	pene		2.55 U	4-Chlorotolue	ne		2.55 U
Carbon tetrachlo	oride		2.55 U	tert-Butylbenz	ene		2.55 U
Benzene			2.55 U	1,2,4-Trimethy	ylbenzene		2.55 U
1,2-Dichloroetha	ane		2.55 U	sec-Butylbenz	ene		2.55 U
Trichloroethene			2.55 U 🖑	1,3-Dichlorob	enzene		2.55 U
1,2-Dichloropro	pane		2.55 U	p-Isopropyltol	uene		6.37 U
Dibromomethan	e		2.55 U J	1,4-Dichlorob	enzene		2.55 U
Bromodichloron	nethane		2.55 U J	n-Butylbenzer	ie		6.37 U
2-Chloroethylvin	nyl ether		6.37 U J	1,2-Dichlorob	enzene		2.55 U
Methyl isobutyl	ketone (MIBK)		2.55 U	1,2-Dibromo-3	3-chloropropane		2.55 U
cis-1,3-Dichloro	propene		-2.55 U R	1,2,4-Trichlor	obenzene		2.55 U
Toluene			2.55 U J	Hexachlorobu	tadiene		2.55 U
trans-1,3-Dichlo	ropropene		2.55 U R	Naphthalene			6.37 U
1,1,2-Trichloroe			2.55 U J	1,2,3-Trichlor	obenzene		2.55 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	87	70-130
1,2-Dichloroethane-d4	75	70-130
Toluene-d8	89	70-130
4-Bromofluorobenzene	80	70-130

N/A - Not Applicable

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:17



ARCADIS **Tidewater MGP**

SC47-S1

SDG: N/A

Sediment

N/A

Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-25 Associated Blank: VS072408B15 **Concentration Units:** µg/Kg

-

Date CollectedDate ReceivedDate AnalyzedPercent SolidAmount (g)Volume (ml)FactorAnalyst07/11/0807/11/0807/24/0883.27.5051BASParameterResultParameterResultParameterResultDichlorodiftuoromethane1.60U2-Hexanone4.01UChloromethane1.60U1.3-Dichloropropane1.60UDiromoethane1.60U1.3-Dichloropropane1.60UChlorotethane1.60U1.2-Dibromoethane1.60UTrichlorofluoromethane1.60UChlorobenzane1.60UAcetone14.2 $B \cup \Im$ 1.1.1.2-Tetrachloroethane1.60UCarbon disulfide1.71p/m-Xylene3.21UMethylene chloride4.01UStyrene1.60UYinyl acetae1.60UStyrene1.60UYinyl acetae1.60UStyrene1.60U2.3-Dichloroethene1.60U1.2.3-Trichloropane1.60U1.3-Dichloroethene1.60U1.2.3-Trichloropane1.60U2.3-Dichloroethene1.60U1.2.3-Trichloropane1.60U1.3-Dichloroethene1.60U1.2.3-Trinchloropane1.60U1.3-Dichloroethene1.60U1.2.3-Trinchloroethane1.60U1.3-Dichloroethene1.60U1.2.3-Trinchloroethane<	Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	et
ParameterResultParameterResultDichlorodifluoromethane1.60U2-Hexanone4.01UChloromethane1.60UTetrachloroethene1.60UBromomethane1.60UJ.3-Dichloropropane1.60UBromomethane1.60UDibromochloromethane1.60UChloroethane1.60UDibromochloromethane1.60UTrichlorofhoromethane1.60UChlorobenzene1.60UAcetone14.2 $B' \cup J'$ 1,1,1,2-Tetrachloroethane1.60UQarbon disulfide1.71p/m-Xylene3.21UMethylene chloride4.01Uo-Xylene1.60UYinyl acetate1.60UBromoform1.60Utrans-1,2-Dichloroethene1.60UStyrene1.60Uyinyl acetate1.60UIsopropylbenzene1.60U2-Butanone (MEK)1.60U1,1,2-Tetrachloroethane1.60U2-Butanone (MEK)1.60U1,3,5-Tinlerdylbenzene1.60U2-Dichloroethane1.60U1,1,2-Tetrachloroethane1.60U1,1-Dichloroethane1.60U1,3,2-Tietholroethane1.60U2-Dichloroppane1.60U1,3,2-Tietholroethane1.60U2-Dichloroppane1.60U1,3,2-Tietholroethane1.60U1,1-Dichloroethane1.60U1,2,2-T								enfo more e e estado	
$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	Parameter	L	Ĭ	Result	Parameter	I			
Vinyl chloride1.60U1.3-Dichloropropane1.60UBromomethane1.60UDibromochloromethane1.60UChloroethane1.60U1.2-Dibromoethane1.60UAcetone1.60UChlorobenzene1.60U1.1-Dichloroethane1.60UChlorobenzene1.60U2.1-Dichloroethane1.60UChlorobenzene1.60U2.1-Dichloroethane1.60UChlorobenzene1.60U2.1-Dichloroethane1.60Uethylenzene3.21UMethylene chloride4.01Uo-Xylene3.21UMethylene chloride4.01Uo-Xylene1.60U1.1-Dichloroethane1.60UStyrene1.60U1.1-Dichloroethane1.60UStyrene1.60U1.1-Dichloroethane1.60U1.1,2,2-Tetrachloroethane1.60U2.2-Dichloroethane1.60U1.2,3-Trinchloropropane1.60U2.2-Dichloroethane1.60U1.2,3-Trinchloropropane1.60U1.1-Dichloroethane1.60U2.2-Chlorotoluene1.60U1.2-Dichloroethane1.60U2.2-Chlorotoluene1.60U1.1-Dichloroethane1.60U1.3,5-Trinchlybenzene1.60U1.1-Dichloroethane1.60U1.2,2-Trinchlybenzene1.60U1.2-Dichloroethane1.60	Dichlorodifluoro	omethane		1.60 U	2-Hexanone			4.01	U
Vinyl chloride1.60U1.3-Dichloropropane1.60UBromomethane1.60UDibromochloromethane1.60UChloroethane1.60U1.2-Dibromoethane1.60UTrichlorofluoromethane1.60UChlorobenzene1.60UAcetone1.42 $B' \cup J'$ 1.1.1.2-Tetrachloroethane1.60U1.1-Dichloroethene1.60UEthylbenzene1.60UCarbon disulfide1.71p/m-Xylene3.21UMethylene chloride4.01Uo-Xylene1.60UMethyl tert-butyl ether (MTBE)1.60UStyrene1.60U1.1-Dichloroethane1.60UBromoform1.60U1.1-Dichloroethane1.60UIntras-1,2-Dichloroethane1.60U1.1-Dichloroethane1.60U1.1,2,2-Tetrachloroethane1.60U2Butanone (MEK)1.60U1.2,3-Trichloropropane1.60U2.2-Dichloroethane1.60U1.2,3-Trichloropropane1.60U1.1-Dichloroethane1.60U1.2,3-Trichloropropane1.60U1.1-Dichloroethane1.60U1.2,3-Trichloropropane1.60U1.1-Dichloroethane1.60U1.2,4-Trimethylbenzene1.60U1.1-Dichloroethane1.60U1.2,4-Trimethylbenzene1.60U1.1-Dichloroethane1.60U1.2,4-Trimethylbenzene1.60	Chloromethane			1.60 U	Tetrachloroeth	ene		1.60	U
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	Vinyl chloride			1.60 U	1,3-Dichlorop	ropane	and an and a set of the set of th		U
$\begin{array}{c c c c c c c c c c c c c c c c c c c $	Bromomethane			1.60 U J	Dibromochlor	omethane		1.60	U
Acetone14.2 $B'\cup$ 1,1,1,2-Tetrachloroethane1.60U1,1-Dichloroethane1.60UEthylbenzene1.60UCarbon disulfide1.71 p/m -Xylene3.21UMethylene chloride4.01Uo-Xylene3.21UMethylene chloride4.01Uo-Xylene1.60UItrans-1,2-Dichloroethene1.60UStyrene1.60U1,1-Dichloroethene1.60UBromoform1.60U1,1,2-2-Tetrachloroethane1.60UIsopropylenzene1.60U2-Butanone (MEK)1.60UIsopropylenzene1.60U2,2-Dichloroethene1.60U1,2,2-Tetrachloroethane1.60U2,2-Dichloropropane1.60U1,2,3-Trichloropropane1.60U1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U1,2,4-Trimethylbenzene1.60U1,1-Dichloropropene1.60U1,2,4-Trimethylbenzene1.60U1,1-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,2-Titmethylbenzene1.60U1,1-Dichloropropane1.60U1,2,4-Trimethylbenzene1.60U1,1-Dichloropropane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1	Chloroethane			1.60 U	1,2-Dibromoet	hane		1.60	U
1,1-Dichloroethene1.60UEthylbenzene1.60UCarbon disulfide1.71 p/m -Xylene3.21UMethylene chloride4.01Uo-Xylene1.60UMethyl tert-butyl ether (MTBE)1.60UStyrene1.60U1.1-Dichloroethene1.60UBromoform1.60U1.1-Dichloroethene1.60UIsopropylbenzene1.60U2-Butanone (MEK)1.60UJ.2.3-Trichloroethane1.60U2-Butanone (MEK)1.60UJ.2.3-Trichloroptopane1.60U2,2-Dichloroethene1.60U1.2.3-Trinethylbenzene1.60U2,2-Dichloroptopane1.60U1.2.3-Trinethylbenzene1.60U1,1-Trichloroethane1.60U2-Chlorotohene1.60U1,1-Dichloroptopene1.60U1.3.5-Trimethylbenzene1.60U1,1-Dichloroptopene1.60U1.2,4-Trimethylbenzene1.60U1,1-Dichloroptopene1.60U1.2,4-Trimethylbenzene1.60U1,2-Dichloroptopene1.60U1.2,4-Trimethylbenzene1.60U1,2-Dichloroptopene1.60U1.2-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1.2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1.2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1.2-Dichlorobenzene<	Trichlorofluoror	nethane				2		1.60	U
Carbon disulfide1.71 $p/m-Xylene$ 3.21UMethylene chloride4.01U $o-Xylene$ 1.60UMethyl tert-butyl ether (MTBE)1.60UStyrene1.60Utrans-1,2-Dichloroethene1.60UBromoform1.60U1,1-Dichloroethane1.60UIsopropylbenzene1.60U2-Butanone (MEK)1.60UI,2,2-Tetrachloroethane1.60U2-Butanone (MEK)1.60UI,2,3-Trichloroptopane1.60U2,2-Dichloroethane1.60U1,2,3-Trichloroptopane1.60U2,2-Dichloroptopane1.60U2-Chlorotohene1.60U1,1-Trichloroethane1.60U2-Chlorotohene1.60U1,1-Trichloroptopane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloroptopene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,2-A-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,2-Dichlorobenzen				<u>14.2 BU</u> J	1,1,1,2-Tetracl	nloroethane		1.60	U
Methylene chloride4.01Uo-Xylene1.60UMethyl tert-butyl ether (MTBE)1.60UStyrene1.60Utrans-1,2-Dichloroethane1.60UBromoform1.60U1,1-Dichloroethane1.60UIsopropylbenzene1.60U2-Butanone (MEK)1.60UJBromobenzene1.60U2,2-Dichloroethane1.60U1,2,3-Trichloropropane1.60U2,2-Dichloroptopane1.60U1,2,3-Trichloropropane1.60U2,2-Dichloroptopane1.60U1,3,5-Trimethylbenzene1.60U1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloroptopane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloroptopane1.60U1,2,4-Trimethylbenzene1.60U1,1-Dichloroptopane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroptopane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroptopane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopane1.	1,1-Dichloroethe	ene		1.60 U	Ethylbenzene			1.60	U
Methyl tert-butyl ether (MTBE)1.60UStyrene1.60Utrans-1,2-Dichloroethene1.60UBromoform1.60U1,1-Dichloroethene1.60UIsopropylbenzene1.60UVinyl acetate1.60U1,1,2,2-Tetrachloroethane1.60U2-Butanone (MEK)1.60U1,2,3-Trichloroethane1.60U2-Butanone (MEK)1.60U1,2,3-Trichloroptopane1.60U2,2-Dichloropropane1.60Un-Propylbenzene1.60U2,2-Dichloroptopane1.60U1,3,5-Trimethylbenzene1.60U1,1-Trichloroptopene1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloroptopene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroptopene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroethane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroethane1.60<	Carbon disulfide	<u>}</u>	41 M A # 7 B #	1.71	p/m-Xylene			3.21	U
trans-1,2-Dichloroethene1.60UBromoform1.60U1,1-Dichloroethane1.60UIsopropylbenzene1.60UYinyl acetate1.60U1,1,2,2-Tetrachloroethane1.60U2-Butanone (MEK)1.60UBromobenzene1.60U2-Dichloroethene1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60U1,2,3-Trichloroptolene1.60U2,2-Dichloropropane1.60U2-Chlorotoluene1.60U1,1-Dichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,1-Dichloroethane1.60U4-Chlorotoluene1.60U1,2-Dichloropropene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene8.17Trichloropropane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloromethane1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropengene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropengene1.60U1,2-Dichlorobenz				4.01 U	o-Xylene			1.60	U
1,1-Dichloroethane1.60UIsopropylbenzene1.60UVinyl acetate1.60U1,1,2,2-Tetrachloroethane1.60U2-Butanone (MEK)1.60UBromobenzene1.60Ucis-1,2-Dichloroethene1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60Un-Propylbenzene1.60U2,2-Dichloropropane1.60U2-Chlorotoluene1.60U1,1-Trichloroethane1.60U2-Chlorotoluene1.60U1,1-Trichloropropene1.60U4-Chlorotoluene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroptopane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U				1.60 U	Styrene			1.60	U
Vinyl acetate1.60U1,1,2,2-Tetrachloroethane1.60U2-Butanone (MEK)1.60UJBromobenzene1.60Ucis-1,2-Dichloroethene1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60U1,2,3-Trichloropropane1.60U1,1-Trichloroethane1.60U2-Chlorotoluene1.60U1,1-Dichloropropene1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60U1,2-Dichlorobenzene1.60U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene1.60U01,4-Dichlorobenzene1.60U1,2-Dichlorobenzene1.60U1,4-Dichlorobe	trans-1,2-Dichlo	roethene		1.60 U	Bromoform			1.60	U
2-Butanone (MEK)1.60UBromobenzene1.60Ucis-1,2-Dichloroethene1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60Un-Propylbenzene1.60UChloroform1.60U2-Chlorotoluene1.60U1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,2-Dichloroethane1.60Utert-Butylbenzene1.60U1,2-Dichloroethane1.60Usec-Butylbenzene1.60U1,2-Dichloroethane1.60Usec-Butylbenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene4.01U1,2-Dichloropengene1.60U1,2-Dichlorobenzene1.60U1,2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60U1,2-Dichloropengene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropengene1.60U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2		ane		1.60 U	Isopropylbenz	ene		1.60	U
2-Butanone (MEK)1.60UBromobenzene1.60Ucis-1,2-Dichloroethene1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60Un-Propylbenzene1.60UChloroform1.60U2-Chlorotoluene1.60U1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,2-Dichloropropene1.60Utert-Butylbenzene1.60UBenzene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroptopane1.60Usec-Butylbenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloroptopane1.60U1,4-Dichlorobenzene4.01U1,2-Chloroethylvinyl ether4.01U1,2-Dichloroptopane1.60U1,2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Di					1,1,2,2-Tetracl	loroethane			U
cis-1,2-Dichloroethene1.60U1,2,3-Trichloropropane1.60U2,2-Dichloropropane1.60Un-Propylbenzene1.60UChloroform1.60U2-Chlorotoluene1.60U1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60U1,2-Dichloroethane1.60U4-Chlorotoluene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene8.171,2-Dichloroptopane1.60U1,2-Dichlorobenzene8.171,2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloroptopene1.60U1,2-Dichloroptopane1.60U1,2-Dichloroptopene1.60U1,2-Dichloroptopane1.60U1,2-Dichloroptopene1.60U1,2-Dichloroptopane1.60U1,3-Dichloroptopene1.60U1,2-Dichloroptopane1.60U1,3-Dichloroptopene1.60U1,2-Dichlorop	2-Butanone (ME	EK)		1.60 U J	Bromobenzene	;		the second second second	
2,2-Dichloropropane1.60Un-Propylbenzene1.60UChloroform1.60U2-Chlorotoluene1.60U1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60UCarbon tetrachloride1.60U4-Chlorotoluene1.60UBenzene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene8.17Dibromomethane1.60U1,2-Dichlorobenzene1.60U2-Chloroethylvinyl ether4.01U1,2-Dichloropropane1.60U1,3-Dichloropropene1.60U1,2-Dichloropropane1.60U1,2-Dichloropenzene1.60U1,2-Dichloropenzene1.60U1,3-Dichloropropene1.60U1,2-Dichloropropane1.60U1,2-Dichloropropene1.60U1,2-Dichloropropane1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene	cis-1,2-Dichloro	ethene		1.60 U	1,2,3-Trichloro	propane			
1,1,1-Trichloroethane1.60U1,3,5-Trimethylbenzene1.60U1,1-Dichloropropene1.60U4-Chlorotoluene1.60UCarbon tetrachloride1.60U4-Chlorotoluene1.60UBenzene1.60Utert-Butylbenzene1.60U1,2-Dichloroethane1.60Usec-Butylbenzene1.60U1,2-Dichloroethane1.60Usec-Butylbenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60U1,4-Dichlorobenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,4-Dichlorobenzene1.60U1,4-Dichlorobenzene6.0U1,2-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropropene1.60U1,2-A-Trichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-A-Trichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-A-Trichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-A-Trichlorobenzene1.60U1,3-Dichloropropene1.60U<		pane		1.60 U	n-Propylbenze	ne			
1,1-Dichloropropene1.60U4-Chlorotoluene1.60UCarbon tetrachloride1.60Utert-Butylbenzene1.60UBenzene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60Usec-Butylbenzene1.60UTrichloroethane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloromethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60U1,2-Dichlorobenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,2-Dichloropropene1.60U1,2-Dichlorobenzene1.60U1,3-Dichloropropene1.60U1,2-Dichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60UTans-1,3-Dichloropropene1.60UNaphthalene4.01U	Chloroform			1.60 U	2-Chlorotoluer	1e		1.60	U
Carbon tetrachloride 1.60 Utert-Butylbenzene 1.60 UBenzene 1.60 U $1,2,4$ -Trimethylbenzene 1.60 U $1,2$ -Dichloroethane 1.60 U $1,2,4$ -Trimethylbenzene 1.60 UTrichloroethene 1.60 U $1,3$ -Dichlorobenzene 1.60 U $1,2$ -Dichloropropane 1.60 U $1,3$ -Dichlorobenzene 1.60 U $1,2$ -Dichloropropane 1.60 U $1,3$ -Dichlorobenzene 4.01 UDibromomethane 1.60 U $1,4$ -Dichlorobenzene 8.17 Bromodichloromethane 1.60 U $1,2$ -Dichlorobenzene 4.01 U 2 -Chloroethylvinyl ether 4.01 U $1,2$ -Dichlorobenzene 1.60 U $1,3$ -Dichloropropene 1.60 U $1,2$ -Dichlorobenzene 1.60 U $1,2$ -Dichlorobenzene 1.60 U $1,2$ -Dichlorobenzene 1.60 U $1,3$ -Dichloropropene 1.60 U $1,2$ -Dichlorobenzene 1.60 U $1,3$ -Dichloropropene 1.60 U $1,2$ -Trichlorobenzene 1.60 U $1,3$ -Dichloropropene 1.60 </td <td>1,1,1-Trichloroe</td> <td>thane</td> <td></td> <td>1.60 U</td> <td>1,3,5-Trimethy</td> <td>lbenzene</td> <td></td> <td>1.60</td> <td>Ų</td>	1,1,1-Trichloroe	thane		1.60 U	1,3,5-Trimethy	lbenzene		1.60	Ų
Benzene1.60U1,2,4-Trimethylbenzene1.60U1,2-Dichloroethane1.60Usec-Butylbenzene1.60UTrichloroethene1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60Up-Isopropyltoluene4.01UDibromomethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60Un-Butylbenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dibromo-3-chloropropane1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U				1.60 U	4-Chlorotoluer	ne		1.60	U
1,2-Dichloroethane1.60Usec-Butylbenzene1.60UTrichloroethene1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60Up-Isopropyltoluene4.01UDibromomethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60Un-Butylbenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dibromo-3-chloropropane1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U	Carbon tetrachlo	oride		1.60 U	tert-Butylbenz	ene		1.60	U
Trichloroethene1.60U1,3-Dichlorobenzene1.60U1,2-Dichloropropane1.60Up-Isopropyltoluene4.01UDibromomethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60Un-Butylbenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dichlorobenzene1.60Ucis-1,3-Dichloropropene1.60U1,2,4-Trichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U	Benzene			1.60 U	1,2,4-Trimethy	lbenzene		1.60	U
1,2-Dichloropropane1.60Up-Isopropyltoluene4.01UDibromomethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60U1,2-Dichlorobenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dibromo-3-chloropropane1.60Ucis-1,3-Dichloropropene1.60U1,2,4-Trichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U		ane		<u>1.60 U</u>	sec-Butylbenze	ene		1.60	U
Dibromomethane1.60U1,4-Dichlorobenzene8.17Bromodichloromethane1.60Un-Butylbenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dibromo-3-chloropropane1.60Ucis-1,3-Dichloropropene1.60U1,2,4-Trichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U	Here's a second s			1.60 U	1,3-Dichlorobe	enzene		1.60	U
Bromodichloromethane1.60Un-Butylbenzene4.01U2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dibromo-3-chloropropane1.60Ucis-1,3-Dichloropropene1.60U1,2,4-Trichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U	1,2-Dichloropro	pane		1.60 U	p-Isopropyltol	uene		4.01	U
2-Chloroethylvinyl ether4.01U1,2-Dichlorobenzene1.60UMethyl isobutyl ketone (MIBK)1.60U1,2-Dibromo-3-chloropropane1.60Ucis-1,3-Dichloropropene1.60U1,2,4-Trichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U		dian statute and the statement of the		<u>1.60 U</u>	1,4-Dichlorobe	enzene		8.17	
Methyl isobutyl ketone (MIBK)1.60U1.2-Dibromo-3-chloropropane1.60Ucis-1,3-Dichloropropene1.60U1,2-Dibromo-3-chloropropane1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U				1.60 U	n-Butylbenzen	e		4.01	U
cis-1,3-Dichloropropene1.60U1,2,4-Trichlorobenzene1.60UToluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U								1.60	U
Toluene6.84BHexachlorobutadiene1.60Utrans-1,3-Dichloropropene1.60UNaphthalene4.01U								1.60	U
trans-1,3-Dichloropropene 1.60 U Naphthalene 4.01 U		is-1,3-Dichloropropene 1.60 U 1,2,4-Trichlorobenzene						1.60	U
					Hexachlorobutadiene			1.60	U
1,1,2-Trichloroethane 1.60 U 1,2,3-Trichlorobenzene 1.60 U				1.60 U	Naphthalene			4.01	U
	1,1,2-Trichloroe	thane		1.60 U	1,2,3-Trichloro	benzene		1.60	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	104	70-130
1,2-Dichloroethane-d4	106	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	81	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:25



ARCADIS Tidewater MGP

SC32-S1

N/A

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-26 Associated Blank: VS072308B15 Concentration Units: ug/Kg

			0. N	5261			Associ	aleu Dialik.	Y 30/2300D
		Matrix:	Sediment				Conce	ntration Units	: µg/Kg
Date	Date	Date	Percent	Methanol	Sample	Extract	Final	Dilution	
Collected	Received	Analyzed	Solid	Volume (ml) Amount (g)	Volume (ml)	Volume (ml)	Factor	Analyst
07/11/08	07/11/08	07/23/08	60.1	5	7.34	0.1	5	1	BAS
Parameter			Re	sult	Parameter	•		R	esult
Dichlorodiflu	uoromethane			180 U	2-Hexanon	e			450 U
Chlorometha				180 U	Tetrachloro	oethene			<u>180 U</u>
Vinyl chloric	le			180 U	1,3-Dichlo				180 U
Bromometha			180 .	<u>140 JB ()</u> .	Dibromoch	loromethane			<u>180 U</u>
Chloroethane	e			180 U	1,2-Dibron				180 U
Trichlorofluc	oromethane			180 U	Chloroben				<u>180 U</u>
Acetone				<u>627 B</u> J	<u>1,1,1,2-Tet</u>	rachloroethar	ıe		180 U
1,1-Dichloro	ethene			180 U	Ethylbenze				180 U
Carbon disul	fide			180 U	p/m-Xylen	e			360 U
Methylene cl	hloride			<u>450 U</u>	o-Xylene				180 U
Methyl tert-b	outyl ether (M	ITBE)		<u>180 U</u>	Styrene	- 14			180 U
trans-1,2-Dic	chloroethene			<u>180 U</u>	Bromoform				180 U
1,1-Dichloro	ethane			180 U	Isopropylb				180 U
Vinyl acetate				180 U		rachloroethau	ne		180 U
2-Butanone	(MEK)			180 U J					180 U
cis-1,2-Dich	loroethene			180 U		loropropane			180 U
2,2-Dichloro	propane			<u>180 U</u>	n-Propylbe			L L L L	<u>180 U</u>
Chloroform				180 U	2-Chloroto	· · · · · · · · · · · · · · · · · · ·			180 U
1,1,1-Trichle	oroethane			180 U		ethylbenzene			180 U
1,1-Dichloro				180 U	<u>4-Chlorotc</u>				180 U
Carbon tetra	chloride			<u>180 U</u>		enzene			180 U
Benzene				132 J		ethylbenzene			180 U
1,2-Dichloro	oethane			<u>180 U</u>	sec-Butylb				180 U
Trichloroeth	ene			180 U		robenzene			180 U
1,2-Dichlore	propane			180 U	p-Isopropy				450 U
Dibromomet	thane			180 U		robenzene			180 U
Bromodichle	oromethane			180 U	n-Butylber				450 U
2-Chloroethy	ylvinyl ether			450 U		robenzene			180 U
Methyl isobu	utyl ketone (N	MIBK)		180 U		no-3-chlorop			180 U
cis-1,3-Dich				180 U		nlorobenzene			180 U
Toluene				180 U		obutadiene			180 U
trans-1,3-Die	chloropropen	e		180 U	Naphthale				2540
1,1,2-Trichle				180 U	1,2,3-Tricl	hlorobenzene		· · · · · ·	180 U

		Acceptance
Surrogate	% Recovery	Range (%)
2-Bromo-1-chloropropane	81	70-130
1-Chloro-2-fluorobenzene	80	70-130
1,4-Dichlorobutane	82	70-130
Dibromofluoromethane	99	70-130
Toluene-d8	95	70-130
4-Bromofluorobenzene	97	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SC33-S1

Sediment

N/A

Matrix:

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-27 Associated Blank: VS072408B15 Concentration Units: µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	zst
07/11/08	07/11/08	07/24/08	31.1	6.57	5	1	BAS	
Parameter	0//11/00		Result	Parameter		·······	Result	
Dichlorodifluoro	methane		4.90 U	2-Hexanone			12.3	U
Chloromethane	71110 (114110		4.90 U	Tetrachloroeth	iene		4.90	U
Vinyl chloride			4.90 U	1,3-Dichlorop		A CONTRACTOR OF A CONTRACTOR A CO	4.90	U
Bromomethane			4.90 U J	Dibromochlor			4.90	U
Chloroethane			4.90 U	1,2-Dibromoe	thane	· · · · · · · · · · · · · · · · · · ·	4.90	U
Trichlorofluoror	nethane		4.90 U	Chlorobenzen			4.90	U
Acetone			368 B J	1.1.1.2-Tetrac		a dana da waran da	4.90	U
1.1-Dichloroeth	ene		4.90 U	Ethylbenzene			4.90	U
Carbon disulfide			42.5	p/m-Xylene			9.80	U
Methylene chlor			12.3 U	o-Xylene			4.90	U
Methyl tert-buty			4.90 U	Styrene			4,90	U
trans-1,2-Dichlo		19 889 ann martairtean ann a' d anna ann ann ann	4.90 U	Bromoform			4.90	U
1,1-Dichloroeth			4.90 U	Isopropylbenz	ene		4.90	U
Vinyl acetate			4.90 U	1,1,2,2-Tetrac			4.90	U
2-Butanone (MI	EK)	1	85.4 J	Bromobenzen	e		4.90	U
cis-1,2-Dichloro			4.90 U	1,2,3-Trichlor	opropane		4.90	U
2,2-Dichloropro			4.90 U	n-Propylbenze	ene		4.90	U
Chloroform			4.90 U	2-Chlorotolue	ne		4.90	U
1,1,1-Trichloroe	thane		4.90 U	1,3,5-Trimeth	ylbenzene		4.90	U
1,1-Dichloropro			4.90 U	4-Chlorotolue	ne		4.90	U
Carbon tetrachle			4.90 U	tert-Butylbenz	zene		4.90	U
Benzene			4.90 U	1,2,4-Trimeth	ylbenzene		4.90	U
1,2-Dichloroeth	ane		4.90 U	sec-Butylbenz	ene		4.90	U
Trichloroethene			4.90 U	1,3-Dichlorob	enzene		4.90	U
1,2-Dichloropro			4.90 U	p-Isopropylto	luene		12.3	U
Dibromomethan			4.90 U	1,4-Dichlorob	enzene		4.90	U
Bromodichloror	nethane		4.90 U	n-Butylbenzer	ne		12.3	U
2-Chloroethylvi	nyl ether		12.3 U	1,2-Dichlorob	enzene		4.90	U
Methyl isobutyl		•	4.90 U		3-chloropropane		4.90	U
cis-1,3-Dichlord			4.90 U	1,2,4-Trichlor			4.90	U
Toluene			4.90 U	Hexachlorobu	itadiene		4.90	U
trans-1,3-Dichle	propropene		4.90 U	Naphthalene		,	12.3	U
1,1,2-Trichloroe			4.90 U	1,2,3-Trichlor	robenzene		4.90	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	101	70-130
1,2-Dichloroethane-d4	103	70-130
Toluene-d8	88	70-130
4-Bromofluorobenzene	78	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:26

	PHA	Client: Project: Client ID: Case: Matrix:	ARCADI Tidewate: SW2-S1 N/A Sediment	r MGP	N/A		ETR: Lab II Assoc	ode: MA00 0807059 D: 0807059- iated Blank: entration Unit	28 VS07:	2308B15 /Kg
Date	Date	Date	Percent	Methanol	Sample	Extract	Final	Dilution Factor	Anal	livet
Collected	Received	Analyzed	Solid				Volume (ml)	1	BA	
07/11/08	07/11/08	07/23/08	39.6	5	5.39	0.1	5			
Parameter			Re	sult	Parameter			F	Result	
Dichlorodiflu	oromethane			387 U	2-Hexanon				967	U
Chlorometha				387 U	Tetrachloro	ethene			387	<u> </u>
Vinyl chlorid	······	· · · · · · · · · · · · · · · · · · ·		387 U	1,3-Dichlor	ropropane			387	<u>U</u>
Bromometha			387 -	321 JB -U乙	Dibromoch	loromethane			387	<u> </u>
Chloroethane				387 U	1,2-Dibrom				387	<u> </u>
Trichlorofluc				387 U	Chlorobenz				387	U
Acetone			1	480 B 🔾 🕽		rachloroetha	ne		387	<u>U</u>
1,1-Dichloro	ethene			<u>387 U</u>	Ethylbenze	And the second s			334	<u> J </u>
Carbon disul				731	p/m-Xylen	e			<u>998</u>	
Methylene ch			and the second s	967 U	o-Xylene				619	
Methyl tert-b	outyl ether (M	ITBE)		<u>387 U</u>	Styrene	······································			387	<u>U</u>
trans-1,2-Dic				<u>387 U</u>	Bromoforn				387	<u> </u>
1,1-Dichloro	ethane			<u>387 U</u>	Isopropylb					<u> </u>
Vinyl acetate				<u>387 U</u>		trachloroetha	ne		387	<u> </u>
2-Butanone ((MEK)			387 U J	Bromoben				387	U
cis-1,2-Dich	loroethene			<u>387 U</u>		loropropane			387	U
2,2-Dichloro	propane			387 U	n-Propylbe				414	
Chloroform				387 U	2-Chloroto			······································	387	<u>U</u>
1,1,1-Trichle	oroethane			387 U		ethylbenzen)		3840	¥ ¥
1,1-Dichloro	propene			<u>387 U</u>	4-Chloroto				387	<u> </u>
Carbon tetra	chloride	·		387 U	tert-Butylb				387	<u> </u>
Benzene				234 J		ethylbenzen	<u> </u>		7800	
1,2-Dichloro				387 U	sec-Butylb				<u>365</u> 387	
Trichloroeth	iene			<u>387 U</u>		orobenzene			<u> </u>	
1,2-Dichloro	opropane			387 U	p-Isopropy				387	<u>т</u> т
Dibromome				387 U		orobenzene	• •!		2680	0
Bromodichle				387 U	n-Butylber				387	U
2-Chloroeth				967 U		probenzene			387	<u> </u>
	utyl ketone (I	MIBK)		387 U		mo-3-chlorop		<u></u>	<u> </u>	
cis-1,3-Dich	loropropene	······		387 U		niorobenzene obutadiene	5		387	<u> </u>
Toluene				265 J	<u>Hexachlor</u> Naphthale			34200		
	chloropropen	ie		387 U 387 U		hlorobenzene		a prior	387	
<u>1,1,2-Trichl</u>	oroethane			387 U	1,2,3-110	moroucuzen			207	

Surrogate	% Recovery	Acceptance Range (%)
2-Bromo-1-chloropropane	81	70-130
1-Chloro-2-fluorobenzene	83	70-130
1,4-Dichlorobutane	85	70-130
Dibromofluoromethane	97	70-130
Toluene-d8	98	70-130
4-Bromofluorobenzene	99	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

E - Estimated value, exceeds the upper limit of calibration.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 15:46

	PHA	Client: Project: Client ID: Case: Matrix:	ARCADI Tidewate SW2-S1 N/A Sediment		N/A		Lab Code: MA ETR: 0807059 Lab ID: 08070 Associated Blan Concentration U	59-28E k: VS07	72408B16 2/Kg
Date	Date	Date	Percent	Methano	l Sample	Extract Fi	nal Dilution		<u>yng</u>
Collected	Received	Analyzed	Solid	1		Volume (ml) Volum			lyst
07/11/08	07/11/08	07/24/08	39.6	5	5.39		5 2	BA	
Parameter	l		Res	ult	Parameter	۰ ۲		Result	
Dichlorodiflu	oromethane			773 U	2-Hexanor	ie.		1930	U
Chlorometha				773 U	Tetrachlor		······	773	U
Vinyl chlorid				73 U	1,3-Dichlo			773	U
Bromometha				534 JB		loromethane		773	<u>U</u>
Chloroethane				773 U	1,2-Dibron	and a second sec		773	Ŭ
Trichlorofluc	Contraction of the second seco			73 U	Chloroben			773	U
Acetone				570 JB		rachloroethane		773	U
1,1-Dichloro	ethene			773 U	Ethylbenze			773	U
Carbon disul				580 J	p/m-Xylen				J
Methylene ch				930 U	o-Xylene			541	J
Methyl tert-b	utyl ether (M	TBE)		773 U	Styrene			773	U
trans-1,2-Dic	hloroethene			773 U	Bromoforr	n		773	U
1,1-Dichloro	ethane			773 U	lsopropylb	enzene		773	U
Vinyl acetate	;			173 U	1,1,2,2-Te	rachloroethane		773	U
2-Butanone (173 U	Bromoben			773	U
cis-1,2-Dichl				<u>773 U</u>		loropropane		773	U
2,2-Dichloro	propane			77 <u>3</u> U	n-Propylbe			773	U
Chloroform				773 U	2-Chloroto			773	<u>U</u>
<u>1,1,1-Trichlo</u>			•	773 U		ethylbenzene		3710	/
1,1-Dichloro				773 U	4-Chloroto			773	<u>U</u>
Carbon tetrac	chloride			773 U	tert-Butylb			773	U
Benzene		•****		77 <u>3</u> U		ethylbenzene		7180	
1,2-Dichloro				773 U	sec-Butylb			773	<u>U</u>
Trichloroethe				773 U	1,3-Dichlo			773	U
1,2-Dichloro	· · · · · · · · · · · · · · · · · · ·			773 U	p-Isopropy		<u></u>	1310	J
Dibromomet			· · · · · · · · · · · · · · · · · · ·	773 U	1,4-Dichlo			773	<u> </u>
Bromodichlo				773 U	n-Butylber			2410	
2-Chloroethy				930 U	1,2-Dichlo			773	
Methyl isobu		IIBK)		773 U		no-3-chloropropane		773	U
cis-1,3-Dichl	oropropene			773 U		lorobenzene		773	<u>U</u>
Toluene		······································		773 U		obutadiene		773	U
trans-1,3-Dic				773 U	Naphthale			34200	
1,1,2-Trichlo	roethane			773 U	<u>1,2,3-Tricl</u>	lorobenzene	Pro	773	<u> </u>

		Acceptance
Surrogate	% Recovery	Range (%)
2-Bromo-1-chloropropane	73	70-130
1-Chloro-2-fluorobenzene	79	70-130
1,4-Dichlorobutane	77	70-130
Dibromofluoromethane	92	70-130
Toluene-d8	94	70-130
4-Bromofluorobenzene	96	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

07/31/08 15:46



ARCADIS **Tidewater MGP**

SC39-S1

SDG: N/A

N/A Sediment Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-29 Associated Blank: VS072408B15 Concentration Units: μg/Kg

				Sample	Final	Dilution		<u>, 126</u>
Date Collected			Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/11/08	07/11/08	07/24/08	23.7	8.39	5	1	BAS	S
Parameter		F	Result	Parameter			Result	
Dichlorodifluoro	methane	· · · · · · · · · · · · · · · · · · ·	5.04 U	2-Hexanone			12.6	
Chloromethane			5.04 U	Tetrachloroeth			5.04	
Vinyl chloride			5.04 U	1,3-Dichlorop			5.04	
Bromomethane			5.04 U J	Dibromochlor	omethane			U
Chloroethane			5.04 U	1,2-Dibromoe			5.04	
Frichlorofluoror	nethane		5.04 U	Chlorobenzen			5.04	
Acetone			466 B J	1,1,1,2-Tetrac	hloroethane		5.04	
1,1-Dichloroethe	ene		5.04 <u>U</u>	Ethylbenzene			5.04	U
Carbon disulfide)		92.0	p/m-Xylene			10.1	U
Methylene chlor	ide		12.6 U	o-Xylene			5.04	U
Methyl tert-buty	l ether (MTBE)		5.04 U	Styrene			5,04	<u>U</u>
rans-1,2-Dichlo	roethene		5,04 U	Bromoform			5.04	U
1,1-Dichloroetha	ane		5.04 U	Isopropylbenz	ene		5.04	U
Vinyl acetate			5.04 U	1,1,2,2-Tetrac	hloroethane		5.04	U
2-Butanone (MI			119 J	Bromobenzen	e .		5.04	U
cis-1,2-Dichloro			5.04 U	1,2,3-Trichlor	opropane		5.04	U
2,2-Dichloropro			5.04 U	n-Propylbenze	ene		5.04	U
Chloroform	<u>.</u>		5.04 U	2-Chlorotolue	ne		5.04	U
1,1,1-Trichloroe	thane		5.04 U	1,3,5-Trimethy	ylbenzene		5.04	U
1,1-Dichloropro			5.04 U	4-Chlorotolue	ne		5.04	U
Carbon tetrachlo			5.04 U	tert-Butylbenz	ene		5.04	U
Benzene			5.04 U	1,2,4-Trimeth	ylbenzene		5.04	U
1,2-Dichloroeth	ane		5.04 U	sec-Butylbenz	ene		5.04	U
Trichloroethene	<u> </u>		5.04 U	1,3-Dichlorob	enzene		5.04	U
1,2-Dichloropro	pane		5.04 U	p-Isopropyltol	uene		12.6	U
Dibromomethan			5.04 U	1,4-Dichlorob			5.04	U
	omodichloromethane 5.04 U n-Butylbenzene					12.6		
	Chloroethylvinyl ether 12.6 U 1,2-Dichlorobenzene					5,04	U	
	butyl ketone (MIBK) 5.04 U 1,2-Dibromo-3-chloropropane					5.04		
cis-1,3-Dichloro			5.04 U	1,2,4-Trichlor			5.04	
Toluene			5.04 U	Hexachlorobu			5.04	
trans-1,3-Dichle	ropropene		5.04 U	Naphthalene			12.6	
1,1,2-Trichloroe			5.04 U	1,2,3-Trichlor	obenzene		5.04	Ū

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	102	70-130
Toluene-d8	88	70-130
4-Bromofluorobenzene	80	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:26



ARCADIS **Tidewater MGP**

SC38-S1

SDG: N/A

N/A Sediment

Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-30 Associated Blank: VS072408B15 **Concentration Units:**

µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	
07/11/08	07/11/08	07/24/08	34.5	7.59	5	1	BAS	
Parameter	• · · · · · · · · · · · · · · · · · · ·	ľ	Result	Parameter	I		Result	
Dichlorodifluoro	methane		3.82 U	2-Hexanone			9.55	U
Chloromethane			3.82 U	Tetrachloroeth	iene			Ū
Vinyl chloride			3.82 U	1,3-Dichlorop	ropane		3.82	
Bromomethane			3.82 U J	Dibromochlor			3.82	
Chloroethane			3.82 U	1,2-Dibromoet	thane		3.82	
Trichlorofluoron	nethane		3.82 U	Chlorobenzena	3		3.82	
Acetone			206 B J	1,1,1,2-Tetracl	hloroethane		3.82	
1,1-Dichloroethe	ne		3.82 U	Ethylbenzene			3.82	
Carbon disulfide		· · · · · · · · · · · · · · · · · · ·	43.9	p/m-Xylene			7.64	
Methylene chlor			9.55 U	o-Xylene			3.82	
Methyl tert-buty	l ether (MTBE)		3.82 U	Sturene		·····	3.82	
trans-1,2-Dichlo	roethene		3.82 U	Bromoform			3.82	*** *******
1,1-Dichloroetha	ine		3.82 U	Isopropylbenz	ene		2 02	
Vinyl acetate			3.82 U	1,1,2,2-Tetracl			3.82	
2-Butanone (ME	K)		48.8 J	Bromobenzene			3.82	
cis-1,2-Dichloro	ethene		3.82 U	1,2,3-Trichloro	opropane		3.82	
2,2-Dichloroprop	bane		3.82 U	n-Propylbenzene			3.82	
Chloroform			3.82 U	2-Chlorotoluene			3.82	
1,1,1-Trichloroe	thane		3.82 U	1,3,5-Trimethylbenzene			3.82	
1,1-Dichloroprop	pene		3.82 U	4-Chlorotoluer	ne		3.82	
Carbon tetrachlo	ride		3,82 U	tert-Butylbenz	ene	······································	3.82	
Benzene			3.82 U	1,2,4-Trimethy	lbenzene		3.82	
1,2-Dichloroetha	ne		3.82 U	sec-Butylbenze	ene		3.82	
Trichloroethene			3.82 U	1,3-Dichlorobe			3.82	
1,2-Dichloroprop	oane		3.82 U	p-Isopropyltol	uene		9.55	
Dibromomethan			3.82 U	1,4-Dichlorobe			3.82	
Bromodichlorom	lethane		3.82 U	n-Butylbenzen	e		9.55	
2-Chloroethylvir			9.55 U	1,2-Dichlorobe			2 02	***
Methyl isobutyl ketone (MIBK) 3.82 U				1,2-Dibromo-3-chloropropane			3.82	
cis-1,3-Dichloro	propene		3.82 U	1,2,4-Trichlorobenzene			3.82	
Foluene			3.82 U	Hexachlorobutadiene			3.82	
rans-1,3-Dichlor			3.82 U	Naphthalene			9.55	****
1,1,2-Trichloroe	hane		3.82 U	1,2,3-Trichlord	obenzene		3.82	

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	99	70-130
1,2-Dichloroethane-d4	103	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	86	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:26



ARCADIS Tidewater MGP

SC34-S1

SDG: N/A

Matrix: Sediment

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-31 Associated Blank: VS072408B15 Concentration Units: µg/Kg

				Sample	Final	Dilution		
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	/st
07/11/08	07/11/08	07/24/08	53.9	6.03	5	1	BAS	3
Parameter		I	Result	Parameter			Result	
Dichlorodifluoro	methane		3.08 U	2-Hexanone			7.69	U
Chloromethane			3.08 U	Tetrachloroeth			3.08	Ŭ
Vinyl chloride			3.08 U	1,3-Dichloropi	ropane	· • · · · · · · · · · · · · · · · · · ·	3.08	U
Bromomethane			3.08 U J	Dibromochlor	omethane		3.08	U
Chloroethane			3.08 U	1,2-Dibromoet	thane		3.08	U
Trichlorofluoron	nethane		3.08 U	Chlorobenzene	>		3.08	U
Acetone			251 B J	1,1,1,2-Tetracl	nloroethane		3.08	U
1,1-Dichloroethe	ene		3.08 U	Ethylbenzene			3.08	U
Carbon disulfide			45.7	p/m-Xylene			6.15	U
Methylene chlor	ide		7.69 U	o-Xylene			3.08	U
Methyl tert-buty	l ether (MTBE)		3.08 U	Styrene			3.08	U
trans-1,2-Dichlo	roethene		3.08 U	Bromoform			3.08	U
1,1-Dichloroetha	ine		3.08 U	Isopropylbenz	ene		3.08	U
Vinyl acetate			3.08 U	1,1,2,2-Tetracl	nloroethane		3.08	U
2-Butanone (ME	K)		61.7 5	Bromobenzene	3		3.08	U
cis-1,2-Dichloro	ethene		3.08 U	1,2,3-Trichlord	opropane		3.08	U
2,2-Dichloropro	oane		3.08 U	n-Propylbenze			3.08	U
Chloroform			3.08 U	2-Chlorotoluer	ne		3.08	U
1,1,1-Trichloroe	thane		3.08 U	1,3,5-Trimethy	/lbenzene		3.08	U
1,1-Dichloropro	oene		3.08 U	4-Chlorotoluer			3.08	U
Carbon tetrachlo	ride		3.08 U	tert-Butylbenz	ene		3.08	U
Benzene		-	3.08 U	1,2,4-Trimethy	/lbenzene		3.08	U
1,2-Dichloroetha	ine		3.08 U	sec-Butylbenz	ene		3.08	Ŭ
Trichloroethene			3.08 U	1,3-Dichlorobe	enzene		3.08	U
1,2-Dichloropro			3.08 U	p-Isopropyltol	uene		7.69	
Dibromomethan			3.08 U	1,4-Dichlorobe			3.08	*****
Bromodichloron	nethane		3.08 U	n-Butylbenzen	e		7.69	
2-Chloroethylvir	nyl ether		7.69 U	1,2-Dichlorobe			3.08	
Methyl isobutyl			3.08 U	ministra	-chloropropane		3.08	
cis-1,3-Dichloro			3.08 U	1,2,4-Trichloro			3.08	
Toluene			3.08 U	Hexachlorobut	the second s		3.08	Ū
trans-1,3-Dichlo	ropropene		3.08 U	Naphthalene			7.69	Ū
1,1,2-Trichloroe			3.08 U	1,2,3-Trichlord	benzene			Ū

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	100	70-130
1,2-Dichloroethane-d4	97	70-130
Toluene-d8	86	70-130
4-Bromofluorobenzene	77	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:27

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

Page 40 / 277

N/A



ARCADIS

Tidewater MGP SC35-S1

SDG;

Matrix: Sediment

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-32 Associated Blank: VS072408B15 Concentration Units: µg/Kg

Date Collected	Data Passived	Date Analyzed	Percent Solid	Sample	Final	Dilution		0
07/11/08				Amount (g)	Volume (ml)	Factor	Anal	
	07/11/08	07/24/08	55.1	7.88	5	1	BA	S
Parameter		ł	Result	Parameter			Result	
Dichlorodifluoro	methane		2.30 U	2-Hexanone			5.75	U
Chloromethane			2.30 U	Tetrachloroeth	ene		2.30	Ų
Vinyl chloride			2.30 U	1,3-Dichlorop	ropane		2.30	U
Bromomethane			2.30 U J	Dibromochlore	omethane		2.30	U
Chloroethane			<u>2.30 U</u>	1,2-Dibromoet	thane		2.30	U
Trichlorofluoron	nethane		2.30 U	Chlorobenzene	•		2.30	U
Acetone			<u>44.2 B ()</u> J	1,1,1,2-Tetracl	hloroethane		2.30	U
1,1-Dichloroethe			<u>2.30 U</u>	Ethylbenzene			2.30	U
Carbon disulfide			3.26	p/m-Xylene			4.60	U
Methylene chlori			5.75 U	o-Xylene			2.30	U
Methyl tert-butyl			2.30 U	Styrene			2.30	U
trans-1,2-Dichlor			<u>2.30 U</u>	Bromoform			2.30	U
1,1-Dichloroetha	ne		2.30 U	Isopropylbenze	ene		2.30	U
Vinyl acetate			2.30 U	1,1,2,2-Tetracl	hloroethane		2.30	U
2-Butanone (ME	te in Conservation is a substantial databased in a substantial datab		<u>6.34</u> T	Bromobenzene	3		2.30	U
cis-1,2-Dichloroe			2.30 U	1,2,3-Trichlord	opropane		2.30	U
2,2-Dichloroprop	oane		2.30 U	n-Propylbenze	ne		2.30	U
Chloroform			<u>2.30 U</u>	2-Chlorotoluer	1¢		2.30	U
1,1,1-Trichloroet			2.30 U	1,3,5-Trimethy	lbenzene		2.30	U
1,1-Dichloroprop	bene		2.30 U	4-Chlorotoluer	ne		2.30	
Carbon tetrachlor	ride		2.30 U	tert-Butylbenze	ene		2.30	U
Benzene			<u>2.30 U</u>	1,2,4-Trimethy	lbenzene		2.30	U
1,2-Dichloroetha	ne		2.30 U	sec-Butylbenze	ene		2.30	U
Trichloroethene	·		2.30 U	1,3-Dichlorobe	enzene		2.30	U
1,2-Dichloroprop			2.30 U	p-Isopropyltoh	uene		5.75	U
Dibromomethane			2.30 U	1,4-Dichlorobe	enzene		2.30	U
Bromodichlorom			2.30 U	n-Butylbenzen	e			U
2-Chloroethylvin			5.75 U	1,2-Dichlorobe	enzene		2.30	···· ·
Methyl isobutyl l			2.30 U	1,2-Dibromo-3	-chloropropane		2.30	
cis-1,3-Dichlorop	propene		2.30 U	1,2,4-Trichlord			2.30	
Toluene	· · · · · · · · · · · · · · · · · · ·		2.30 U	Hexachlorobut	adiene		2.30	
trans-1,3-Dichlor	opropene		2.30 U	Naphthalene			4,34	 J
1,1,2-Trichloroet	hane		2.30 U	1,2,3-Trichlord	benzene		2.30	

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	103	70-130
1,2-Dichloroethane-d4	106	70-130
Toluene-d8	89	70-130
4-Bromofluorobenzene	91	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



ARCADIS

Tidewater MGP

Trip Blank SDG: N/A

Matrix: Sediment Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-33 Associated Blank: VS072408B15 Concentration Units: µg/Kg

				Sample	Final	Dilution	<u>υπτο, μ</u> ε/.	***6
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst	t l
07/02/08	07/11/08	07/24/08	100	5.00	5	1	BAS	
Parameter	l	F	Result	Parameter			Result	;
Dichlorodifluoro	methane		2.00 U J	2-Hexanone			5.00 U	υJ
Chloromethane			2.00 U	Tetrachloroeth	iene			U (
Vinyl chloride			2.00 U	1,3-Dichlorop	ropane		2.00 0	U
Bromomethane			2.00 U	Dibromochlor	omethane		2.00 1	
Chloroethane			2.00 U	1,2-Dibromoe	thane		2.00 1	υŤ
Trichlorofluoron	nethane		2.00 U	Chlorobenzen	¢		2.00 1	
Acetone		······	9.42 B	1,1,1,2-Tetrac			2.00 1	
1,1-Dichloroethe	ene		2.00 U	Ethylbenzene			2.00 0	
Carbon disulfide			2.00 U	p/m-Xylene			4.00 T	
Methylene chlor	ide		5.00 U	o-Xylene		· · · · · · · · · · · · · · · · · · ·		υŤ
Methyl tert-buty	l ether (MTBE)		2.00 U	Styrene			2.00 T	σΤ
trans-1,2-Dichlo	roethene		2.00 U	Bromoform				U
1,1-Dichloroetha	me		2.00 U	Isopropylbenz	ene		2.00 1	U
Vinyl acetate			2.00 U	1,1,2,2-Tetrac	hloroethane		2.00 1	U
2-Butanone (ME	CK)		2.00 U	Bromobenzen	e		2.00 1	U
cis-1,2-Dichloro	ethene		2.00 U	1,2,3-Trichloropropane				U
2,2-Dichloroprop	oane		2.00 U	n-Propylbenze	ne		2.00 U	U
Chloroform			2.00 U	2-Chlorotolue	ne		2.00 t	U
1,1,1-Trichloroe	thane		2.00 U	1,3,5-Trimethy	ylbenzene			U
1,1-Dichloroprop	oene		2.00 U	4-Chlorotolue	ne		2.00 1	U
Carbon tetrachlo	ride		2.00 U	tert-Butylbenz	ene		2.00 U	U
Benzene			2.00 U	1,2,4-Trimethy	ylbenzene		2.00 T	U
1,2-Dichloroetha	ine		2.00 U	sec-Butylbenz	ene		2.00 (U
Trichloroethene			2.00 U	1,3-Dichlorob	enzene		2.00 T	U
1,2-Dichloroprop	bane		2.00 U	p-Isopropyltol	uene		5.00 U	U
Dibromomethan	e		2.00 U	1,4-Dichlorob	enzene		2.00 t	υT
Bromodichloron	nethane		2.00 U	n-Butylbenzer	ie		5.00 U	U
2-Chloroethylvir	yl ether		5.00 U	1,2-Dichlorob	enzene		2.00 l	IJ
Methyl isobutyl	ketone (MIBK)		2.00 U	1,2-Dibromo-	3-chloropropane		2.00 t	
cis-1,3-Dichloro	propene		2.00 U	1,2,4-Trichlor			2.00 (
Toluene			2.00 U	Hexachlorobu	tadiene		2.00 (υl
trans-1,3-Dichlo	ropropene		2.00 U	Naphthalene			5.00 T	U
1,1,2-Trichloroe	thane		2.00 U 🖞	1,2,3-Trichlor	2.00 T	U V		

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	104	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	83	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 11:24

	Client: Project: Client ID: Case:	ARCADI Tidewate Trip Blank N/A	r MGP	N/A		ETR: Lab II	Code: MA0 0807059 D: 0807059 siated Blank;	-33E	2308B15
	Matrix:	Sediment				Conce	entration Uni	te. III	g/Kg
Date Date	Date	Percent	Methanol	Sample	Extract	Final	Dilution	<u>, , , , , , , , , , , , , , , , , , , </u>	
Collected Received	Analyzed	Solid				Volume (ml)		Ana	lvet
07/02/08 07/11/08	07/23/08	100	5	5.00	0.1	5	1	BA	
Parameter		Res	ult	Parameter		L	L	Result	<u></u>
Dichlorodifluoromethane			100 U T	2-Hexanon			-		¥7 ~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~
Chloromethane			100 U	Tetrachloro				250	
Vinyl chloride			100 U	1,3-Dichlor				100	
Bromomethane			106 B		loromethane			100	f
Chloroethane			100 U	1,2-Dibron				100	
Trichlorofluoromethane	· · · • ······		100 U	Chloroben:				<u>100</u> 100	يؤد به معمد ت
Acetone	• • • • • • • • • • • • • • • • • • • •		232 JB		rachloroethar			100	UU
1,1-Dichloroethene		the second balance and the second	100 U	Ethylbenze		10		100	
O-1-1-101			100 U	p/m-Xylen				200	UU
Methylene chloride			250 U	o-Xylene				100	
Methyl tert-butyl ether (M	TBE)		100 U	Styrene				100	
trans-1,2-Dichloroethene	()		00 U	Bromoforn				100	
11 101 11 1			100 U	Isopropylb			·····	100	U
Vinyl acetate			100 U		rachloroethar			100	
2-Butanone (MEK)			100 U	Bromobenz				100	
cis-1,2-Dichloroethene			100 U		loropropane			100	
2,2-Dichloropropane			100 U	n-Propylbe	nzene			100	U
Chloroform			100 U	2-Chloroto				100	
1,1,1-Trichloroethane			00 U		ethylbenzene			100	U
1,1-Dichloropropene			00 U	4-Chloroto				100	
Carbon tetrachloride			00 U	tert-Butylb				100	cane ander
Benzene]	00 U		ethylbenzene			100	U
1,2-Dichloroethane			00 U	sec-Butylb				*************	UT
Trichloroethene]	00 U	1,3-Dichlor				100	U
1,2-Dichloropropane			00 U	p-Isopropy	ltoluene			250	Ū
Dibromomethane]	.00 U	1,4-Dichlor	robenzene			100	····
Bromodichloromethane]	00 U	n-Butylben					Ŭ
2-Chloroethylvinyl ether			250 U	1,2-Dichlor				100	
Methyl isobutyl ketone (M	IBK)]	00 U	1,2-Dibron	no-3-chloropr	opane		100	
cis-1,3-Dichloropropene			00 U		lorobenzene			100	
Toluene			00 U	Hexachloro	butadiene			100	U
trans-1,3-Dichloropropene	,	<u>[</u>]	00 U	Naphthaler				250	U
1,1,2-Trichloroethane			100 U 🦞	1,2,3-Trich	lorobenzene			100	UV

Surrogate	% Recovery	Acceptance Range (%)
2-Bromo-1-chloropropane	101	70-130
1-Chloro-2-fluorobenzene	111	70-130
1,4-Dichlorobutane	96	70-130
Dibromofluoromethane	101	70-130
Toluene-d8	93	70-130
4-Bromofluorobenzene	95	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

WOODS	CADIS lewater MG 3-S1 SDG: iment		ET La As	b Code: MA0 R: 0807059 b ID: 0807059 sociated Blank: oncentration Un	9-01 : SS071808B06			
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/18/08	07/26/08	59.3	5.41	10	1	SEO
	Para	ameter			F	Result	<u>.</u>	
	Napl	hthalene				281		
	Acer	naphthylene		268				
	Ace	naphthene				199		
	Fluo					267		
		anthrene				2710		
		iracene		- ~ 11 × 11 Part # 1 +		640		
		ranthene		5030				
	Pyre		·····			4330		
		z[a]anthracene			** ***	1950		
	······································	sene				2390		
Benzo[b]fluoranthene						1990	· · · · ·	
Benzo[k]fluoranthene						1860		
Benzo[a]pyrene					2110			
		no[1,2,3-cd]pyr				1380		
		nz[a,h]anthrace				331		
	Benz	o[g,h,i]perylen	<u>e</u>			1430		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	95	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	96	30-150

N/A - Not Applicable

Client: ARCADIS Project: Tidewater MGP Client ID: SC-24-S1 N/A SDG: N/A Matrix: Sediment						ET Lat Ass Cor	ncentration Ur	9-02 : SS071808B06
Date	Date	Date	Date	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
Collected	Received 07/11/08	Extracted 07/18/08	Analyzed	82.6	30.02	20	1	SEO
07/08/08			01121100		F	Result		,
	Parameter Naphthalene Acenaphthylene Acenaphthene Fluorene Phenanthrene Anthracene Fluoranthene Pyrene Benz[a]anthracene Chrysene Benzo[b]fluoranthene Benzo[b]fluoranthene Benzo[k]fluoranthene Benzo[a]pyrene Indeno[1,2,3-cd]pyrene Dibenz[a,h]anthracene Benzo[g,h,i]perylene					461 283 3 454 514 514 4620 2060 6000 5080 2730 2780 1780 1980 2720 1330 344 1160 1160		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	98	30-150
Pyrene-d10	113	30-150
Benzo[b]fluoranthene-d12	107	30-150

N/A - Not Applicable

WOODS A	PHA P	roject: Tid lient ID: SC- ase: N/A				ET La As	b Code: MA0 TR: 0807059 b ID: 0807059 ssociated Blank pncentration Un	9-03 : SS071808B06
Date	Date	Date	Date	D (C.K.I	Sample	Final	Dilution Factor	Analyst
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	1	SEO
07/08/08	07/11/08	07/18/08	07/27/08	42.5	7.07	10	1	SEU
	Para	ameter			F	Result		
	Nap	hthalene				399	August 111	
		naphthylene			615			
		naphthene			261			
	Fluc	rene				367	<u> </u>	
	Pher	nanthrene				3520		
	Antl	hracene				1030		
	Fluc	oranthene				8150		
	Pyre	ene			7400			
	·	z[a]anthracene			3590			
	Chrysene					4140		
	Benzo[b]fluoranthene					3950		
	Benzo[k]fluoranthene					3810		
	Benzo[a]pyrene					4090		
Indeno[1,2,3-cd]pyrene					3040			
		enz[a,h]anthrac		<u> </u>		713		
	Ben	zo[g,h,i]peryle	ne			2870		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	96	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	99	30-150

N/A - Not Applicable

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Page 70 / 277

Client: WOODSHOLELABSCLIENTID: WOODSHOLELABSCLIENTID: Case: Matrix: Sediment ARCADIS Tidewater MGP SC-10-S1 N/A SDG: N/A						ET La As	b Code: MAC R: 0807059 b ID: 080705 sociated Blank ncentration Ur	9-04 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/18/08	07/27/08	82.4	30.15	11.11	1	SEO
L	Para	ameter			F	lesult		
	Nap	hthalene				438		
		naphthylene		224				
		naphthene			465			
		orene				546		
	Phe	nanthrene				4430		
	Ant	hracene				1460		
	Fluc	oranthene				7110		
	Pyre					6060		
		z[a]anthracene				3910		
		ysene				3760		
	Benzo[b]fluoranthene					3540		
Benzo[k]fluoranthene					3350	· · · · · · · · · · · · · · · · · · ·		
Benzojajpyrene					4430			
Indeno[1,2,3-cd]pyrene			,		<u>2560</u> 725			
Dibenz[a,h]anthracene					2250			
	Ben	zo[g,h,i]peryle				<u> </u>		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	97	30-150
Pyrene-d10	110	30-150
Benzo[b]fluoranthene-d12	103	30-150

N/A - Not Applicable

08/04/08 01:39

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Client: ARCADIS Client: Droject: Tidewater MGP SC-12-S1 N/A SDG: N/A Matrix: Sediment						ET Lal As	b Code: MA 0 R: 0807059 b ID: 080705 sociated Blank ncentration Un	9-05 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/18/08	07/27/08	86.0	30.40	5	1	SEO
L	Para	ameter			F	Result		
	Nap	hthalene				57.7		
		naphthylene		173				
		naphthene			24.2			
		rene				34.9		
	Pher	nanthrene			480			
	Anth	nracene				235		
	Fluo	oranthene		1380				
	Pyre	ne				1170	20.1.FTTTT	
	Ben	z[a]anthracene			697			
	Chrysene				630			
	Benzo[b]fluoranthene					491		
	Benzo[k]fluoranthene					487		
	Benzo[a]pyrene					616		
Indeno[1,2,3-cd]pyrene						360		
		enz[a,h]anthrac				88.2		
Benzo[g,h,i]perylene						307		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	92	30-150
Pyrene-d10	111	30-150
Benzo[b]fluoranthene-d12	100	30-150

N/A - Not Applicable

WOODS	PHA P	Project: Tid Client ID: SC- Case: N/A				ET Lai As	b Code: MA(R: 0807059 b ID: 080705 sociated Blank	9-06 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	10 0
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/18/08	07/27/08	76.7	20.93	20	1	SEO
	Par	ameter			F	Result		
	Nap	hthalene				3780		
		naphthylene		1100				
	Ace	naphthene				2530		
	Fluc	orene		<u>3140</u> 24800				
		nanthrene						
		nracene			4	8130		
	<u>Fluc</u>	ranthene			24400 3	2 8800 E 🗋	<u>></u>	
	Pyre			23300				
Benz[a]anthracene						2400		
Chrysene]	1000			
Benzo[b]fluoranthene					8790			
Benzo[k]fluoranthene Benzo[a]pyrene			8440					
]	1900			
		no[1,2,3-cd]py				6780		
		enz[a,h]anthrac				1710		
	Ben	zo[g,h,i]perylen	<u>e</u>			5850		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	93	30-150
Pyrene-d10	105	30-150
Benzo[b]fluoranthene-d12	98	30-150

N/A - Not Applicable

E - Estimated value, exceeds the upper limit of calibration.

WOODS ,	PHA P	roject: Tic lient ID: SC- ase: N/A				ET Lal Ass	b Code: MA R: 0807059 b ID: 080705 sociated Blank ncentration Ur	9-06E :: SS071808B06
Date	Date	Date	Date	D . 0 111	Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/08/08	07/11/08	07/18/08	08/02/08	76.7	20.93	20	10	SEO
	Par	ameter			F	Result		
	Nap	hthalene			125 U			
		naphthylene			125 U			
		naphthene			125 U			
		orene				125 U		
	Pher	nanthrene				125 U		
	Antl	nracene			125 U			
	Fluc	oranthene			24400			
	Pyre				125 U			
		z[a]anthracene				125 U		
		ysene				125 U		
	Benzo[b]fluoranthene					<u>125 U</u>		
	Benzo[k]fluoranthene					<u>125 U</u>		
Benzo[a]pyrene				125 U 125 U				
	Indeno[1,2,3-cd]pyrene				125 U 125 U	,, .		
	Dibenz[a,h]anthracene Benzo[g,h,i]perylene				125 U			
		2913,1,1,1,001 y 101						

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	60	30-150
Pyrene-d10	90	30-150
Benzo[b]fluoranthene-d12	74	30-150

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

WOODS ,	PHA P	Project: Tid Client ID: SC3 Case: N/A				e ET Lal As	b Code: MAC R: 0807059 b ID: 080705 sociated Blank ncentration Un	9-07 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/18/08	08/02/08	80.2	30.44	66.67	10	AC
		ameter			F	Result		
		hthalene		2530				
	·······	naphthylene				<u>6060</u> 4190		
	ww.w	naphthene			1	<u>4190</u> 12100		
		nanthrene				20000		
	2	racene				52300		
		ranthene				12000		
	Pyre	······)8000	· · · · · · ·	
		z[a]anthracene				54900		
	Chrysene				4	45000		
	Benzo[b]fluoranthene					29500		
Benzo[k]fluoranthene					32200			
Benzo[a]pyrene					43300			
	Indeno[1,2,3-cd]pyrene					23000		
	-	enz[a,h]anthrac			,	5870	at 4400 - 48 100 - 47 47 47 47 47	
Benzo[g,h,i]perylene					•	18000		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	63	30-150
Pyrene-d10	83	30-150
Benzo[b]fluoranthene-d12	123	30-150

N/A - Not Applicable

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08/04/08 14:33

Client: ARCADIS Project: Tidewater MGP Client ID: SC41-S1 N/A SDG: N/A Matrix: Sediment						ET La As	b Code: MA0 R: 0807059 b ID: 0807059 sociated Blank: ncentration Uni	-08 SS071808B06
Date	Date	Date	Date	D	Sample	Final	Dilution	A
Collected 07/09/08	Received 07/11/08	Extracted 07/18/08	Analyzed 07/27/08	Percent Solid 81.3	Amount (g) 30.90	Volume (ml) 5	Factor	Analyst SEO
07709708		ameter	0//2//08	01.5		Result	<u> </u>	510
					r			
		hthalene naphthylene			······································	<u>42.9</u> 113		
		naphthene				26.6		
		rene				48.3		
	Pher	nanthrene				537		
	Antl	nracene				201		
	Fluc	oranthene				1080		
	Pyre	me				874		
	Ben	z[a]anthracene				557		
	Chrysene					524	· · · · · · · · · · · · · · · · · · ·	
Benzo[b]fluoranthene					4.5/ - .	444		
Benzo[k]fluoranthene						434		
Benzo[a]pyrene					540			
Indeno[1,2,3-cd]pyrene			······································		335			
		enz[a,h]anthrac				89.9		
Benzo[g,h,i]perylene						296		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	107	30-150
Benzo[b]fluoranthene-d12	100	30-150

N/A - Not Applicable

08/04/08 01:40

Client: ARCADIS Project: Tidewater MGP Client ID: Client ID: Case: N/A SDG: N/A Matrix: Sediment						ET Lal As	o Code: MA0 R: 0807059 b ID: 080705 sociated Blank ncentration Un	9-09 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/18/08	07/27/08	84.4	30.84	5	1	SEO
· · · · · · · · · · · · · · · · · · ·	Para	ameter			F	Result		
	Ace	hthalene naphthylene		<u>41.7</u> <u>89.3</u>				
		naphthene				20.5	<u> </u>	
		anthrene				<u>39.7</u> 527		
		nracene				<u> </u>		
		ranthene				1090		
	Pyre					870	·····	
		z[a]anthracene				522		
	Chrysene					538		
	Benzo[b]fluoranthene					449		
	Benzo[k]fluoranthene					461		
	Benzo[a]pyrene					532		
Indeno[1,2,3-cd]pyrene						359		
		enz[a,h]anthrac				93.4		
	Ben	zo[g,h,i]peryler	1ê			311		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	93	30-150
Pyrene-d10	112	30-150
Benzo[b]fluoranthene-d12	103	30-150

N/A - Not Applicable

WOODS	CADIS lewater MG -S1 SDG: iment		ET La As	b Code: MA0 R: 0807059 b ID: 0807059 sociated Blank: ncentration Un	9-10 : SS071808B06			
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/18/08	07/27/08	89.3	30.33	33.33	1	SEO
	Par	ameter			F	Result		
		hthalene		651				
		naphthylene		917				
		naphthene				565		
		orene				508		
		nanthrene		5410				
	÷	hracene				1940		
		oranthene				9230		
	Pyre			8290				
		z[a]anthracene			A. 17	3800		
	Chrysene					3590		
	Benzo[b]fluoranthene Benzo[k]fluoranthene				······	2770	and the second second of the second	
	Benzo[a]pyrene					<u>2840</u> 4450		
						2130		
	Indeno[1,2,3-cd]pyrene Dibenz[a,h]anthracene					<u>2130</u> 540		
		zo[g,h,i]perylen				<u> </u>		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	93	30-150
Pyrene-d10	112	30-150
Benzo[b]fluoranthene-d12	106	30-150

N/A - Not Applicable

WOODS	PHA P	Project: Tid Client ID: SC2 Case: N/A				ET Lab Ass	o Code: MA(R: 0807059 o ID: 080705 sociated Blank ncentration Ur	9-11 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/18/08	07/27/08	80.2	30,90	6.67	1	SEO
i	Para	ameter			F	Result		
	Ace Ace	hthalene naphthylene naphthene				50.1 204 38.0		
		nanthrene				<u>71.9</u> 880		
		hracene		326				
	Aug	oranthene				1620		
	Pyre	ene	f+			1480		
		z[a]anthracene				788		
		ysene				742	z.j.,	
	Benzo[b]fluoranthene					582		
Benzo[k]fluoranthene					561			
Benzo[a]pyrene					744			
	Indeno[1,2,3-cd]pyrene					456		
		enz[a,h]anthrac				118		
	Ben	zo[g,h,i]peryle	ne			418		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	96	30-150
Pyrene-d10	113	30-150
Benzo[b]fluoranthene-d12	106	30-150

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N/A - Not Applicable

Client: ARCADIS Project: Tidewater MGP SC5-S1 N/A SDG: N/A Matrix: Sediment						ETI Lab Ass	Code: MA R: 0807059 DD: 080705 sociated Blank incentration Un	9-12 :: SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/18/08	07/27/08	81.5	30.14	20	1	SEO
<u></u>	Para	ameter			F	Result		
		hthalene			226			
	Ace	naphthylene			252			
	Ace	naphthene			87.2			
		orene			126			
		nanthrene			711			
		hracene				340		
		oranthene			1340			
	Pyre				1410			
		z[a]anthracene			733			
	Chrysene				715			
	Benzo[b]fluoranthene				536			
	Benzo[k]fluoranthene					550		
	Benzo[a]pyrene					828		
	Indeno[1,2,3-cd]pyrene					462		
		enz[a,h]anthrac				120		
	Benzo[g,h,i]perylene					606		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	94	30-150
Pyrene-d10	112	30-150
Benzo[b]fluoranthene-d12	106	30-150

N/A - Not Applicable

Alkylated Polynuclear Aromatic Hydrocarbons

WOODS	PHA P	Project: Ti Client ID: SC Case: N/	RCADIS dewater MG 6-S1 A SDG: liment		Lab ID: 0807059-13			
Date	Date	Date	Date		Sample	Final	Dilution	
Collected 07/09/08	Received 07/11/08	Extracted 07/18/08	Analyzed 07/27/08	Percent Solid 30.9	Amount (g) 6.12	Volume (ml) 6.67	Factor	Analyst SEO
07/09/00		ameter	0//2//08	50.9		Result *	1	32.0
					Г			
		hthalene Naphthalenes				<u>636</u> 376		
		Naphthalenes				392		
		Naphthalenes		·····		290		
		Naphthalenes				215		
	Ace	naphthylene				962		
		naphthene				235		
		orene		308				
		Fluorenes Fluorenes		<u> </u>				
		Fluorenes		765				
		anthrene		3180				
		nracene		1080				
	C1-I	Phenanthrenes	Anthracenes	1790				
		Phenanthrenes		1270				
		Phenanthrenes				758		
		Phenanthrenes	Anthracenes			<u>394</u>		
	<u>Fluo</u> Pyre	oranthene				<u>10600</u>		
		Fluoranthenes/	Pyrenes	4290				
		z[a]anthracene		····		4020	·····	
		ysene		5420				
		Chrysenes		2620				
		Chrysenes		1580				
	<u>C3-Chrysenes</u>				1340			
		Chrysenes				793		
		zo[b]fluoranth zo[k]fluoranth			······································	<u>5740</u> 4890		
		zo[a]pyrene	<u></u>			5460		
	·	lene				1360		
		zo[e]pyrene				4330		
		no[1,2,3-cd]p				4400		
		enz[a,h]anthra			······································	1010		
	Ben:	zo[g,h,i]peryle	ne			4160		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	95	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	103	30-150

N/A - Not Applicable

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WOODS	PHAP HOLE LABS	Project: Tid Client ID: SC2 Case: N/A				ET Lat Ass	D Code: MA0 R: 0807059 D ID: 080705 sociated Blank ncentration Un	9-14 : SS071808B06
Date	Date	Date	Date		Sample .	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/09/08	07/11/08	07/18/08	07/28/08	77.6	10.03	6.67	1	SEO
	Para	ameter			F	Result		
	Ace	hthalene naphthylene naphthene				358 189 746		
		brene				835		
		nanthrene		·····		6160		
	Ant	hracene				2170		
	Fluc	oranthene				7030		
	Pyre					5620		
	Ben	z[a]anthracene				3120		
		ysene				2870		
	Ben	zo[b]fluoranthe	ene			2050		
	Ben	zo[k]fluoranthe	ene			2230		
		zo[a]pyrene				2830		
		eno[1,2,3-cd]py				1500		
		enz[a,h]anthrac				396		
	Ben	zo[g,h,i]peryle	ne			1270	//=	

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	92	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

WOODS A	PHA P	roject: Tid lient ID: SC2 ase: N/A				ET Lal As	b Code: MA0 R: 0807059 b ID: 0807055 sociated Blank: ncentration Un	9-15 SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	83.5	21.64	10	1	SEO
	Para	ımeter			F	lesult		
	Ace	hthalene naphthylene				52.6 134		
		naphthene				44.7		
		rene				75.4		
		nanthrene				1220	· · · · · · · · · · · · · · · · · · ·	
		nracene				379	<u> </u>	
		ranthene				2680		
	Pyre		~~~~~			2140 1160		
		z[a]anthracene /sene			a	1170		
		zo[b]fluoranthe				954		
Benzo[k]fluoranthene					913			
Benzo[a]pyrene				······································	1100			
		no[1,2,3-cd]py	rene			703		
		enz[a,h]anthrac				178		
		zo[g,h,i]peryler				624		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	94	30-150
Pyrene-d10	114	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

Alkylated Polynuclear Aromatic Hydrocarbons

Client: ARCADIS Project: Tidewater MGP Client ID: SC19-S1 Case: N/A SDG: N/A Matrix: Sediment						Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-16 Associated Blank: SS071808 Concentration Units: μg/Kg		
Date Collected	Date Received	Date Extracted	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	35.7	6.34	12.5	1	SEO
07/10/00		imeter	01720/00			Result		
	Napl C1-N C2-N C3-N C4-N Accer Accer Fluo C1-F C2-F C3-F Pher Anth C1-F C2-F C3-F Pher Anth C1-F C2-F C3-F Pher Anth C1-F C2-F C3-F Pher C1-F C2-F C3-F	hthalene Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthene rene Fluorenes Fluorenes Fluorenes Pluorenes Pluorenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes	/Anthracenes /Anthracenes /Anthracenes Pyrenes			1570 787 1770 2640 2640 1160 907 766 882 1660 2200 4550 2770 5380 4820 3890 3780 8310 8940 5900 4310 4920 3310 2230 1980 1160		
	Ben: Ben Ben	zo[b]fluoranth zo[k]fluoranth zo[a]pyrene /lene				3310 3530 4480 989	······································	
	Ben Inde	zo[e]pyrene mo[1,2,3-cd]p enz[a,h]anthra				3080 2660 731 2710		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	96	30-150
Pyrene-d10	107	30-150
Benzo[b]fluoranthene-d12	103	30-150

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N/A - Not Applicable

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08/04/08 13:58

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WOODS A	PHA P	roject: Tid lient ID: SC4 ase: N/A	CADIS lewater MG I3-S1 SDG: iment			ET La As	b Code: MA0 TR: 0807059 b ID: 0807059 sociated Blank oncentration Un	9-17 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	A
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	62.0	10.14	10	1	SEO
	Para	ımeter			F	lesult		
		hthalene naphthylene				308 373		
	Ace	naphthene				207		
	Fluo	rene				251		
	Pher	anthrene			2460			
	Anth	racene			726			
	<u>Fluo</u>	ranthene			4520			
	Pyre				3770			
		z[a]anthracene			1940			
		/sene			2030			
	Benzo[b]fluoranthene					1780		
Benzo[k]fluoranthene				1860				
	Benzo[a]pyrene				· · · · · · · · · · · · · · · · · · ·	2130		
		no[1,2,3-cd]py				1400	. <u> </u>	
		enz[a,h]anthrac				341		
	Ben	zo[g,h,i]perylei	<u>1e</u>			1290		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	92	30-150
Pyrene-d10	107	30-150
Benzo[b]fluoranthene-d12	102	30-150

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N/A - Not Applicable

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

WOODS,	PHA P	roject: Tid lient ID: SC4 lase: N/A				ET Lat Ass	D Code: MA0 R: 0807059 D ID: 080705 sociated Blank ncentration Un	9-18 : SS071808B00
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	77.3	30.45	33.33	1	SEO
	Para	ameter			F	Result		
	Acer Acer	hthalene naphthylene naphthene prene				442 338 154 187		
	1.000	anthrene		- · · · · · · · · · · · · · · · · · · ·		1870		
		nracene				603		
		ranthene				3640		
	Pyre	ne				3550		
	Benz	z[a]anthracene				1710		
	Chry	/sene				1750		
	Benz	zo[b]fluoranthe	ene			1270		
	Benzo[k]fluoranthene					1390		
		zo[a]pyrene				1970		
		no[1,2,3-cd]py				1060		
		enz[a,h]anthrac				268		
	Benz	zo[g,h,i]peryler	1e			1070		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	88	30-150
Pyrene-d10	105	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

Alkylated Polynuclear Aromatic Hydrocarbons

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WOODS HOLE LABSClient ID: SC9-S1 Case: N/A SDG: N/A Matrix: Sediment				ETR: 0807059 Lab ID: 0807059-19 Associated Blank: SS071808				
Date	Date	Date	Date		Sample	Cor Final	ncentration Un Dilution	its: µg/Kg
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	36.6	7.51	10	1	SEO
	Para	ameter		J	F	Result		
	Napl	hthalene				541		
		Naphthalenes		·····		366		
		Naphthalenes				449		
		Vaphthalenes				358		
		Naphthalenes				270		
		naphthylene				1020 ×		
		naphthene				286		
		Rhorenes				<u>370</u> 235		
C1-Fluorenes C2-Fluorenes						362		
C3-Fluorenes				,	781			
Phenanthrene					3450			
Anthracene						1310		
	**************************************	henanthrenes/			2090			
		henanthrenes/			1420			
		Phenanthrenes/				888		
		Phenanthrenes/	Anthracenes	····		549		
		ranthene				9520		
	Pyre C1-I	luoranthenes/	Duranac			8450 4370		
	•	z[a]anthracene				4170		
		/sene	•	······································		4850		
		Chrysenes				2640		
C2-Chrysenes					1490			
C3-Chrysenes					1260			
C4-Chrysenes					815			
Benzo[b]fluoranthene					4720			
Benzo[k]fluoranthene					4280			
Benzo[a]pyrene Perylene				*********	<u>4900</u> 1610	<u> </u>		
		zo[e]pyrene				3730	· · · · · · · · · · · · · · · · · · ·	
		no[1,2,3-cd]py	/rene			3560		
		enz[a,h]anthrac				881		
		zo[g,h,i]peryle			· · · · · · · · · · · · · · · · · · ·	3460		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	91	30-150
Pyrene-d10	108	30-150
Benzo[b]fluoranthene-d12	103	30-150

N/A - Not Applicable

08/04/08 13:59

Alkylated Polynuclear Aromatic Hydrocarbons

Client: ARCADIS Project: Tidewater MGP DUP-1 N/A SDG: N/A Matrix: Sediment						e El La As	b Code: MAC R: 0807059 b ID: 080705 sociated Blank oncentration Un	9-20 : SS071808B06
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	36.2	6.33	6.67	1	SEO
	Para	ameter			F	Result		
	Nap	hthalene				615		
		Naphthalenes				382		
		Naphthalenes				425		
		Naphthalenes				334		
		Naphthalenes				263		
		naphthylene				<u>943</u> 305		
		naphthene				382		
		Fluorenes				246		
		Fluorenes		·····		348		
		Fluorenes				916		
	Phenanthrene					3580		
		hracene				1350		
		Phenanthrenes						
		Phenanthrenes,				1350		
		Phenanthrenes,	······································			804		
		Phenanthrenes,	Anthracenes					
		oranthene		<u>9600</u> 8200 ^c				
	Pyre	Fluoranthenes/	Duranag			4210		
		z[a]anthracene				4090		
		ysene				4790		
		Chrysenes		······		2370		
		Chrysenes				1240		
	C3-Chrysenes					1110		
C4-Chrysenes					724			
		zo[b]fluoranth				4530		
		zo[k]fluoranth	ene			4260		
	Benzo[a]pyrene					4780		
		/lene zo[e]pyrene				<u>1530</u> 3620		
		zo[e]pyrene eno[1,2,3-cd]p	vrene			3550		
		enz[a,h]anthra				861		
		zo[g,h,i]peryle				3380		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	92	30-150
Pyrene-d10	108	30-150
Benzo[b]fluoranthene-d12	103	30-150

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N/A - Not Applicable

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08/04/08 14:00

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Page 91 / 277

Client: ARCADIS Project: Tidewater MGP Client ID: SC29-S1 Case: N/A SDG: N/A Matrix: Sediment						ET Lat As	b Code: MA00 R: 0807059 b ID: 0807059 sociated Blank: ncentration Unit	21 SS071808B07
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	42.1	5.42	6.67	1	SEO
	Para	ameter	in the second		F	Result		
	Ace	hthalene naphthylene				1010 1310		
		naphthene				224		
		orene		· · · · · · · · · · · · · · · · · · ·		344		
	<u></u>	nanthrene				3300		
	Providence -	nracene				1310		
		oranthene	and the second			7850		
	Pyre	·····				8570		
		z[a]anthracene				3980	·····	
	Chrysene					4870		
	Benzo[b]fluoranthene				J	4640		
	Benzo[k]fluoranthene					4550		
	Benzo[a]pyrene					<u>4470</u> 3660		
		eno[1,2,3-cd]py				926		
		enz[a,h]anthrac zo[g,h,i]peryle				3540		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	91	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	100	30-150

N/A - Not Applicable

08/04/08 01:43

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

Client: ARCADIS Project: Client ID: Client ID: Case: N/A SDG: Matrix: Sediment						ET Lal As	o Code: MA(R: 0807059 o ID: 080705 sociated Blank ncentration Ur	9-22 :: SS071808B07
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/28/08	52.9	5.57	6.67	1	SEO
	Par	ameter		-	F	Result		
		hthalene naphthylene				288 253		
	Ace	naphthene				59.9		
		orene				88.8		
		nanthrene				821		
		hracene			·····	310		
		oranthene			,	2730		
	Pyre					2340		
		z[a]anthracene				1080		
	Chrysene					1380		
	Benzo[b]fluoranthene					1400		
	Benzo[k]fluoranthene					1320		
	Benzo[a]pyrene					1210		
	Indeno[1,2,3-cd]pyrene					1050		
		enz[a,h]anthrac zo[g,h,i]peryler	•/•			<u>252</u> 1020		
	Den	Loight, porytor				1040		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	108	30-150
Benzo[b]fluoranthene-d12	101	30-150

N/A - Not Applicable

WOODS ,		roject: Tid lient ID: SC2 lase: N/A				ET La As	b Code: MA00 R: 0807059 b ID: 0807059- sociated Blank: ncentration Unit	23 SS071808B07
Date Collected	Date	Date	Date		Sample	Final	Dilution	
07/10/08	Received 07/11/08	Extracted	Analyzed	Percent Solid	(8)	Volume (ml)	Factor	Analyst
07/10/08		07/18/08	07/28/08	34.6	5.25	6.67		SEO
Parameter Naphthalene Acenaphthylene Acenaphthene Fluorene Phenanthrene Anthracene Fluoranthene Pyrene Benz[a]anthracene Chrysene Benzo[b]fluoranthene Benzo[k]fluoranthene Benzo[k]fluoranthene Benzo[k]fluoranthene					tesult 534 834 162 241 2400 942 6810 6820 3440 4220 4040 3970 3910			
	Dibe	no[1,2,3-cd]py nz[a,h]anthraco co[g,h,i]perylen	ene			3190 804 3070		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	89	30-150
Pyrene-d10	107	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

08/04/08 01:42

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

WOODS F	CADIS lewater MG 0-S1 . SDG: iment		ET Lal As	o Code: MA(R: 0807059 b ID: 080705 sociated Blank ncentration Ur	9-24 :: SS071808B07			
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/10/08	07/11/08	07/18/08	07/29/08	53.8	5.07	10	1	SEO
	Para	ameter			F	Result		
	Nap	hthalene		4320				
	Ace	naphthylene		5670				
	Ace	naphthene				757		
	Fluc	orene				1340		
	Pher	anthrene				9510		
	Ant	iracene				4620		
	Fluc	oranthene				19800		
	Pyre	ne				21500		
	Ben	z[a]anthracene				11600		
	Chrysene					11300		
	Benzo[b]fluoranthene					9000		
	Benzo[k]fluoranthene					9570		
Benzo[a]pyrene					11000			
Indeno[1,2,3-cd]pyrene					7690			
	· · · · · · · · · · · · · · · · · · ·	enz[a,h]anthrac				2170 J		
	Ben	zo[g,h,i]peryler	ne			7480		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	106	30-150
Benzo[b]fluoranthene-d12	100	30-150

N/A - Not Applicable

WOODS	PHA P	Project: Tid Client ID: SC4 Case: N/A				ET Lal As	b Code: MAC R: 0807059 b ID: 080705 sociated Blank ncentration Ur	9-25 : SS071808B07
Date	Date	Date	Date		Sample	Final	Dilution	Aurahunt
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/11/08	07/11/08	07/18/08	07/29/08	83.2	10.23	2.86	1	SEO
	Par	ameter			F	Result		
		hthalene naphthylene		<u>215</u> 425				
	Ace	naphthene			51.7			
	Fluc	orene			95.9			
	Phe	nanthrene			1460			
	Ant	hracene			976			
	Fluc	oranthene				6630		
	Pyre					5960		
		z[a]anthracene			3380			
		ysene		,,	2970			
Benzo[b]fluoranthene					2520			
Benzo[k]fluoranthene						2390		
Benzo[a]pyrene						2670		
		ono[1,2,3-cd]py				2070		
		enz[a,h]anthrac zo[g,h,i]peryler				<u>527</u> 1790		
	Den	zol g'u'i lhei Alei				1//		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	87	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

WOODS h	PHA P	Project: Tic lient ID: SC: Case: N/A	CADIS lewater MG 32-S1 SDG: liment			ET Lak Ass	o Code: MA(R: 0807059 o ID: 080705 sociated Blank ncentration Ur	9-26 :: SS071808B07
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/11/08	07/11/08	07/18/08	07/29/08	60.1	20.88	40	1	SEO
() () () () () () () () (Para	ameter			F	Result		
		hthalene				16500		
		naphthylene			8590			
		naphthene			· · · · · · · · · · · · · · · · · · ·	2470		
		brene				5780		
		nanthrene				<u>32800</u>		
		hracene				<u>13000</u> 29500		
		oranthene	, , , , , , , , , , , , , , , , ,			34500		
	<u>Pyre</u> Don	z[a]anthracene				<u>19200</u>		
		vsene		20200				
	··········	zo[b]fluoranthe	ene			11700		
Benzo[k]fluoranthene				· · · · · · · · · · · · · · · · · · ·	13000			
Benzo[a]pyrene					16800	Abr		
Indeno[1,2,3-cd]pyrene					10200	<u> </u>		
		enz[a,h]anthrac				3180		
		zo[g,h,i]peryle				10300		

	Acceptance
% Recovery	Range (%)
92	30-150
112	30-150
104	30-150
	92 112

N/A - Not Applicable

08/04/08 01:42

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Alkylated Polynuclear Aromatic Hydrocarbons

Client: ARCADIS Project: Tidewater MGP SCIII ID: Case: N/A SDG: N/A Matrix: Sediment						E' La A	ab Code: MA(FR: 0807059 ab ID: 080705 ssociated Blank oncentration Ur	9-27 :: SS071808B07
Date	Date Date	Date Extracted	Date	Demont Calle	Sample	Final	Dilution	A
Collected 07/11/08	Received 07/11/08	07/18/08	Analyzed 07/29/08	Percent Solid 31.1	Amount (g) 5.91	Volume (ml) 6.67	Factor 1	Analyst SEO
0771708	L		01123108	51.1			1	SEO
		imeter			r	lesult 4		
		hthalene				528		
		Naphthalenes				288		
		Naphthalenes Naphthalenes				<u>337</u> 268		
		Vaphthalenes				194	************	
		haphthylene				685		
		naphthene	and the second s	222				
		rene		279				
	<u>C1-</u> F	luorenes		173 [°]				
	C2-Fluorenes				308			
	C3-Fluorenes					631		
	Phenanthrene					2710		
	Anthracene					892		
		henanthrenes/		1370				
		henanthrenes/.		<u> </u>				
		henanthrenes/						
		henanthrenes/.	Anunracenes	<u> </u>				
	Pyre			6460				
		Fluoranthenes/I	vrenes			3120		
		z[a]anthracene				3000		
		/sene				3990		
		Chrysenes				1920		
	<u>C2-C</u>	Chrysenes				1130		
	C3-Chrysenes					1060		
	C4-Chrysenes				· · · · · · · · · · · · · · · · · · ·	684		
Benzo[b]fluoranthene					3940			
	Benzo[k]fluoranthene					3600		
	Benzo[a]pyrene Perylene					3620		
		zo[e]pyrene			······································	996 3080		
		no[1,2,3-cd]py	rene	<u></u>		2990		
		no[1,2,5-cu py nz[a,h]anthrac				720		
		zo[g,h,i]peryler		· · · · · · · · · · · · · · · · · · ·		2850		
		· · · · · · · · · · · · · · · · · · ·						

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	89	30-150
Pyrene-d10	106	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

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08/04/08 14:00

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Client: ARCADIS Project: Tidewater MGP WOODS HOLE LABS Client ID: Client ID: Client ID: Case: N/A SDG: N/A Matrix: Sediment						ET La As	b Code: MA0 R: 0807059 b ID: 0807059 sociated Blank: ncentration Uni	9-28 SS071808B06
Date Collected	Date Received	Date Extracted	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
07/11/08	07/11/08	07/18/08	08/02/08	39.6	5.13	200	10	Analyst
	Para	ameter		<u> </u>	F	lesult	↓L	······································
	Naphthalene Acenaphthylene Acenaphthene				42)4000 29000 53000		
		rene				38000 38000		
		nanthrene		2350000				
		iracene				20000		
		ranthene		3460009				
	<u>Pyre</u> Born	ne z[a]anthracene		<u> </u>				
		/sene				35000 35000		
	Benzo[b]fluoranthene)4000		
Benzo[k]fluoranthene					57000			
Benzo[a]pyrene				8'	72000			
Indeno[1,2,3-cd]pyrene					48000			
		enz[a,h]anthrac				1000	······································	
Benzo[g,h,i]perylene					4;	54000		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	117	30-150
Pyrene-d10	105	30-150
Benzo[b]fluoranthene-d12	107	30-150

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N/A - Not Applicable

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08/04/08 14:35

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320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

WOODS ,	PHA P	roject: Tid lient ID: SC3 ase: N/A				ET. Lat Ass	Code: MA(R: 0807059 DD: 080705 sociated Blank ncentration Un	9-29 :: SS071808B07
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/11/08	07/11/08	07/18/08	07/29/08	23.7	5.08	6.67	1	SEO
	Para	ameter			F	lesult		
	Acer	hthalene naphthylene				514 840		
		naphthene		176				
	Fluo					241		
	A. W	anthrene		1900				
	••••••••••••••••••••••••••••••••••••••	racene						
		ranthene		6200				
	Pyre					5730	<u> </u>	
		z[a]anthracene		2700				
	Chrysene					3390		
Benzo[b]fluoranthene						3680		
Benzo[k]fluoranthene					3220			
Benzo[a]pyrene					3040			
Indeno[1,2,3-cd]pyrene					2720			
		mz[a,h]anthrac zo[g,h,i]peryler				<u>682</u> 2660		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	105	30-150
Benzo[b]fluoranthene-d12	102	30-150

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N/A - Not Applicable

WOODS	PHA P	Project: Tid Client ID: SC: Case: N/A				ET Lal As	b Code: MA R: 0807059 b ID: 080705 sociated Blank ncentration Ur	9-30 :: SS071808B07
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/11/08	07/11/08	07/18/08	07/29/08	34.5	5.21	6.67	1	SEO
	Para	ameter			F	Result		· · · · · · · · · · · · · · · · · · ·
		hthalene naphthylene		<u> </u>				
	Acer	naphthene				314		
	Fluo	rene				481		
	Pher	anthrene		v - 17,1 = 14 - 14 - 14 - 14 - 14 - 14 - 14 - 1		4010		
	Anth	iracene				1840		
	Fluo	ranthene				8880		
	Pyre	ne				9410		
	Benz	z[a]anthracene				4780		
		/sene			5390			
Benzo[b]fluoranthene					5170			
Benzo[k]fluoranthene					5070			
Benzo[a]pyrene						5030		
Indeno[1,2,3-cd]pyrene						4260		
		nz[a,h]anthrac				1110		
	Benz	zo[g,h,i]perylen	le			4070		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	103	30-150
Benzo[b]fluoranthene-d12	100	30-150

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N/A - Not Applicable

Date Date	Project: Ti Client ID: SC Case: N/A Matrix: Sec	A SDG: liment			ETI Lab Ass	Code: MA R: 0807059 DD: 080705 sociated Blank accentration Ur	9-31 :: SS071808B07
Date Date Collected Received	Date Extracted	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution	
07/11/08 07/11/08	07/18/08	07/29/08	53.9	5,44	4	Factor	Analyst SEO
N A A F P A F P B C B B C B B C B B C B B C B B C B B C B B C B B C D D	arameter aphthalene .cenaphthylene .cenaphthene luorene henanthrene nthracene luoranthene yrene enz[a]anthracene hrysene enzo[b]fluoranthe enzo[k]fluoranthe enzo[a]pyrene deno[1,2,3-cd]py ibenz[a,h]anthrac	rene ene		F	Result 131 246 50.5 79.0 836 295 2390 2090 954 1240 1140 1050 946 239 911		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	106	30-150
Benzo[b]fluoranthene-d12	102	30-150

N/A - Not Applicable

08/04/08 01:43

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WOODS	PHA P	Project: Tic lient ID: SC: Case: N/A	CADIS lewater MG 55-S1 SDG: iment			ET La As	b Code: MA0 TR: 0807059 b ID: 0807059 ssociated Blank: oncentration Uni	0-32 SS071808B07	
Date	Date	Date	Date		Sample	Final	Dilution		
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst	
07/11/08	07/11/08	07/18/08	07/29/08	55.1	10.73	20	1	SEO	
	Parameter				F	Result			
	Nap	hthalene				4290			
	Acenaphthylene					4920			
	Ace	naphthene		650					
		rene		1750					
		nanthrene		18200					
		nracene	P.t.=1	6770					
		ranthene				19300			
	Pyre					23400			
	Statement and all	z[a]anthracene				12300			
	No.	/sene			1	12900			
		zo[b]fluoranthe				7840			
	Benzo[k]fluoranthene					8810			
	Benzo[a]pyrene				1	1400			
	Indeno[1,2,3-cd]pyrene					6980			
		nz[a,h]anthrac				2140			
	Benz	zo[g,h,i]peryler	e			6940			

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	90	30-150
Pyrene-d10	102	30-150
Benzo[b]fluoranthene-d12	99	30-150

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N/A - Not Applicable

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08/04/08 01:43

Page 105 / 277



ARCADIS : Tidewater MGP

SDG: N/A

Client ID: SC23-S1 Matrix: Sediment Percent Solid: 59.3

N/A

Lab Code: MIA00030 ETR: 0807059 Lab ID: 0807059-01 Concentration Units: mg/Kg Date Collected: 07/08/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.484	J	0.119	5	08/05/08	07/29/08	6020A	LCP
Arsenic	8.39	5	0.122	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.445	T	0.0488	5	08/01/08	07/29/08	6020A	LMR
Cadmium	5.87	T	0.0488	5	08/01/08	07/29/08	6020A	LMR
Chromium	85.0	J	0.488	5	08/01/08	07/29/08	6020A	LMR
Copper	173		0.244	5	08/01/08	07/29/08	6020A	LMR
Lead	178	J	0.122	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.374	T	0.0187	5	07/31/08	07/29/08	7474	LCP
Nickel	24.9		0.244	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.918		0.244	5	08/01/08	07/29/08	6020A	LMR
Silver	1.94		0.119	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.132		0.0488	5	08/01/08	07/29/08	6020A	LMR
Zinc	320		1.22	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable



ARCADIS t: Tidewater MGP

SDG: N/A

'n

Matrix: Sediment Percent Solid: 82.6

N/A

SC-24-S1

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-02 Concentration Units: mg/Kg Date Collected: 07/08/08 Date Received: 07/11/08

Parameter	Result	Oualifier	Reporting Limit	Dilution	Date	Date	Analytical	
A		Yuumitei	L.11010C	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.206	<u> </u>	0.0958	5	08/05/08	07/29/08	6020A	LCP
Arsenic	2.91		0.0971	5	08/01/08	07/29/08	6020A	LMR
Beryllium	1.11	T	0.0388	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.619	J	0.0388	5	08/01/08	07/29/08	6020A	LMR LMR
Chromium	10.4	J	0.388	5	08/01/08	07/29/08	6020A	LMR
Copper	45.6		0.194	5	08/01/08	07/29/08	6020A	
Lead	60.4	June	0.0971	5	08/01/08	07/29/08	6020A	
Mercury	0.0252	J	0.0139	5	07/31/08	07/29/08	7474	LIVIR
Nickel	13.4		0.194	5	08/01/08	07/29/08	6020A	LCP LMR
Selenium	0.318		0.194	5	08/01/08	07/29/08	6020A	LMR
Silver	0.120		0.0958	5	08/05/08	07/29/08	6020A	
Thallium	0.0388	U	0.0388	5	08/01/08	07/29/08		LCP
Zinc	105		0.971		08/01/08		6020A	LMR
					00/01/08	07/29/08	6020A	LMR

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

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08/07/08 06:50

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Page 114 / 277



ARCADIS : Tidewater MGP

SDG: N/A

Client ID: SC-21-S1 Matrix: Sediment Percent Solid: 42.5

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-03 Concentration Units: mg/Kg Date Collected: 07/08/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.654	Jun	0.157	5	08/05/08	07/29/08	6020A	LCP
Arsenic	12.7	J	0.168	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.827	J	0.0672	5	08/01/08	07/29/08	6020A	LMR
Cadmium	11.0	7	0.0672	5	08/01/08	07/29/08	6020A	LMR
Chromium	130	J	0.672	5	08/01/08	07/29/08	6020A	LMR
Copper	249		0.336	5	08/01/08	07/29/08	6020A	LMR
Lead	253		0.168	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.635	J	0.0265	5	07/31/08	07/29/08	7474	LCP
Nickel	38.4		0.336	5	08/01/08	07/29/08	6020A	LMR
Selenium	1.77		0.336	5	08/01/08	07/29/08	6020A	LMR
Silver	4.00		0.157	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.247		0.0672	5	08/01/08	07/29/08	6020A	LMR
Zinc	408		1.68	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable



ARCADIS Tidewater MGP

SDG: N/A

Client ID: SC-10-S1 Matrix: Sediment Percent Solid: 82.4

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-04 Concentration Units: mg/Kg Date Collected: 07/08/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.345	T	0.0960	5	08/05/08	07/29/08	6020A	LCP
Arsenic	7.07	T	0.0985	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.460	5	0.0394	5	08/01/08	07/29/08	6020A	LMR
Cadmium	1.37	<u> </u>	0.0394	5	08/01/08	07/29/08	6020A	LMR
Chromium	18.9	<u> </u>	0.394	5	08/01/08	07/29/08	6020A	LMR
Copper	336		0.197	5	08/01/08	07/29/08	6020A	LMR
Lead	170	J	0.0985	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0529		0.0146	5	07/31/08	07/29/08	7474	LCP
Nickel	15.7		0.197	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.241		0.197	5	08/01/08	07/29/08	6020A	LMR
Silver	0.272		0.0960	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0494		0.0394	5	08/01/08	07/29/08	6020A	LMR
Zinc	158		0.985	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable

N/A



ARCADIS **Tidewater MGP**

SDG: N/A Client ID: SC-12-S1 Sediment Percent Solid: 86.0

Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-05 Concentration Units: mg/Kg Date Collected: 07/08/08 Date Received: 07/11/08

Parameter	Result	Oualifier	Reporting Limit	Dilution	Date Analyzed	Date Prepared	Analytical Method	Analyst
Antimony	0.154	5	0.0944	5	08/05/08	07/29/08	6020A	LCP
Arsenic	3.41	J	0.0881	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.212	T	0.0352	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.555	7	0.0352	5	08/01/08	07/29/08	6020A	LMR
Chromium	10.1	5	0.352	5	08/01/08	07/29/08	6020A	LMR
Copper	25.8		0.176	5	08/01/08	07/29/08	6020A	LMR
Lead	40.1	T	0.0881	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0141	υJ	0.0141	5	07/31/08	07/29/08	7474	LCP
Nickel	11.7		0.176	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.176	U	0.176	5	08/01/08	07/29/08	6020A	LMR
Silver	0.0951		0.0944	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0352	U	0.0352	5	08/01/08	07/29/08	6020A	LMR
Zinc	64.9		0.881	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



: ARCADIS t: Tidewater MGP

> N/A SDG: SC-25-S1 Sediment

Matrix: Sedime Percent Solid: 76.7 Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-06 Concentration Units: mg/Kg Date Collected: 07/08/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	5.13	T	0.107	5	08/05/08	07/29/08	6020A	LCP
Arsenic	5.39	5	0.104	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.283	T	0.0418	5	08/01/08	07/29/08	6020A	LMR
Cadmium	2.57	5	0.0418	5	08/01/08	07/29/08	6020A	LMR
Chromium	28.7	<u> </u>	0.418	5	08/01/08	07/29/08	6020A	LMR
Copper	74.5		0.209	5	08/01/08	07/29/08	6020A	LMR
Lead	124	T	0.104	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0819	J	0.0154	5	07/31/08	07/29/08	7474	LCP
Nickel	20.6	*	0.209	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.277		0.209	5	08/01/08	07/29/08	6020A	LMR
Silver	0.658		0.107	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0549		0.0418	5	08/01/08	07/29/08	6020A	LMR
Zinc	131		1.04	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable



Matrix:

Percent Solid: 80.2

:: ARCADIS et: Tidewater MGP

> N/A SDG: SC3-S1 Sediment

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-07 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.107	Ţ	0.101	5	08/05/08	07/29/08	6020A	LCP
Arsenic	4.03	T	0.0987	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.200	J	0.0395	5	08/01/08	07/29/08	6020A	LMR
Cadmium	1.15	J.	0.0395	5	08/01/08	07/29/08	6020A	LMR
Chromium	16.2	<u>J</u>	0.395	5	08/01/08	07/29/08	6020A	LMR
Copper	24.8		0.197	5	08/01/08	07/29/08	6020A	LMR
Lead	71.0	T	0.0987	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0591	J	0.0151	5	07/31/08	07/29/08	7474	LCP
Nickel	10.1		0.197	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.230		0.197	5	08/01/08	07/29/08	6020A	LMR
Silver	0.224		0.101	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0531		0.0395	5	08/01/08	07/29/08	6020A	LMR
Zinc	170		0.987	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable

N/A



ARCADIS **Tidewater MGP**

Matrix:

SDG: N/A SC41-S1 Sediment Percent Solid: 81.3

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-08 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.102	2	0.102	5	08/05/08	07/29/08	6020A	LCP
Arsenic	2.28	J	0.0961	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.184	J	0.0384	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.752	5	0.0384	5	08/01/08	07/29/08	6020A	LMR
Chromium	9.88	Ţ	0.384	5	08/01/08	07/29/08	6020A	LMR
Copper	80.1		0.192	5	08/01/08	07/29/08	6020A	LMR
Lead	28.3	J	0.0961	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0264	J	0.0146	5	07/31/08	07/29/08	7474	LCP
Nickel	7.99		0.192	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.192	U	0.192	5	08/01/08	07/29/08	6020A	LMR
Silver	0.167		0.102	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0385		0.0384	5	08/01/08	07/29/08	6020A	LMR
Zinc	64.4		0.961	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



Matrix:

ARCADIS **Tidewater MGP**

SDG: N/A Client ID: SC40-S1 Sediment Percent Solid: 84.4

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-09 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

			Reporting		Date Analyzed	Date	Analytical	Analyst
Parameter	Result	Qualifier	Limit	Dilution		Prepared	Method	
Antimony	0.208		0.0903	5	08/05/08	07/29/08	6020A	LCP
Arsenic	2.39	5	0.0937	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.165	5	0.0375	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.673	7"	0.0375	5	08/01/08	07/29/08	6020A	LMR
Chromium	8.96		0.375	5	08/01/08	07/29/08	6020A	LMR
Copper	26.6		0.187	5	08/01/08	07/29/08	6020A	LMR
Lead	40,9	J.	0.0937	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0264	J	0.0140	5	07/31/08	07/29/08	7474	LCP
Nickel	7.41		0.187	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.187	U	0.187	5	08/01/08	07/29/08	6020A	LMR
Silver	0.109		0.0903	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0375	U	0.0375	5	08/01/08	07/29/08	6020A	LMR
Zinc	55.2		0.937	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



Client:ARCADISProject:Tidewater MGPCase:N/ASDG:

Client ID: SC4-S1 Matrix: Sediment Percent Solid: 89.3 Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-10 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

		Reporting		Date	Date	Analytical	
Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
0.180	J.	0.0903	5	08/05/08	07/29/08	6020A	LCP
5.32	J	0.0881	5	08/01/08	07/29/08	6020A	LMR
0.333	T	0.0352	5	08/01/08	07/29/08	6020A	LMR
1.07	J	0.0352	5	08/01/08	07/29/08	6020A	LMR
15.7	5	0.352	5	08/01/08	07/29/08	6020A	LMR
44.7		0.176	5	08/01/08	07/29/08	6020A	LMR
79.2		0.0881	5	08/01/08	07/29/08	6020A	LMR
0.0312	J	0.0131	5	07/31/08	07/29/08	7474	LCP
13.1		0,176	5	08/01/08	07/29/08	6020A	LMR
0.448		0.176	5	08/01/08	07/29/08	6020A	LMR
0.225		0.0903	5	08/05/08	07/29/08	6020A	LCP
0.0530		0.0352	5	08/01/08	07/29/08	6020A	LMR
95.1		0.881	5	08/01/08	07/29/08	6020A	LMR
	0.180 5.32 0.333 1.07 15.7 44.7 79.2 0.0312 13.1 0.448 0.225 0.0530	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	Result Qualifier Limit 0.180 J 0.0903 5.32 J 0.0881 0.333 J 0.0352 1.07 J 0.0352 15.7 J 0.352 44.7 0.176 79.2 J 0.0881 0.0312 J 0.0131 13.1 0.176 0.225 0.0903 0.0530 0.0352	Result Qualifier Limit Dilution 0.180 J 0.0903 5 5.32 J 0.0881 5 0.333 J 0.0352 5 1.07 J 0.0352 5 15.7 J 0.352 5 44.7 0.176 5 79.2 J 0.0881 5 0.0312 J 0.0131 5 13.1 0.176 5 5 0.225 0.0903 5 0.0903 5 0.0530 0.0352 5 5	Result Qualifier Limit Dilution Analyzed 0.180 J 0.0903 5 08/05/08 5.32 J 0.0881 5 08/01/08 0.333 J 0.0352 5 08/01/08 1.07 J 0.0352 5 08/01/08 15.7 J 0.352 5 08/01/08 44.7 0.176 5 08/01/08 79.2 J 0.0881 5 08/01/08 13.1 0.176 5 08/01/08 0.448 0.176 5 08/01/08 0.225 0.0903 5 08/05/08 0.0352 5 08/01/08 0.0352 5	Result Qualifier Limit Dilution Analyzed Prepared 0.180 J 0.0903 5 08/05/08 07/29/08 5.32 J 0.0881 5 08/01/08 07/29/08 0.333 J 0.0352 5 08/01/08 07/29/08 1.07 J 0.0352 5 08/01/08 07/29/08 15.7 J 0.352 5 08/01/08 07/29/08 44.7 0.176 5 08/01/08 07/29/08 79.2 J 0.0881 5 08/01/08 07/29/08 0.0312 J 0.0131 5 07/31/08 07/29/08 13.1 0.176 5 08/01/08 07/29/08 0.448 0.176 5 08/01/08 07/29/08 0.225 0.0903 5 08/05/08 07/29/08 0.0530 0.0352 5 08/01/08 07/29/08	Result Qualifier Limit Dilution Analyzed Prepared Method 0.180 J 0.0903 5 08/05/08 07/29/08 6020A 5.32 J 0.0881 5 08/01/08 07/29/08 6020A 0.333 J 0.0352 5 08/01/08 07/29/08 6020A 1.07 J 0.0352 5 08/01/08 07/29/08 6020A 1.07 J 0.0352 5 08/01/08 07/29/08 6020A 15.7 J 0.352 5 08/01/08 07/29/08 6020A 44.7 0.176 5 08/01/08 07/29/08 6020A 79.2 J 0.0881 5 08/01/08 07/29/08 6020A 0.0312 J 0.0131 5 07/31/08 07/29/08 6020A 0.448 0.176 5 08/01/08 07/29/08 6020A 0.225 0.0903 5 08/05/08

N/A - Not Applicable

N/A



Matrix:

ARCADIS **Tidewater MGP**

SDG: N/A SC22-S1 Sediment Percent Solid: 80.2

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-11 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.16	J	0.0951	5	08/05/08	07/29/08	6020A	LCP
Arsenic	3.07	J	0.0963	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.184	are and a second	0.0385	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.541	J	0.0385	5	08/01/08	07/29/08	6020A	LMR
Chromium	11.6	J	0.385	5	08/01/08	07/29/08	6020A	LMR
Copper	68.3		0.193	5	08/01/08	07/29/08	6020A	LMR
Lead	43.6	J	0.0963	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0492	J	0.0151	5	07/31/08	07/29/08	7474	LCP
Nickel	10.4		0.193	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.193	U	0.193	5	08/01/08	07/29/08	6020A	LMR
Silver	0.129		0.0951	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0385	U	0.0385	5	08/01/08	07/29/08	6020A	LMR
Zinc	69.8		0.963	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



Matrix:

ARCADIS **Tidewater MGP**

SDG: N/A SC5-S1 Sediment Percent Solid: 81.5

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-12 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

		Date	Date	Analytical				
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0:0952	H R	0.0952	5	08/05/08	07/29/08	6020A	LCP
Arsenic	2.58	J	0.0996	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.385	- Jam	0.0398	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.823	J.	0.0398	5	08/01/08	07/29/08	6020A	LMR
Chromium	10.4	J	0.398	5	08/01/08	07/29/08	6020A	LMR
Copper	29.0		0.199	5	08/01/08	07/29/08	6020A	LMR
Lead	56.0	3	0.0996	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0650	J	0.0150	5	07/31/08	07/29/08	7474	LCP
Nickel	10.1		0.199	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.199	U	0.199	5	08/01/08	07/29/08	6020A	LMR
Silver	0.115		0.0952	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0398	U	0.0398	5	08/01/08	07/29/08	6020A	LMR
Zinc	89.9		0.996	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable $\ensuremath{\mathbb{U}}$ - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



:: ARCADIS ct: Tidewater MGP N/A SDG:

Case:N/ASClient ID:SC6-S1Matrix:SedimentPercent Solid:30.9

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-13 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.633	J	0.231	5	08/05/08	07/29/08	6020A	LCP
Arsenic	16.8	J	0.219	5	08/01/08	07/29/08	6020A	LMR
Beryllium	1.11	J	0.0875	5	08/01/08	07/29/08	6020A	LMR
Cadmium	8.45	5	0.0875	5	08/01/08	07/29/08	6020A	LMR
Chromium	140	J	0.875	5	08/01/08	07/29/08	6020A	LMR
Copper	266		0.438	5	08/01/08	07/29/08	6020A	LMR
Lead	242	J	0.219	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.718	T	0.0382	5	07/31/08	07/29/08	7474	LCP
Nickel	35.9		0.438	5	08/01/08	07/29/08	6020A	LMR
Selenium	2.39		0.438	5	08/01/08	07/29/08	6020A	LMR
Silver	3.46		0.231	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.224	.,	0.0875	5	08/01/08	07/29/08	6020A	LMR
Zinc	439		2.19	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable



Matrix:

Percent Solid: 77.6

ARCADIS Tidewater MGP

> N/A SDG: N/A SC20-S1 Sediment

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-14 Concentration Units: mg/Kg Date Collected: 07/09/08 Date Received: 07/11/08

		Date	Date	Analytical				
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.395	J.	0.0976	5	08/05/08	07/29/08	6020A	LCP
Arsenic	3.96	7	0.0994	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.212	J	0.0398	5	08/01/08	07/29/08	6020A	LMR
Cadmium	1.58	J	0.0398	5	08/01/08	07/29/08	6020A	LMR
Chromium	18.3	J	0.398	5	08/01/08	07/29/08	6020A	LMR
Copper	54.4		0.199	5	08/01/08	07/29/08	6020A	LMR
Lead	128	T	0.0994	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.182	J	0.0156	5	07/31/08	07/29/08	7474	LCP
Nickel	9.71		0.199	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.271		0.199	5	08/01/08	07/29/08	6020A	LMR
Silver	0.436		0.0976	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0398	U	0.0398	5	08/01/08	07/29/08	6020A	LMR
Zinc	93.6		0.994	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable

U - The analyte was analyzed for but not detected at the sample specific level reported.

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Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: N/A SC26-S1 Sediment Percent Solid: 83.5

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-15 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.372	J	0.0992	5	08/05/08	07/29/08	6020A	LCP
Arsenic	1.79	J	0.0936	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.143	5	0.0374	5	08/01/08	07/29/08	6020A	LMR
Cadmium	0.551	J	0.0374	5	08/01/08	07/29/08	6020A	LMR
Chromium	7.68	\mathcal{T}	0.374	5	08/01/08	07/29/08	6020A	LMR
Copper	22.5		0.187	5	08/01/08	07/29/08	6020A	LMR
Lead	28.4	2	0.0936	5	08/01/08	07/29/08	6020A	LMR
Mercury	0.0300	J	0.0145	5	07/31/08	07/29/08	7474	LCP
Nickel	4.91		0.187	5	08/01/08	07/29/08	6020A	LMR
Selenium	0.264		0.187	5	08/01/08	07/29/08	6020A	LMR
Silver	0.124		0.0992	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.0374	U	0.0374	5	08/01/08	07/29/08	6020A	LMR
Zinc	58.2		0.936	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: N/A Client ID: SC19-S1 Sediment Percent Solid: 35.7

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-16 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

	Reporting				Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.981	J	0.200	5	08/05/08	07/29/08	6020A	LCP
Arsenic	17.6	J	0.198	5	08/01/08	07/29/08	6020A	LMR
Beryllium	0.706	T	0.0792	5	08/01/08	07/29/08	6020A	LMR
Cadmium	19.7	J	0.0792	5	08/01/08	07/29/08	6020A	LMR
Chromium	280	J	0.792	5	08/01/08	07/29/08	6020A	LMR
Copper	773		0.396	5	08/01/08	07/29/08	6020A	LMR
Lead	383	J	0.198	5	08/01/08	07/29/08	6020A	LMR
Mercury	1.39	J	0.0331	5	07/31/08	07/29/08	7474	LCP
Nickel	49.6		0.396	5	08/01/08	07/29/08	6020A	LMR
Selenium	1.77		0.396	5	08/01/08	07/29/08	6020A	LMR
Silver	5.25		0.200	5	08/05/08	07/29/08	6020A	LCP
Thallium	0.200		0.0792	5	08/01/08	07/29/08	6020A	LMR
Zinc	792		1.98	5	08/01/08	07/29/08	6020A	LMR

N/A - Not Applicable

N/A



Client: ARCADIS Project: Tidewater MGP Case: N/A SDG: Client ID: SC43-S1

Matrix: Sediment Percent Solid: 62.0 Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-17 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.334	J	0.117	5	08/05/08	07/31/08	6020A	LCP
Arsenic	4,93	T	0.110	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.408	J	0.0438	5	08/01/08	07/31/08	6020A	LMR
Cadmium	2.95	J	0.0438	5	08/01/08	07/31/08	6020A	LMR
Chromium	46.5	T	0.438	5	08/01/08	07/31/08	6020A	LMR
Copper	117		0.219	5	08/01/08	07/31/08	6020A	LMR
Lead	112	J	0.110	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.292	J	0.0178	5	07/31/08	07/29/08	7474	LCP
Nickel	16.3		0.219	5	08/01/08	07/31/08	6020A	LMR
Selenium	0.707		0.219	5	08/01/08	07/31/08	6020A	LMR
Silver	1.71		0.117	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.115		0.0438	5	08/01/08	07/31/08	6020A	LMR
Zinc	176		1.10	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable

N/A



ARCADIS **Tidewater MGP**

SDG: N/A Client ID: SC42-S1 Sediment Percent Solid: 77.3

Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-18 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.139		0.0934	5	08/05/08	07/31/08	6020A	LCP
Arsenic	3.87	Ĵ	0.0934	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.237	J	0.0374	5	08/01/08	07/31/08	6020A	LMR
Cadmium	2.76	5	0.0374	5	08/01/08	07/31/08	6020A	LMR
Chromium	15.2		0.374	5	08/01/08	07/31/08	6020A	LMR
Copper	44.4		0.187	5	08/01/08	07/31/08	6020A	LMR
Lead	83.1	5	0.0934	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.0542	J	0.0152	5	07/31/08	07/29/08	7474	LCP
Nickel	10.8		0.187	5	08/01/08	07/31/08	6020A	LMR
Selenium	0.346		0.187	5	08/01/08	07/31/08	6020A	LMR
Silver	0.783		0.0934	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.0441		0.0374	5	08/01/08	07/31/08	6020A	LMR
Zinc	992		1.87	10	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS **Tidewater MGP**

N/A

Sediment

Matrix:

Percent Solid: 36.6

SDG: N/A Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-19 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.09	5	0.212	5	08/05/08	07/31/08	6020A	LCP
Arsenic	15.7	2	0.192	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.982	5	0.0767	5	08/01/08	07/31/08	6020A	LMR
Cadmium	10.4	5	0.0767	5	08/01/08	07/31/08	6020A	LMR
Chromium	154	5	0.767	5	08/01/08	07/31/08	6020A	LMR
Copper	322		0.384	5	08/01/08	07/31/08	6020A	LMR
Lead	308	J	0.192	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.903	5	0.0292	5	07/31/08	07/29/08	7474	LCP
Nickel	41.0	0.5	0.384	5	08/01/08	07/31/08	6020A	LMR
Selenium	1.99		0.384	5	08/01/08	07/31/08	6020A	LMR
Silver	5.80		0.212	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.180		0.0767	5	08/01/08	07/31/08	6020A	LMR
Zinc	455		1.92	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



: ARCADIS t: Tidewater MGP

N/A

Matrix:

Percent Solid: 36.2

DUP-1

Sediment

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-20 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.25	J	0.207	5	08/05/08	07/31/08	6020A	LCP
Arsenic	15.9	T	0.208	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.04	T	0.0833	5	08/01/08	07/31/08	6020A	LMR
Cadmium	11.1	J.	0.0833	5	08/01/08	07/31/08	6020A	LMR
Chromium	158	5	0.833	5	08/01/08	07/31/08	6020A	LMR
Copper	314		0.416	5	08/01/08	07/31/08	6020A	LMR
Lead	302		0.208	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.942		0.0295	5	07/31/08	07/29/08	7474	LCP
Nickel	41.0		0.416	5	08/01/08	07/31/08	6020A	LMR
Selenium	1.88		0.416	5	08/01/08	07/31/08	6020A	LMR
Silver	6.10		0.207	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.183		0.0833	5	08/01/08	07/31/08	6020A	LMR
Zinc	446		2.08	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable

N/A



Matrix:

ARCADIS **Tidewater MGP**

SDG: N/A Client ID: SC29-S1 Sediment Percent Solid: 42.1

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-21 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.848	J	0.174	5	08/05/08	07/31/08	6020A	LCP
Arsenic	13.2	J	0.172	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.04	alere a	0.0686	5	08/01/08	07/31/08	6020A	LMR
Cadmium	10.8	J	0.0686	5	08/01/08	07/31/08	6020A	LMR
Chromium	132	3	0.686	5	08/01/08	07/31/08	6020A	LMR
Copper	245		0.343	5	08/01/08	07/31/08	6020A	LMR
Lead	257	J	0.172	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.678	J	0.0277	5	07/31/08	07/29/08	7474	LCP
Nickel	38.6		0.343	5	08/01/08	07/31/08	6020A	LMR
Selenium	2.13		0.343	5	08/01/08	07/31/08	6020A	LMR
Silver	4.05		0.174	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.192		0.0686	5	08/01/08	07/31/08	6020A	LMR
Zinc	384		1.72	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS Project: **Tidewater MGP**

> SDG: N/A SC17-S1

Matrix: Sediment Percent Solid: 52.9

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-22 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.19	and the second s	0.141	5	08/05/08	07/31/08	6020A	LCP
Arsenic	6.94	J	0.130	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.512	T	0.0522	5	08/01/08	07/31/08	6020A	LMR
Cadmium	4.68	J	0.0522	5	08/01/08	07/31/08	6020A	LMR
Chromium	67.9	J	0.522	5	08/01/08	07/31/08	6020A	LMR
Copper	158		0.261	5	08/01/08	07/31/08	6020A	LMR
Lead	158	T	0.130	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.426	5	0.0223	5	07/31/08	07/29/08	7474	LCP
Nickel	20.4		0.261	5	08/01/08	07/31/08	6020A	LMR
Selenium	0.996		0.261	5	08/01/08	07/31/08	6020A	LMR
Silver	2.76		0.141	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.0986		0.0522	5	08/01/08	07/31/08	6020A	LMR
Zinc	206		1.30	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable

N/A



ARCADIS **Tidewater MGP**

> N/A SDG: SC28-S1

Sediment Percent Solid: 34.6

Matrix:

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-23 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.928	5	0.198	5	08/05/08	07/31/08	6020A	LCP
Arsenic	14.7	T	0.222	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.09	T	0.0886	5	08/01/08	07/31/08	6020A	LMR
Cadmium	11.0	J	0.0886	5	08/01/08	07/31/08	6020A	LMR
Chromium	158	T	0.886	5	08/01/08	07/31/08	6020A	LMR
Copper	320		0.443	5	08/01/08	07/31/08	6020A	LMR
Lead	306	J	0.222	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.875	T	0.0337	5	07/31/08	07/29/08	7474	LCP
Nickel	39.8		0.443	5	08/01/08	07/31/08	6020A	LMR
Selenium	2.14		0.443	5	08/01/08	07/31/08	6020A	LMR
Silver	5.69		0.198	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.214		0.0886	5	08/01/08	07/31/08	6020A	LMR
Zinc	446		2.22	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS **Tidewater MGP**

Matrix:

SDG: N/A N/A Client ID: SC30-S1 Sediment Percent Solid: 53.8

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-24 Concentration Units: mg/Kg Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.40	T	0.136	5	08/05/08	07/31/08	6020A	LCP
Arsenic	13.5	7	0.136	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.35	5	0.0543	5	08/01/08	07/31/08	6020A	LMR
Cadmium	6.46	J	0.0543	5	08/01/08	07/31/08	6020A	LMR
Chromium	143	J	0.543	5	08/01/08	07/31/08	6020A	LMR
Copper	271		0.272	5	08/01/08	07/31/08	6020A	LMR
Lead	246	5	0.136	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.882	J	0.0213	5	07/31/08	07/29/08	7474	LCP
Nickel	61.7		0.272	5	08/01/08	07/31/08	6020A	LMR
Selenium	2.62		0.272	5	08/01/08	07/31/08	6020A	LMR
Silver	3.35		0.136	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.179		0.0543	5	08/01/08	07/31/08	6020A	LMR
Zinc	496		1.36	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS **Tidewater MGP**

> SDG: N/A SC47-S1

Matrix: Sediment Percent Solid: 83.2

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-25 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.264	T	0.0964	5	08/05/08	07/31/08	6020A	LCP
Arsenic	3.84	angere .	0.0964	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.254	J	0.0385	5	08/01/08	07/31/08	6020A	LMR
Cadmium	0.847	J	0.0385	5	08/01/08	07/31/08	6020A	LMR
Chromium	12.4	5	0.385	5	08/01/08	07/31/08	6020A	LMR
Copper	30.3		0.193	5	08/01/08	07/31/08	6020A	LMR
Lead	40.6	2	0.0964	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.0210	5	0.0140	5	07/31/08	07/29/08	7474	LCP
Nickel	12.3		0.193	5	08/01/08	07/31/08	6020A	LMR
Selenium	0.232		0.193	5	08/01/08	07/31/08	6020A	LMR
Silver	0.175		0.0964	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.0731		0.0385	5	08/01/08	07/31/08	6020A	LMR
Zinc	79.9		0.964	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS : Tidewater MGP

SDG: N/A

Client ID: SC32-S1 Matrix: Sediment Percent Solid: 60.1

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-26 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.66	T	0.120	5	08/05/08	07/31/08	6020A	LCP
Arsenic	11.7	5	0.120	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.53	5	0.0478	5	08/01/08	07/31/08	6020A	LMR
Cadmium	6.93	J	0.0478	5	08/01/08	07/31/08	6020A	LMR
Chromium	138	J	0.478	5	08/01/08	07/31/08	6020A	LMR
Copper	295		0.239	5	08/01/08	07/31/08	6020A	LMR
Lead	239	Ţ	0.120	5	08/01/08	07/31/08	6020A	LMR
Mercury	1.23	- Landon - L	0.0184	5	07/31/08	07/29/08	7474	LCP
Nickel	33.7		0.239	5	08/01/08	07/31/08	6020A	LMR
Selenium	1.70		0.239	5	08/01/08	07/31/08	6020A	LMR
Silver	2.74		0.120	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.148		0.0478	5	08/01/08	07/31/08	6020A	LMR
Zinc	391		1.20	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS **Tidewater MGP**

> SDG: N/A SC33-S1

Matrix: Sediment Percent Solid: 31.1

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-27 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

Parameter	Result	Qualifier	Reporting Limit	Dilution	Date Analyzed	Date Prepared	Analytical Method	Analyst
Antimony	1.38	enterna and	0.224	5	08/05/08	07/31/08	6020A	LCP
Arsenic	17.1	T	0.221	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.18	J	0.0885	5	08/01/08	07/31/08	6020A	LMR
Cadmium	10.0	T	0.0885	5	08/01/08	07/31/08	6020A	LMR
Chromium	172	J	0.884	5	08/01/08	07/31/08	6020A	LMR
Copper	357		0.442	5	08/01/08	07/31/08	6020A	LMR
Lead	311	and the second	0.221	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.922	T	0.0373	5	07/31/08	07/29/08	7474	LCP
Nickel	39.1		0.442	5	08/01/08	07/31/08	6020A	LMR
Selenium	2.45		0.442	5	08/01/08	07/31/08	6020A	LMR
Silver	6.72	b.#/3	0.224	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.240		0.0885	5	08/01/08	07/31/08	6020A	LMR
Zinc	465		2.21	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



Client: ARCADIS Project: Tidewater MGP

> N/A SDG: N/A SW2-S1

Matrix: Sediment Percent Solid: 39.6 Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-28 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	3.22	J	0.188	5	08/05/08	07/31/08	6020A	LCP
Arsenic	10.4	7	0.180	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.144	7	0.0721	5	08/01/08	07/31/08	6020A	LMR
Cadmium	1.88	5	0.0721	5	08/01/08	07/31/08	6020A	LMR
Chromium	108	J	0.721	5	08/01/08	07/31/08	6020A	LMR
Copper	283		0.361	5	08/01/08	07/31/08	6020A	LMR
Lead	318	T	0.180	5	08/01/08	07/31/08	6020A	LMR
Mercury	1,48	J	0.0301	5	07/31/08	07/29/08	7474	LCP
Nickel	15.0		0.361	5	08/01/08	07/31/08	6020A	LMR
Selenium	2,52		0.361	5	08/01/08	07/31/08	6020A	LMR
Silver	0.697		0.188	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.375		0.0721	5	08/01/08	07/31/08	6020A	LMR
Zinc	206		1.80	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS **Tidewater MGP**

SDG: N/A SC39-S1

Matrix: Sediment Percent Solid: 23.7

N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-29 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

Parameter	Result	Qualifier	Reporting Limit	Dilution	Date Analyzed	Date Prepared	Analytical Method	Analyst
Antimony	1.52	7	0.285	5	08/05/08	07/31/08	6020A	LCP
Arsenic	16.8	T	0.311	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.34	Ţ	0.124	5	08/01/08	07/31/08	6020A	LMR
Cadmium	16.1	J	0.124	5	08/01/08	07/31/08	6020A	LMR
Chromium	232	5	1.24	5	08/01/08	07/31/08	6020A	LMR
Copper	516		0.621	5	08/01/08	07/31/08	6020A	LMR
Lead	413	J	0.311	5	08/01/08	07/31/08	6020A	LMR
Mercury	1.35	J	0.0444	5	07/31/08	07/29/08	7474	LCP
Nickel	55.0		0.621	5	08/01/08	07/31/08	6020A	LMR
Selenium	2.90		0.621	5	08/01/08	07/31/08	6020A	LMR
Silver	11.0		0.285	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.283		0.124	5	08/01/08	07/31/08	6020A	LMR
Zinc	583		3.11	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



Client: ARCADIS Project: Tidewater MGP

N/A SDG: N/A SC38-S1

Matrix: Sediment Percent Solid: 34.5 Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-30 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

		Reporting					Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.06	Ţ	0.225	5	08/05/08	07/31/08	6020A	LCP
Arsenic	17.5	5	0.211	5	08/01/08	07/31/08	6020A	LMR
Beryllium	1.58	5	0.0843	5	08/01/08	07/31/08	6020A	LMR
Cadmium	18.2	J	0.0843	5	08/01/08	07/31/08	6020A	LMR
Chromium	243	J	0.843	5	08/01/08	07/31/08	6020A	LMR
Copper	446		0.421	5	08/01/08	07/31/08	6020A	LMR
Lead	450	5	0.211	5	08/01/08	07/31/08	6020A	LMR
Mercury	1.58	J	0.0336	5	07/31/08	07/29/08	7474	LCP
Nickel	71.7		0.421	5	08/01/08	07/31/08	6020A	LMR
Selenium	2.17		0.421	5	08/01/08	07/31/08	6020A	LMR
Silver	9.50		0.225	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.243		0.0843	5	08/01/08	07/31/08	6020A	LMR
Zinc	593		2.11	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable

N/A



: ARCADIS t: Tidewater MGP

Case: N/A SDG: Client ID: SC34-S1 Matrix: Sediment

Matrix: Sedimer Percent Solid: 53.9 Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-31 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

		Reporting				Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.482	and the second s	0.136	5	08/05/08	07/31/08	6020A	LCP
Arsenic	7.07	T	0.132	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.526	J	0.0527	5	08/01/08	07/31/08	6020A	LMR
Cadmium	4.50	J	0.0527	5	08/01/08	07/31/08	6020A	LMR
Chromium	69.1	J	0.527	5	08/01/08	07/31/08	6020A	LMR
Copper	184		0.263	5	08/01/08	07/31/08	6020A	LMR
Lead	153	T	0.132	5	08/01/08	07/31/08	6020A	LMR
Mercury	0.617	3	0.0227	5	07/31/08	07/29/08	7474	LCP
Nickel	21.8		0.263	5	08/01/08	07/31/08	6020A	LMR
Selenium	0.957		0.263	5	08/01/08	07/31/08	6020A	LMR
Silver	2.85		0.136	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.112		0.0527	5	08/01/08	07/31/08	6020A	LMR
Zinc	200		1.32	5	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



ARCADIS **Tidewater MGP**

N/A

Matrix:

Percent Solid: 55.1

SC35-S1

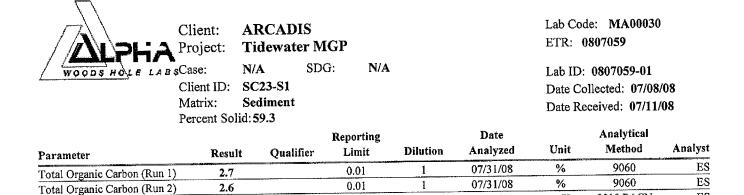
Sediment

SDG: N/A

Lab Code: MA00030 ETR: 0807059 Lab ID: 0807059-32 Concentration Units: mg/Kg Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date	Date Prepared	Analytical Method	Analyst
Parameter	Result	Qualifier	Limit	Dilution	Analyzed			
Antimony	1.19	T	0.132	5	08/05/08	07/31/08	6020A	LCP
Arsenic	20.1	and an	0.141	5	08/01/08	07/31/08	6020A	LMR
Beryllium	0.610	J	0.0563	5	08/01/08	07/31/08	6020A	LMR
Cadmium	10.4	J.	0.0563	5	08/01/08	07/31/08	6020A	LMR
Chromium	210	5	0.563	5	08/01/08	07/31/08	6020A	LMR
Copper	348		0.282	5	08/01/08	07/31/08	6020A	LMR
Lead	553		0.141	· 5	08/01/08	07/31/08	6020A	LMR
Mercury	0.914	T	0.0222	5	07/31/08	07/29/08	7474	LCP
Nickel	50.3		0.282	5	08/01/08	07/31/08	6020A	LMR
Selenium	1.78		0.282	5	08/01/08	07/31/08	6020A	LMR
Silver	3.37		0.132	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.174		0.0563	5	08/01/08	07/31/08	6020A	LMR
Zinc	1440		2.82	10	08/01/08	07/31/08	6020A	LMR

N/A - Not Applicable



0.14

U

0.14

Physiologically Available Cyanide

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:49

Page 168 / 277

9010 PACN

mg/Kg

07/21/08

1

ES

	Client: Project:	ARCADIS Tidewater						de: MA00030 0807059	I
WOODS HOLE LABS	Case: Client ID: Matrix: Percent So	SC-24-S1 Sediment	DG: 1	N/A			Date C	: 0807059-02 ollected: 07/08 eceived: 07/11/	
i.			Reporti	ıg		Date		Analytical	
Parameter	Resu	lt Qualifier	r Limit	D	ilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.77	す	0.01		1	07/31/08	%	9060	ES
Total Organic Carbon (Run 2)		J	0.01		1	07/31/08	%	9060	ES

0.093

U

1

07/21/08

mg/Kg

9010 PACN

ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

Physiologically Available Cyanide

0.093



ARCADIS : Tidewater MGP

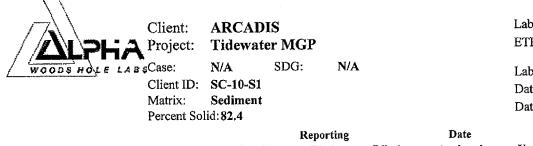
N/A SDG: N/A

Client ID: SC-21-S1 Matrix: Sediment Percent Solid: 42.5 Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-03 Date Collected: 07/08/08 Date Received: 07/11/08

		Reporting			Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	2.9		0.01	1	07/31/08	%	9060	ES
Total Organic Carbon (Run 2)	2.9		0.01	1	07/31/08	%	9060	ES
Physiologically Available Cyanide	0.15	U	0.15	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-04 Date Collected: 07/08/08 Date Received: 07/11/08

			Reporting		Date		Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	0.56		0.01	1	07/31/08	%	9060	ES	
Total Organic Carbon (Run 2)	0.60		0.01	1	07/31/08	%	9060	ES	
Physiologically Available Cyanide	0.075	U	0.075	1	07/21/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SDG: N/A

SC-12-S1

Matrix: Sediment Percent Solid: 86.0

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-05 Date Collected: 07/08/08 Date Received: 07/11/08

		Reporting			Date	Analytical			
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	0.10		0.01	1	07/31/08	%	9060	ES	
Total Organic Carbon (Run 2)	0.14		0.01	1	07/31/08	%	9060	ES	
Physiologically Available Cyanide	0.076	U	0.076	1	07/21/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SDG: N/A

SC-25-S1

Sediment

Matrix: Sediment Percent Solid: 76.7

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-06 Date Collected: 07/08/08 Date Received: 07/11/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.3		0.01	1	07/31/08	%	9060	ES
Total Organic Carbon (Run 2)	1.6		0.01	1	07/31/08	%	9060	ES
Physiologically Available Cyanide	0.089	U	0.089	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS **Tidewater MGP**

> SDG: N/A

Matrix: Sediment Percent Solid: 80.2

N/A

Lab Code: MA00030 ETR: 0807059

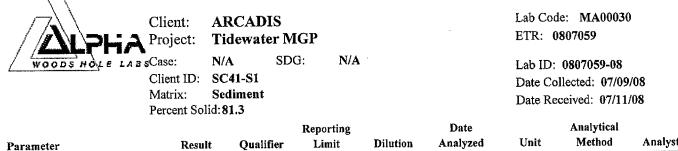
Lab ID: 0807059-07 Date Collected: 07/09/08 Date Received: 07/11/08

		Reporting			Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.0		0.01	1	07/31/08	%	9060	ES
Total Organic Carbon (Run 2)	1.0		0.01	1	07/31/08	%	9060	ES
Physiologically Available Cyanide	0.091	U	0.091	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:49

Page 176 / 277

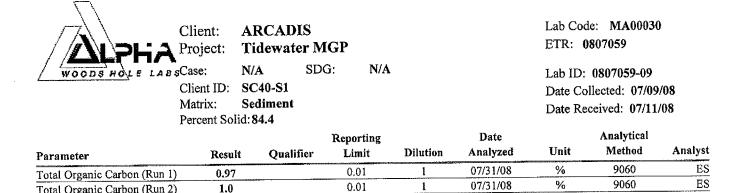


Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.1		0.01	1	07/31/08	%	9060	ES
Total Organic Carbon (Run 2)	1.3		0.01	1	07/31/08	%	9060	ES
Physiologically Available Cyanide	0.076	U	0.076	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:50

Page 177 / 277



0.01

0.089

U

1

1

07/21/08

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

1.0

0.089

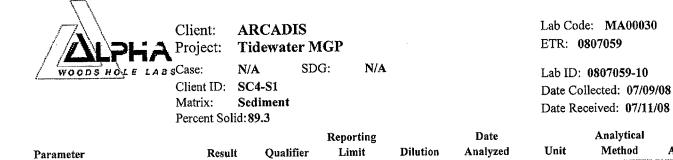
Total Organic Carbon (Run 2)

Physiologically Available Cyanide

9010 PACN

mg/Kg

ES



U

1.4

1.9

0.092

Total Organic Carbon (Run 1)

Total Organic Carbon (Run 2)

Physiologically Available Cyanide

0.01

0.01

0.092

1 07/21/08 mg/Kg

07/31/08

07/31/08

1

1

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:50

Analyst

ES

ES

ES

9060

9060

9010 PACN

%

%



ARCADIS **Tidewater MGP**

> SDG: N/A

Matrix: Sediment Percent Solid: 80.2

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-11 Date Collected: 07/09/08 Date Received: 07/11/08

		Reporting			Date	Analytical			
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	0.33		0.01	1	08/01/08	%	9060	ES	
Total Organic Carbon (Run 2)	0.28		0.01	1	08/01/08	%	9060	ES	
Physiologically Available Cyanide	0.089	U	0.089	1	07/21/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

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ARCADIS **Tidewater MGP**

> SDG: N/A N/A

Sediment Percent Solid: 81.5

Matrix:

ETR: 0807059 Lab ID: 0807059-12

> Date Collected: 07/09/08 Date Received: 07/11/08

Lab Code: MA00030

		Reporting			Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.55		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	0.50		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.099	U	0.099	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS **Tidewater MGP**

N/A

Matrix:

Percent Solid: 30.9

Sediment

SDG: N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-13 Date Collected: 07/09/08 Date Received: 07/11/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.8		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	4.8		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.25	υ	0.25	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

	Client: Project:	ARCADIS Tidewater N	AGP				ode: MA00030 0807059)
]	Case: Client ID: Matrix: Percent Sol	SC20-S1 Sediment	DG: N/2	A		Date C	: 0807059-14 ollected: 07/09 eceived: 07/11/	
			Reporting		Date		Analytical	
Parameter	Resu	lt Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.1		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	0.85		0.01	1	08/01/08	%	9060	ES
					* - ** * ** *			

0.080

U

1

07/21/08

mg/Kg

9010 PACN

ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

Physiologically Available Cyanide

0.080

N/A



ARCADIS :: Tidewater MGP

N/A SDG:

Client ID: SC26-S1 Matrix: Sediment

Percent Solid: 83.5

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-15 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.37		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	0.39		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.092	U	0.092	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

		ARCADIS Tidewater					de: MA0003 0807059	0
WOODS HOLE LAB	sCase: Client ID:		DG: N/A	L			: 0807059-16 ollected: 07/10)/08
	Matrix: Percent So	Sediment lid:35.7				Date R	eceived: 07/11	/08
			Reporting		Date		Analytical	
Parameter	Resu	lt Qualifier	· Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)) 5.2		0.01	1	08/01/08	%	9060	ES

0.01

0.17

υ

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

Total Organic Carbon (Run 2)

Physiologically Available Cyanide

4.0

0.17

08/07/08 16:50

ES

ES

9060

9010 PACN

%

mg/Kg

08/01/08

07/21/08

1

1

N/A



ARCADIS **Tidewater MGP**

N/A

Matrix:

Percent Solid: 62.0

Sediment

SDG:

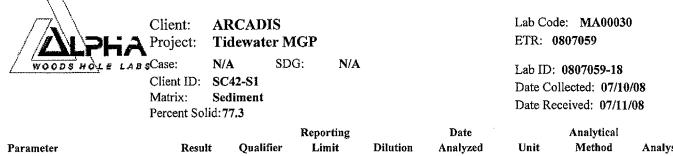
Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-17 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.7		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	1.8		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.11	U	0.11	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

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			Reporting		Date		renaryticat	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.92		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	0.81		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.10	U	0.10	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



ARCADIS : Tidewater MGP

N/A SDG:

Client ID: SC9-S1

Matrix: Sediment Percent Solid: 36.6 Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-19 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date		Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	4.1		0.01	1	08/01/08	%	9060	ES	
Total Organic Carbon (Run 2)	4.2		0.01	1	08/01/08	%	9060	ES	
Physiologically Available Cyanide	0.18	U	0.18	1	07/21/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



ARCADIS t: Tidewater MGP

N/A SDG:

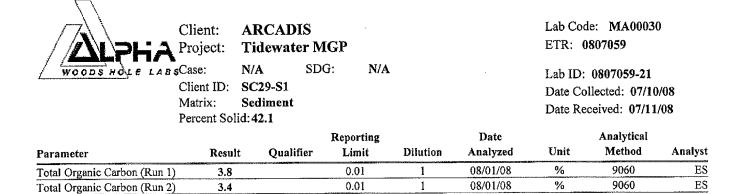
Client ID: **DUP-1** Matrix: **Sediment**

Matrix: Sediment Percent Solid: 36.2 Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-20 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	3.9	a an	0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	3.6	······	0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.22	U	0.22	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



0.17

1

U

0.17

mg/Kg

07/21/08

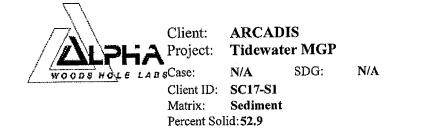
9010 PACN

ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

Physiologically Available Cyanide

08/07/08 16:50



Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-22 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	2.3		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	2.4		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.13	U	0.13	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SDG: N/A

SC28-S1

Matrix: Sediment Percent Solid: 34.6

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-23 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.8		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	3.7		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.23	U	0.23	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS : Tidewater MGP

SC30-S1

Sediment

Matrix:

Percent Solid: 53.8

N/A SDG: N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-24 Date Collected: 07/10/08 Date Received: 07/11/08

			Reporting		Date	Analytical			
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	9.2	J.	0.01	1	08/01/08	%	9060	ES	
Total Organic Carbon (Run 2)	10	J	0.01	1	08/01/08	%	9060	ES	
Physiologically Available Cyanide	0.15	U	0.15	1	07/21/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:50



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: ARCADIS t: Tidewater MGP

SDG: N/A

Client ID: SC47-S1

Matrix: Sediment Percent Solid:83.2

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-25 Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.2		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	1.4		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.091	U	0.091	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:50



ARCADIS Tidewater MGP

N/A SDG: N/A

Client ID: SC32-S1 Matrix: Sediment Percent Solid:60.1 Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-26 Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	8.3		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	10		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.13	U	0.13	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



Matrix: Sediment Percent Solid: 31.1 Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-27 Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	5.1		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	4.8		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.25	U	0.25	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SDG: N/A

Client ID: SW2-S1 Matrix: Sediment

N/A

Percent Solid: 39.6

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-28 Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	34		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	28		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.19	U	0.19	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 16:50



ARCADIS : Tidewater MGP

SDG: N/A

SC39-S1

Matrix: Sediment Percent Solid:23.7

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-29 Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	6.5		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	4.7		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.32	υ	0.32	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



: ARCADIS t: Tidewater MGP

SDG: N/A

SC38-S1

Matrix: Sediment Percent Solid: 34.5

N/A

Lab Code: MA00030 ETR: 0807059

Lab ID: 0807059-30 Date Collected: 07/11/08 Date Received: 07/11/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.1		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	4.7		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.21	U	0.21	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

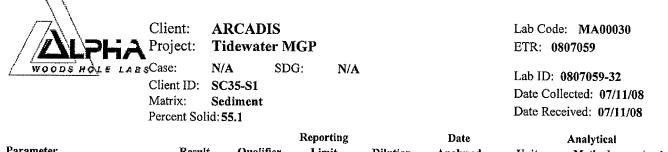
Client: ARCADIS Lab Code: MA00030 Project: **Tidewater MGP** ETR: 0807059 WOODS HOLE LAB Case: N/A SDG: N/A Lab ID: 0807059-31 Client ID: SC34-S1 Date Collected: 07/11/08 Matrix: Sediment Date Received: 07/11/08 Percent Solid: 53.9 Reporting Date Analytical

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Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.7		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	1.6		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.13	U	0.13	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

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08/07/08 16:50



Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.8		0.01	1	08/01/08	%	9060	ES
Total Organic Carbon (Run 2)	5,4		0.01	1	08/01/08	%	9060	ES
Physiologically Available Cyanide	0.12	U	0.12	1	07/21/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

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08/07/08 16:50

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-01 SC23-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe	ers: 1-Glass	

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	55	ક	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	4880	mg/kg	72.7	1 9071B	0718 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 4 of 39

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-02	Date Collected: 08-JUL-2008 11:40			
	SC-24-S1	Date Received : 16-JUL-2008			
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008			
Condition of Sample:	Satisfactory	Field Prep: None			
Number & Type of Containers: 2-Glass					

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	82	ę	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	2260	mg/kg	48.8	1 9071B	0710 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 5 of 39

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	SC-21-S1 .	Date Received :	
Sample Matrix: Condition of Sample:	SOIL Satisfactory	Date Reported :	28-JUL-2008
Number & Type of Containe:	-	agener trup.	None

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	41	₽	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	3690	mg/kg	97.6	1 9071B	0718 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 6 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAC00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-04 SC-10-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe:	rs: l-Glass	

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	72	ę	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	1980	mg/kg	55.6	1 9071B	0718 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 7 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-05 SC-12-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None
Number & Type of Container	cs: 1-Glass		

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	90	8	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	406	mg/kg	44.4	1 9071B	0718 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 8 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-06 SC-25-S1	Date Collected Date Received	: 08-JUL-2008 11:40 : 16-JUL-2008		
Sample Matrix:	SOIL	Date Reported	28-JUL-2008		
Condition of Sample:	Satisfactory	Field Prep:	None		
Number & Type of Containers: 1-Glass					

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	72.	8	0.10	30 2540g	0717 14:23 SD
TPH, HEM-SGT	1620	mg/kg	50.0	1 9071B	0718 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 9 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-07 SC3-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008		
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008		
Condition of Sample:	Satisfactory	Field Prep:	None		
Number & Type of Containers: 1-Glass					

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	83	8	0.10	30 2540G	0717 14;23 SD
TPH, HEM-SGT	1790	mg/kg	48.2	1 9071B	0718 11:15 0721 15:00 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 10 of 39

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-08 SC41-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None
Number & Type of Containe	rs: 1-Glass		

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID	
					PREP ANAL	
Solids, Total	78	¥	0.10	30 2540G	0717 14:23 SD	
TPH, HEM-SGT	729	mg/kg	46.2	1 9071B	0718 11:15 0721 15:00 AT	

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 11 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-09 SC40-S1 SOIL	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008 Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	date id Prep Anal
Solids, Total	82	8	0,10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	538	mg/kg	48.8	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 12 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-10 SC4-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	date id Prep Anal
Solids, Total	89	8	0.10	30 2540g	0717 14:23 SD
TPH, HEM-SGT	980	mg/kg	44.9	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 13 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:L0810512-11Date Collected:08-JUL-200811:40SC22-S1Date Received :16-JUL-2008Sample Matrix:SOILDate Reported :28-JUL-2008Condition of Sample:SatisfactoryField Prep:None

Number & Type of Containers: 1-Glass

Parameter	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	78	Se a constantino de la constan	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	518	mg/kg	46.2	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 14 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-12 SC5~S1 SOIL	Date Collected: Date Received : Date Reported :	
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	85	8	0.10	30 2540g	0717 14:23 SD
TPH, HEM-SGT	611	mg/kg	47.0	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 15 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:		Date Collected: 08-JUL-2008 11:40
	SC6-S1	Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

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PARAMETER	RESULT	UNITS	RDI.	REF METHOD	DATE ID PREP ANAL
Solids, Total	31	8	0.10	30 2540g	0717 14:23 SD
TPH, HEM-SGT	7280	mg/kg	129	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 16 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-14 SC20-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	78	z	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	924	mg/kg	51.3	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 17 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAC000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-15 SC26-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	81	ę	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	594	mg/kg	49.4	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 18 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-16 SC19-S1 SOIL	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008 Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	35	8	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	13600	mg/kg	114	1 9071B	0721 12:45 0722 10:15 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 19 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-17 SC43-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Glass

Parameter	RESULT	UNITS	RDL	REF METHOD	date id Prep anal
Solids, Total	62	8	0.10	30 2540g	0717 14:23 SD
TPH, HEM-SGT	3150	mg/kg	64.5	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07260813:27 Page 20 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-18 SC42-S1 SOIL	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008 Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	78	<u>0</u>	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	1480	mg/kg	51.3	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 21 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-19 SC9-S1 SOIL	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008 Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

Parameter	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	35	ę	0.10	30 2540G	0717 14:23 5D
TPH, HEM-SGT	5820	mg/kg	103	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 22 of 39

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:		Date Collected: 08-JUL-2008 11:40
	DUP-1	Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	35	9	0.10	30 2540G	0717 14:23 SD
TPH, HEM-SGT	5410	mg/kg	114	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 23 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-21 SC29-S1 SOIL	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008 Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	43	8	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	3870	mg/kg	93.0	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I 07280013:27 Page 24 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-22 SC17-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: N	Jone

Number & Type of Containers: 1-Glass

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	54	ę	0,10	30 2540g	0717 15:07 SD
TPH, HEM-SGT	4330	mg/kg	74.1	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 25 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810512-23 SC28-S1 SOIL	Date Collected: 0 Date Received : 1 Date Reported : 2	6-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: No	one

Number & Type of Containers: 1-Glass

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	36	8	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	7770	mg/kg	100	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 26 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-24 SC30-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 2-Glass

Parameter	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	54	8	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	6090	mg/kg	74.1	1 9071B	0722 14:15 0723 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 27 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-25 SC47-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Glass

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	83	ę	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	1380	mg/kg	48.2	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 28 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-26 SC32-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	date Id Prep Anal
Solids, Total	57	ę	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	6900	mg/kg	70.2	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 29 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-27 SC33-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	29	ę	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	9080	mg/kg	138	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 30 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-28 SW2-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	53	Þ	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	87700	mg/kg	75.5	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 31 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-29 SC39-S1	Date Collected: Date Received :	08-JUL-2008 11:40 16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL.	REF METHOD	date id Prep Anal
Solids, Total	23	8	0.10	30 2540g	0717 15:07 SD
TPH, HEM-SGT	12400	mg/kg	156	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 32 of 39

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-30 SC38-S1	Date Collected: 08-JUL-2008 11:40 Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	36	9	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	7300	mg/kg	100	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 33 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-31	Date Collected:	08-JUL-2008 11:40
	SC34-S1	Date Received :	16-JUL-2008
Sample Matrix:	SOIL	Date Reported :	28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Glass

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Parameter	RESULT	UNITS	RDL	REF METHOD	DATE ID PRE P ANAL
Solids, Total	49	ę	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	3480	mg/kg	81.6	1 9071B	0723 14:45.0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 34 of 39

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810512-32	Date Collected: 08-JUL-2008 11:40
	SC35-S1	Date Received : 16-JUL-2008
Sample Matrix:	SOIL	Date Reported : 28-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Glass

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Parameter	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	52	₽ Ŏ	0.10	30 2540G	0717 15:07 SD
TPH, HEM-SGT	6600	mg/kg	76.9	1 9071B	0723 14:45 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07280813:27 Page 35 of 39

CHAIN OF CUSTODY

CHAIN OF CUSTO	CUSTODY PAGE 1 OF M.	Date Recid in Lab:	ALPHA Job#: 0807039
MANSFIELD, MA	Project Information	Report Information - Data Deliverables	Billing Information
TEL: 508-822-9300 FAX: 508-822-3288	Project Name: Tidewoter MGP		X Same as Client into PO #:
Client Information	Project Location: DOWTH ICKET RI	XADEx C Add'l Deliverables	
client: Arcadis	{	'ements/Repo	
Address: 100 CUMmings Ctr, 135P	Project Manager: MOLK MOLDOR	State /Fed Program Criteria	
Beverly, MA OTHER	ALPHA Quote #:	MA MCP PRESUMPTIVE CERTAINTY CT REASONABLE CONFIDENCE PROTOCOLS	EASONABLE CONFIDENCE PROTOCOLS
Õ	Turn-Around Time		Required?
Fax: 978-921-0939		D Yes D No Are CT RCP (Reasonable Co	Are CT RCP (Reasonable Confidence Protocols) Required?
Email:	L HUSH (mit)		/cm/ / / /
These samples have been previously analyzed by Alpha	Date Due: Time:	-	
Other Project Specific Requirements/Comments/Detection L	ents/Detection Limits:		
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Page 265 / 277

ALRHAJob#: 0807059 Billing Information Same as Client info		MAMCP PRESUMIPTIVECERTAINTYCTREASONABLECONFIDENCE PROTOCOLS	Filtration		Plaate specify below 1 Sample Specific Comments 'S	Archive 1		Archive 1	Henve 1	Achive 1	1 20/112/14	Hrhive 1	Arche VC]	CREasephrinteleath; regibly and	Date/Time difference of the randound und clock without standound under a clock of 15 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2
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CUSTODY PAGE 2 OF M Project Information Project Name: Ticlewerker M6P	1357 Project Location: Low tyckst, BL Project #: 36697 1357 Project Manager: Mark Mahoney	ALPHA quote #: Turn-Around Time Kstandard DRISH intercenting it measured	bate Due: Time: nts/Detection Limits:	(ronlers)	ction Sample Time Matrix	1402 Sed	1453 Sed	718108 1454 Sed KEUBT	1537 Sed	1538	1013 Sec.	llois Sed	1017 Sed	Container Type Preservative	Relinquistred By: Date/Time
CHAIN OF CHAIN OF SFIELD, MA SOF 822, 3288	5 yumings Ctr. 135P	Phone: 978 - 921 - 0442 01915 Fax: 978 - 921 - 0442 01915	Email: Date Due: These samples have been previously analyzed by Apha D These samples have been previously analyzed by Apha Other Project Specific Requirements/Comments/Detection Limits:		(Lab.Use Only). Sample ID	5C-21-55	-4 X-0-51	<u> SC-10-52</u> <u> </u>	-5 SC-12-51	<u>86-12-53</u>	<u> </u>	Sc - X - X	SC - 35 - 53	PLEASE ANSWER QUESTIONS ABOVE!	IS YOUR PROJECT MA MCP or CT RCP? 3.2

Page 266 / 277

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CHAIN OF	CHAIN OF CUSTODY ME S of M	Date Recid in Lab	
ncal. Mansfield, Ma	Project Information	Report Information - Data Deliverables	ables Billing Information
TEL: 508-898-9220 TEL: 508-822-3300 FAX: 508-898-9133 FAX: 508-822-3288	Project Name: Tidewater MGP	-	□ Same as Client info PO #:
Client Information	Project Location: Powtucket, RI	ADEx D Add'i Deliverables	5
Client: Arcadis	Project #: 36697	ements/Repo	
Address: 100 Cummings Ctr., #135P	Project Manager: MGrK MghOney		Uniteria
Beverly, MA GOUGIS	ALPHA Quote #:	MAMCP PRESUMPTIVE CERTAINT	MAMCP PRESUMPTIVE CERTAINTY CT REASONABLE CONFIDENCE PROTOCOLS
Phone: 978-921-0442	Turn-Around Time	٩ N	Are MCP Analytical Methods Required?
Fax: 978-921-0939		C Yes CI No Are CT RCP (Reas	Are CT RCP (Reasonable Confidence Protocols) Required?
Email:	Timo.		
These samples have been previously analyzed by Alpha			
Other Project Specific Requirements/Comments/Detection Limits:		24 29	
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Page 267 / 277

CHAIN OF CUSTO	F CUSTODY PAGE 4 OF Date Recontinuation	ALPHA JOD#: 0807059
	Project Information	s Billing Information
TEL: 508-998-9220 TEL: 508-822-9300 FAX: 508-998-9193 FAX: 508-822-3288	-	Same as Glient Info PO #:
atic	Project Location: POW + UCKC+, RI ADEx D Add'i Deliverables	
Client: Arcadis	rements/Repo	
Address: 100 Cummings Ctr. #135P	Project Manager: MOrK MOhoney	
Beveriy, MA OZHEZ 01915	ALPHA Quote #:	MA MCP PRESUMPTIVE CERTAINTY CT REASONABLE CONFIDENCE PROTOCOLS
Phone: 978-921-0442	Turn-Around Time Date Date D No Are MCP Analytical Methods Required?	hods Required?
Fax: 978-921-0939		Are CT RCP (Reasonable Confidênce Protocols) Required?
Email:		
These samples have been previously analyzed by Alpha		
Other Project Specific Requirements/Comments/Detection Limits:	20	
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PLEASE ANSWER QUESTIONS ABOVE!	Container Type G G G G G	Ribasefoljing deaty Neology and
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MA MCP or CT RCP?	Relinquished By: Date/Time Repeived By:	not start unt lived. All so
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CHAIN (CHAIN (CHAIN OF CUSTODY	Date Rec(d)h Lab. 5 h	ALPHA Job#:
WESTBORO, MA WESTBORO, MA TEL: 500 0000 0001 TEL: 500 000 0000	Project Information	Report Information - Data Deliverables	Billing Information
	Project Name: Tidewoter MGP	D FAX XEMAIL	Came as Client info PO #:
Client Information	Project Location: POWTUCKET, RI	ADEX Dadd' Deliverables	
client: Arcadis		ements/Repo	
Address: 100 CUMMings Ctr, #135P	5P Project Manager: MGrK MQhOney	State / Fug Frogram	
Beveriv. M.A. Olgis	ALPHA Quote #:	MAMCPPRESUMPTIVE CERTAINTY CTREASONABLE CONFIDENCE PROTOCOLS	FEASONABLE CONFIDENCE PROTOCOLS
Phone: 978 - 921 - 0442	Turm-Around Time	Yes U No Are MCP Analytical Methods:Bernitred?	Bernined?
Fax: 978-921-0939		U Yes D No Are CT RCP (Reasonable Co	Are CT RCP (Reasonable Confidence Protocols) Required?
Email:	L RUSH (anty continued if pro-approved)		
These samples have been previously analyzed by Alpha	Date Due:		
Other Project Specific Requirements/Comments/Detection Li	mits:	x / 2	
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DATA REVIEW FOR

TIDEWATER MGP

PAWTUCKET, RHODE ISLAND

SDG #0807065

VOLATILE, SEMIVOLATILE, METALS, AND MISCELLANEOUS ANALYSES

Analyses performed by:

Alpha Woods Hole Labs Mansfield, Massachusetts

Review performed by:



Syracuse, New York Report #8850

Summary

The following is an assessment of the data package for sample delivery group (SDG) #0807059 for sampling from the Tidewater MGP Site. Included with this assessment are the corrected sample results and chain of custody. Analyses were performed on the following samples:

Sample ID	Lab ID	Matrix	Sample			Analy	/sis		
			Date	VOC	SVOC	ТРН	MET	PCBs	MISC
SC27-S1	0807065-01	Sediment	7/14/2008	Х	Х		Х		Х
SC44-S1	0807065-02	Sediment	7/14/2008	Х	Х		Х		Х
DUP-2	0807065-03	Sediment	7/14/2008	Х	Х		Х		Х
SC45-S1	0807065-04	Sediment	7/14/2008	Х	Х		Х		Х
SC46-S1	0807065-05	Sediment	7/14/2008	Х	Х		Х		Х
SC18-S1	0807065-06	Sediment	7/14/2008	Х	Х		Х		Х
SC7-S1	0807065-07	Sediment	7/14/2008	Х	Х		Х		Х
SC8-S1	0807065-08	Sediment	7/14/2008	Х	Х		Х		Х
SC31-S1	0807065-09	Sediment	7/14/2008	Х	Х		Х		Х
SC36-S1	0807065-10	Sediment	7/14/2008	Х	Х		Х		Х
SC37-S1	0807065-11	Sediment	7/14/2008	Х	Х		Х		Х
Trip Blank	0807065-12	Water	7/14/2008	Х					
SC1-S1	0807065-13	Sediment	7/15/2008	Х	Х		Х		Х
SC2-S1	0807065-14	Sediment	7/15/2008	Х	Х		Х		Х
Dup-3	0807065-15	Sediment	7/15/2008	Х	Х		Х		Х
SC11-S1	0807065-16	Sediment	7/15/2008	Х	Х		Х		Х
SC13-S1	0807065-17	Sediment	7/15/2008	Х	Х		Х		Х
SC14-S1	0807065-18	Sediment	7/15/2008	Х	Х		Х		Х
SC15-S1	0807065-19	Sediment	7/16/2008	Х	Х		Х		Х
SC16-S1	0807065-20	Sediment	7/16/2008	Х	Х		Х		Х
Rinsate-1	0807065-21	Water	7/16/2008	Х	Х		Х		Х

Notes:

1. Miscellaneous parameters include total organic carbon, physiologically available cyanide, and oil and grease.

2. Sample location DUP-2 is the field duplicate of parent sample location SC44-S1.

3. Sample location DUP-3 is the field duplicate of parent sample location SC2-S1.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

Introduction

Analyses were performed according to (United Stated Environmental Protection Agency) USEPA SW-846 Method 8260. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999. The quality indicators of this limited data review included holding times, associated blanks, matrix spike/matrix spike duplicate (MS/MSD) analysis, field duplicates, laboratory control sample and surrogate recoveries.

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 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
	Water	14 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
SW-846 8260	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
Trip Blank	Analysis Completed	26 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.

	Quali	fication
Criteria	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
SC31-S1	Acetone	Detected sample results <rl <bal<="" and="" td=""><td>"U" at the PQL</td></rl>	"U" at the PQL
SC27-S1 SC44-S1 SC45-S1 SC46-S1 SC7-S1 SC8-S1 SC37-S1 SC1-S1 SC2-S1 Dup-3 SC13-S1	Acetone	Detected sample results >RL and <bal< td=""><td>"U" at detected sample concentration</td></bal<>	"U" at detected sample concentration

RL = reporting limit

3. 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
	1,2-Dichloroethane-d4	AC
DUP-2	4-Bromofluorobenzene	< LL but > 10%
DUP-2	Dibromofluoromethane	AC
	Toluene-d8	AC
Lower control limit (LL)		

Lower control limit (LL) 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	Non-detect	No Action

Control Limit	Sample Result	Qualification
calibration curve due to the high concentration of a target compounds	Detect	

4. 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 were the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

Almost all compounds associated with sample location SC37-S1 had one or both MS/MSD recoveries below the lower control limit and above 10%, and all of the RPDs between the MS/MSD recoveries did not meet acceptance limits; therefore the associated sample result was qualified as estimated (J).

Sample Locations	Compound	MS Recovery	MSD Recovery
	Bromomethane	<ll but=""> 10%</ll>	<ll but=""> 10%</ll>
	Vinyl acetate	AC	<ll but=""> 10%</ll>
	p-Isopropyltoluene	AC	<ll but=""> 10%</ll>
SC1-S1	n-Butylbenzene	AC	<ll but=""> 10%</ll>
	1,2,4-Trichlorobenzene	AC	<ll but=""> 10%</ll>
	Hexachlorobutadiene	<ll but=""> 10%</ll>	<ll but=""> 10%</ll>
	1,2,3-Trichlorobenzene	AC	<ll but=""> 10%</ll>

Sample locations associated with the MS/MSD exhibiting recoveries outside of the control limits are presented in the following table.

AC = Acceptable

The criteria used to evaluate the MS/MSD recoveries are presented in the following table. In the case of an MS/MSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper control limit (III.)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (LL) but > 10%	Non-detect	UJ
	Detect	J
< 10%	Non-detect	R
< 10%	Detect	J
Parent sample concentration > four times the MS/MSD	Detect	No Action
spiking solution concentration (D).	Non-detect	NU ACIION

Sample locations associated with MS/MSD recoveries exhibiting an RPD greater than of the control limit presented in the following table.

Sample Locations	Compound
	Vinyl acetate
	sec-Butylbenzene
	p-Isopropyltoluene
SC1-S1	n-Butylbenzene
	1,2,4-Trichlorobenzene
	Hexachlorobutadiene
	1,2,3-Trichlorobenzene

The criteria used to evaluate the RPD between the MS/MSD recoveries are presented in the following table. In the case of an RPD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
S10	Non-detect	J
> UL	Detect	J

5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/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.

Sample locations associated with LCS/LCSD analysis exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Compound	LCS Recovery	LCSD Recovery
Rinsate-1	Dichlorodifluoromethane	>UL	AC
SC27-S1 SC44-S1 SC45-S1 SC46-S1 SC18-S1 SC7-S1	Bromomethane	<ll but=""> 10%</ll>	<ll but=""> 10%</ll>
SC8-S1 SC31-S1 SC36-S1 SC37-S1 Trip Blank	Acetone	AC	>UL
DUP-2 SC1-S1 SC2-S1 Dup-3	Bromomethane	<ll but=""> 10%</ll>	AC

Sample Locations	Compound	LCS Recovery	LCSD Recovery
SC11-S1			
SC13-S1			
SC14-S1	Acetone	AC	>UL
SC15-S1			
SC16-S1			

The criteria used to evaluate the LCS/LCSD recoveries are presented in the following table. In the case of an LCS deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper central limit (LLL)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (1.1.) but > 100/	Non-detect	J
< the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
< 10%	Detect	J

6. 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 and 100% for soil 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
	Acetone	ND(3.87)	97.3	NC
SC44-S1/DUP-2	Carbon disulfide	5.3	8.9	50.7%
	2-Butanone (MEK)	7.0	15	72.7%
SC2-S1/Dup-3	Carbon disulfide	5.7	4.6	21.4%

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

NC = Not compliant.

The compound acetone associated with samples SC44-S1 and DUP-2 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.

7. 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.

8. 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.

SEMI-VOLATILE ORGANIC COMPOUND (SVOC) ANALYSES

Introduction

Analyses were performed according to (United Stated Environmental Protection Agency) USEPA SW-846 Method 8270 selective ion monitoring (SIM). Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999. The quality indicators of this limited data review included holding times, associated blanks, matrix spike/matrix spike duplicate (MS/MSD) analysis, field duplicates, laboratory control sample and surrogate recoveries.

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 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 8270 SIM	Water	7 days from collection to extraction and 40 days from extraction to analysis	Cooled @ 4 °C
3W-040 0270 3IW	Soil	14 days from collection to extraction and 40 days from extraction to analysis	Cooled @ 4 °C

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance blanks (i.e., method 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. 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. 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. SVOC analysis requires that two of the three SVOC surrogate compounds within each fraction exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

4. Matrix Spike/Matrix Spike Duplicate (MS/MSD)/Laboratory Duplicate 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 were the compounds concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

Sample locations associated with the MS/MSD exhibiting recoveries outside of the control limits are presented in the following table.

Sample Locations	Compound	MS Recovery	MSD Recovery
	Fluorene	> UL	AC
SC1-S1	Anthracene	> UL	AC
301-31	Indeno(1,2,3-cd)pyrene	> UL	AC
	Benzo(g,h,i)perylene	> UL	AC

AC = Acceptable

The criteria used to evaluate the MS/MSD recoveries are presented in the following table. In the case of an MS/MSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper central limit (LLL)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (1.1.) but > 10%	Non-detect	J
< the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
< 10%	Detect	J

Sample locations associated with MS/MSD recoveries exhibiting an RPD greater than of the control limit presented in the following table.

Sample Locations	Compound
SC1-S1	Acenaphthene Fluorene Anthracene
	Indeno(1,2,3-cd)pyrene Benzo(g,h,i)perylene

The criteria used to evaluate the RPD between the MS/MSD recoveries are presented in the following table. In the case of an RPD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> UL	Non-detect	J
2 0L	Detect	J

5. 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.

6. 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 and 100% for soil matrices is applied to the RPD between the parent sample and the field duplicate.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
SC44-S1/DUP-2	Acenaphthene	36.6	56	41.9%
	Acenaphthylene	205	125	48.5%
	Anthracene	208	191	8.5%
	Benz[a]anthracene	841	739	12.9%
	Benzo[a]pyrene	688	606	12.7%
	Benzo[b]fluoranthene	701	635	9.9%
	Benzo[g,h,i]perylene	509	464	9.2%
	Benzo[k]fluoranthene	750	703	6.5%
	Chrysene	961	817	16.2%
	Dibenz[a,h]anthracene	140	124	12.1%
	Fluoranthene	1760	1670	5.2%
	Fluorene	66.4	67.1	1.0%
	Indeno[1,2,3-cd]pyrene	493	481	2.5%
	Naphthalene	58	96	49.4%
	Phenanthrene	608	721	17.0%
	Pyrene	1610	1420	12.5%
	Acenaphthene	296	50.4	141.8%
	Acenaphthylene	323	179	57.4%
	Anthracene	3200	282	167.6%
SC2-S1/Dup-3	Benz[a]anthracene	3390	929	114.0%
	Benzo[a]pyrene	2570	798	105.2%
	Benzo[b]fluoranthene	2070	790	89.5%
	Benzo[g,h,i]perylene	1250	444	95.2%
	Benzo[k]fluoranthene	2120	691	101.7%
	Chrysene	3320	1040	104.6%
	Dibenz[a,h]anthracene	406	127	104.7%
	Fluoranthene	7870	2090	116.1%
	Fluorene	523	69.4	153.1%
	Indeno[1,2,3-cd]pyrene	1440	478	100.3%
	Naphthalene	483	93.7	135.0%
	Phenanthrene	6160	1050	141.7%
	Pyrene	6570	2230	98.6%

Results for duplicate samples are summarized in the following table.

- ND = Not detected.
- AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

All compounds except acenaphthylene, benzo[b]fluoranthene, benzo[g,h,i]perylene, and pyrene associated with samples SC2-S1 and Dup-3 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.

7. 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
	Phenanthrene	112000 E	105000 D	105000 D
SC36-S1	Fluoranthene	73600 E	73300 D	73300 D
	Pyrene	95200 E	94000 D	94000 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

8. 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.

METALS ANALYSES

Introduction

Analyses were performed according to United States Environmental Protection Agency (USEPA) SW-846 Method 6000/7000. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1994. The quality indicators of this limited data review included holding times, associated blanks, matrix spike (MS) and duplicate analysis, serial dilution analysis, field duplicate, and laboratory control sample recoveries.

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 that it was already 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 the USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The analyte was analyzed for but not detected. The associated value is the analyte instrument detection limit.
 - B The reported value was obtained from a reading less than the contract-required detection limit (CRDL), but greater than or equal to the instrument detection limit (IDL).
- Quantitation (Q) Qualifiers
 - E The reported value is estimated due to the presence of interference.
 - N Spiked sample recovery is not within control limits.
 - * Duplicate analysis is not within control limits.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - 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 6020	Water	180 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
	Soil	180 days from collection to analysis	Cooled @ 4 °C.
SW-846 7470	Water	28 days from collection to analysis	Cooled @ 4 °C; preserved to a pH of less than 2.
SW-846 7471	Soil	28 days from collection to analysis	Cooled @ 4 °C.

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method or rinse blanks), are prepared to identify any contamination that may have been introduced into the samples during sample preparation or field activity. Method blanks (including initial and continuing calibration blanks, and preparation blanks) measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected analyte in an associated blank is calculated for QA blanks containing concentrations greater than the IDL. The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No analytes were detected above the reporting limit in the associated blanks.

3. Matrix Spike (MS)/Laboratory Duplicate Analysis

MS and laboratory duplicate data are used to assess the precision and accuracy of the analytical method.

3.1 MS Analysis

All metal analytes must exhibit a percent recovery within the established acceptance limits of 75% to 125%. The MS recovery control limits do not apply for MS performed on sample locations where the analyte's concentration detected in the parent sample exceeds the MS concentration by a factor of four or greater. In instance where this is true, the data will not be qualified even if the percent recovery does not meet the control limits and the laboratory

qualifier "N" will be removed.

All analytes associated with MS recoveries were within control limits with the exception of the following analytes present in the table below.

Sample Location	Analytes	MS Recovery
SC1-S1	Antimony	64%
0.037.04	Antimony	33%
SC37-S1	Mercury	127%

The criteria used to evaluate MS recoveries are presented in the following table. In the case of an MS deviation, the sample results are qualified. The qualifications are applied to all sample results associated with this SDG.

Control limit	Sample Result	Qualification
MS percent	Non-detect	UJ
recovery 30% to 74%	Detect	J
MS percent	Non-detect	R
recovery <30%	Detect	J
MS percent	Non-detect	No Action
recovery >125%	Detect	J

3.2 Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the CRDL. A control limit of 20% for water matrices and 35% for soil matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the CRDL, a control limit of one times the CRDL is applied for water matrices and two times the CRDL for soil matrices.

All analytes associated with laboratory duplicate RPD were within the control limit, with the exception of the analytes presented in the following table.

Sample Location	Analytes	Laboratory RPD
SC37-S1	Selenium	26%
	Beryllium	42%
	Chromium	37%
SC1-S1	Lead	23%
	Nickel	56%
	Zinc	32%

The criteria used to evaluate laboratory duplicate RPD are presented in the following table. In the case of a laboratory duplicate RPD deviation, the sample results are qualified. The qualifications are applied to the all sample results associated with this SDG.

Sample Concentration	Control Limit	Sample Result	Qualification
Parent sample and laboratory sample concentration >5 times	Water 20%	Non-detect	UJ
CRDL	Soil 35%	Detect	J
Parent sample and/or laboratory duplicate sample result ≤ five	Water one times RL	Non-detect	UJ
times the RL and difference between samples >RL	Soil two times RL	Detect	J

4. 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 and 100% for soil 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
	Antimony	0.172	0.155	10.3 %
	Arsenic	3.04	2.79	8.5 %
	Beryllium	0.233	0.227	2.6 %
	Cadmium	1.26	1.28	1.5 %
	Chromium	18.2	17.9	1.6 %
	Copper	39.5	39.8	0.7 %
SC44-S1/DUP-2	Lead	48.9	48.3	1.2 %
	Nickel	8.87	8.57	3.4 %
	Selenium	0.34	0.333	2.0 %
	Silver	0.447	0.444	0.6 %
	Thallium	0.0566	0.0562	0.7 %
	Zinc	89.3	84.4	5.6 %
	Mercury	0.104	0.101	2.9 %
	Antimony	0.169	0.273	47.0 %
SC2-S1/Dup-3	Arsenic	5.18	3.54	37.6 %
	Beryllium	0.364	0.224	47.6 %

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
	Cadmium	0.816	0.716	13.0 %
	Chromium	15.5	16.4	5.6 %
	Copper	35.8	29	20.9 %
	Lead	55.9	47.9	15.4 %
	Nickel	11.9	9.73	20.0 %
	Silver	0.105	0.0997	5.1 %
	Thallium	0.0645	ND(0.0377)	AC
	Zinc	75.6	76.1	0.6 %
	Mercury	0.0311	0.03	3.6 %

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two 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.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences.

All LCS recoveries were within control limits, with the exception of the analytes associated with sample locations, as presented in the following table.

Sample Location	Analytes	LCS Recovery
Rinsate-1	Nickel	122%
Rinsale-1	Zinc	121%

The criteria used to evaluate LCS recoveries are presented in the following table. In the case of an LCS deviation, the sample results are qualified.

Control limit	Sample Result	Qualification
LCS (water) percept receivery E0% to 70%	Non-detect	UJ
LCS (water) percent recovery 50% to 79%	Detect	J
LCS (water) percent recovery <50%	Non-detect	R
ECS (water) percent recovery <50%	Detect	J
LCS (water) percent receiver (>120%	Non-detect	No Action
LCS (water) percent recovery >120%	Detect	J
LCC (apil) percent receiver (< lower limit	Non-detect	J
LCS (soil) percent recovery < lower limit	Detect	J

Control limit	Sample Result	Qualification
	Non-detect	No Action
LCS (soil) percent recovery > upper limit	Detect	J

6. Serial Dilution

The serial dilution analysis is used to assess if a significant physical or chemical interference exists due to sample matrix. Analytes exhibiting concentrations greater than 50 times the MDL in the undiluted sample are evaluated to determine if matrix interference exists. These analytes are required to have less than a 10% difference (%D) between sample results from the undiluted (parent) sample and results associated with the same sample analyzed with a five-fold dilution.

All serial dilutions were within control limits.

7. 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.

MISCELLANEOUS ANALYSES

Introduction

Analyses were performed according to United States Environmental Protection Agency (USEPA) SW-846 methods 9010, 9060, and 9071b. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1994. The quality indicators of this limited data review included holding times, associated blanks, matrix spike (MS) and laboratory duplicate analysis, field duplicate, and laboratory control sample recoveries.

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 that it was already 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 the USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The analyte was analyzed for but not detected. The associated value is the analyte instrument detection limit.
 - B The reported value was obtained from a reading less than the contract-required detection limit (CRDL), but greater than or equal to the instrument detection limit (IDL).
- Quantitation (Q) Qualifiers
 - N Spiked sample recovery is not within control limits.
 - * Duplicate analysis is not within control limits.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - 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 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
Physiologically Available Cyanide by SW-846 9010	Soil	14 days from collection to analysis	Cooled @ 4 °C
Total Organic Carbon by 9060	Soil	28 days from collection to analysis	Cooled @ 4 °C
Oil and Grease by 9071b	Soil	28 days from collection to analysis	Cooled @ 4 °C

All samples were analyzed within the specified holding times.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method or rinse blanks), are prepared to identify any contamination that may have been introduced into the samples during sample preparation or field activity. Method blanks (including initial and continuing calibration blanks, and preparation blanks) measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected analyte in an associated blank is calculated for QA blanks containing concentrations greater than the MDL. The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

No analytes were detected above the reporting limit in the associated blanks.

3. Matrix Spike (MS)/Laboratory Duplicate Analysis

MS and laboratory duplicate data are used to assess the precision and accuracy of the analytical method.

3.1 MS Analysis

All analyses must exhibit a percent recovery within the established acceptance limits of 75% to 125%. The MS recovery control limits do not apply for MS performed on sample locations where the analyte's concentration detected in the parent sample exceeds the MS concentration by a factor of four or greater

All analytes associated with MS recoveries were within control limits with the exception of the following analytes present in the table below.

Sample Location	Analytes	MS Recovery
SC37-S1	TOC	132%
SC1-S1	TOC	215%

The criteria used to evaluate MS recoveries are presented in the following table. In the case of an MS deviation, the sample results are qualified. The qualifications are applied to all sample results associated with this SDG.

Control limit	Sample Result	Qualification
MS percent recovery 30% to	Non-detect	UJ
74%	Detect	J
MS percent	Non-detect	R
recovery <30%	Detect	J
MS percent	Non-detect	No Action
recovery >125%	Detect	J

3.2 Laboratory Duplicate Analysis

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the CRDL. A control limit of 20% for water matrices and 35% for soil matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the CRDL, a control limit of one times the CRDL is applied for water matrices and two times the CRDL for soil matrices.

All analytes associated with laboratory duplicate RPD were within the control limit, with the exception of the analytes presented in the following table.

Sample Location	Analytes	Laboratory RPD	
SC37-S1	TOC	35%	

The criteria used to evaluate laboratory duplicate RPD are presented in the following table. In the case of a laboratory duplicate RPD deviation, the sample results are qualified. The qualifications are applied to the all sample results associated with this SDG.

Sample Concentration	Control Limit	Sample Result	Qualification
Parent sample and laboratory sample concentration >5 times	Soil 35%	Non-detect	UJ
CRDL	3011 33 76	Detect	J
Parent sample and/or laboratory duplicate sample result ≤ five	Soil two times RL	Non-detect	UJ
times the RL and difference between samples >RL		Detect	J

4. 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 and 100% for soil matrices is applied to the RPD between the parent sample and the field duplicate.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
	Total Organic Carbon (Run 1)	1.2	1.2	0%
SC44-S1/DUP-2	Total Organic Carbon (Run 2)	1.2	1.4	15.4%
	TPH, HEM-SGT	958	1260	27.2%
	Physiologically Available Cyanide	ND	ND	AC
	Total Organic Carbon (Run 1)	2.2	0.4	139%
	Total Organic Carbon (Run 2)	1.9	0.22	159%
SC2-S1/Dup-3	TPH, HEM-SGT	520	573	9.7%
	Physiologically Available Cyanide	ND	ND	AC

Results for duplicate samples are summarized in the following table.

ND = Not detected.

AC = The field duplicate RPD is acceptable when the RPD between parent sample and field duplicate sample is less than one/or/two times the RL and where the parent sample and/or duplicate concentration is less than five times the RL.

The compound total organic carbon associated with samples SC2-S1 and Dup-3 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.

5. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences.

The LCS analysis exhibited recoveries within the control limits.

6. System Performance and Overall Assessment

Upon validation of the TOC analysis performed within this SDG it was discovered that all samples were analyzed in duplicate. Technical compliance of method SW-846 9060 requires analysis to be performed in quadruplicate. The purpose of the quadruplicate analysis is to establish data precision. Since the laboratory performed the sample analysis in duplicate for each sample location, for data qualification and usability purposes the sample analysis precision of the two reported runs for each sample location were evaluated against laboratory duplicate criterion:

The laboratory duplicate relative percent difference (RPD) criterion is applied when parent and duplicate sample concentrations are greater than or equal to 5 times the reporting limit (RL). A control

limit of 20% for water matrices and 35% for soil matrices is applied when the criteria above is true. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of one times the RL is applied for water matrices and two times the RL for soil matrices.

All analytes associated with laboratory duplicate RPD were within the control limit, with the exception of the analytes presented in the following table.

Sample Location	Analytes	Laboratory RPD	
SC11-S1	Total Organic Carbon	40.0%	
SC13-S1	Total Organic Carbon	42.9%	

The criteria used to evaluate laboratory duplicate RPD (in this case the two sample runs) are presented in the following table. In the case of a laboratory duplicate RPD deviation, the sample results are qualified.

Sample Concentration	Control Limit	Sample Result	Qualification
Parent sample and laboratory	Water 20%	Non-detect	UJ
sample concentration >5 times RL	Soil 35%	Detect	J
Parent sample and/or laboratory		Non-detect	UJ
duplicate sample result ≤ five times the RL and difference between samples >RL	Water one times RL Soil two times RL	Detect	J

The duplicate sample results exhibited RPD within the control limit for all sample locations within this SDG; therefore no results were qualified due to the method deviation.

CORRECTED SAMPLE ANALYSIS DATA SHEETS



ARCADIS

Tidewater MGP SC27-S1

SDG: N/A

Case: N/A Matrix: Sediment Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-01

Associated Blank: VS072508B04

Concentration Units:	μg/Kg
Dilution	

Dete Callestad	Data Bassivad	Data Analyzad	Percent Solid	Sample	Final Volume (mt)	Dilution Factor		5/122
		Date Analyzed	· · · · · · · · · · · · · · · · · · ·	Amount (g)	Volume (ml)		Analy	
07/14/08	07/16/08	07/25/08	50.8	7.74	5	1	BAS	5
Parameter		F	Result	Parameter	Parameter		Result	
Dichlorodifluoro	methane		<u>2.54 U</u>	2-Hexanone			6.36	
Chloromethane			<u>2.54 U</u>	Tetrachloroeth	ene		2.54	U
Vinyl chloride			2.54 U	1,3-Dichlorop			2.54	
Bromomethane			<u>2.54 U J</u>	Dibromochlor			2.54	U
		2.54 U	1,2-Dibromoe			2.54	Ŭ	
Frichlorofluoromethane		2.54 U	Chlorobenzen	3		2.54	U	
Acetone			32.3 B U	1,1,1,2-Tetrac	hloroethane		2.54	U
1,1-Dichloroethe	ne		<u>2.54 U</u>	Ethylbenzene			2.54	U
Carbon disulfide			2.99	p/m-Xylene			5.09	U
Methylene chlor	ide		6.36 U	o-Xylene			2.54	U
Methyl tert-buty	l ether (MTBE)		2.54 U	Styrene			2.54	U
rans-1,2-Dichlo	roethene		2.54 U	Bromoform			2.54	U
1,1-Dichloroethane		2.54 U	Isopropylbenzene			2.54	U	
Vinyl acetate		2.54 U	1,1,2,2-Tetrachloroethane			2,54	U	
		7.28	Bromobenzene			2.54	U	
cis-1,2-Dichloro	ethene		2.54 Ŭ	1,2,3-Trichloropropane			2.54	U
2,2-Dichloropro	pane		2.54 U	n-Propylbenzene			2.54	U
Chloroform			2.54 U	2-Chlorotoluene			2.54	U
1,1,1-Trichloroe	thane		2.54 U	1,3,5-Trimethy	ylbenzene		2.54	U
1,1-Dichloropro	pene		2.54 U	4-Chlorotolue	ne		2.54	U
Carbon tetrachlo	ride		2.54 U	tert-Butylbenz	ene		2.54	
Benzene			2.54 U	1,2,4-Trimethy	ylbenzene		2.54	
1,2-Dichloroetha	me		2.54 U	sec-Butylbenz	ene		2,54	U
Trichloroethene			2,54 U	1,3-Dichlorob	enzene		2.54	U
1,2-Dichloropro	pane		2.54 U	p-Isopropyltol	uene		6.36	U
Dibromomethan			2.54 U	1,4-Dichlorob	enzene		2.54	
Bromodichloron			2.54 U	n-Butylbenzen			6.36	
2-Chloroethylvii	nyl ether		6.36 U	1,2-Dichlorob			2.54	
	ketone (MIBK)		2.54 Ŭ	1,2-Dibromo-3	3-chloropropane		2.54	
cis-1,3-Dichloro			2.54 U	1,2,4-Trichlor		<u></u>	2.54	ch free as well.
Toluene			2.54 U	Hexachlorobu			2.54	
trans-1,3-Dichlo	ropropene		2.54 U	Naphthalene			6.36	
1,1,2-Trichloroe			2.54 U	1,2,3-Trichlor	obenzene		2.54	

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	111	70-130
1,2-Dichloroethane-d4	115	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	87	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

Sample



ARCADIS

Tidewater MGP SC44-S1

SDG: N/A

Case: N/A Matrix: Sediment Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-02 Associated Blank: VS072508B04

Concentration Units: µg/Kg Dilution

Final

Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/14/08	07/16/08	07/25/08	76.0	8.49	5	1	BAS
Parameter		I	Result	Parameter	Parameter		
Dichlorodifluoro	methane		1.55 U	2-Hexanone			3.87 U
Chloromethane		and the second	1.55 U	Tetrachloroeth	ene		1.55 U
Vinyl chloride			1.55 U	1,3-Dichlorop			1.55 U
Bromomethane	(_) _ (<i>_</i>) _ _ (<i>_</i>) _ _ (<i>_</i>) _ (<i>_</i>) _ (<i>_</i>) _ _ (<i>_</i>) _ (<i>_</i>) _ (<i>_</i>) _ _ (<i>_</i>) _ (<i>_</i>) _ _ (<i>_</i>) _		1.55 U J	Dibromochlor	omethane		1.55 U
		1.55 U	1,2-Dibromoe	thane		1.55 U	
		1.55 U	Chlorobenzene	e		1,55 U	
Acetone			43.8 BUJ	1,1,1,2-Tetrac	hloroethane		1.55 U
	me		1.55 U	Ethylbenzene			1.55 U
Carbon disulfide			5.25	p/m-Xylene			3.10 U
Methylene chlor			3.87 U	×2 1			1.55 U
			1.55 U	Styrene			1.55 U
		1.55 U	Bromoform			1.55 U	
			1.55 U	Isopropylbenz			1.55 U
		1.55 U	1,1,2,2-Tetrac	1,1,2,2-Tetrachloroethane			
		7.04	Bromobenzen			1.55 U	
cis-1.2-Dichloro	ethene		1.55 U	1,2,3-Trichlor	opropane		1.55 U
2.2-Dichloropro	pane		1.55 U	n-Propylbenze	n-Propylbenzene		1.55 U
Chloroform			1.55 U	2-Chlorotolue	2-Chlorotoluene		
	thane		1.55 U	1,3,5-Trimeth	1,3,5-Trimethylbenzene		
	pene		1.55 U	4-Chlorotolue	4-Chlorotoluene		
Carbon tetrachlo			1.55 U	tert-Butylbenzene			1.55 U
Benzene			1.55 U	1,2,4-Trimeth	1,2,4-Trimethylbenzene		
	ane		1.55 U	sec-Butylbenz			1.55 U
Trichloroethene			1.55 U	1,3-Dichlorob	enzene		1.55 U
			1.55 U	p-Isopropyltol	p-Isopropyltoluene		<u>3.87 U</u>
Dibromomethan			1.55 U	1,4-Dichlorob	enzene		<u>1.55 Ŭ</u>
		1.55 U	1,4-Dichlorobenzene n-Butylbenzene				
2-Chloroethylvinyl ether 3.87 U			1,2-Dichlorobenzene		1 66 11		
Methyl isobutyl ketone (MIBK) 1.55 U		1,2-Dibromo-3-chloropropane			<u>1.55 U</u>		
cis-1,3-Dichloropropene 1.55 U		1,2,4-Trichlorobenzene			1.55 U		
Toluene	4		1,55 Ŭ	Hexachlorobu	itadiene		1.55 U
trans-1,3-Dichlo	ropropene		1.55 U	Naphthalene			3.87 U
1.1.2-Trichloroe			1.55 U	1,2,3-Trichlor	obenzene		1.55 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1.2-Dichloroethane-d4	101	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	80	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.



ARCADIS Tidewater MGP

N/A

Matrix:

DUP-2

SDG: N/A Sediment

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-03

Associated Blank: VS072608B03

µg/Kg

Concentration Units:

				Sample	Final	Dilution			
[Date Received		Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	~~~~·	
07/14/08	07/16/08	07/26/08	73.4	7.40	5	11	BAS	3	
Parameter		ľ	Result	Parameter			Result		
Dichlorodifluoro	omethane		<u>1.84 U J</u>	2-Hexanone	••••••••••••••••••••••••••••••••••••••		4.60	UJ	
Chloromethane			<u>1.84 U</u>	<u>Tetrachloroeth</u>	ene		1.84	U	
Vinyl chloride 1.84 U		1,3-Dichlorop	ropane		1.84	U			
Bromomethane	**		<u>1.84 U</u>	Dibromochlor	omethane		1.84	U	
Chloroethane			1.84 U	1,2-Dibromoe	hane		1.84	U	
Trichlorofluoror	nethane		1.84 U	Chlorobenzen	2		1.84	<u> </u>	
Acetone			97.3 B	1,1,1,2-Tetrac	loroethane		1,84	U	
1,1-Dichloroethe	ene		1.84 U	Ethylbenzene			1.84	<u> </u>	
Carbon disulfide			8.92	p/m-Xylene			3.68	U	
Methylene chlor	ide		4.60 U	o-Xylene			1.84	U	
Methyl tert-buty	l ether (MTBE)		1.84 U	Styrene			1.84	U	
trans-1,2-Dichlo	roethene		1.84 U	Bromoform			1.84	U	
1,1-Dichloroethane 1.84		1.84 U	Isopropylbenz	Isopropylbenzene			U		
Vinyl acetate 1.84		1.84 U	1,1,2,2-Tetrac	1,1,2,2-Tetrachloroethane			U		
2-Butanone (ME	and the set of the set		14.6	Bromobenzene			1.84	U	
cis-1,2-Dichloro	ethene		1.84 U	1,2,3-Trichloropropane			1.84	U	
2,2-Dichloropro	pane		1.84 U	n-Propylbenzene			1.84	U	
Chloroform			1.84 U	2-Chlorotoluene			1.84	U	
1,1,1-Trichloroe	thane		1.84 U	1,3,5-Trimethy	/lbenzene			U	
1,1-Dichloropro	pene		1.84 U	4-Chlorotolue	ne		1.84	U	
Carbon tetrachic	oride		1.84 U	tert-Butylbenz	1.84				
Benzene			1.84 U	1,2,4-Trimethy	/lbenzene		1.84	U	
1,2-Dichloroeth	ane		1.84 U	sec-Butylbenz	ene		1.84	U	
Trichloroethene			1.84 U	1,3-Dichlorob	enzene		1.84	U	
1,2-Dichloropro	pane		1.84 U	p-Isopropyltoluene			4.60	U	
Dibromomethan			1.84 U	1,4-Dichlorobenzene			1.84	U	
Bromodichloron	nethane		1.84 U	n-Butylbenzene			4.60	U	
2-Chloroethylvi	2-Chloroethylvinyl ether 4.60 U		4.60 U	1,2-Dichlorobenzene			1.84	U	
Methyl isobutyl ketone (MIBK) 1.84 U		1,2-Dibromo-3-chloropropane			1.84	U			
cis-1,3-Dichloropropene 1.84 U			1.84 U	1,2,4-Trichlor	obenzene		1.84	U	
Toluene		11 8 4 76 4 - 4	1.84 U	Hexachlorobu	tadiene		1.84	U	
trans-1,3-Dichlo	oropropene		1.84 U	Naphthalene		.,	4.60	U	
1,1,2-Trichloroe			1.84 U 🌡	1,2,3-Trichlor	obenzene		1.84	<u>U</u> -/	

Surrogate	% Recovery		Acceptance Range (%)
Dibromofluoromethane	109		70-130
1,2-Dichloroethane-d4	105		70-130
Toluene-d8	86		70-130
4-Bromofluorobenzene	65	§	70-130

N/A - Not Applicable

§ - Surrogate value outside of acceptable range.

B - Found in associated blank as well as sample.

N/A



ARCADIS **Tidewater MGP**

N/A SDG:

SC45-S1

Sediment Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-04 Associated Blank: VS072508B04

						Concentration	Juns. µg/ng
				Sample	Final	Dilution	
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/14/08	07/16/08	07/25/08	77.8	8.43	5	1	BAS
Parameter		I	Result	Parameter			Result
Dichlorodifluor	omethane		1.52 U	2-Hexanone			<u>3.81 U</u>
Chloromethane			1.52 U	Tetrachloroeth			<u>1.52 U</u>
Vinyl chloride			1,52 U	1,3-Dichlorop			1.52 U
Bromomethane			1.52 U J	Dibromochlor	omethane		<u>1.52 U</u>
Chloroethane			<u>1.52 U</u>	1,2-Dibromoet			1.52 U
Trichlorofluoror	nethane		1.52 U	Chlorobenzen			<u>1.52 U</u>
Acetone			13.0 BU	1,1,1,2-Tetrac	hloroethane		1.52 U
1,1-Dichloroeth	ene		<u>1.52 Ű</u>	Ethylbenzene			1.52 U
Carbon disulfide	•		1.90	p/m-Xylene			3.05 U
Methylene chlor	ide		3.81 U	o-Xylene			<u>1.52 U</u>
Methyl tert-buty			1.52 U	Styrene			1.52 U
trans-1,2-Dichlo	roethene		1.52 U	Bromoform			<u>1.52 U</u>
1,1-Dichloroeth			1.52 U	Isopropylbenzene			1.52 U
Vinyl acetate			1,1,2,2-Tetrachloroethane			<u>1.52</u> U	
2-Butanone (MI	EK)		1.20 J	Bromobenzene			1.52 U
cis-1,2-Dichloro	ethene		1.52 U	1,2,3-Trichloropropane			1.52 Ŭ
2,2-Dichloropro	pane		1.52 U	n-Propylbenzene			1.52 U
Chloroform			1.52 U	2-Chlorotoluene			1.52 U
1,1,1-Trichloroe	thane		1.52 U	1,3,5-Trimethylbenzene			1.52 U
1,1-Dichloropro	pene		1.52 U	4-Chlorotoluene			1.52 U
Carbon tetrachlo	oride		1.52 U	tert-Butylbenzene			1.52 U
Benzene			1.52 U	1,2,4-Trimethylbenzene			1.52 U
1,2-Dichloroeth	ane		1.52 U	sec-Butylbenzene			<u>1.52 U</u>
Trichloroethene			1.52 U	1,3-Dichlorobenzene		1.52 U	
	pane		1.52 U	p-Isopropyltoluene			
Dibromomethar			1.52 U	1,4-Dichlorobenzene			1.52 U
Bromodichloror			1.52 U	n-Butylbenzene			<u>3.81 U</u>
2-Chloroethylvi			3,81 U	1,2-Dichlorobenzene			1.52 U
	Methyl isobutyl ketone (MIBK) 1.52 U		1,2-Dibromo-3-chloropropane			1.52 U	
cis-1,3-Dichloro			1.52 U	1,2,4-Trichlorobenzene			1.52 U
Toluene	••••••••••••••••••••••••••••••••••••••		1.52 U	Hexachlorobutadiene			1.52 U
trans-1,3-Dichlo	ropropene		1.52 U	Naphthalene			3.81 U
1,1,2-Trichloroethane 1.52 U		1,2,3-Trichlor	ohanzana		1.52 U		

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	103	70-130
1.2-Dichloroethane-d4	102	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	79	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

N/A



ARCADIS **Tidewater MGP**

SC18-S1 SDG: N/A

Sediment Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-06

Associated Blank: VS072508B04 Concentration Units:

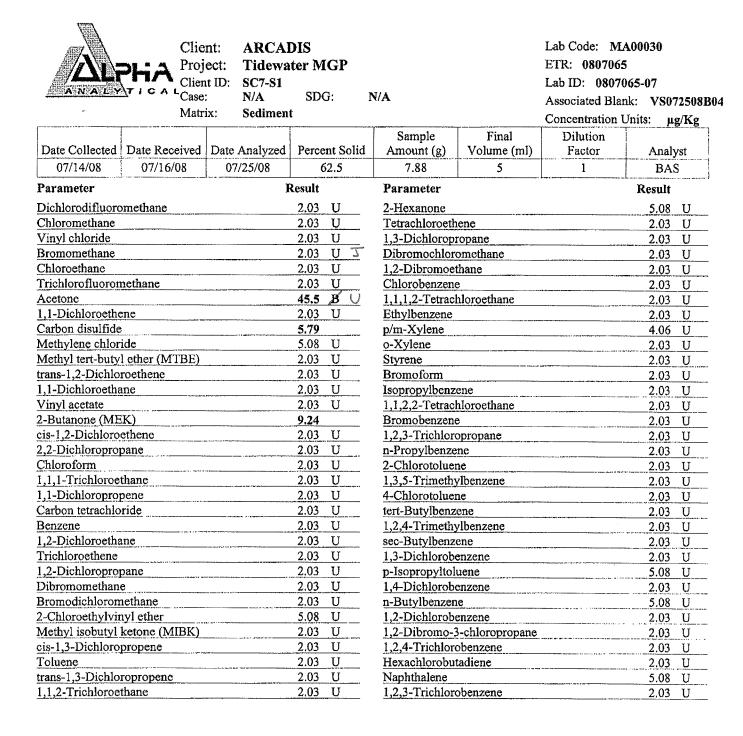
µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	vst
07/14/08	07/16/08	07/25/08	24.3	7.32	5	1	BAS	
Parameter			Result	Parameter			Result	
Dichlorodifluoro	omethane		5.63 U	2-Hexanone			14.1	U
Chloromethane	anna - afa fi tangana		5.63 U	Tetrachloroeth	iene		5.63	U
Vinyl chloride			5.63 U	1,3-Dichlorop	ropane		5.63	U
Bromomethane			5.63 U J	Dibromochlor	omethane		5.63	U
Chloroethane	· · · · · · · · · · · · · · · · · · ·		5.63 U	1,2-Dibromoe	thane		5.63	U
Trichlorofluoror	nethane		5.63 U	Chlorobenzen	8		5.63	U
Acetone		4 H m	252 B J	1,1,1,2-Tetrac	hloroethane		5.63	U
1.1-Dichloroethe	ene		5.63 U	Ethylbenzene			5.63	U
Carbon disulfide)		32.2	p/m-Xylene			11.2	U
Methylene chlor			14.1 U	o-Xylene			5.63	U
Methyl tert-buty			5.63 U	Styrene			5.63	U
trans-1,2-Dichlo			5.63 U	Bromoform			5.63	U
1,1-Dichloroetha			5.63 U	Isopropylbenz	ene		5.63	U
Vinyl acetate			5.63 U	1,1,2,2-Tetrac			5,63	
2-Butanone (MI	EK)		52.2	Bromobenzen	e		5.63	U
cis-1.2-Dichloro	ethene		5.63 U	1,2,3-Trichlor	opropane		5.63	Ŭ
2,2-Dichloropro	pane		5.63 U	n-Propylbenze	ene		5.63	U
Chloroform			5.63 U	2-Chlorotolue	ne		5,63	U
1,1,1-Trichloroe	thane		5.63 U	1,3,5-Trimeth	ylbenzene		5.63	U
1,1-Dichloropro			5.63 U	4-Chlorotolue	ne		5.63	U
Carbon tetrachlo	oride		5.63 U	tert-Butylbenz	zene		5.63	U
Benzene			5.63 U	1,2,4-Trimeth	ylbenzene		5.63	U
1,2-Dichloroeth	ane		5.63 U	sec-Butylbenz	ene		5.63	U
Trichloroethene			5.63 U	1,3-Dichlorob	enzene		5.63	U
1,2-Dichloropro	pane		5.63 U	p-Isopropyltol	uene		14.1	U
Dibromomethan			5.63 U	1,4-Dichlorob	enzene		5.63	U
Bromodichloror	nethane		5.63 U	n-Butylbenzer	ne		14.1	U
2-Chloroethylvi	nyl ether		14.1 U	1,2-Dichlorob	enzene		5.63	U
Methyl isobutyl			5.63 U	1,2-Dibromo-	3-chloropropane		5.63	U
cis-1,3-Dichloro			5.63 U	1,2,4-Trichlor	obenzene		5.63	U
Toluene			5.63 U	Hexachlorobu	tadiene		5.63	U
trans-1,3-Dichlo	ropropene		5.63 U	Naphthalene		,	14.1	U
1,1,2-Trichloroe			5.63 U	1,2,3-Trichlor	obenzene		5.63	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	105	70-130
1,2-Dichloroethane-d4	104	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	82	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.



Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	101	70-130
1,2-Dichloroethane-d4	105	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	88	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

U - The analyte was analyzed for but not detected at the sample specific level reported.

13 of 208 17:43



ARCADIS **Tidewater MGP**

SC8-S1 SDG: N/A

N/A Sediment Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-08 Associated Blank: VS072508B04

Concentration Units: µg/Kg 1

F		,				Concentration	<u> </u>	
				Sample	Final	Dilution	41	
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/14/08	07/16/08	07/25/08	52.7	7.88	5	1	BAS	,
Parameter		F	Result	Parameter			Result	
Dichlorodifluoro	methane		2.41 U	2-Hexanone	L		6.02	
Chloromethane			2.41 U	Tetrachloroeth		· · · · · · · · · · · · · · · · · · ·		U
Vinyl chloride			<u>2.41 U</u>	1,3-Dichlorop				U
Bromomethane			2.41 U J	Dibromochlor				U
Chloroethane			2.41 U	1,2-Dibromoe				U
Trichlorofluoror	nethane		<u>2.41 U</u>	Chlorobenzen				U
Acetone			63.2 BU	1,1,1,2-Tetrac	hloroethane			U
1,1-Dichloroethe	ene		<u>2.41 U</u>	Ethylbenzene				U
Carbon disulfide	,		9.04	p/m-Xylene				<u> </u>
Methylene chlor	ide		6.02 U	o-Xylene	مرب به			U
Methyl tert-buty	l ether (MTBE)		2.41 U	Styrene			2.41	
trans-1,2-Dichlo	roethene		2.41 U	Bromoform			2.41	U
1,1-Dichloroeth	ane		2.41 U	Isopropylbenzene			3.39	
Vinyl acetate	invl acetate 2.41 U		1,1,2,2-Tetrachloroethane				U	
2-Butanone (MI	EK)		12.0	Bromobenzen	e		2.41	
cis-1,2-Dichloro	ethene		2,41 U	1,2,3-Trichloropropane			2.41	
2,2-Dichloropro	pane		2.41 U	n-Propylbenzene			2.41	U
Chloroform			2.41 U	2-Chlorotoluene			2.41	U
1,1,1-Trichloroe	thane		2.41 U	1,3,5-Trimethylbenzene				U
1,1-Dichloropro	pene		2.41 U	4-Chlorotoluene			2,41	U
Carbon tetrachlo	oride		2.41 U	tert-Butylbenz	ene			U
Benzene			2.41 U	1,2,4-Trimeth	ylbenzene		2.41	
1,2-Dichloroeth	ane		2.41 U	sec-Butylbenz		199 (1 1 1	2.41	
Trichloroethene			2.41 U	1,3-Dichlorobenzene		2.41		
1,2-Dichloropro	pane		2.41 U	p-Isopropyltoluene		6.02		
Dibromomethan			2.41 U	1,4-Dichlorobenzene		2.41	U	
Bromodichloror			2.41 U	n-Butylbenzene			6.02	
2-Chloroethylvi	nyl ether		6.02 U	1,2-Dichlorobenzene			2.41	
Methyl isobutyl			2.41 U	1,2-Dibromo-3-chloropropane				
cis-1,3-Dichlord			2.41 U	1,2,4-Trichlorobenzene			2.41	U
Toluene			2.41 U	Hexachlorobutadiene			2.41	U
trans-1,3-Dichlo	propropene		2.41 U	Naphthalene			1.74	J
1,1,2-Trichloroe			2.41 U	1,2,3-Trichlor	obenzene		2,41	Ų

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	100	70-130
1,2-Dichloroethane-d4	101	70-130
Toluene-d8	89	70-130
4-Bromofluorobenzene	92	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

	PHA	Client: Project: Client ID: Case: Matrix:	ARCADI Tidewate SC31-S1 N/A Sediment	r MGP	N/ A		ETR: Lab II Assoc		09 VS081108B10
r	· · · · · · · · · · · · · · · · · · ·	r	T					entration Units	s: µg/Kg
Date	Date	Date	Percent	Methanol	Sample	Extract Volume (ml)	Final Volumo (ml)	Dilution Factor	Analyst
Collected	Received	Analyzed	Solid					1	
07/14/08	07/16/08	07/26/08	75.4	5	7.72	0.1	5		BAS
Parameter			Res		Parameter			R	esult
Dichlorodiflu	uoromethane			118 U	2-Hexanon				<u>296 U</u>
Chlorometha	ine			118 U	Tetrachloro	the second se			<u>118 U</u>
Vinyl chloric	1e			118 U	1,3-Dichlor				<u>118 U</u>
Bromometha	ine			<u>118 U T</u>		loromethane			118 U
Chloroethane	<u>e</u>	119 }		118 U	1,2-Dibron				118 U
Trichlorofluc	oromethane			118 U	Chloroben:				<u>118 U</u>
Acetone								118 U	
1,1-Dichloro				118 U Ethylbenzene				<u>118 U</u>	
Carbon disul			118 U p/m-Xylene					<u>126 J</u>	
Methylene cl			1 weeks	296 U	o-Xylene				96.6 J
	outyl ether (M						118 U		
trans-1,2-Dic	chloroethene			118 U Bromoform					118 U
1,1-Dichloro				118 U	Isopropylb				118 U
Vinyl acetate	8			<u>118 U</u>	and the second sec	rachloroetha	10		118 U
2-Butanone				118 U	Bromoben				118 U
	loroethene			118 U		loropropane			118 U
2,2-Dichloro	propane			<u>118 U</u>	n-Propylbe				118 U
Chloroform				118 U	2-Chloroto				118 U
	oroethane			118 U		ethylbenzene			74.1 J
1,1-Dichloro				118 U	4-Chloroto				<u>118 U</u>
Carbon tetra	chloride			118 U	tert-Butylb		······		118 U
Benzene	and a state of the			139		ethylbenzene			232
1,2-Dichlore			Accel & development and the second	118_U	sec-Butylb				118 U
Trichloroeth	ene			118 U		robenzene			118 U
1,2-Dichloro				118 U	p-Isopropy			.	296 U
Dibromomet	thane			118 U	water and the second	robenzene			<u>118 U</u>
Bromodichle				118 U n-Butylbenzene			<u>296 U</u>		
	ylvinyl ether			296 U 1,2-Dichlorobenzene				<u>118 U</u>	
	utyl ketone (N	AIBK)		118 U		no-3-chlorop			<u>118 U</u>
cis-1,3-Dich	loropropene			118 U		hlorobenzene			<u>69.9 J</u>
Toluene				59.8 J	and an article statement of the statemen	obutadiene			118 U
	chloropropen	e		<u>118 U</u>	Naphthale				<u>9530</u>
1,1,2-Trichle	oroethane			118 U	1,2,3-Tricl	hlorobenzene			118 U

		Acceptance
Surrogate	% Recovery	Range (%)
2-Bromo-1-chloropropane	98	70-130
1-Chloro-2-fluorobenzene	110	70-130
1,4-Dichlorobutane	92	70-130
Dibromofluoromethane	80	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	87	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

N/A



ARCADIS Tidewater MGP

SC36-S1 N/A SDG:

N/A SDC Sediment Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-10 Associated Blank: VS072508B04 Concentration Units: µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analy	
07/14/08	07/16/08	07/25/08	63.3	7.16	5	1	BAS	
Parameter	01110/00		Result	Parameter	ll		Result	
Dichlorodifluoro	methane		2.21 U	2-Hexanone			5.51	U
Chloromethane			2.21 U	Tetrachloroeth	ene		2,21	U
Vinyl chloride	······································		2,21 U	1,3-Dichlorop	ropane		2.21	U
Bromomethane			2.21 U J	Dibromochlor	omethane		2,21	U
Chloroethane			2.21 U	1,2-Dibromoet	thane		2.21	U
Trichlorofluoron	nethane		2.21 U	Chlorobenzen	2		2.21	U
Acatono			63.3 B J	1,1,1,2-Tetrac	hloroethane		2.21	U
1,1-Dichloroethe			2.21 U	Ethylbenzene			2.21	U
Carbon disulfide			3.37	p/m-Xylene			1.32	J
Methylene chlor	ide		5.51 U	o-Xylene			2.21	U
Methyl tert-buty			2.21 U	Styrene			2.21	U
trans-1,2-Dichlo			2.21 U	Bromoform		•	2.21	U
1,1-Dichloroetha				2.21	U			
Vinyl acetate		1411 40000	2.21 U	1,1,2,2-Tetrachloroethane			2.21	U
2-Butanone (ME	EK)		13.3	Bromobenzene			2.21	U
cis-1,2-Dichloro			2.21 U	1,2,3-Trichloropropane			2,21	υ
2,2-Dichloropro			2.21 U	n-Propylbenze	ne		2.21	U
Chloroform			2.21 U	2-Chlorotolue			2.21	U
1,1,1-Trichloroe	thane		2.21 U	1,3,5-Trimethy	lbenzene		2.21	U
1,1-Dichloropro	pene		2.21 U	4-Chlorotolue	ne		2.21	U
Carbon tetrachlo			2.21 U	tert-Butylbenzene		2.21	U	
Benzene			1.26 J	1,2,4-Trimethy	ylbenzene		1.97	J
1,2-Dichloroetha	ane		2.21 U	sec-Butylbenzene			2.21	U
Trichloroethene			2.21 U	1,3-Dichlorobenzene			2.21	U
1,2-Dichloropro	pane		2.21 U	p-Isopropyltoluene			5.51	U
Dibromomethan			2.21 U	1,4-Dichlorobenzene		2.21	U	
Bromodichloron	nethane		2.21 U	n-Butylbenzene		5.51	Ŭ.	
2-Chloroethylvin	nyl ether		5.51 U	1,2-Dichlorobenzene			2.21	U
Methyl isobutyl	ketone (MIBK)		2.21 U	1,2-Dibromo-3-chloropropane			2.21	U
cis-1,3-Dichloro			2.21 U	1,2,4-Trichlorobenzene			2.21	U
Toluene			2.21 U	Hexachlorobutadiene			2.21	U
trans-1,3-Dichlo	ropropene		2.21 U	Naphthalene			21.8	
1,1,2-Trichloroe			2.21 U	1,2,3-Trichlor			2.21	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	93	70-130
1,2-Dichloroethane-d4	88	70-130
Toluene-d8	91	70-130
4-Bromofluorobenzene	88	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

N/A



ARCADIS

Tidewater MGP

SC37-S1 SDG: N/A

Sediment

Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-11

Associated Blank: VS072508B04

Concentration Units: µg/Kg

		1			Sample	Final	Dilution		
Date Collected	Date Received	Date Analyzed	Percent	Solid	Amount (g)	Volume (ml)	Factor	Analy	'st
07/14/08	07/16/08	07/26/08	52.	0	8.03	5	1	BAS	}
Parameter		J	Result		Parameter	J		Result	
Dichlorodifluoro	methane		2.39 U	JJ	2-Hexanone			5.98	บป
Chloromethane			2.39 L	J	Tetrachloroeth	nene		2.39	U
Vinyl chloride			2.39 L	J	1,3-Dichlorop	ropane		2.39	U
Bromomethane	Jure-		2.39 L	J	Dibromochlor	omethane		2.39	U
Chloroethane			2.39 L	J	1,2-Dibromoe	thane		2.39	U
Trichlorofluoror	nethane		2.39 U	j 🚽	Chlorobenzen	e		2.39	U
Acetone			52.9 E	805	1,1,1,2-Tetrac	hloroethane		2.39	U
1,1-Dichloroethe	ene		2.39 L	JJ	Ethylbenzene			2.39	U
Carbon disulfide			14.6	1	p/m-Xylene			4.79	U
Methylene chlor			5.98 C	J	o-Xylene			2.39	U
Methyl tert-buty			2.39 L	J	Styrene			2.39	U
trans-1,2-Dichlo		1	2.39 U	J	Bromoform	and a second second second second		2,39	U
1,1-Dichloroetha	A cally all provide a second and a second and a second as		2.39 L	J	Isopropylbenzene			3.22	
Vinyl acetate	, , , , , , , , , , , , , , , , ,	, , , , , , , , , , , , , , , , , , ,	2.39 L	J	1,1,2,2-Tetrac	hloroethane		2,39	U
2-Butanone (ME	EK)		11.8		Bromobenzen	e		2.39	U
cis-1,2-Dichloro			2.39 L	J	1,2,3-Trichlor	opropane		2.39	U
2,2-Dichloropro			2.39 U	J	n-Propylbenze			2.39	U
Chloroform			2.39 T	J	2-Chlorotolue	ne		2.39	U
1,1,1-Trichloroe	thane	,, ,	2.39 T	J	1,3,5-Trimeth	ylbenzene		2.39	U
1,1-Dichloropro			2.39 U	J	4-Chlorotolue	ne		2.39	U
Carbon tetrachic			2.39 U	J	tert-Butylbenz	zene		2.39	U
Benzene			2.39 U	J	1,2,4-Trimeth			2.39	υ
1,2-Dichloroetha	ane		2.39 1	J	sec-Butylbenz	ene		2.39	U
Trichloroethene			2.39 U	J	1,3-Dichlorob	enzene		2.39	U
1,2-Dichloropro			2.39 (J	p-Isopropylto	luene		5.98	U
Dibromomethan			2.39 I	J	1,4-Dichlorob			2.39	U
Bromodichloror	• • • • • • • • • • • • • • • • • • •		2.39 T	J	n-Butylbenzer	ne		5,98	Ŭ
2-Chloroethylvi			5.98 T	J	1,2-Dichlorob		2.39	U	
Methyl isobutyl			2.39 1	J	1,2-Dibromo-	3-chloropropane		2.39	U
cis-1,3-Dichloro			2.39	υT	1,2,4-Trichlor	obenzene		2.39	U
Toluene			2.39 T	υT	Hexachlorobu	itadiene		2.39	U
trans-1,3-Dichlo	ropropene		2.39 T	J	Naphthalene			1.24	J
1,1,2-Trichloroe			2.39		1,2,3-Trichlor	obenzene		2.39	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	86	70-130
1.2-Dichloroethane-d4	76	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	85	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

N/A



ARCADIS Tidewater MGP

Trip Blank N/A SDG:

Sediment

Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-12

Associated Blank: VS072508B04

× · · · · ·

07/02/08 07/16/08 07/25/08 100 5.00 5 1 BAS Parameter Result Parameter Result Parameter Result Dichlorodifluoromethane 2.00 U 2-Hexanonc 5.00 U Choromethane 2.00 U 1,3-Dichloropropane 2.00 U Sromomethane 2.00 U 1,3-Dichloropropane 2.00 U Sromomethane 2.00 U 1,3-Dichloropropane 2.00 U Chorobethane 2.00 U 1,2-Dichoropropane 2.00 U Cyctone 10.7 B 1,1,1,2-Tetrachloroethane 2.00 U Actoro disulfide 2.00 U p/m-Xylene 4.00 U Carbon disulfide 2.00 U p/m-Xylene 2.00 U Intristice 2.00 U Bromoform 2.00 U Intristice 2.00 U Bromoform 2.00 U 1,2,3-Trichloroptopane<	Date Collected	Date Received	Date Analyzed	Percent Sol	d Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
Name2.00U \checkmark 2-Hexanonc5.00UChloromethane2.00UTetrachloroethene2.00UAinorothane2.00U1,3-Dichloropropane2.00UBromomethane2.00U1,2-Dibromochloromethane2.00UChloroethane2.00U1,2-Dibromochloromethane2.00UChloroethane2.00U1,2-Dibromochloromethane2.00UAcetone10.7B'1,1,1,2-Tetrachloroethane2.00UCarbon disulfide2.00Up/m-Xylene2.00UCarbon disulfide2.00Up/m-Xylene2.00UAethylere chloride5.00Uo-Xylene2.00UMethyl etr-thyl ether (MTBE)2.00UStyrene2.00UI,1-Dichloroethane2.00UIspropylbenzene2.00UJ,2-Dichloroethane2.00UIspropylbenzene2.00UJ,2-Dichloroethane2.00UIspropylbenzene2.00UJ,2-Dichloroethane2.00U1,1,2,2-Tetrachloroethane2.00UJ,2-Dichloropropane2.00UIspropylbenzene2.00UJ,2-Dichloropropane2.00Un-Propylbenzene2.00UJ,2-Dichloroethane2.00Un-Propylbenzene2.00UJ,1-Dichloropropane2.00Un-Propylbenzene2.00UJ,1-Dichloroethane <t< td=""><td></td><td>1</td><td></td><td>100</td><td></td><td></td><td></td><td></td></t<>		1		100				
Chloromethane 2.00 U Tetrachloroethene 2.00 U Vinyl chloride 2.00 U 1,3-Dichloropropane 2.00 U Bromomethane 2.00 U Dibromochloromethane 2.00 U Chloroethane 2.00 U Chloroethane 2.00 U Acetone 10.7 B' 1,1,1,2-Tetrachloroethane 2.00 U Acetone 10.7 B' 1,1,1,2-Tetrachloroethane 2.00 U Chlorobenzene 2.00 U Ethylbenzene 2.00 U Acetone 1.07 B' 1,1,1,2-Tetrachloroethane 2.00 U Chlorobenzene 2.00 U Styrene 2.00 U Methyl ter-thoryl ether (MTBE) 2.00 U Styrene 2.00 U I,1-Dichloroethane 2.00 U Ispropylbenzene 2.00 U I,1-Dichloroethane 2.00 U Ispropylbenzene 2.00 U I,1-Dichlorop	Parameter		1	Result	Parameter			Result
Vinyl chloride 2.00 U 1,3-Dichloropropane 2.00 U Aromomethane 2.00 U Dibromochloromethane 2.00 U Chloroethane 2.00 U 1,2-Dibromochloromethane 2.00 U Chlorobethane 2.00 U Chlorobenzene 2.00 U Acetone 10.7 B' 1,1,1,2-Tetrachloroethane 2.00 U Carbon disulfide 2.00 U pm-Xylene 4.00 U Methylen chloride 5.00 U o-Xylene 2.00 U Methyl tert-butyl ether (MTBE) 2.00 U Styrene 2.00 U JDichloroethane 2.00 U Isopropylbenzene 2.00 U JDichloroethane 2.00 U Isopropylbenzene 2.00 U JDichloroethane 2.00 U 1,2,3-Trichloropthane 2.00 U JDichloroethane 2.00 U 1,2,3-Trichloroptopane 2.00 U	Dichlorodifluoro	omethane		2.00 U	2-Hexanone			
Bromomethane 2.00 U Dibromochloromethane 2.00 U Chloroethane 2.00 U 1,2-Dibromochlane 2.00 U Chloroothane 2.00 U 1,2-Dibromochlane 2.00 U Acotone 10.7 B 1,1,1.2-Tetrachloroothane 2.00 U Acotone 10.7 B 1,1,1.2-Tetrachloroothane 2.00 U Acotone 2.00 U Ethylenzene 2.00 U Arbon disulfide 2.00 U p/m-Xylene 4.00 U Arbon disulfide 2.00 U Styrene 2.00 U Methyl ent-butyl ether (MTBE) 2.00 U Bromoform 2.00 U 1.1-Dichloroothane 2.00 U Isopropylbenzene 2.00 U 2.1-Dichloroothane 2.00 U Isopropylbenzene 2.00 U 2.2-Dichloroothane 2.00 U 1,2.3-Trichloroothane 2.00 U 1,1-Tichlorooth	Chloromethane			2.00 U	Tetrachloroe	thene		
Chloroethane2.00U1.2-Dibromoethane2.00UChlorofluoromethane2.00UChlorobenzene2.00UAcetone10.7B1,1,1,2-Tetrachloroethane2.00U1.1-Dichloroethene2.00UEthylbenzene2.00UCarbon disulfide2.00Up/m-Xylene4.00UMethylene chloride5.00Uo-Xylene2.00UMethylene chloride2.00UBromoform2.00UMethyl tert-butyl ether (MTBE)2.00UBromoform2.00U1.1-Dichloroethane2.00UBromoform2.00U1.1-Dichloroethane2.00U1,1,2,2-Tetrachloroethane2.00U2.1-Dichloroethene2.00U1,1,2,2-Tetrachloroethane2.00U2.2-Dichloroethene2.00U1,1,2,2-Tetrachloroethane2.00U2.2-Dichloroethene2.00U1,2,3-Trichloroethane2.00U2.2-Dichloroethane2.00U1,2,3-Trinethylbenzene2.00U1,1-Trichloroethane2.00U1,2,4-Trimethylbenzene2.00U1,1-Trichloroethane2.00U1,2,4-Trimethylbenzene2.00U1,1-Trichloroethane2.00U1,2,4-Trimethylbenzene2.00U1,1-Trichloroethane2.00U1,2,4-Trimethylbenzene2.00U1,1-Dichloroethane2.00U1,2,4-Trimethylbenzene <td>Vinyl chloride</td> <td></td> <td></td> <td>2.00 U</td> <td></td> <td></td> <td></td> <td><u>2.00 U</u></td>	Vinyl chloride			2.00 U				<u>2.00 U</u>
$\begin{array}{c c c c c c c c c c c c c c c c c c c $	Bromomethane				Dibromochl	oromethane		
Acctone 10.7 B 1,1,1,2-Tetrachloroethane 2,00 U 1,1-Dichloroethene 2.00 U Ethylbenzene 2.00 U Acton disulfide 2.00 U p/m-Xylene 4.00 U Methyl tert-butyl ether (MTBE) 2.00 U Styrene 2.00 U Arbon disulfide 2.00 U Styrene 2.00 U Arbon disulfide 2.00 U Bromoform 2.00 U Arbon disulfide 2.00 U Ispropylbenzene 2.00 U Isprichloroptopane	Chloroethane							
1.1-Dichloroethene2.00UEthylbenzene2.00UCarbon disulfide2.00U $p'm$ -Xylene4.00UMethylenc chloride5.00U o -Xylene2.00UMethyl tert-butyl ether (MTBE)2.00UStyrene2.00UJans-1,2-Dichloroethene2.00UBromoform2.00U1,1-Dichloroethene2.00UIsopropylbenzene2.00U1,1-Dichloroethene2.00UIsopropylbenzene2.00U2-Dichloroethene2.00U1,1,2,2-Tetrachloroethane2.00U2-Butanone (MEK)2.00UBromoform2.00U2-Dichloropropane2.00U1,2,3-Trichloropropane2.00U2,2-Dichloropropane2.00U1,3,5-Trimethylbenzene2.00U2,1-Dichloropropane2.00U1,3,5-Trimethylbenzene2.00U1,1-Trichloropenene2.00U4-Chlorotoluene2.00U2,2-Dichloropropane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloropropane2.00U1,2,4-Trimethylbenzene2.00U2,2-Dichloropropane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloropropane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloropropane2.00U1,4-Dichlorobenzene2.00U1,2-Dichloropropane2.00U1,2-Dichlorobenzen	Trichlorofluoror	nethane		2.00 U				
Larbon disulfide2.00 Up/m-Xylene4.00 UMethylene chloride5.00 Uo-Xylene2.00 UMethyl tert-butyl ether (MTBE)2.00 UStyrene2.00 UIrans-1,2-Dichloroethene2.00 UBromoform2.00 UJDichloroethane2.00 UIsopropylbenzene2.00 UVinyl acetate2.00 U1,1,2,2-Tetrachloroethane2.00 U2-Dichloroethene2.00 U1,1,2,2-Tetrachloroethane2.00 U2-Dichloroethene2.00 U1,2,3-Trichloropropane2.00 U2-Dichloropropane2.00 U1,2,3-Trichloropropane2.00 U2.1-Dichloroethane2.00 U1,3,5-Trimethylbenzene2.00 U1,1-Dichloropropane2.00 U1,3,5-Trimethylbenzene2.00 U1,1-Dichloropropane2.00 U1,3,5-Trimethylbenzene2.00 U1,1-Dichloropropane2.00 U1,2,4-Trimethylbenzene2.00 U1,1-Dichloropropane2.00 U1,2,4-Trimethylbenzene2.00 U1,1-Dichloropropane2.00 U1,2,4-Trimethylbenzene2.00 U1,2-Dichloroethane2.00 U1,3-Dichlorobenzene2.00 U2,2-Dichloroethane2.00 U1,2,2-Tetraethylbenzene2.00 U1,2-Dichloropropane2.00 U1,2,4-Trimethylbenzene2.00 U2,2-Dichloroethane2.00 U1,2-Dichlorobenzene2.00 U2,2-Dichloroethane2.00 U1,2-Dichlorobenzene2.00 U2,2-Dichloroethane2.00 U1,2-Dichlorobenzene2.00 U2,2-Dichloroethane2.00 U1,2-	Acetone			10.7 B				
Methylene chloride 5.00 Uo-Xylene 2.00 UMethyl tert-butyl ether (MTBE) 2.00 UStyrene 2.00 Urans-1,2-Dichloroethene 2.00 UBromoform 2.00 U $1,1-Dichloroethane$ 2.00 UBromoform 2.00 U $1,1-Dichloroethane$ 2.00 UIsopropylbenzene 2.00 U 2.01 u $1,1,2,2$ -Tetrachloroethane 2.00 U 2.02 -Dichloroethene 2.00 U $1,2,3$ -Trichloropropane 2.00 U 2.2 -Dichloroethene 2.00 U $1,2,3$ -Trichloropropane 2.00 U 2.2 -Dichloroethane 2.00 U $1,2,3$ -Trichloropropane 2.00 U 2.2 -Dichloropropane 2.00 U $1,2,3$ -Trimethylbenzene 2.00 U $1,1$ -Trichloropropene 2.00 U $1,3,5$ -Trimethylbenzene 2.00 U $1,1$ -Dichloropropene 2.00 U $1,2,4$ -Trimethylbenzene 2.00 U $2.arbon tetrachloride2.00U1,2,4-Trimethylbenzene2.00U2.2-Dichloroethane2.00U1,2,4-Trimethylbenzene2.00U2.2-Dichloroethane2.00U1,2,4-Trimethylbenzene2.00U2.2-Dichloroethane2.00U1,2,4-Trimethylbenzene2.00U2.2-Dichloroethane2.00U1,2,4-Trimethylbenzene2.00U2.2-Dichloroethane2.00U1,2,$	1,1-Dichloroethe	ene		2.00 U	Ethylbenzen	e		2.00 U
Methyl tert-butyl ether (MTBE)2.00UStyrene2.00Urans-1,2-Dichloroethene2.00UBromoform2.00U1,1-Dichloroethane2.00UIsopropylbenzene2.00UVinyl acetate2.00U1,1,2,2-Tetrachloroethane2.00U2-Butanone (MEK)2.00UBromobenzene2.00U2-Dichloroptopane2.00U1,2,3-Trichloroptopane2.00U2,2-Dichloroptopane2.00U1,2,3-Trichloroptopane2.00U1,1-Dichloroptopane2.00U1,3,5-Trimethylbenzene2.00U1,1-Dichloroptopene2.00U1,3,5-Trimethylbenzene2.00U1,1-Dichloroptopene2.00U1,2,4-Trimethylbenzene2.00U2,2-Dichloroptopene2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloroptopane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,2-Dichlorobenzene2.00U1,2-Dichloroethane2.00 <td< td=""><td>Carbon disulfide</td><td>9</td><td></td><td>2.00 U</td><td>p/m-Xylene</td><td></td><td></td><td></td></td<>	Carbon disulfide	9		2.00 U	p/m-Xylene			
rans-1,2-Dichloroethene2.00UBromoform2.00U1,1-Dichloroethane2.00UIsopropylbenzene2.00U1,1,2,2-Tetrachloroethane2.00U1,1,2,2-Tetrachloroethane2.00U1-Butanone (MEK)2.00UBromobenzene2.00U1,2-Dichloroptopane2.00U1,2,3-Trichloroptopane2.00U2,2-Dichloroptopane2.00Un-Propylbenzene2.00U1,1-Trichloroethane2.00U1,3,5-Trimethylbenzene2.00U1,1-Dichloroptopane2.00U1,3,5-Trimethylbenzene2.00U1,1-Dichloroptopene2.00U1,3,5-Trimethylbenzene2.00U2,2-Dichloroptopene2.00U4-Chlorotoluene2.00U1,1-Dichloroptopene2.00U1,3,5-Trimethylbenzene2.00U2,2-Dichloroptopene2.00U1,2,4-Trimethylbenzene2.00U2,2-Dichloroptopene2.00U1,2,4-Trimethylbenzene2.00U2,2-Dichloroptopane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloroptopane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloroptopane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloroptopane2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloroptopane2.00U1,2,0UU1,2-Dichloroptopene2.00 </td <td>Methylene chlor</td> <td>ride</td> <td></td> <td>5.00 U</td> <td>o-Xylene</td> <td></td> <td></td> <td>2.00 U</td>	Methylene chlor	ride		5.00 U	o-Xylene			2.00 U
1.1-Dichloroethane2.00UIsopropylbenzene2.00UVinyl acetate2.00U $1,1,2,2$ -Tetrachloroethane2.00U2-Butanone (MEK)2.00UBromobenzene2.00U1.2,2-Dichloroethene2.00U $1,2,3$ -Trichloropropane2.00U2,2-Dichloroptropane2.00U n -Propylbenzene2.00U2,2-Dichloroptropane2.00U n -Propylbenzene2.00U1,1-Trichloroethane2.00U $1,3,5$ -Trimethylbenzene2.00U1,1-Dichloroptropene2.00U 4 -Chlorotoluene2.00U2arbon tetrachloride2.00U 4 -Chlorotoluene2.00U2arbon tetrachloride2.00U $1,2,4$ -Trimethylbenzene2.00U3enzene2.00U $1,2,4$ -Trimethylbenzene2.00U1,2-Dichloroethane2.00U $1,2,4$ -Trimethylbenzene2.00U1,2-Dichloroptropane2.00U $1,2,4$ -Trimethylbenzene2.00U1,2-Dichloroptropane2.00U $1,2-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,4-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,4-Dichlorobenzene2.00U1,2-Dichloroethane2.00U1,2-Dichlorobenzene2.00U2-Dichloroethane2.00U1,2-Dichlorobenzene2.00U2-Dichloroethane$	Methyl tert-buty	l ether (MTBE)		2.00 U	Styrene			<u>2.00 U</u>
Vinyl acetate2.00U $1,1,2,2$ -Tetrachloroethane2.00U2-Butanone (MEK)2.00UBromobenzene2.00U2-Dichloroethene2.00U $1,2,3$ -Trichloropropane2.00U2,2-Dichloropropane2.00U $1,2,3$ -Trichloropropane2.00U2,2-Dichloropropane2.00U $1,2,3$ -Trichloropropane2.00U2,2-Dichloropropane2.00U 2 -Chlorotoluene2.00U1,1-Trichloroethane2.00U $1,3,5$ -Trimethylbenzene2.00U1,1-Dichloropropene2.00U 4 -Chlorotoluene2.00U2arbon tetrachloride2.00U 4 -Chlorotoluene2.00U2arbon tetrachloride2.00U $1,2,4$ -Trimethylbenzene2.00U3enzene2.00U $1,2,4$ -Trimethylbenzene2.00U1,2-Dichloroethane2.00U $1,3,5$ -Dichlorobenzene2.00U1,2-Dichloropropane2.00U $1,3,5$ -Dichlorobenzene2.00U1,2-Dichloropropane2.00U $1,4,5,4$ -Trimethylbenzene2.00U1,2-Dichloropropane2.00U $1,4,5,4$ -Trimethylbenzene2.00U1,2-Dichloropropane2.00U $1,4,5,4$ -Trimethylbenzene2.00U1,2-Dichloropropane2.00U $1,4,5,4$ -Trimethylbenzene2.00U2-Chloroethylvinyl ether5.00U $1,2,2,5,6,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0$	trans-1,2-Dichlo	oroethene		2.00 U	Bromoform			2.00 U
Vinyl acetate2.00U $1,1,2,2$ -Tetrachloroethane2.00U2-Butanone (MEK)2.00UBromobenzene2.00U 2 -Dichloroethene2.00U $1,2,3$ -Trichloropropane2.00U $2,2$ -Dichloropropane2.00U $1,2,3$ -Trichloropropane2.00U $2,2$ -Dichloropropane2.00U $1,2,3$ -Trichloropropane2.00U $2,2$ -Dichloropropane2.00U $1,2,3$ -Trichloropropane2.00U $2,2$ -Dichloropropane2.00U 2 -Chlorotoluene2.00U $1,1$ -Trichloroethane2.00U $1,3,5$ -Trimethylbenzene2.00U $1,1$ -Dichloropropene2.00U 4 -Chlorotoluene2.00UCarbon tetrachloride2.00U $1,2,4$ -Trimethylbenzene2.00UBenzene2.00U $1,2,4$ -Trimethylbenzene2.00U $1,2$ -Dichloroethane2.00U $1,3,4$ -Trimethylbenzene2.00U $1,2$ -Dichloropropane2.00U $1,3,4$ -Trimethylbenzene2.00U $1,2$ -Dichloropropane2.00U $1,3,4$ -Trimethylbenzene2.00U $1,2$ -Dichloropropane2.00U $1,4$ -Dichlorobenzene2.00U $1,2$ -Dichloropropane2.00U $1,4$ -Dichlorobenzene2.00U $2,2$ -Chloroethylvinyl ether5.00U $1,2$ -Dichlorobenzene2.00U $2,2$ -Chloropropene2.00U $1,2$ -Dichlorobenzene <td< td=""><td>1,1-Dichloroetha</td><td>ane</td><td></td><td>2.00 U</td><td>Isopropylbe</td><td>nzene</td><td></td><td>2.00 U</td></td<>	1,1-Dichloroetha	ane		2.00 U	Isopropylbe	nzene		2.00 U
is-1,2-Dichloroethene2.00U1,2,3-Trichloropropane2.00U $2,2$ -Dichloropropane2.00Un-Propylbenzene2.00UChloroform2.00U2-Chlorotoluene2.00U $1,1,1$ -Trichloroethane2.00U1,3,5-Trimethylbenzene2.00U $1,1$ -Dichloropropene2.00U4-Chlorotoluene2.00UCarbon tetrachloride2.00U4-Chlorotoluene2.00UCarbon tetrachloride2.00U1,2,4-Trimethylbenzene2.00U3enzene2.00U1,2,4-Trimethylbenzene2.00U $1,2$ -Dichloroethane2.00U1,3-Dichlorobenzene2.00U $1,2$ -Dichloropropane2.00U1,3-Dichlorobenzene2.00U $1,2$ -Dichloropropane2.00U1,4-Dichlorobenzene2.00U $1,2$ -Dichloromethane2.00U1,2-Dichlorobenzene2.00U $1,2$ -Dichloropropane2.00U1,2-Dichlorobenzene2.00U $1,2$ -Dichloroethane2.00U1,2-Dichlorobenzene2.00U 2.00 U1,2-Dichlorobenzene2.00UU $1,2$ -Dichloropropene2.00U1,2-Dichlorobenzene2.00U 2.00 U1,2-Dichlorobenzene2.00UU 2.00 U1,2-Dichlorobenzene2.00U 2.00 U1,2-Dibromo-3-chloropropane2.00U<	Vinyl acetate			2.00 U	1,1,2,2-Tetr	achloroethane		2.00 U
2,2-Dichloropropane 2.00 Un-Propylbenzene 2.00 UChloroform 2.00 U 2 -Chlorotoluene 2.00 U $1,1$ -Trichloroethane 2.00 U $1,3,5$ -Trimethylbenzene 2.00 U $1,1$ -Dichloropropene 2.00 U 4 -Chlorotoluene 2.00 UCarbon tetrachloride 2.00 U 4 -Chlorotoluene 2.00 UCarbon tetrachloride 2.00 U $1,2,4$ -Trimethylbenzene 2.00 UBenzene 2.00 U $1,2,4$ -Trimethylbenzene 2.00 U $1,2$ -Dichloroethane 2.00 U $1,3$ -Dichlorobenzene 2.00 U $1,2$ -Dichloropropane 2.00 U $1,3$ -Dichlorobenzene 2.00 U $1,2$ -Dichloropropane 2.00 U $1,4$ -Dichlorobenzene 2.00 U $1,2$ -Dichloromethane 2.00 U $1,4$ -Dichlorobenzene 2.00 U 2.00 U $1,4$ -Dichlorobenzene 2.00 U 2.00 U $1,2$ -Dichlorobenzene 2.00 U 2.00 U $1,2,4$ -Trichlorobenzene 2.00 U 2.00 U $1,2,4$ -Trichloro	2-Butanone (MI	EK)		2.00 U	Bromobenze	ene		2.00 U
Chloroform2.00U2-Chlorotoluene2.00U $1,1$ -Trichloroethane2.00U $1,3,5$ -Trimethylbenzene2.00U $1,1$ -Dichloropropene2.00U 4 -Chlorotoluene2.00UCarbon tetrachloride2.00U 4 -Chlorotoluene2.00UBenzene2.00U $1,2,4$ -Trimethylbenzene2.00U $1,2$ -Dichloroethane2.00U $1,2,4$ -Trimethylbenzene2.00U $1,2$ -Dichloroethane2.00U $1,3$ -Dichlorobenzene2.00U $1,2$ -Dichloropropane2.00U $1,3$ -Dichlorobenzene2.00U $1,2$ -Dichloromethane2.00U $1,4$ -Dichlorobenzene2.00UDibromomethane2.00U $1,4$ -Dichlorobenzene2.00U 2 -Chloroethylvinyl ether5.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloroethylvinyl ether5.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloroethylvinyl ether2.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloroethylvinyl ether2.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloropropene2.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloroethylvinyl ether2.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloropropene2.00U $1,2$ -Dichlorobenzene2.00U 2 -Chloropropene2.00U $1,2$ -Dichlorobenzene2.00U <td>cis-1,2-Dichloro</td> <td>ethene</td> <td></td> <td>2.00 U</td> <td>1,2,3-Trichl</td> <td>oropropane</td> <td></td> <td>2.00 U</td>	cis-1,2-Dichloro	ethene		2.00 U	1,2,3-Trichl	oropropane		2.00 U
1,1-Trichloroethane2.00U1,3,5-Trimethylbenzene2.00U1,1-Dichloropropene2.00U4-Chlorotoluene2.00UCarbon tetrachloride2.00U4-Chlorotoluene2.00UBenzene2.00U1,2,4-Trimethylbenzene2.00U1,2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloropropane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloropropane2.00U1,3-Dichlorobenzene2.00U1,2-Dichloromethane2.00U1,4-Dichlorobenzene2.00U2.01U1,4-Dichlorobenzene2.00U2.02U1,2-Dichlorobenzene2.00U1,2-Dichloropropane2.00U1,2-Dichlorobenzene2.00U2.01U1,2-Dichlorobenzene2.00U2.02U1,2-Dichlorobenzene2.00U2.03U1,2-Dichlorobenzene2.00U2.04U1,2-Dichlorobenzene2.00U2.05U1,2-Dichlorobenzene2.00U2.06U1,2-Dichlorobenzene2.00U2.07U1,2-Dichlorobenzene2.00U2.08U1,2-Dichlorobenzene2.00U2.09U1,2-Dichlorobenzene2.00U2.00U1,2-A-Trichlorobenzene2.00U2.08U1,2-Dichlorobenzene <td>2,2-Dichloropro</td> <td>pane</td> <td></td> <td>2.00 U</td> <td>n-Propylben</td> <td>zene</td> <td>_</td> <td>2.00 U</td>	2,2-Dichloropro	pane		2.00 U	n-Propylben	zene	_	2.00 U
1.1-Dichloropropene2.00U4-Chlorotoluene2.00UCarbon tetrachloride2.00Utert-Butylbenzene2.00UBenzene2.00U1,2,4-Trimethylbenzene2.00U3.2-Dichloroethane2.00Usec-Butylbenzene2.00U1.2-Dichloroethane2.00U1,3-Dichlorobenzene2.00U1.2-Dichloropropane2.00U1,3-Dichlorobenzene2.00U1.2-Dichloropropane2.00U1,4-Dichlorobenzene2.00UDibromomethane2.00U1,4-Dichlorobenzene2.00UBromodichloromethane2.00U1,2-Dichlorobenzene2.00U2-Chloroethylvinyl ether5.00U1,2-Dichlorobenzene2.00U1.2-Dichloropropane2.00U1,2-Dichlorobenzene2.00U1.3-Dichloropenethane2.00U1,2-Dichlorobenzene2.00U2-Chloroethylvinyl ether5.00U1,2-Dichlorobenzene2.00U1.3-Dichloropropene2.00U1,2-Dibromo-3-chloropropane2.00U1.3-Dichloropropene2.00U1,2,4-Trichlorobenzene2.00U1.3-Dichloropropene2.00U1,2,4-Trichlorobenzene2.00U1.3-Dichloropropene2.00U1,2,4-Trichlorobenzene2.00U1.3-Dichloropropene2.00U1,2,4-Trichlorobenzene2.00U1.3-Dichloropropene <td>Chloroform</td> <td>-</td> <td></td> <td>2.00 U</td> <td>2-Chlorotol</td> <td>uene</td> <td></td> <td>2.00 U</td>	Chloroform	-		2.00 U	2-Chlorotol	uene		2.00 U
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	Toluene			2.00 U	Hexachloro	butadiene		2.00 U
	trans-1,3-Dichlo	propropene		2.00 U	Naphthalene)		5.00 U
				2.00 U 🕔	1,2,3-Trichl	orobenzene		2.00 U 🗸

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	107	70-130
1,2-Dichloroethane-d4	113	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	81	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

Concentration Units: µg/Kg

Date Collected Received AnalyzedNotinal Motume (m) Notume (m)Amount (g) Notume (ml) Volume (ml) Volume (ml)Factor FactorAnalyst BAS Farameter Result Parameter Result Parameter ResultDichlorodifluoromethane100U 2 2-Hexanone250UChoromethane100U 2 2-Hexanone100UBromomethane100U1,3-Dichloropropane100UBromomethane167BDibromochloromethane100UTrichlorofluoromethane100U1,2-Dibromochane100UTrichlorofluoromethane100UChlorobenzzne100UAcetone244JR1,1,2-Tetrachloroethane100UI.1-Dichloroethene100UEthylenzene100UI.1-Dichloroethene100UStyrene200UAcetone244JR1,1,2-Tetrachloroethane100UI.1-Dichloroethene100UStyrene100UI.1-Dichloroethene100UStyrene100UI.1-Dichloroethane100UStyrene100UI.1-Dichloroethane100UStyrene100UI.1-Dichloroethane100UStyrene100UI.1-Dichloroethane100UStyrene100UI.1-Dichloroethane100UIsoropylbenzene100<		PHA	Client: Project: Client ID: Case: Matrix:	ARCADI Tidewate Trip Blank N/A Sediment	er MGP	N/A		ETR: Lab II Associ	ntration Unit	-12E VS072808B14
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		hloroproper	A						.,	
			<u>.</u>		100 U					

		Acceptance
Surrogate	% Recovery	Range (%)
2-Bromo-1-chloropropane	99	70-130
1-Chloro-2-fluorobenzene	108	70-130
1,4-Dichlorobutane	98	70-130
Dibromofluoromethane	97	70-130
Toluene-d8	94	70-130
4-Bromofluorobenzene	96	70-130

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N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.



ARCADIS

Tidewater MGP SC1-S1

SDG: N/A

Matrix: Sediment

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-13 Associated Blank: VS072608B03

Concentration Units: µg/Kg

					,		<u>, , , , , , , , , , , , , , , , , , , </u>	
				Sample	Final	Dilution		
Date Collected	Date Received	Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/15/08	07/16/08	07/26/08	82.7	6.97	5	1	BAS	
Parameter		I	Result	Parameter			Result	
Dichlorodifluoro	methane		1.73 U	2-Hexanone		·····	4.34	
Chloromethane			1.73 U	Tetrachloroeth	iene		1.73	
Vinyl chloride			1.73 U	1,3-Dichlorop	ropane		1.73	
Bromomethane			1.73 U J	Dibromochlor	omethane		1.73	
Chloroethane		1	1.73 U	1,2-Dibromoe	thane		1.73	
Trichlorofluoror	nethane		1.73 U	Chlorobenzen	e			Ŭ
Acetone			23.7 BU	1,1,1,2-Tetrac	hloroethane			U
1,1-Dichloroeth	ene		1.73 U	Ethylbenzene				Ŭ
Carbon disulfide			7.58	p/m-Xylene				U
Methylene chlor			4.34 U	o-Xylene				U
Methyl tert-buty		2.10 ⁻¹⁷	1.73 U	Styrene			1.73	
trans-1,2-Dichlo			1.73 U	Bromoform			and the second se	U
1,1-Dichloroeth			1.73 U	Isopropylbenz	ene	<u></u>	1.73	
Vinyl acetate			1.73 U J	1,1,2,2-Tetrac	hloroethane			U
2-Butanone (MI	EK)		2.26	Bromobenzen	e			U
cis-1,2-Dichlore			1.73 U	1,2,3-Trichloropropane				U
2,2-Dichloropro			1.73 U	n-Propylbenzene				U
Chloroform	<u>P4117</u>		1.73 U	2-Chlorotolue	ne		1.73	
1,1,1-Trichloroe	ethane		1.73 U	1,3,5-Trimeth	ylbenzene		1.73	
1,1-Dichloropro			1.73 U	4-Chlorotolue	ne		1.73	
Carbon tetrachle			1.73 U	tert-Butylben:	zene		1.73	
Benzene			1.73 U	1,2,4-Trimeth	ylbenzene		1.73	
1,2-Dichloroeth	ane		1.73 U	sec-Butylben:	zene		1.73	
Trichloroethene			1.73 U	1,3-Dichlorot	enzene		1.73	
1,2-Dichloropro			1.73 U	p-Isopropylto	luene		4.34	
Dibromometha			1.73 U	1,4-Dichlorob	oenzene		1.36	
Bromodichloro			1.73 U	n-Butylbenze	ne		4.34	
2-Chloroethylv			4.34 U	1,2-Dichlorol	oenzene	an	1.73	
	l ketone (MIBK)		1.73 U	1,2-Dibromo-	-3-chloropropane		1.73	
	opropene		1.73 U	1,2,4-Trichlo	robenzene		1.73	
			0.87 J	Hexachlorob	utadiene		1.73	
trans-1 3-Dichl	oropropene		1.73 U	Naphthalene			4.34	
1,1,2-Trichloro			1.73 U	1,2,3-Trichlo	robenzene		1.73	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	93	70-130
1,2-Dichloroethane-d4	93	70-130
Toluene-d8	92	70-130
4-Bromofluorobenzene	85	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.



ARCADIS

Tidewater MGP SC2-S1

SDG: N/A

Matrix: Sediment

N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-14

Associated Blank: VS072608B03 Concentration Units: µg/Kg

	D . D . 1	D	D	Sample	Final Volume (ml)	Dilution Factor	Analyst
	Date Received	Date Analyzed	Percent Solid	Amount (g)	· · · · · · · · · · · · · · · · · · ·		
07/15/08	07/16/08	07/26/08	83.0	7.40	5	1	BAS
Parameter		F	Result	Parameter			Result
Dichlorodifluoro	omethane		1.63 U	2-Hexanone			4.07 U
Chloromethane			<u>1.63 U</u>	Tetrachloroeth			1.63 U
Vinyl chloride			<u>1.63 U</u>	1,3-Dichlorop			1.63 U
Bromomethane			1.63 U J	Dibromochlor			1.63 U
Chloroethane			1.63 U	1,2-Dibromoe	· · · · · · · · · · · · · · · · · · ·		<u>1.63 U</u>
Trichlorofluoror	nethane		1.63 U	Chlorobenzen			1.63 U
Acetone			<u>12.4 BU</u>	1,1,1,2-Tetrac	hloroethane		1.63 U
1,1-Dichloroethe	ene		1.63 U	Ethylbenzene			<u>1.63 U</u>
Carbon disulfide	2		5.69	p/m-Xylene			3.26 U
Methylene chlor	ride		4.07 U	o-Xylene			<u>1.63 U</u>
Methyl tert-buty	l ether (MTBE)		1.63 U	Styrene			<u> </u>
trans-1,2-Dichlo	roethene		1.63 U	Bromoform			<u>1.63 U</u>
1,1-Dichloroeth	ane		1.63 U	Isopropylbenz	ene		<u>1.63 U</u>
Vinyl acetate			1.63 U	1,1,2,2-Tetrac	hloroethane		<u>1.63 U</u>
2-Butanone (MI			1.63 U	Bromobenzen	e		<u>1.63 U</u>
cis-1,2-Dichloro	ethene		1.63 U	1,2,3-Trichlor	opropane		<u>1.63 U</u>
2,2-Dichloropro	pane		1.63 U	n-Propylbenze	ene		<u>1.63 U</u>
Chloroform			1.63 U	2-Chlorotolue	ne		1.63 U
1,1,1-Trichloroe	thane		1.63 U	1,3,5-Trimeth	ylbenzene		1.63 U
1,1-Dichloropro	pene		1.63 U	4-Chlorotolue	ne		1.63 U
Carbon tetrachlo			1.63 U	tert-Butylbenz	zene		1.63 U
Benzene			1.63 U	1,2,4-Trimeth	ylbenzene		1.63 U
1,2-Dichloroeth	ane		1.63 U	sec-Butylbenz	zene		1.63 U
Trichloroethene			1.63 U	1,3-Dichlorob	enzene		1.63 U
1,2-Dichloropro	pane		1.63 U	p-Isopropylto	luene		<u>4.07</u> U
Dibromomethan			1.63 U	1,4-Dichlorob	enzene		<u>1.63 U</u>
Bromodichloror	nethane		1.63 U	n-Butylbenzer	ne		4.07 U
2-Chloroethylvi	nyl ether		4.07 U	1,2-Dichlorob	enzene		<u>1.63 U</u>
Methyl isobutyl			1.63 U	1,2-Dibromo-	3-chloropropane		1.63 U
cis-1,3-Dichloro			1.63 U	1,2,4-Trichlor	robenzene		1.63 U
Toluene	<u></u>		1.63 U	Hexachlorobu	itadiene		1.63 U
trans-1,3-Dichlo	propropene		1.63 U	Naphthalene			4.07 U
1,1,2-Trichloroe			1.63 U	1,2,3-Trichlor	robenzene		1.63 U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	106	70-130
1,2-Dichloroethane-d4	112	70-130
Toluene-d8	90	70-130
4-Bromofluorobenzene	85	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

N/A



ARCADIS Tidewater MGP

Dup-3 N/A SDG:

Sediment

Matrix:

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-15 Associated Blank: VS072608B03

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Concentration Units: µg/Kg

			D . 0 . 11	Sample	Final	Dilution		
Date Collected		Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/15/08	07/16/08	07/26/08	84.0	7.06	5	1	BAS	S
Parameter		I	Result	Parameter			Result	•
Dichlorodifluoro	methane		1.69 U	2-Hexanone			4.22	** ********
Chloromethane			1.69 U	Tetrachloroeth			1.69	
Vinyl chloride			<u>1.69 U</u>	1,3-Dichloropi	ropane		1.69	**********
Bromomethane			1.69 U J	Dibromochloro			1.69	
Chloroethane			1.69 U	1,2-Dibromoet	thane		1.69	U
Trichlorofluoron	nethane		<u>1.69 U</u>	Chlorobenzene	3		1.69	
Acetone			<u>11,5 B U</u>	1,1,1,2-Tetracl	hloroethane		1.69	
1,1-Dichloroethe	ene		1.69 Ű	Ethylbenzene			1.69	U
Carbon disulfide	;		4.64	p/m-Xylene			3.37	U
Methylene chlor	ide		4.22 U	o-Xylene			1.69	U
Methyl tert-buty	l ether (MTBE)		1.69 U	Styrene			1.69	U
trans-1,2-Dichlo			1,69 U	Bromoform			1.69	U
1,1-Dichloroetha	me		1.69 U	Isopropylbenz	ene		1.69	U
Vinyl acetate			1.69 U	1,1,2,2-Tetracl	hloroethane		1.69	U
2-Butanone (ME	EK)		1.69 U	Bromobenzene	3		1.69	U
	ethene		1,69 U	1,2,3-Trichloro	opropane		1.69	U
2,2-Dichloropro			1.69 U	n-Propylbenzene			1.69	U
Chloroform			1.69 U	2-Chlorotoluer	ne		1.69	U
1,1,1-Trichloroe			1.69 U	1,3,5-Trimethy	lbenzene		1.69	Ŭ
1,1-Dichloropro			1.69 U	4-Chlorotoluer	ne		1.69	U
Carbon tetrachlo	oride		1.69 U	tert-Butylbenz	ene		1.69	U
Benzene			1.69 U	1,2,4-Trimethy	ylbenzene		1.69	U
1,2-Dichloroetha	ane		1.69 U	sec-Butylbenz	ene		1.69	U
Trichloroethene			1.69 U	1,3-Dichlorob	enzene		1.69	U
1,2-Dichloropro	pane		1.69 U	p-Isopropyltol	uene		4.22	U
Dibromomethan		· · · · · · · · · · · · · · · · · · ·	1.69 U	1,4-Dichlorob			1,69	
Bromodichloron			1.69 U	n-Butylbenzen			4.22	
2-Chloroethylvii			4.22 U	1,2-Dichlorob			1.69	
	ketone (MIBK)		1.69 U	1,2-Dibromo-3	3-chloropropane		1.69	U
cis-1,3-Dichloro			1.69 U	1,2,4-Trichlor			1.69	
Toluene	······		1.19 J	Hexachlorobu		<u> </u>	1.69	
trans-1,3-Dichlo	ropropene		1.69 U	Naphthalene			4.22	
1,1,2-Trichloroe			1.69 U	1,2,3-Trichlor	obenzene		1.69	Ū

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	106	70-130
Toluene-d8	91	70-130
4-Bromofluorobenzene	82	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.



ARCADIS

Tidewater MGP SC11-S1

SDG: N/A

Sediment

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-16 Associated Blank: VS072608B03

Concentration Units: µg/Kg

	,			Sample	Final	Dilution	
Date Collected		Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/15/08	07/16/08	07/26/08	74.7	8.39	5	1	BAS
Parameter		I	Result	Parameter			Result
Dichlorodifluoro	omethane		1.60 U	2-Hexanone			3.99 U
Chloromethane		4.64 F 10 ⁻¹⁰	<u>1.60 U</u>	Tetrachloroeth			<u>1.60 U</u>
Vinyl chloride			1.60 U	1,3-Dichloropi			<u>1.60 U</u>
Bromomethane			1.60 U J	Dibromochlor	omethane		<u>1.60 U</u>
Chloroethane			1.60 U	1,2-Dibromoet	thane		<u>1.60 U</u>
Trichlorofluoror	nethane		1.60 U	Chlorobenzene	0		<u>1.60 U</u>
Acetone			126 B J	1,1,1,2-Tetrac	hloroethane		<u>1.60 U</u>
1,1-Dichloroethe	ene		1.60 U	Ethylbenzene		·······	<u>1.60 U</u>
Carbon disulfide	;		14.6	p/m-Xylene			<u>3.19 U</u>
Methylene chlor	ide		3.99 U	o-Xylene			1.60 U
Methyl tert-buty	l ether (MTBE)		1.60 U	Styrene			1.60 U
trans-1,2-Dichlo	roethene		1.60 U	Bromoform			<u>1.60</u> U
1,1-Dichloroetha	1-Dichloroethane 1.60 U Isopropylbenzene					4.77	
Vinyl acetate			1.60 U	1,1,2,2-Tetrac	hloroethane		<u>1.60 U</u>
2-Butanone (ME	EK)		26.9	Bromobenzene			<u>1.60 U</u>
cis-1,2-Dichloro	2-Dichloroethene 1.60 U 1,2,3-Trichloropropane				<u>1.60 U</u>		
2,2-Dichloropro	pane		1.60 U	n-Propylbenze	me		<u>1.60 U</u>
Chloroform			1.60 <u>U</u>	2-Chlorotolue	ne		<u>1.60 U</u>
1,1,1-Trichloroe	thane		1.60 U	1,3,5-Trimethy	ylbenzene		<u>1.60 U</u>
1,1-Dichloropro	pene		1.60 U	4-Chlorotolue	ne		<u>1.60 U</u>
Carbon tetrachlo	oride		1.60 U	tert-Butylbenz	zene		<u>1.60 U</u>
Benzene			1.00 J	1,2,4-Trimethy	ylbenzene		1.60 U
1,2-Dichloroeth	ane		1.60 U	sec-Butylbenz	ene		2.48
Trichloroethene			1.60 U	1,3-Dichlorob	1,3-Dichlorobenzene		<u>1.60 U</u>
1,2-Dichloropro	pane		1.60 U	p-Isopropyltol	uene		3.99 U
Dibromomethan	ie		1.60 U	1,4-Dichlorob	enzene		3.61
Bromodichloror	nethane		1.60 U	n-Butylbenzer			<u>3.99 U</u>
2-Chloroethylvi	nyl ether		3.99 U	1,2-Dichlorob	enzene		1.07 J
Methyl isobutyl			1.60 U	1,2-Dibromo-	3-chloropropane		1.60 U
cis-1,3-Dichloro			1.60 U	1,2,4-Trichlor	obenzene		1.60 U
Toluene			1.55 J	Hexachlorobu	tadiene		1.60 U
trans-1,3-Dichlo	propropene		1.60 U	Naphthalene			3.71 J
1,1,2-Trichloroe			1.60 U	1,2,3-Trichlor	1.60 U		

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	102	70-130
1,2-Dichloroethane-d4	103	70-130
Toluene-d8	84	70-130
4-Bromofluorobenzene	79	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

-23 of 208/11/08 17:21



ARCADIS

Tidewater MGP SC13-S1

SDG: N/A

Sediment

N/A

Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-17

Associated Blank: VS072608B03

Concentration Units: µg/Kg

		D ()]]	D	Sample	Final	Dilution Factor	Analyz	. *
Date Collected		Date Analyzed	Percent Solid	Amount (g)	Volume (ml)		Analys	
07/15/08	07/16/08	07/26/08	80.9	8.16	5	1	BAS	
Parameter		I	Result	Parameter			Result	
Dichlorodifluoro	methane		1.52 U	2-Hexanone			3:79	
Chloromethane			<u>1.52 U</u>	Tetrachloroeth	and all a balance by traper to the second		1.52	
Vinyl chloride			1.52 U	1,3-Dichlorop				<u>U</u>
Bromomethane			1.52 U J	Dibromochlor				U
Chloroethane			1.52 U	1,2-Dibromoe				U
Trichlorofluoron	nethane		1.52 U	Chlorobenzen				U
Acetone			16.8 BU	1,1,1,2-Tetrac	hloroethane	,,,,,,,		U
1,1-Dichloroethe	ene		1.52 U	Ethylbenzene			1.52	
Carbon disulfide			4.11	p/m-Xylene				U
Methylene chlor	ide		3.79 U	o-Xylene			1.52	
Methyl tert-buty	l ether (MTBE)		1.52 U	Styrene			1.52	
trans-1,2-Dichlo			1.52 U	Bromoform	1.52	U		
1,1-Dichloroetha	ine		0.98 J	Isopropylbenz			1.52	U
Vinyl acetate			1.52 U	1,1,2,2-Tetrac	hloroethane		1.52	U
2-Butanone (ME	ΣK)		3.36	Bromobenzen	e		1.52	U
cis-1,2-Dichloro			1.52 U	1,2,3-Trichloropropane			1.52	U
2,2-Dichloropro	pane		1,52 U	n-Propylbenze	1.52	U		
Chloroform			1.52 U	2-Chlorotolue	ne		1.52	U
1,1,1-Trichloroe	thane		1.52 U	1,3,5-Trimethy	ylbenzene		1.52	U
1,1-Dichloropro			1.52 U	4-Chlorotolue	ne		1.52	U
Carbon tetrachlo	ride		1.52 U	tert-Butylbenz	ene		1.52	U
Benzene			1.52 U	1,2,4-Trimethy	ylbenzene		1.52	U
1,2-Dichloroetha	ane		1.52 U	sec-Butylbenz	ene		1.52	U
Trichloroethene			1.52 U	1,3-Dichlorob	enzene		1.52	U
1,2-Dichloropro	pane		1.52 U	p-Isopropyltol	uene		3.79	U
Dibromomethan			1.52 U	1,4-Dichlorob	enzene		3.44	
Bromodichloron	nethane		1.52 U	n-Butylbenzer	ıe		3.79	U
2-Chloroethylvin			3.79 U	1,2-Dichlorob	enzene		1.52	U
Methyl isobutyl			1.52 U	1,2-Dibromo-	3-chloropropane		1.52	U
cis-1,3-Dichloro			1.52 U	1,2,4-Trichlor			1.52	U
Toluene	K indiana		1.52 U	Hexachlorobu	tadiene		1,52	U
trans-1,3-Dichlo	ropropene		1.52 U	Naphthalene			3.79	U
1,1,2-Trichloroe			1.52 U	1.2.3-Trichlor	obenzene			U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	94	70-130
1,2-Dichloroethane-d4	85	70-130
Toluene-d8	88	70-130
4-Bromofluorobenzene	78	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.



Tidewater MGP

SDG: N/A

Sediment Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-18

Concentration Units:

Associated Blank: VS072608B03

µg/Kg

Date Collected	Date Received	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst		
07/15/08	07/16/08	07/26/08	30.5	7.97	5	1	BAS		
Parameter		l	Result	Parameter	<u>.</u>		Result		
Dichlorodifluoro	omethane		4.11 U	2-Hexanone			10.3 U		
Chloromethane			4.11 U	Tetrachloroeth	iene		4.11 U		
Vinyl chloride			4.11 U	1,3-Dichlorop	ropane		4.11 U		
Bromomethane			4.11 U J	Dibromochlor	omethane		4.11 U		
Chloroethane			4.11 U	1,2-Dibromoe	thane		4.11 U		
Trichlorofluoror	nethane		4.11 U	Chlorobenzen	e		4.11 U		
Acetone			241 B	1,1,1,2-Tetrac	hloroethane		4.11 U		
1,1-Dichloroethe	ene		4.11 U	Ethylbenzene			4.11 U		
Carbon disulfide	;		27.4	p/m-Xylene			8.22 U		
Methylene chlor			10.3 U	o-Xylene			4.11 U		
Methyl tert-buty			4.11 U	Styrene					
trans-1,2-Dichlo	roethene		4.11 U	Bromoform		4.11 U			
1,1-Dichloroetha			4.11 U	Isopropylbenz	ene		4.11 U		
Vinyl acetate			4.11 U	1,1,2,2-Tetrac			4.11 U		
2-Butanone (ME	EK)		67.7	Bromobenzen	e		4.11 U		
cis-1,2-Dichloro			4.11 U	1,2,3-Trichlor	opropane		4.11 U		
2,2-Dichloropro			4.11 U	n-Propylbenze		4.11 U			
Chloroform	,		4.11 U	2-Chlorotolue	ne		4.11 U		
1,1,1-Trichloroe	thane		4.11 U	1,3,5-Trimeth	ylbenzene		4.11 U		
1,1-Dichloropro	pene		4.11 U	4-Chlorotolue	ne		4.11 U		
Carbon tetrachlo	oride		4.11 U	tert-Butylbenzene			4.11 U		
Benzene			4.11 U	1,2,4-Trimeth	ylbenzene		4.11 U		
1,2-Dichloroetha	ane		4.11 U	sec-Butylbenz	ene		4.11 U		
Trichloroethene			4.11 U	1,3-Dichlorob	enzene		4.11 U		
1,2-Dichloropro	pane		4.11 U	p-Isopropyltol	p-Isopropyltoluene				
Dibromomethan	e		4,11 U	1,4-Dichlorobenzene			4.11 U		
Bromodichloron	nethane		4,11 U	n-Butylbenzer	ıe		10.3 U		
2-Chloroethylvin	nyl ether		10.3 U	1,2-Dichlorob	enzene		4.11 U		
Methyl isobutyl	ketone (MIBK)		4.11 U	1,2-Dibromo-	3-chloropropane		4.11 U		
cis-1,3-Dichloro	propene		4.11 U	1,2,4-Trichlor	obenzene		4.11 U		
Toluene			2.14 J	Hexachlorobu	tadiene		4.11 U		
trans-1,3-Dichlo	ropropene		4.11 U	Naphthalene			11.8		
1,1,2-Trichloroe			4.11 U	1,2,3-Trichlor	obenzene		4.11 U		

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	90	70-130
1,2-Dichloroethane-d4	85	70-130
Toluene-d8	87	70-130
4-Bromofluorobenzene	77	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.

U - The analyte was analyzed for but not detected at the sample specific level reported.

08/11/08 17:21

25 of 208

N/A



ARCADIS Tidewater MGP

SC15-S1 N/A SD

Matrix:

N/A SDG: Sediment Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-19 Associated Blank: VS072608B03

Concentration Units: µg/Kg

				Sample	Final	Dilution		
Date Collected		Date Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analy	
07/16/08	07/16/08	07/26/08	30.6	6.88	5	1	BAS	5
Parameter		1	Result	Parameter			Result	
Dichlorodifluoro	omethane		4.75 U	2-Hexanone			11.9	U
Chloromethane			4.75 U	Tetrachloroeth	ene		4.75	U
Vinyl chloride			<u>4.75 U</u>	1,3-Dichlorop	ropane		4.75	U
Bromomethane			4.75 U J	Dibromochlor	omethane		4.75	U
Chloroethane			4.75 U	1,2-Dibromoet	hane		4,75	Ų
Trichlorofluoror	nethane		4.75 U	Chlorobenzene	3		4.75	U
Acetone			269 B J	1,1,1,2-Tetracl	nloroethane		4.75	U
1,1-Dichloroethe	ene		4.75 U	Ethylbenzene			4.75	U
Carbon disulfide)		40.2	p/m-Xylene			9.50	U
Methylene chlor	ide		11.9 U	o-Xylene			4.75	U
Methyl tert-buty	l ether (MTBE)		4.75 U	Styrene			4.75	U
trans-1,2-Dichlo	roethene		4.75 U	Bromoform			4.75	U
1,1-Dichloroetha	me		4.75 U	Isopropylbenz	ene		4.75	U
Vinyl acetate			4.75 U	1,1,2,2-Tetracl	nloroethane		4.75	U
2-Butanone (MF	K)		73.3	Bromobenzene	3		4.75	U
cis-1,2-Dichloro	ethene		4.75 U	1,2,3-Trichloro	opropane		4.75	U
2,2-Dichloropro	pane		4.75 U	n-Propylbenze	ne		4.75	U
Chloroform			4.75 U	2-Chlorotolue	1e		4.75	U
1,1,1-Trichloroe	thane		4.75 U	1,3,5-Trimethy	/lbenzene		4.75	U
1,1-Dichloropro	pene		4.75 U	4-Chlorotoluer	ne		4.75	U
Carbon tetrachlo	ride		4.75 U	tert-Butylbenz	ene		4.75	U
Benzene			4.75 U	1,2,4-Trimethy	/lbenzene		4,75	U
1,2-Dichloroetha	ine		4.75 U	sec-Butylbenz	ene		4.75	U
Trichloroethene			4.75 U	1,3-Dichlorobe	enzene		4,75	U
1,2-Dichloropro	pane		4.75 U	p-Isopropyltol	uene		11.9	
Dibromomethan	e		4.75 U	1,4-Dichlorobe	enzene		4.75	
Bromodichloron	nethane		4.75 U	n-Butylbenzen	e		11,9	U
2-Chloroethylvii	nyl ether		11.9 U	1,2-Dichlorobe	enzene		4.75	**********
Methyl isobutyl	ketone (MIBK)		4.75 U	1,2-Dibromo-3	-chloropropane		4.75	
cis-1,3-Dichloro	propene		4.75 U	1,2,4-Trichlord	obenzene		4.75	
Toluene			4.75 U	Hexachlorobut	tadiene		4.75	
trans-1,3-Dichlo	ropropene		4.75 U	Naphthalene			11.9	
1,1,2-Trichloroe		·····	4.75 U	1,2,3-Trichlord	henzene	······································	4.75	

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	95	70-130
1,2-Dichloroethane-d4	89	70-130
Toluene-d8	88	70-130
4-Bromofluorobenzene	80	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.



ARCADIS

Tidewater MGP SC16-S1

SDG: N/A

Matrix: Sediment

N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-20

Associated Blank: VS072608B03

Concentration Units: µg/Kg

Dilution Sample Final Date Collected | Date Received | Date Analyzed | Percent Solid Amount (g) Volume (ml) Factor Analyst 07/16/08 07/16/08 07/26/08 38.3 6.50 5 1 BAS Parameter Result Result Parameter 10.0 U 4.01 U 2-Hexanone Dichlorodifluoromethane Tetrachloroethene 4.01 U 4.01 U Chloromethane 4.01 4.01 U 1,3-Dichloropropane U Vinyl chloride 4.01 U J Dibromochloromethane 4.01 U Bromomethane 4.01 U Chloroethane 1,2-Dibromoethane 4.01 U 4.01 U____ 4.01 U Chlorobenzene Trichlorofluoromethane 3 4.01 U 1,1,1,2-Tetrachloroethane 189 B Acetone 4.01 U 4.01 U Ethylbenzene 1.1-Dichloroethene 8.03 U p/m-Xylene Carbon disulfide 25.6 Methylene chloride 10.0 U o-Xylene 4.01 U Methyl tert-butyl ether (MTBE) 4.01 U Styrene 4.01 U Bromoform 4.01 U trans-1,2-Dichloroethene 4.01 U 4.01 U Isopropylbenzene 4.01 U 1,1-Dichloroethane 1,1,2,2-Tetrachloroethane 4.01 U 4.01 U Vinyl acetate 4.01 Bromobenzene U 2-Butanone (MEK) 46.0 4.01 U 4.01 U 1,2,3-Trichloropropane cis-1,2-Dichloroethene n-Propylbenzene 4.01 U 4.01 U 2.2-Dichloropropane 2-Chlorotoluene 4.01 U 4.01 U Chloroform 1,1,1-Trichloroethane 4.01 U 1,3,5-Trimethylbenzene 4.01 U 4-Chlorotoluene 4.01 U 1,1-Dichloropropene 4.01 U 4.01 U tert-Butylbenzene 4.01 U Carbon tetrachloride 4.01 U 4.01 U 1,2,4-Trimethylbenzene Benzene 4.01 U sec-Butylbenzene 4.01 U 1,2-Dichloroethane 1.3-Dichlorobenzene 4.01 U 4.01 U Trichloroethene 10.0 U p-Isopropyltoluene 1.2-Dichloropropane 4.01 U Dibromomethane 4.01 U 1,4-Dichlorobenzene 4.01 U 10.0 U n-Butylbenzene 4.01 U Bromodichloromethane 10.0 U 1,2-Dichlorobenzene 4.01 U 2-Chloroethylvinyl ether 4.01 U 1,2-Dibromo-3-chloropropane Methyl isobutyl ketone (MIBK) 4.01 U 1,2,4-Trichlorobenzene 4.01 U 4.01 U cis-1,3-Dichloropropene 4.01 U 4.01 U Hexachlorobutadiene Toluene Naphthalene 3.15 J trans-1,3-Dichloropropene 4.01 U 4.01 U 1,1,2-Trichloroethane 4.01 U 1,2,3-Trichlorobenzene

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	94	70-130
1.2-Dichloroethane-d4	90	70-130
Toluene-d8	88	70-130
4-Bromofluorobenzene	84	70-130

N/A - Not Applicable

B - Found in associated blank as well as sample.

J - Estimated value, below quantitation limit.



ARCADIS

Tidewater MGP Rinsate-1

SDG: N/A

Matrix: Water

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-21 Associated Blank: VW072208B18

μg/L

Concentration Units:

		T	San		Final			
Date Collected	Date Received	Date Analyzed			Volume (ml)	Dilution Factor	Analys	St
07/16/08	07/16/08	07/23/08		5	5	1	MLR	
Parameter		Result		Parame	eter		Result	
Dichlorodifluorom	ethane	2.00	U	2-Hexa			2.00	U
Chloromethane		2.00	U		loroethene	·····	2.00	U
Vinyl chloride		2.00	U	<u>1,3-Dic</u>	hloropropane		2.00	U
Bromomethane		2.00	U	Dibrom	ochloromethane	······································	2.00	U
Chloroethane		2.00	U	<u>1,</u> 2-Dib	romoethane		2.00	U
Trichlorofluorome	thane	2.00	U	<u>Chlorol</u>			2.00	U
Acetone		5.00	U	1,1,1,2-	Tetrachloroethane		2.00	U
1,1-Dichloroethene	>	2.00	U	Ethylbe	nzene		2.00	U
Carbon disulfide		2.00	U	p/m-Xy	lene		4.00	U
Methylene chloride	5	5.00	U	o-Xyler	ne		2.00	U
Methyl tert-butyl e	ther (MTBE)	2.00	U	Styrene			2.00	U
trans-1,2-Dichloro	ethene	2.00	U	Bromot	form		2.00	U
1,1-Dichloroethane	•	2.00	U	Isoprop	ylbenzene		2.00	U
Vinyl acetate		2,00	U	1,1,2,2-	Tetrachloroethane		2.00	U
2-Butanone (MEK)	2.00	U	Bromol	penzene		2.00	U
cis-1,2-Dichloroetl	nene	2.00	<u>U</u>	<u>1,2,3-T</u>	richloropropane		2.00	U
2,2-Dichloropropa	ne	2.00	U	n-Propy	/lbenzene		2.00	Ŭ
Chloroform		2.00	U	2-Chlor	otoluene		2.00	U
1,1,1-Trichloroetha	ane	2.00	U	1,3,5-T	rimethylbenzene		2.00	U
1,1-Dichloroprope	ne	2.00	U	4-Chlor	otoluene		2.00	U
Carbon tetrachlorid	de	2.00	U	tert-But	ylbenzene		2.00	U
Benzene		2.00	U	<u>1,2,4-T</u>	rimethylbenzene		2.00	U
1,2-Dichloroethane	9	2.00	U	sec-But	ylbenzene		2,00	U
Trichloroethene		2.00	U	<u>1,3-Dic</u>	hlorobenzene		2.00	U
1,2-Dichloropropa	ne	2.00	U	p-Isopr	opyltoluene		2.00	U
Dibromomethane		2.00	U	<u>1,4-Dic</u>	hlorobenzene		2.00	U
Bromodichloromet	thane	2.00	U	n-Butyl	benzene		2.00	U
2-Chloroethylviny	l ether	2.00	U	<u>1,2-Dic</u>	hlorobenzene		2.00	U
Methyl isobutyl ke	tone (MIBK)	2.00	U		romo-3-chloroproj	pane	2.00	U
cis-1,3-Dichloropr	opene	2.00	U	1,2,4-T	richlorobenzene		2.00	U
Toluene		2.00	U	Hexach	lorobutadiene		2,00	U
trans-1,3-Dichloro	propene	2.00	U	Naphth	alene		2.00	U
1,1,2-Trichloroeth		2.00	U	1,2,3-T	richlorobenzene		2.00	U

Surrogate	% Recovery	Acceptance Range (%)
Dibromofluoromethane	106	70-130
1,2-Dichloroethane-d4	109	70-130
Toluene-d8	100	70-130
4-Bromofluorobenzene	101	70-130

N/A - Not Applicable

Semi-Volatile Organics by 8270 - SIM

WOODS,	PHA P	roject: Tid lient ID: SC2 case: N/A				ET Lal As	b Code: MA0 R: 0807065 b ID: 080706 sociated Blank ncentration Un	5-01 : SS072208B0
Date	Date	Date	Date	Demonst Calib	Sample	Final	Dilution Factor	Analyst
Collected	Received 07/16/08	Extracted 07/22/08	Analyzed 07/30/08	Percent Solid 50.8	5.77	Volume (ml) 11.11	1	Analyst SEO
07/14/08		07/22/08	07/30/08	50.8		Result		310
	Naphthalene Acenaphthylene Acenaphthene Fluorene					7960 3210 10900 3880		
	Antl	nanthrene hracene pranthene		<u>15600</u> 6760 23400				
	Pyrene Benz[a]anthracene Chrysene					27000 12500 13400		
Benzo[b]fluoranthene Benzo[k]fluoranthene Benzo[a]pyrene Indeno[1,2,3-cd]pyrene Dibenz[a,h]anthracene Benzo[g,h,i]perylene			7640 8200 10300					
					6000 1630 5950			

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	86	30-150
Pyrene-d10	104	30-150
Benzo[b]fluoranthene-d12	97	30-150

N/A - Not Applicable

52 of 208^{3/04/08 01:49}

WOODS A	PHA P	roject: Tid lient ID: SC4 ase: N/A				ET La As	b Code: MA(TR: 0807065 b ID: 080706 ssociated Blank oncentration Ur	5-02 :: SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/14/08	07/16/08	07/22/08	07/30/08	76.0	6.18	11.11	1	SEO
Las	Para	ameter			F	Result		
	Naphthalene Acenaphthylene					58.0 205		
	to be a second stress of the s	naphthene			,	<u>36.6</u> 66.4		
		nanthrene				608		
		iracene	······			208		
	·	ranthene				1760		
	Pvre					1610		
		z[a]anthracene				841		
		sene				961		
Benzo[b]fluoranthene						701		
Benzo[k]fluoranthene				\$**	750			
Benzo[a]pyrene					688			
Indeno[1,2,3-cd]pyrene						493		
		enz[a,h]anthrac				140		
	Ben	zo[g,h,i]perylei	ne			509		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	73	30-150
Pyrene-d10	111	30-150
Benzo[b]fluoranthene-d12	95	30-150

53 of 208

WOODS ,	PHA P	Project: Tid Client ID: DU Case: N/A				ET La As	b Code: MA0 R: 0807065 b ID: 0807063 sociated Blank: incentration Un	5-03 : SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	4 1
Collected	Received	Extracted	Analyzed	Percent Solid		Volume (ml)	Factor	Analyst
07/14/08	07/16/08	07/22/08	07/30/08	73.4	6.28	4.44	1	SEO
	Para	ameter			F	lesult		
	Naphthalene Acenaphthylene Acenaphthene					96.0 125 56.0		
		orene				67.1		
	Pher	nanthrene				721		
		hracene				191		
	Fluc	oranthene				1670		
	Pyre	ene				1420		
	Ben	z[a]anthracene				739		
		ysene				817		
Benzo[b]fluoranthene						635		
Benzo[k]fluoranthene						703		
Benzo[a]pyrene						606		
Indeno[1,2,3-cd]pyrene				· · · · <u></u>		481		
Dibenz[a,h]anthracene						124		
	Ben	zo[g,h,i]peryle	ne			464	y	

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	77	30-150
Pyrene-d10	107	30-150
Benzo[b]fluoranthene-d12	96	30-150

Client: ARCADIS Project: Tidewater MGP Client ID: Client ID: Case: N/A SDG: N/A Matrix: Sediment						ET Lai As	o Code: MA(R: 0807065 b ID: 080706 sociated Blank ncentration Ur	5-04 :: SS072208B08
Date	Date	Date	Date	n	Sample	Final Valuma (ml)	Dilution Factor	Analyst
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	1 1 1	
07/14/08	07/16/08	07/22/08	07/30/08	77.8	5.79	4.44	1	SEO
·	Para	ameter			F	Result		
	Ace	hthalene naphthylene			<u>612</u> <u>590</u>			
		naphthene				<u> 540 </u>		
		orene				7050		
		nanthrene				1360		
	· · · · · · · · · · · · · · · · · · ·	hracene				10500		
	••••••			······································		9310		
	Pyre					4240		
	Benz[a]anthraceneChrysene					4440		
Benzo[b]fluoranthene						3330		
Benzo[k]fluoranthene						3060		
Benzo[a]pyrene						3000		
Indeno[1,2,3-cd]pyrene						2510		
Dibenz[a,h]anthracene					()	626		
		zo[g,h,i]peryle				2500	// // =	

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	81	30-150
Pyrene-d10	107	30-150
Benzo[b]fluoranthene-d12	92	30-150

N/A - Not Applicable

55 of 2088/04/08 01:49

WOODS M		roject: Tid lient ID: SC4 ase: N/A				ET Lat Ass	o Code: MAC R: 0807065 o ID: 080706 sociated Blank ncentration Un	5-05 : SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution Factor	Amaliyat
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	ractor	Analyst
07/14/08	07/16/08	07/22/08	07/30/08	78.4	6.41	4.44	<u> </u>	SEO
v	Para	ameter			F	Result		
	Nap	hthalene				52.4		
		naphthylene			101			
		naphthene			68.8			
		orene				75.2		
	Pher	nanthrene			1030			
	Antl	hracene		, ,		223		
	Fluc	oranthene			2240			
	Pyre	ene			1790			
	Ben	z[a]anthracene			762			
		ysene				924		
	Benzo[b]fluoranthene					672		
Benzo[k]fluoranthene						691		
Benzo[a]pyrene						569		
	Indeno[1,2,3-cd]pyrene					470		
		enz[a,h]anthrac				125		
	Ben	zo[g,h,i]peryle	ne			466		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	77	30-150
Pyrene-d10	113	30-150
Benzo[b]fluoranthene-d12	94	30-150

N/A - Not Applicable

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WOODS A	PHA P	Project: Tid Client ID: SC Case: N/A				ET La As	b Code: MA0 R: 0807065 b ID: 0807065 sociated Blank: oncentration Uni	5-06 SS072208B0
Date	Date	Date	Date	Deveent Calit	Sample	Final	Dilution	Austral
Collected 07/14/08	Received	Extracted 07/22/08	Analyzed 07/30/08	Percent Solid 24.3	Amount (g) 5.09	Volume (ml) 11.11	Factor	Analyst SEO
07/14/08		ameter	07730708	24.3		Result	ļ 1	320
Naphthalene Acenaphthylene Acenaphthene Fluorene						470 781 257 331 3270		
Phenanthrene Anthracene Fluoranthene						911 7280		
Pyrene Benz[a]anthracene Chrysene						6470 2860 3850		
Benzo[b]fluoranthene Benzo[k]fluoranthene						3450 3300		
	Inde	zo[a]pyrene mo[1,2,3-cd]py enz[a,h]anthrac				3340 2510 636		
		zo[g,h,i]peryler				2590		

	Acceptance
% Recovery	Range (%)
81	30-150
107	30-150
98	30-150
	81 107

N/A - Not Applicable

57 of 20804/08 01:50

WOODS A	PHA P	roject: Tid lient ID: SC7 ase: N/A				ET. Lat As: Co	ncentration Ur	5-07 : SS072208B08
Date	Date	Date	Date	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
Collected 07/14/08	Received 07/16/08	Extracted 07/22/08	Analyzed 07/30/08	62.5	5.39	11.11	1	SEO
0//14/08			01150100	0210	v	Result		L
Parameter Naphthalene Acenaphthylene Acenaphthene Fluorene Phenanthrene Anthracene Fluoranthene Pyrene Benz[a]anthracene Chrysene Benzo[b]fluoranthene Benzo[k]fluoranthene Benzo[a]pyrene Indeno[1,2,3-cd]pyrene Dibenz[a,h]anthracene						336 612 140 181 1300 578 3260 3600 1730 2030 2120 1950 1880 1360 373 1380		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	83	30-150
Pyrene-d10	108	30-150
Benzo[b]fluoranthene-d12	101	30-150

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 N/A - Not Applicable

58 of 208,8/04/08 01:50

Alkylated Polynuclear Aromatic Hydrocarbons

		roject: Tie lient ID: SC ase: N/A				ET) Lat	Code: MA R: 0807065 DD: 080706 sociated Blank	5-08 :: SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/14/08	07/16/08	07/22/08	07/30/08	52.7	5.30	11.11	1	SEO
L	Para	ameter			F	Result		
	Nan	hthalene				2310		
		Naphthalenes				1080		
		Naphthalenes		· · · · · · · · · · · · · · · · · · ·		1480		
		Naphthalenes				2290		
		Naphthalenes				1560		
		naphthylene				<u>1980</u> 1350	<u></u>	
		naphthene						
		orene				647	······	
C1-Fluorenes						975		
C2-Fluorenes			<u> </u>					
C3-Fluorenes Phenanthrene			6330					
		hracene		· ····		3520		
		Phenanthrenes	/Anthracenes			7000	<u></u>	
		Phenanthrenes				5860		
	and the second se	Phenanthrenes				3380		
		Phenanthrenes				1720		
		oranthene				17900		
	Pyre					17600		
	C1-	Fluoranthenes	/Pyrenes			11200		
	Ben	z[a]anthracene	2			8320	·····	
		ysene				8810		
		Chrysenes				5860		
		Chrysenes			<u></u>	3380		
		Chrysenes				2570		
		Chrysenes				<u>1240</u> 6200		
		nzo[b]fluoranth nzo[k]fluoranth				6740 ⁽	<u></u>	
		nzo[a]pyrene	10110			6900		
		vlene				1760		
		nzo[e]pyrene				5120		
		eno[1,2,3-cd]p	yrene			4630		
		enz[a,h]anthra				1260		
	Ber	nzo[g,h,i]peryl	ene			4450	· ·····	

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	84	30-150
Pyrene-d10	106	30-150
Benzo[b]fluoranthene-d12	100	30-150

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N/A - Not Applicable

59 of 208 13:05

Alkylated Polynuclear Aromatic Hydrocarbons

Date	PHA P	Project: Tic Client ID: SC Case: N/A Matrix: Sed Date	liment Date	N/A	Sample	ET Lat As: Co Final	ncentration U	5-09 κ: SS072208B08 nits: μg/Kg
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml) 11.11	Factor	Analyst SEO
07/14/08	07/16/08	07/22/08	07/30/08	75.4	6.60	L	1	360
	Nap C1- C2- C3- C4- Ace Ace Fluc C1- C2- C3- Phe Ant C1- C2- C3- C4- Flu C1- C2- C3- C4- Flu C1- C2- C3- C4- Elu C1- C2- C3- C4- Ace Ber Ber Ber Ber Ber Per	ameter hthalene Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Phaphthene orene Fluorenes Fluorenes Fluorenes Fluorenes Fluorenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Chrysenes -Chrysenes	/Anthracenes /Anthracenes /Anthracenes /Pyrenes			Result 9680 2500 2650 1800 * 1000 3830 990 2580 1210 1020 963 5480 2910 6030 3980 1760 632 12000 11000 8020 5990 6200 * 4510 2460 1730 844 4890 4750 4890 1270 3690 *		
	Dit	leno[1,2,3-cd]p benz[a,h]anthra	cene			<u>3660</u> 999		
	Ber	nzo[g,h,i]peryle	ene			3420		

	A())	Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	87	30-150
Pyrene-d10	104	30-150
Benzo[b]fluoranthene-d12	95	30-150

N/A - Not Applicable

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_61 of 208/04/08 13:06

WOODS	PHA P	roject: Tid lient ID: SC3 ase: N/A	CADIS ewater MG 6-S1 SDG: iment			ET Lat Ass	D Code: MA0 R: 0807065 D ID: 080706 sociated Blank ncentration Un	5-10 : SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/14/08	. 07/16/08	07/22/08	07/30/08	63.3	5.85	11.11	1	SEO
<u> </u>	Para	ameter			F	lesult		
		hthalene naphthylene				2 <u>1200</u> 30500		
		naphthene				3220		
	Fluc	orene				22900		
	Pher	nanthrene		· · · · · · · · · · · · · · · · · · ·		12000 E 1	2	
	Antl	nracene				33700		
	Fluc	oranthene	1			73600 E I	>	
	Pyre	me				95200 E	D	
	Ben	z[a]anthracene				50500	·····	
	Chr	ysene		P E		52700		
	Ben	zo[b]fluoranthe	me			24200		
	Ben	zo[k]fluoranthe	ene			27500		
		zo[a]pyrene				37600		
		eno[1,2,3-cd]py				20700		
		enz[a,h]anthrac				7230		
	Ben	zo[g,h,i]peryle	ne			19600	······································	

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	85	30-150
Pyrene-d10	105	30-150
Benzo[b]fluoranthene-d12	98	30-150

N/A - Not Applicable E - Estimated value, exceeds the upper limit of calibration.

63 of 208,8/04/08 01:50

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Lab Code: MA00030 ARCADIS Client: ETR: 0807065 **Tidewater MGP** Project: Lab ID: 0807065-10E Client ID: SC36-S1 WOODS HOL SDG: N/A Associated Blank: SS072208B08 N/A Case: Matrix: Sediment Concentration Units: µg/Kg Dilution Final Sample Date Date Date Date Factor Analyst Volume (ml) Percent Solid Amount (g) Analyzed Extracted Collected Received 10 SEO 11.11 5.85 08/02/08 63.3 07/22/08 07/14/08 07/16/08 Result Parameter 300 U Naphthalene 300 U Acenaphthylene 300 U Acenaphthene U 300 Fluorene 105000 Phenanthrene U 300 Anthracene 73300 Fluoranthene 94000 Pyrene 300 υ Benz[a]anthracene U 300 Chrysene 300 U Benzo[b]fluoranthene Ŭ 300 Benzo[k]fluoranthene U 300 Benzo[a]pyrene 300 U Indeno[1,2,3-cd]pyrene 300 U Dibenz[a,h]anthracene \U 300 Benzo[g,h,i]perylene

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	66	30-150
Pyrene-d10	98	30-150
Benzo[b]fluoranthene-d12	78	30-150

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

64 of 208 8/04/08 01:52

Semi-Volatile Organics by 8270 - SIM

Client: ARCADIS Project: Tidewater MGP Client ID: SC37-S1 N/A SDG: N/A Matrix: Sediment Date Date Date Date				ETI Lat Ass	sociated Blank	5-11 : SS072208B08
Date			Sample	Final		
Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/22/08	07/30/08	52.0	5.74	11.11	1	SEO
meter			F	Result		
NaphthaleneAcenaphthyleneAcenaphtheneFluorenePhenanthreneAnthraceneFluoranthenePyreneBenz[a]anthraceneChryseneBenzo[b]fluorantheneBenzo[k]fluorantheneBenzo[a]pyreneIndeno[1,2,3-cd]pyrene						
	ient ID: SC3 ise: N/A atrix: Sedi Date Extracted 07/22/08 meter thalene aphthylene aphthylene aphthene ene anthrene racene racene anthrene racene anthrene ie [a]anthracene sene so[b]fluoranthe so[a]pyrene no[1,2,3-cd]py nz[a,h]anthracene	ient ID: SC37-S1 ise: N/A SDG: atrix: Sediment Date Date Extracted Analyzed 07/22/08 07/30/08 meter thalene aphthylene aphthylene aphthene rene anthrene racene anthrene ie [a]anthracene sene o[b]fluoranthene io[a]pyrene	ient ID: SC37-S1 ise: N/A SDG: N/A atrix: Sediment Date Date Extracted Analyzed Percent Solid 07/22/08 07/30/08 52.0 meter thalene aphthylene aphthylene aphthene ene anthrene racene racene anthrene racene anthrene iene anthrene iene anthrene iene anthrene iene anthrene iene anthrene iene iene anthrene iene ialanthracene sene o[b]fluoranthene io[a]pyrene no[1,2,3-cd]pyrene nz[a,h]anthracene	ient ID: SC37-S1 ise: N/A SDG: N/A atrix: Sediment Date Date Percent Solid Amount (g) 07/22/08 07/30/08 52.0 5.74 meter F thalene aphthylene aphthylene aphthene ene anthrene racene racene anthrene racene anthrene sene sene solb]fluoranthene io[1,2,3-cd]pyrene nz[a,h]anthracene	Lak Lak se:Lak Ass ass:N/ASDG:N/AAss Ass atrix:Date ExtractedSample Percent SolidFinal Volume (ml)07/22/0807/30/0852.05.7411.11meterResultthalene2510aphthylene2200anthrene1210ene1210anthrene12100anthrene5760sene60440colspan="2">S760sene6040colspan="2">5200colspan="2">Ar60no[1,2,3-cd]pyrene4760nz[a,h]anthracene1180	Jain ID:SC37-S1Lab ID:0807063sse:N/ASDG:N/AAssociated Blankatrix:SedimentConcentration UnDateDateSampleFinalDilutionExtractedAnalyzedPercent SolidAmount (g)Volume (ml)Factor07/22/0807/30/0852.05.7411.111meterResultthalene2510aphthylene3220aphthene1210ene1220anthrene5430racene2480anthene11100ne12100sene6040sole5200sole6040sole5200sole6040sole4760no[1,2,3-cd]pyrene4020nz[a,h]anthracene1180

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	88	30-150
Pyrene-d10	105	30-150
Benzo[b]fluoranthene-d12	97	30-150

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N/A - Not Applicable

65 of 20808/04/08 01:50

Alkylated Polynuclear Aromatic Hydrocarbons

Date CollectedDate ReceivedDate ExtractedDate AnalyzedSample Percent SolidFinal Mmount (g)Dilution FactorAnalyst07/15/0807/16/0807/22/0807/31/0882.75.924.441SEOParameterResultNaphthalene298C1-Naphthalenes91.3C2-Naphthalenes125C3-Naphthalenes157C4-Naphthalenes95.4Acenaphthylene351Acenaphthylene153C1-Fluorenes158C2-Fluorenes146C3-Fluorenes146C3-Fluorenes100Anthracene101 * JC1-Phenanthrenes/Anthracenes924C3-Phenanthrenes/Anthracenes122C4-Naphthalenes112Phonenthrenes/Anthracenes112		PHA C	Project: Tie lient ID: SC Case: N/A	CADIS lewater MC 1-S1 A SDG: liment			ET Lat Ass Co	ncentration Ur	5-13 :: SS072208B08
Concected Interver Interver					D (0114				Analyst
On 15/03 On 22/03 On 22/03 Parameter Result Naphthalene 298 C1-Naphthalenes 91.3 C2-Naphthalenes 125 C3-Naphthalenes 157 C4-Naphthalenes 95.4 Acenaphthylene 351 Acenaphthylene 351 Fluorene 263 C1-Fluorenes 158 C2-Fluorenes 146 C3-Fluorenes 248 Phenanthrene 1010 Anthracene 1010 C1-Phenanthrenes/Anthracenes 924 C3-Phenanthrenes/Anthracenes 384 C4-Phenanthrenes/Anthracenes 112		.l					and the second s		
Naphthalene298C1-Naphthalenes91.3C2-Naphthalenes125C3-Naphthalenes157C4-Naphthalenes95.4Acenaphthylene351Acenaphthylene153JFluoreneC1-Fluorenes158C2-Fluorenes146C3-Fluorenes248Phenanthrene3100Anthracene1010 * JC1-Phenanthrenes/Anthracenes1680C2-Phenanthrenes/Anthracenes384C4-Phenanthrenes/Anthracenes112	07/15/08			07/31/08	82.1		l	1	560
Pyrene5660 $C1$ -Fluoranthenes/Pyrenes2680Benz[a]anthracene2830Chrysene2850C1-Chrysenes1160C2-Chrysenes516C3-Chrysenes375C4-Chrysenes152Benzo[b]fluoranthene1940Benzo[k]fluoranthene2150Benzo[a]pyrene542Benzo[a]pyrene1500Perylene542Benzo[e]pyrene1500Indeno[1,2,3-cd]pyrene1400Dibenz[a,h]anthracene361Benzo[g,h,i]perylene1240JJ		NaplC1-1C2-1C3-1C4-1AceaFluoC1-1C2-1C3-1C4-1C1-1C2-1C3-1C4-1C1-1C1-1C2-1C3-1C4-1FluoC1-1C1-1C2-1C3-1C4-1BenC1-1C2-1C3-1C4-1BenBenBenBenBenBenDib	hthalene Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes naphthylene naphthene orene Fluorenes Fluorenes Fluorenes Fluorenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes nzo[b]fluoranth nzo[k]fluoranth	/Anthracenes /Anthracenes /Anthracenes /Pyrenes e hene hene hene hene		F	298 91.3 125 157 95.4 351 153 J 263 J 158 146 248 3100 1010 J 1680 924 384 112 6540 5660 2680 2830 2850 1160 516 375 152 1940 2150 2010 542 1500 361		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	81	30-150
Pyrene-d10	114	30-150
Benzo[b]fluoranthene-d12	99	30-150

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N/A - Not Applicable

66 of 208'08 13:07

WOODS A	Client: ARCADIS Project: Tidewater MGP Client ID: SC2-S1 Case: N/A SDG: N/A Matrix: Sediment Date Date Date Date					ET Lat Ass	o Code: MA0 R: 0807065 b ID: 080706 sociated Blank ncentration Un	5-14 : SS072208B08
				_	Sample	Final	Dilution	A
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/15/08	07/16/08	07/22/08	07/31/08	83.0	20.65	11.11	<u> </u>	SEO
	Para	ameter			F	lesult		
	• <u>Ace</u> Ace Fluc Phen Antl Fluc • Pyre					483 J 323 296 523 6160 3200 7870 6570 7300		
	Benz[a]anthracene				3390 J 3320 J			
	Chrysene - Benzo[b]fluoranthene					<u>3320</u> 2070		
	Benzo[k]fluoranthene					<u>2120</u> プ		
	Benzo[a]pyrene					2570		
		eno[1,2,3-cd]py	rene			1440		
	Dib	enz[a,h]anthrac	ene			406		
	- Ben	zo[g,h,i]peryle	ne			1250		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	82	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	98	30-150

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N/A - Not Applicable

68 of 208/04/08 01:51

Client: ARCADIS Project: Client ID: Client ID: Case: N/A SDG: N/A Matrix: Sediment						ET. Lat Ass	Code: MA0 R: 0807065 DD: 080706 sociated Blank ncentration Un	5-15 : SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/15/08	07/16/08	07/22/08	07/31/08	84.0	20.31	11.11	1	SEO
L	Para	ameter			F	lesult		
	Ace Ace	hthalene naphthylene naphthene orene				93.7 J 179 50.4 J 69.4		
		nanthrene		· · · · · · · · · · · · · · · · · · ·		1050		
	Antl	nracene				282		
	Fluc	oranthene				2090		
	Pyre	ene				2230		
	Ben	z[a]anthracene				<u>929 J</u>		
		ysene				<u>1040</u>	·	
	Benzo[b]fluoranthene					790		
	Benzo[k]fluoranthene					<u>691</u> T		
Benzo[a]pyrene						798		
Indeno[1,2,3-cd]pyrene						478		
Dibenz[a,h]anthracene								
	Ben	zo[g,h,i]peryler	ne			444		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	82	30-150
Pyrene-d10	109	30-150
Benzo[b]fluoranthene-d12	99	30-150

N/A - Not Applicable

69 of 208,/04/08 01:51

Alkylated Polynuclear Aromatic Hydrocarbons

Date CollectedDate ReceivedDate ExtractedDate AnalyzedSample Percent SolidFinal Amount (g)Dilution FactorAnalyse07/15/0807/16/0807/16/0874.72.354.441SEOParameterResultNameterResultNameterResultNameterResultNameterResultNameterResultNameterStateAdvantage504C1-Naphthalenes555Accomplithylene476Accomplithylene476Accomplithylene451C1-Fluorenes274C2-Pluorenes388Phenanthrene/Anthracenes27:Pluorenes738Phenanthrene/Anthracenes1780C1-Phenanthrenes/Anthracenes1420C4-Phenanthrenes/Anthracenes1420C4-Phenanthrenes/Anthracenes1420C4-Phenanthrenes/Anthracenes1940C1-Chrysenes3860Princes1940C1-Chrysenes1940C2-Chrysenes1040C1-Chrysenes1940C2-Chrysenes1940C2-Chrysenes1940C2-Chrysenes1940C1-Chrysenes1940C2-Chrysenes1940C1-Chrysenes1940C2-Chrysenes1940C1-Chrysenes </th <th>WOODS ,</th> <th>PHAP DIE LABSC</th> <th>roject: Tid lient ID: SCI ase: N/A</th> <th>CADIS lewater MG 11-S1 SDG: iment</th> <th></th> <th></th> <th>ETF Lab Ass</th> <th></th> <th></th>	WOODS ,	PHAP DIE LABSC	roject: Tid lient ID: SCI ase: N/A	CADIS lewater MG 11-S1 SDG: iment			ETF Lab Ass		
Collected Received Extracted Analyzed Percent Solid Amount (g) Volume (ml) Pactor Analyst 07/15/08 07/16/08 07/22/08 07/31/08 74.7 2.35 Result Parameter Result Naphthalenes 279	Date	Date	Date	Date			Final	Dilution	
07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 07/13/06 Result Naphthalene 504 $(1-Naphthalenes)$ 504 $(2-Naphthalenes)$ 613 C2-Naphthalenes 613 $(2-Naphthalenes)$ 613 $(2-Naphthalenes)$ 613 C4-Naphthalenes 555 $(2-Naphthalenes)$ 555 $(2-Naphthalenes)$ 613 Acenaphthylene 476 $(2-Naphthalenes)$ 398 $(2-Naphthalenes)$ $(2-Naphthalenes)$ $(2-Naphthalenes)$ $(2-Naphthalenes)$ $(2-Naphthalenes)$ $(2-Naphthalenes)$ $(2-Naphthalenes)$ $(2-Naphtalenes)$ $(2-Naphtalen$		Received	Extracted	Analyzed	Percent Solid				
Name 504 C1-Naphthalenes 279 C2-Naphthalenes 470 C3-Naphthalenes 613 C4-Naphthalenes 555 Acenaphthylene 476 Acenaphthylene 398 Fluorene 398 Fluorene 451 C1-Fluorenes 274 C2-Fluorenes 403 C3-Fluorenes 738 Phenanthrene 4830 Anthracene 1780 C1-Phenanthrenes/Anthracenes 2550 C2-Phenanthrenes/Anthracenes 1640 C3-Phenanthrenes/Anthracenes 1660 Fluorantee 8690 Pyrene 7420 C2-Chrysenes 1040 C3-Chrysenes 1040 C1-Chrysenes 3630 Benzofalnthracene 3860 Chrysenes 1040 C3-Chrysenes 1040 C3-Chrysenes 1040 C3-Chrysenes 131 Benzofalnthracene 2560 Benzofalntharcene 2560 Benzofalpyrene 2560	07/15/08	07/16/08	07/22/08	07/31/08	74.7	2.35	4.44	1	SEO
C1-Naphthalenes 279 C2-Naphthalenes 470 C3-Naphthalenes 613 C4-Maphthalenes 555 Acenaphthylene 476 Acenaphthylene 398 Fluorene 451 C1-Fluorenes 274 C2-Fluorenes 403 C3-Fluorenes 738 Phenanthrene 4830 Anthracene 1780 C1-Phenanthrenes/Anthracenes 1840 C3-Phenanthrenes/Anthracenes 1420 C4-Phenanthrenes/Anthracenes 1420 C4-Phenanthrenes/Anthracenes 1660 Fluoranthene 8699 Pyrene 7420 C1-Fluoranthenes/Pyrenes 3630 Benz[a]anthracene 3860 C1-Strusrenes 1040 C2-Chrysenes 1940 C2-Chrysenes 1040 C3-Chrysenes 113 Benzo[k]fluoranthene 2560 Benzo[k]fluoranthene 2560 Benzo[k]fluoranthene 2560 Benzo[k]fluoranthene 2560 Benzo[k]fluoranthene	L	Para	ameter			F	Result		
Benzo[a]pyrene2850Perylene741Benzo[e]pyrene2090Indeno[1,2,3-cd]pyrene1810Dibenz[a,h]anthracene506		$\begin{array}{c} Napl \\ C1-1 \\ C2-1 \\ C3-1 \\ C4-1 \\ Acceler \\ Acceler \\ Acceler \\ C1-1 \\ C2-1 \\ C3-1 \\ C1-1 \\ C1-1 \\ C2-1 \\ C3-1 \\ C4-1 \\ Ben \\ Chr \\ C1-1 \\ C1-1 \\ C2-1 \\ C4-1 \\ Ben \\ Chr \\ C1-1 \\ C2-1 \\ C4-1 \\ Ben \\ Chr \\ C1-1 \\ C2-1 \\ C4-1 \\ Ben \\ Chr \\ C1-1 \\ C2-1 \\ C4-1 \\ Ben \\ Chr \\ C1-1 \\ C2-1 \\ C4-1 \\ Ben \\ C1-1 \\ C2-1 \\ C4-1 \\ Ben \\ C1-1 \\ C2-1 \\ C4-1 $	hthalene Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Naphthalenes Fluorenes Fluorenes Fluorenes Fluorenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Phenanthrenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes Chrysenes	/Anthracenes /Anthracenes /Anthracenes /Pyrenes			504 279 470 613 555 476 398 451 274 403 738 4830 1780 2550 1840 1420 1060 8690 7420 3630 3860 4040 1940 1040 812 413 2560		
Perylene741Benzo[e]pyrene2090Indeno[1,2,3-cd]pyrene1810Dibenz[a,h]anthracene506								·····	
Benzo[e]pyrene2090Indeno[1,2,3-cd]pyrene1810Dibenz[a,h]anthracene506				· · · · · · · · · · · · · · · · · · ·					
Indeno[1,2,3-cd]pyrene1810Dibenz[a,h]anthracene506							2090		
Dibenz[a,h]anthracene 506				yrene					
Benzolghilpervlene 1740		Dib	enz[a,h]anthra	cene					
P. A. (P. A. V. A.		Ber	zo[g,h,i]peryle	ene	<u></u>		1740		

		Acceptance		
Surrogate	% Recovery	Range (%)		
2-Methylnaphthalene-d10	81	30-150		
Pyrene-d10	105	30-150		
Benzo[b]fluoranthene-d12	99	30-150		

N/A - Not Applicable

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70 of 208/08 13:27

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WOODS A	PHA P	roject: Tid lient ID: SC1 case: N/A	CADIS lewater MG 3-S1 SDG: iment			ETI Lab Ass	Code: MA R: 0807065 DD: 080706 sociated Blank accentration U	55-17 <: SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/15/08	07/16/08	07/22/08	07/31/08	80.9	6.17	11.11	1	SEO
<u> </u>	Para	ameter			F	Result		
		hthalene naphthylene		1 g1 <u>ar</u>		<u>515</u> 372		
		naphthene				800		
		orene	· · · · · · · · · · · ·			555		
	Phe	nanthrene		- ()		11600		
	Ant	hracene		and the second		2070		
	Fluc	oranthene				15400		
	Pyre	ene			······································	12400		
	Ben	z[a]anthracene				5770		
	Chr	ysene				6060		
	Benzo[b]fluoranthene					5130		
Benzo[k]fluoranthene						5430		
	Benzo[a]pyrene					5880		
Indeno[1,2,3-cd]pyrene						4930		
		enz[a,h]anthrac				1390		
	Ben	zo[g,h,i]peryle	ne		, , ,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	4950		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	56	30-150
Pyrene-d10	84	30-150
Benzo[b]fluoranthene-d12	91	30-150

N/A - Not Applicable

72 of 208

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WOODS A	P N/A		ET. Lat Ass Coi	ncentration Un	5-18 : SS072208B08			
Date	Date	Date	Date	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
Collected	Received	Extracted 07/22/08	Analyzed 07/31/08	30.5	5.10	11.11	1	SEO
07/15/08			0//51/00			Result		l
Parameter Naphthalene Acenaphthylene Acenaphthene Fluorene Phenanthrene Anthracene Fluoranthene Pyrene Benz[a]anthracene Chrysene Benzo[b]fluoranthene Benzo[k]fluoranthene Benzo[a]pyrene Indeno[1,2,3-cd]pyrene Dibenz[a,h]anthracene						556 1330 330 373 3840 1260 10600 9840 4660 5830 5730 5140 5490 4020 1030 4330		

Surrogate	% Recovery	Acceptance Range (%)
2-Methylnaphthalene-d10	69	30-150
Pyrene-d10	101	30-150
Benzo[b]fluoranthene-d12	94	30-150

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N/A - Not Applicable

73 of 2088/04/08 01:51

WOODS,	PHA P	roject: Tid lient ID: SC1 ase: N/A	CADIS lewater MG 5-S1 SDG: iment			ET Lal As Co	ncentration Un	5-19 : SS072208B08
Date	Date	Date	Date Analyzed	Percent Solid	Sample Amount (g)	Final Volume (ml)	Dilution Factor	Analyst
Collected	Received	Extracted 07/22/08	07/31/08	30.6	5.50	11.11	1	SEO
07/10/08		ameter	07751700	V und	F	Result		······································
			1670 1100 1500 1750 16700 3880 19100 18000 7940 9200 6520 6960 7560 4860 1190 5110					

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	76	30-150
Pvrene-d10	104	30-150
Benzo[b]fluoranthene-d12	98	30-150

N/A - Not Applicable

74 of 208^{8/04/08 01:51}

WOODS,	PHA	Project: Tid Client ID: SCI Case: N/A	CADIS lewater MG 6-S1 SDG: iment			ET Lat Ass	Code: MA0 R: 0807065 DD: 080706 sociated Blank ncentration Un	5-20 : SS072208B08
Date	Date	Date	Date		Sample	Final	Dilution	
Collected	Received	Extracted	Analyzed	Percent Solid	Amount (g)	Volume (ml)	Factor	Analyst
07/16/08	07/16/08	07/22/08	07/31/08	38.3	5.75	11.11	1	SEO
(,)	Par	ameter			F	lesult		
	Nar	ohthalene				793		
		enaphthylene				1080		
		enaphthene				494		
		orene				583		
	Phe	manthrene				5990		
	Ant	thracene				1660		
	Flu	oranthene				11200		
	استاهية	ene				11000		
		nz[a]anthracene				4700		
		rysene	x, ,			5840	• • • • • •	
Benzo[b]fluoranthene						4900		
Benzo[k]fluoranthene						5100		
		nzo[a]pyrene		, , , , , , , , , , , , , , , ,		4810		
Indeno[1,2,3-cd]pyrene						3520		
Dibenz[a,h]anthracene					<u>ــــــــــــــــــــــــــــــــــــ</u>	933		
	Ber	nzo[g,h,i]peryle	ne			3630		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	81	30-150
Pyrene-d10	104	30-150
Benzo[b]fluoranthene-d12	97	30-150

N/A - Not Applicable

75 of 208

WOODS H	Clien Clien Clien Case Matr	ect: Tidewa at ID: Rinsate -1 : N/A	ter MGP I	I/A		Concentration U	65-21 k: SW071808B04
				Sample	Final Volume (ml)	Dilution Factor	Analyst
Date Collected	Date Received	Date Extracted				1	SEO
07/16/08	07/16/08	07/18/08	08/01/08	960	<u> </u>	1	<u> 3EO</u>
L	Paramo	eter			Result		
	Naphth				<u>10.4 U</u> 10.4 U	· · · · · · · · · · · · · · · · · · ·	
		hthylene			10.4 U		
	Acenap				10.4 U	· · · · · · · · · · · · · · · · · · ·	
	Fluoren				10.4 U		
	Phenan				10.4 U		
	Anthra				<u>10.4 U</u> 10.4 U		
	Fluorar	thene			10.4 U		
	Pyrene				<u> </u>	······································	
	Benz[a]anthracene					
	Chryse		· · · · · · · · · · · · · · · · · · ·	· ····································	<u> </u>		
		b]fluoranthene					
		k]fluoranthene			<u> 10.4 U</u> 10.4 U	······································	
		a]pyrene				,,	
		[1,2,3-cd]pyrene			<u> </u>		
		[a,h]anthracene	~		<u> </u>		
	Benzo	[g,h,i]perylene			<u> </u>		

		Acceptance
Surrogate	% Recovery	Range (%)
2-Methylnaphthalene-d10	86	30-150
Pyrene-d10	115	30-150
Benzo[b]fluoranthene-d12	93	30-150

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N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

76 of 20808/04/08 01:52

N/A



Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: Client ID: SC27-S1 Sediment Percent Solid: 50.8

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-01 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	3.56	Ţ	0.135	5	08/05/08	07/31/08	6020A	LCP
Arsenic	25.2		0.138	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.671	J	0.0553	5	08/04/08	07/31/08	6020A	LCP
Cadmium	2.40		0.0553	5	08/04/08	07/31/08	6020A	LCP
Chromium	327	, see	0.553	5	08/04/08	07/31/08	6020A	LCP
Copper	840		0.277	5	08/04/08	07/31/08	6020A	LCP
Lead	14700	T	5.53	100	08/04/08	07/31/08	6020A	LCP
Mercury	4.91		0.0456	10	08/04/08	07/31/08	7474	EYC
Nickel	20.2	25	0.277	5	08/04/08	07/31/08	6020A	LCP
Selenium	1.53	J.	0.277	5	08/04/08	07/31/08	6020A	LCP
Silver	1.67		0.135	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.358		0.0553	5	08/04/08	07/31/08	6020A	LCP
Zinc	689	J	1.38	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable

84 of 208 11:49



Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: N/A Client ID: SC44-S1 Sediment Percent Solid: 76.0

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-02 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

		Reporting		Date	Date	Analytical	
Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
0.172	J	0.103	5	08/05/08	07/31/08	6020A	LCP
3.04		0.0991	5	08/04/08	07/31/08	6020A	LCP
0.233	J.	0.0396	5	08/04/08	07/31/08	6020A	LCP
1.26		0.0396	5	08/04/08	07/31/08	6020A	LCP
18.2		0.396	5	08/04/08	07/31/08	6020A	LCP
39.5		0.198	5	08/04/08	07/31/08	6020A	LCP
48.9	J	0.198	5	08/04/08	07/31/08	6020A	LCP
0.104		0.0322	5	08/04/08	07/31/08	7474	EYC
8.87	J	0.198	5	08/04/08	07/31/08	6020A	LCP
0.340	J	0.198	5	08/04/08	07/31/08	6020A	LCP
0.447		0.103	5	08/05/08	07/31/08	6020A	LCP
0.0566		0.0396	5	08/04/08	07/31/08	6020A	LCP
89.3	J	0.991	5	08/04/08	07/31/08	6020A	LCP
	0.172 3.04 0.233 1.26 18.2 39.5 48.9 0.104 8.87 0.340 0.447 0.0566	0.172 J 3.04 J 0.233 J 1.26 J 18.2 J 39.5 J 0.104 S.87 0.340 J 0.447 J	Result Qualifier Limit 0.172 J 0.103 3.04 0.0991 0.233 J 0.0396 1.26 0.0396 18.2 J 0.396 39.5 0.198 48.9 J 0.198 0.104 0.0322 8.87 J 0.198 0.340 J 0.198 0.447 0.103 0.103	Result Qualifier Limit Dilution 0.172 J 0.103 5 3.04 0.0991 5 0.233 J 0.0396 5 1.26 0.0396 5 39.5 0.198 5 48.9 J 0.198 5 0.104 0.0322 5 8.87 J 0.198 5 0.340 J 0.198 5 0.447 0.103 5 5	Result Qualifier Limit Dilution Analyzed 0.172 J 0.103 5 08/05/08 3.04 0.0991 5 08/04/08 0.233 J 0.0396 5 08/04/08 1.26 0.0396 5 08/04/08 18.2 J 0.396 5 08/04/08 39.5 0.198 5 08/04/08 48.9 J 0.198 5 08/04/08 0.104 0.0322 5 08/04/08 0.340 J 0.198 5 08/04/08 0.340 J 0.198 5 08/04/08 0.340 J 0.198 5 08/04/08 0.447 0.103 5 08/05/08 08/05/08 0.0566 0.0396 5 08/04/08 08/05/08	ResultQualifierLimitDilutionAnalyzedPrepared 0.172 J 0.103 5 $08/05/08$ $07/31/08$ 3.04 0.0991 5 $08/04/08$ $07/31/08$ 0.233 J 0.0396 5 $08/04/08$ $07/31/08$ 1.26 0.0396 5 $08/04/08$ $07/31/08$ 18.2 J 0.396 5 $08/04/08$ $07/31/08$ 39.5 0.198 5 $08/04/08$ $07/31/08$ 48.9 J 0.198 5 $08/04/08$ $07/31/08$ 0.104 0.0322 5 $08/04/08$ $07/31/08$ 8.87 J 0.198 5 $08/04/08$ $07/31/08$ 0.340 J 0.198 5 $08/04/08$ $07/31/08$ 0.447 0.103 5 $08/04/08$ $07/31/08$ 0.0566 0.0396 5 $08/04/08$ $07/31/08$	ResultQualifierLimitDilutionAnalyzedPreparedMethod0.172J0.103508/05/0807/31/086020A3.040.0991508/04/0807/31/086020A0.233J0.0396508/04/0807/31/086020A1.260.0396508/04/0807/31/086020A18.2J0.396508/04/0807/31/086020A39.50.198508/04/0807/31/086020A48.9J0.198508/04/0807/31/086020A0.1040.0322508/04/0807/31/086020A0.340J0.198508/04/0807/31/086020A0.4470.103508/04/0807/31/086020A0.03660.0396508/04/0807/31/086020A

N/A



Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: Client ID: DUP-2 Sediment Percent Solid: 73.4

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-03 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.155	J	0.106	5	08/05/08	07/31/08	6020A	LCP
Arsenic	2.79		0.108	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.227	5	0.0434	5	08/04/08	07/31/08	6020A	LCP
Cadmium	1.28		0.0434	5	08/04/08	07/31/08	6020A	LCP
Chromium	17.9	J	0.434	5	08/04/08	07/31/08	6020A	LCP
Copper	39.8		0.217	5	08/04/08	07/31/08	6020A	LCP
Lead	48.3	T	0.217	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.101		0.0321	5	08/04/08	07/31/08	7474	EYC
Nickel	8.57	J	0.217	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.333	J	0.217	5	08/04/08	07/31/08	6020A	LCP
Silver	0.444		0.106	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.0562		0.0434	5	08/04/08	07/31/08	6020A	LCP
Zinc	84.4	J	1.08	5	08/04/08	07/31/08	6020A	LCP



Matrix:

Percent Solid: 77.8

ARCADIS t: Tidewater MGP

> N/A SDG: N/A SC45-S1 Sediment

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-04 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.103	υΓ	0.103	5	08/05/08	07/31/08	6020A	LCP
Arsenic	2.43		0.100	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.181	super s	0.0401	5	08/04/08	07/31/08	6020A	LCP
Cadmium	1.05		0.0401	5	08/04/08	07/31/08	6020A	LCP
Chromium	12.6	J	0.401	5	08/04/08	07/31/08	6020A	LCP
Copper	26.6		0.201	5	08/04/08	07/31/08	6020A	LCP
Lead	56.3	J	0.201	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.0641		0.0315	5	08/04/08	07/31/08	7474	EYC
Nickel	7.35		0.201	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.201	υJ	0.201	5	08/04/08	07/31/08	6020A	LCP
Silver	0.190		0.103	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.0424		0.0401	5	08/04/08	07/31/08	6020A	LCP
Zinc	70.2	J	1.00	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

87 of 208/07/08 12:02

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288



Matrix:

ARCADIS **Tidewater MGP**

SDG: N/A N/A SC46-S1 Sediment Percent Solid: 78.4

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-05 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.100	U J-	0.100	5	08/05/08	07/31/08	6020A	LCP
Arsenic	1.72		0.0985	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.155	T	0.0394	5	08/04/08	07/31/08	6020A	LCP
Cadmium	0.814		0.0394	5	08/04/08	07/31/08	6020A	LCP
Chromium	9.05	J	0.394	5	08/04/08	07/31/08	6020A	LCP
Copper	21.7		0.197	5	08/04/08	07/31/08	6020A	LCP
Lead	26.4	J.	0.197	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.0577		0.0295	5	08/04/08	07/31/08	7474	EYC
Nickel	5.77	all a	0.197	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.197	υJ	0.197	5	08/04/08	07/31/08	6020A	LCP
Silver	0.141		0.100	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.0394	U	0.0394	5	08/04/08	07/31/08	6020A	LCP
Zinc	63.5	J	0.985	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

-88 of 208/07/08 12:03

320 Forbes Blvd, Mansfield, MA 02048, (508) 822-9300, Fax (508) 822-3288

N/A



ARCADIS **Tidewater MGP**

N/A SDG: Client ID: SC18-S1 Sediment Percent Solid: 24.3

Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-06 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.77	J	0.299	5	08/05/08	07/31/08	6020A	LCP
Arsenic	26.6		0.303	5	08/04/08	07/31/08	6020A	LCP
Beryllium	1.43	T	0.121	5	08/04/08	07/31/08	6020A	LCP
Cadmium	23.5		0.121	5	08/04/08	07/31/08	6020A	LCP
Chromium	345	J	1.21	5	08/04/08	07/31/08	6020A	LCP
Copper	668		0.606	5	08/04/08	07/31/08	6020A	LCP
Lead	428	Te	0.606	5	08/04/08	07/31/08	6020A	LCP
Mercury	1.83		0.0895	5	08/04/08	07/31/08	7474	EYC
Nickel	76.8	7	0.606	5	08/04/08	07/31/08	6020A	LCP
Selenium	2.97	J	0.606	5	08/04/08	07/31/08	6020A	LCP
Silver	10.6		0.299	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.310		0.121	5	08/04/08	07/31/08	6020A	LCP
Zinc	750	J	3.03	5	08/04/08	07/31/08	6020A	LCP



Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: N/A SC7-S1 Sediment Percent Solid: 62.5

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-07 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.817	T	0.122	5	08/05/08	07/31/08	6020A	LCP
Arsenic	7.50		0.123	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.529	J	0.0494	5	08/04/08	07/31/08	6020A	LCP
Cadmium	7.20		0.0494	5	08/04/08	07/31/08	6020A	LCP
Chromium	103	J	0.494	5	08/04/08	07/31/08	6020A	LCP
Copper	244		0.247	5	08/04/08	07/31/08	6020A	LCP
Lead	177	J	0.247	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.404		0.0377	5	08/04/08	07/31/08	7474	EYC
Nickel	34.9	J	0.247	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.667	3	0.247	5	08/04/08	07/31/08	6020A	LCP
Silver	3.01		0.122	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.104		0.0494	5	08/04/08	07/31/08	6020A	LCP
Zinc	272	5	1.24	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable

90 of 208/07/08 12:03



Client: ARCADIS Project: **Tidewater MGP** N/A SDG: Client ID: SC8-S1

Sediment

Matrix:

Percent Solid: 52.7

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-08 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.73	J.	0.147	5	08/05/08	07/31/08	6020A	LCP
Arsenic	16.3		0.145	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.711	7	0.0578	5	08/04/08	07/31/08	6020A	LCP
Cadmium	8.67		0.0578	5	08/04/08	07/31/08	6020A	LCP
Chromium	326	J.	0.578	5	08/04/08	07/31/08	6020A	LCP
Copper	588		0.289	5	08/04/08	07/31/08	6020A	LCP
Lead	376	7	0.289	5	08/04/08	07/31/08	6020A	LCP
Mercury	1.76		0.0451	5	08/04/08	07/31/08	7474	EYC
Nickel	37.0	J.	0.289	5	08/04/08	07/31/08	6020A	LCP
Selenium	1.48	J	0.289	5	08/04/08	07/31/08	6020A	LCP
Silver	2.88	····	0.147	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.222		0.0578	5	08/04/08	07/31/08	6020A	LCP
Zinc	810	J	1.44	5	08/04/08	07/31/08	6020A	LCP



: ARCADIS t: Tidewater MGP

Case: N/A SDG: N/A Client ID: SC31-S1 Matrix: Sediment Percent Solid: 75.4 Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-09 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

		Reporting		Date	Date	Analytical	
Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
0.350	J	0.108	5	08/05/08	07/31/08	6020A	LCP
7.19		0.109	5	08/04/08	07/31/08		LCP
0.376	J	0.0436	5	08/04/08			LCP
0.620		0.0436	5	08/04/08	07/31/08		LCP
28.7	J	0.436	5	08/04/08	07/31/08		LCP
65.8		0.218	5	08/04/08	07/31/08		LCP
77.0	J	0.218	5	08/04/08			LCP
0.240		0.0147	5	08/04/08			EYC
21.6	<i>J</i>	0.218	5	08/04/08	07/31/08		LCP
0.752	J.	0.218	5	08/04/08	07/31/08		LCP
0.250		0.108	5	08/05/08	07/31/08		LCP
0.0577		0.0436	5	08/04/08			LCP
230	<u>J</u>	1.09	5	08/04/08	07/31/08	6020A	LCP
	0.350 7.19 0.376 0.620 28.7 65.8 77.0 0.240 21.6 0.752 0.250 0.0577	0.350 J 7.19 0.376 J 0.620 28.7 J 65.8 77.0 J 0.240 21.6 J 0.752 J 0.250 0.0577	Result Qualifier Limit 0.350 J 0.108 7.19 0.109 0.376 J 0.0436 0.620 0.0436 28.7 J 0.436 65.8 0.218 77.0 J 0.218 0.240 0.0147 21.6 J 0.218 0.250 0.108 0.0577 0.0436	Result Qualifier Limit Dilution 0.350 0.108 5 7.19 0.109 5 0.376 0.0436 5 0.620 0.0436 5 28.7 0.218 5 77.0 0.218 5 0.240 0.0147 5 0.752 0.218 5 0.752 0.218 5 0.250 0.108 5	Result Qualifier Limit Dilution Analyzed 0.350 J 0.108 5 08/05/08 7.19 0.109 5 08/04/08 0.376 J 0.0436 5 08/04/08 0.620 0.0436 5 08/04/08 28.7 J 0.436 5 08/04/08 65.8 0.218 5 08/04/08 77.0 J 0.218 5 08/04/08 0.240 0.0147 5 08/04/08 0.752 J 0.218 5 08/04/08 0.752 J 0.218 5 08/04/08 0.250 0.108 5 08/04/08 0.250 0.108 5 08/05/08 0.0577 0.0436 5 08/04/08	Result Qualifier Limit Dilution Analyzed Prepared 0.350 J 0.108 5 08/05/08 07/31/08 7.19 0.109 5 08/04/08 07/31/08 6.376 J 0.0436 5 08/04/08 07/31/08 0.620 0.0436 5 08/04/08 07/31/08 28.7 J 0.436 5 08/04/08 07/31/08 65.8 0.218 5 08/04/08 07/31/08 77.0 J 0.218 5 08/04/08 07/31/08 0.240 0.0147 5 08/04/08 07/31/08 0.752 J 0.218 5 08/04/08 07/31/08 0.250 0.108 5 08/04/08 07/31/08 0.250 0.108 5 08/04/08 07/31/08	Result Qualifier Limit Dilution Analyzed Prepared Method 0.350 J 0.108 5 08/05/08 07/31/08 6020A 7.19 0.109 5 08/04/08 07/31/08 6020A 0.376 J 0.0436 5 08/04/08 07/31/08 6020A 0.620 0.0436 5 08/04/08 07/31/08 6020A 0.620 0.0436 5 08/04/08 07/31/08 6020A 28.7 J 0.436 5 08/04/08 07/31/08 6020A 65.8 0.218 5 08/04/08 07/31/08 6020A 77.0 J 0.218 5 08/04/08 07/31/08 6020A 0.240 0.0147 5 08/04/08 07/31/08 6020A 0.752 0.218 5 08/04/08 07/31/08 6020A 0.250 0.108 5 08/04/08 07/31/08 6020A 0.2



Matrix:

ARCADIS **Tidewater MGP**

SDG: N/A N/A Client ID: SC36-S1 Sediment Percent Solid: 63.3

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-10 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.14	J	0.117	5	08/05/08	07/31/08	6020A	LCP
Arsenic	19.1		0.110	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.483	T	0.0439	5	08/04/08	07/31/08	6020A	LCP
Cadmium	4.77		0.0439	5	08/04/08	07/31/08	6020A	LCP
Chromium	310	J	0.439	5	08/04/08	07/31/08	6020A	LCP
Copper	219		0.219	5	08/04/08	07/31/08	6020A	LCP
Lead	198	J	0.219	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.673		0.0186	5	08/04/08	07/31/08	7474	EYC
Nickel	23.8	"Ja	0.219	5	08/04/08	07/31/08	6020A	LCP
Selenium	1.72	J	0.219	5	08/04/08	07/31/08	6020A	LCP
Silver	2.33		0.117	5	08/05/08	07/31/08	6020A	LCP
Thallium	0.108		0.0439	5	08/04/08	07/31/08	6020A	LCP
Zinc	680	J	1,10	5	08/04/08	07/31/08	6020A	LCP



ARCADIS **Tidewater MGP**

Matrix:

N/A SDG: N/A SC37-S1 Sediment Percent Solid: 52.0

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-11 Concentration Units: mg/Kg Date Collected: 07/14/08 Date Received: 07/16/08

		Reporting		Date	Date	Analytical	
Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
0.815	T	0.139	5	08/06/08	07/31/08	6020A	LCP
12.6		0.135	5	08/04/08	07/31/08	6020A	LCP
1.50	T	0.0540	5	08/04/08	07/31/08	6020A	LCP
11.1		0.0540	5	08/04/08	07/31/08	6020A	LCP
165	J	0.540	5	08/04/08	07/31/08	6020A	LCP
282		0.270	5	08/04/08	07/31/08	6020A	LCP
255	5	0.270	5	08/04/08	07/31/08		LCP
0.723		0.0213	5	08/04/08	07/31/08		EYC
152	3	0.270	5	08/04/08	07/31/08		LCP
1,44	5	0.270	5	08/04/08	07/31/08		LCP
4.51		0.139	5	08/06/08	07/31/08		LCP
0.140		0.0540	5	08/04/08	07/31/08		LCP
376	7	1.35	5	08/04/08	07/31/08	6020A	LCP
	0.815 12.6 1.50 11.1 165 282 255 0.723 152 1.44 4.51 0.140	$\begin{array}{c c} 0.815 & \mathbb{J}\\ \hline 12.6 \\ \hline 1.50 & \mathbb{J}\\ \hline 11.1 \\ \hline 165 & \mathbb{J}\\ \hline 282 \\ \hline 255 & \mathbb{J}\\ \hline 0.723 \\ \hline 152 & \mathbb{J}\\ \hline 1.44 & \mathbb{J}\\ \hline 4.51 \\ \hline 0.140 \\ \hline \end{array}$	Result Qualifier Limit 0.815 J 0.139 12.6 0.135 1.50 J 0.0540 11.1 0.0540 165 J 0.540 282 0.270 255 J 0.270 0.723 0.0213 152 J 0.270 1.44 J 0.270 4.51 0.139 0.140	Result Qualifier Limit Dilution 0.815 J 0.139 5 12.6 0.135 5 1.50 J 0.0540 5 11.1 0.0540 5 165 J 0.540 5 282 0.270 5 0.723 0.0213 5 152 J 0.270 5 1.44 J 0.270 5 0.139 5 5 5	Result Qualifier Limit Dilution Analyzed 0.815 J 0.139 5 08/06/08 12.6 0.135 5 08/04/08 1.50 J 0.0540 5 08/04/08 11.1 0.0540 5 08/04/08 165 J 0.540 5 08/04/08 282 0.270 5 08/04/08 255 J 0.270 5 08/04/08 152 J 0.270 5 08/04/08 152 J 0.270 5 08/04/08 152 J 0.270 5 08/04/08 1.44 J 0.270 5 08/04/08 4.51 0.139 5 08/06/08 0.140 0.0540 5 08/04/08	Result Qualifier Limit Dilution Analyzed Prepared 0.815 J 0.139 5 08/06/08 07/31/08 12.6 0.135 5 08/04/08 07/31/08 1.50 J 0.0540 5 08/04/08 07/31/08 1.11 0.0540 5 08/04/08 07/31/08 165 J 0.540 5 08/04/08 07/31/08 282 0.270 5 08/04/08 07/31/08 255 J 0.270 5 08/04/08 07/31/08 152 O.270 5 08/04/08 07/31/08 152 O.270 5 08/04/08 07/31/08 1.44 J 0.270 5 08/06/08	Result Qualifier Limit Dilution Analyzed Prepared Method 0.815 J 0.139 5 08/06/08 07/31/08 6020A 12.6 0.135 5 08/04/08 07/31/08 6020A 1.50 J 0.0540 5 08/04/08 07/31/08 6020A 1.11 0.0540 5 08/04/08 07/31/08 6020A 165 J 0.540 5 08/04/08 07/31/08 6020A 282 0.270 5 08/04/08 07/31/08 6020A 255 J 0.270 5 08/04/08 07/31/08 6020A 0.723 0.0213 5 08/04/08 07/31/08 6020A 152 J 0.270 5 08/04/08 07/31/08 6020A 1.44 J 0.270 5 08/04/08 07/31/08 6020A 4.51 0.139 5 08/04/08 07/31/08 6020A </td

N/A - Not Applicable

08/07/08 11.58



Client: ARCADIS Project: Tidewater MGP Case: N/A SDG:

SC1-S1

Matrix:

Percent Solid: 82.7

Sediment

SDG: N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-13 Concentration Units: mg/Kg Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.283	J	0.0900	5	08/06/08	07/31/08	6020A	LCP
Arsenic	4.71		0.0922	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.211	J	0.0369	5	08/04/08	07/31/08	6020A	LCP
Cadmium	0.551	an a	0.0369	5	08/04/08	07/31/08	6020A	LCP
Chromium	12.7	5	0.369	5	08/04/08	07/31/08	6020A	LCP
Copper	79.7		0.184	5	08/04/08	07/31/08	6020A	LCP
Lead	95.8	7	0.184	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.0516		0.0137	5	08/04/08	07/31/08	7474	EYC
Nickel	9.39	5	0.184	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.184	UJ	0.184	5	08/04/08	07/31/08	6020A	LCP
Silver	0.142		0.0900	5	08/06/08	07/31/08	6020A	LCP
Thallium	0.0486		0.0369	5	08/04/08	07/31/08	6020A	LCP
Zinc	83.2	5	0.922	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

96 of 208

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ARCADIS : Tidewater MGP

SDG: N/A

SC2-S1

Matrix: Sediment Percent Solid: 83.0

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-14 Concentration Units: mg/Kg Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.169	J	0.0991	5	08/06/08	07/31/08	6020A	LCP
Arsenic	5.18		0.0972	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.364	J	0.0389	5	08/04/08	07/31/08	6020A	LCP
Cadmium	0.816		0.0389	5	08/04/08	07/31/08	6020A	LCP
Chromium	15.5	2	0.389	5	08/04/08	07/31/08	6020A	LCP
Copper	35.8		0.194	5	08/04/08	07/31/08	6020A	LCP
Lead	55.9	5	0.194	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.0311		0.0146	5	08/04/08	07/31/08	7474	EYC
Nickel	11.9	J	0.194	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.194	υJ	0.194	5	08/04/08	07/31/08	6020A	LCP
Silver	0.105		0.0991	5	08/06/08	07/31/08	6020A	LCP
Thallium	0.0645		0.0389	5	08/04/08	07/31/08	6020A	LCP
Zinc	75.6	T	0.972	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

08/07/08 12:07

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98 of 208



Matrix:

Percent Solid: 84.0

ARCADIS **Tidewater MGP** SDG: N/A Client ID: Dup-3

Sediment

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-15 Concentration Units: mg/Kg Date Collected: 07/15/08 Date Received: 07/16/08

Parameter			Reporting	Date	Date	Analytical		
	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.273	J	0.0918	5	08/06/08	07/31/08	6020A	LCP
Arsenic	3.54		0.0942	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.224	J	0.0377	5	08/04/08	07/31/08	6020A	LCP
Cadmium	0.716		0.0377	5	08/04/08	07/31/08	6020A	LCP
Chromium	16.4	Jac	0.377	5	08/04/08	07/31/08	6020A	LCP
Copper	29.0		0.188	5	08/04/08	07/31/08	6020A	LCP
Lead	47.9	and the second	0.188	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.0300		0.0138	5	08/04/08	07/31/08	7474	EYC
Nickel	9.73	3	0.188	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.188	ŲΓ	0.188	5	08/04/08	07/31/08	6020A	LCP
Silver	0.0997		0.0918	5	08/06/08	· 07/31/08	6020A	LCP
Thallium	0.0377	U	0.0377	5	08/04/08	07/31/08	6020A	LCP
Zinc	76.1	and the second s	0.942	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



ARCADIS Tidewater MGP

Case:N/ASDG:Client ID:SC11-S1Matrix:SedimentPercent Solid:74.7

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-16 Concentration Units: mg/Kg Date Collected: 07/15/08 Date Received: 07/16/08

08/07/08 12:08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	r Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.557	J	0.0990	5	08/06/08	07/31/08	6020A	LCP
Arsenic	7,48		0.103	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.321	2	0.0413	5	08/04/08	07/31/08	6020A	LCP
Cadmium	3.73		0.0413	5	08/04/08	07/31/08	6020A	LCP
Chromium	66.4	T	0.413	5	08/04/08	07/31/08	6020A	LCP
Copper	202		0.207	5	08/04/08	07/31/08	6020A	LCP
Lead	193	T	0.207	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.453		0.0155	5	08/04/08	07/31/08	7474	EYC
Nickel	19.4	J.	0.207	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.473	J	0.207	5	08/04/08	07/31/08	6020A	LCP
Silver	1.39		0.0990	5	08/06/08	07/31/08	6020A	LCP
Thallium	0.0778		0.0413	5	08/04/08	07/31/08	6020A	LCP
Zinc	280	J	1.03	5	08/04/08	07/31/08	6020A	LCP



ARCADIS **Tidewater MGP**

SDG:

N/A

SC13-S1 Sediment

N/A

Percent Solid: 80.9

Matrix:

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-17 Concentration Units: mg/Kg Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	0.271	J	0.0926	5	08/06/08	07/31/08	6020A	LCP
Arsenic	3.71		0.0942	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.248	T	0.0377	5	08/04/08	07/31/08	6020A	LCP
Cadmium	1.44		0.0377	5	08/04/08	07/31/08	6020A	LCP
Chromium	17.8	T	0.377	5	08/04/08	07/31/08	6020A	LCP
Copper	47.2		0.188	5	08/04/08	07/31/08	6020A	LCP
Lead	214	J	0.188	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.0616		0.0146	5	08/04/08	07/31/08	7474	EYC
Nickel	12.4	J	0.188	5	08/04/08	07/31/08	6020A	LCP
Selenium	0.200	T	0.188	5	08/04/08	07/31/08	6020A	LCP
Silver	0.211		0.0926	5	08/06/08	07/31/08	6020A	LCP
Thallium	0.0400		0.0377	5	08/04/08	07/31/08	6020A	LCP
Zinc	225	J	0.942	5	08/04/08	07/31/08	6020A	LCP

Total Metals



Client: ARCADIS Project: Tidewater MGP Case: N/A SDG:

SC14-S1

Sediment

Matrix:

Percent Solid: 30.5

SDG: N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-18 Concentration Units: mg/Kg Date Collected: 07/15/08 Date Received: 07/16/08

102 of 208

Parameter	Result	Qualifier	Reporting Limit	Dilution	Date Analyzed	Date Prepared	Analytical Method	Analyst
Antimony	1.25	J.	0.256	5	08/06/08	07/31/08	6020A	LCP
Arsenic	24.7		0.258	5	08/04/08	07/31/08	6020A	LCP
Beryllium	1.58	intro.	0.103	5	08/04/08	07/31/08	6020A	LCP
Cadmium	13.1		0.103	5	08/04/08	07/31/08	6020A	LCP
Chromium	233	2	1.03	5	08/04/08	07/31/08	6020A	LCP
Copper	430		0.515	5	08/04/08	07/31/08	6020A	LCP
Lead	355	<u> </u>	0.515	5	08/04/08	07/31/08	6020A	LCP
Mercury	1.52		0.0410	5	08/04/08	07/31/08	7474	EYC
Nickel	51.6	T	0.515	5	08/04/08	07/31/08	6020A	LCP
Selenium	3.14	, and the second	0.515	5	08/04/08	07/31/08	6020A	LCP
Silver	6.02	- i	0.256	5	08/06/08	07/31/08	6020A	LCP
Thallium	0.333		0.103	5	08/04/08	07/31/08	6020A	LCP
Zinc	654	J	2.58	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable

1



ARCADIS t: Tidewater MGP

SDG: N/A

Client ID: SC15-S1 Matrix: Sediment Percent Solid: 30.6

N/A

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-19 Concentration Units: mg/Kg Date Collected: 07/16/08 Date Received: 07/16/08

							0111	0,00
Parameter	Result	Qualifier	Reporting Limit	Dilution	Date Analyzed	Date Prepared	Analytical Method	Analyst
Antimony	1.34	J	0.227	5	08/06/08	07/31/08	6020A	LCP
Arsenic	18.0		0.218	5	08/04/08	07/31/08	6020A	LCP
Beryllium	1.16	J	0.0874	5	08/04/08	07/31/08	6020A	
Cadmium	13.8		0.0874	5	08/04/08	07/31/08	6020A	LCP
Chromium	203	J	0.874	5	08/04/08	07/31/08	6020A	LCP
Copper	401		0.437	5	08/04/08	07/31/08	6020A	LCP
Lead	352	J.	0.437	5	08/04/08	07/31/08	6020A	LCP
Mercury	1.22		0.0378	5	08/04/08	07/31/08		LCP
Nickel	50.0	J	0.437	5	08/04/08	07/31/08		EYC
Selenium	2,37	3	0.437	5	08/04/08	07/31/08	6020A	LCP
Silver	6.94		0.227	5	08/06/08		6020A	LCP
Thallium	0.228	**************************************	0.0874			07/31/08	6020A	LCP
Zinc	578			5	08/04/08	07/31/08	6020A	LCP
			2,18	<u> </u>	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable

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103 of 208

Total Metals



Matrix:

ARCADIS **Tidewater MGP**

N/A SDG: N/A SC16-S1 Sediment Percent Solid: 38.3

Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-20 Concentration Units: mg/Kg Date Collected: 07/16/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	1.15	J.	0.193	5	08/06/08	07/31/08	6020A	LCP
Arsenic	15.0		0.198	5	08/04/08	07/31/08	6020A	LCP
Beryllium	0.974	J	0.0790	5	08/04/08	07/31/08	6020A	LCP
Cadmium	9.24		0.0790	5	08/04/08	07/31/08	6020A	LCP
Chromium	136	7.	0.790	5	08/04/08	07/31/08	6020A	LCP
Copper	293		0.395	5	08/04/08	07/31/08	6020A	LCP
Lead	277	J	0.395	5	08/04/08	07/31/08	6020A	LCP
Mercury	0.792		0.0289	5	08/04/08	07/31/08	7474	EYC
Nickel	37.0	T	0.395	5	08/04/08	07/31/08	6020A	LCP
Selenium	1.79	star and a star a st	0.395	5	08/04/08	07/31/08	6020A	LCP
Silver	8.23		0.193	5	08/06/08	07/31/08	6020A	LCP
Thallium	0.219		0.0790	5	08/04/08	07/31/08	6020A	LCP
Zinc	467	T	1.98	5	08/04/08	07/31/08	6020A	LCP

N/A - Not Applicable

Total Metals

N/A



Matrix:

ARCADIS Tidewater MGP

N/A SDG: Rinsate-1 Water Lab Code: MA00030 ETR: 0807065 Lab ID: 0807065-21 Concentration Units: μg/L Date Collected: 07/16/08 Date Received: 07/16/08

			Reporting		Date	Date	Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Prepared	Method	Analyst
Antimony	2.50	U	2.50	5	08/05/08	08/01/08	6020A	LCP
Arsenic	2.50	U	2.50	5	08/04/08	08/01/08	6020A	LCP
Beryllium	2.50	U	2.50	5	08/04/08	08/01/08	6020A	LCP
Cadmium	1.00	U	1.00	5	08/04/08	08/01/08	6020A	LCP
Chromium	2.50	υ	2.50	5	08/04/08	08/01/08	6020A	LCP
Copper	2.50	U	2.50	5	08/04/08	08/01/08	6020A	LCP
Lead	5.00	U	5.00	5	08/04/08	08/01/08	6020A	LCP
Mercury	0.100	U	0.100	1	08/05/08	08/05/08	7470A	EYC
Nickel	2.50	U	2.50	5	08/04/08	08/01/08	6020A	LCP
Selenium	5.00	U	5.00	5	08/04/08	08/01/08	6020A	LCP
Silver	1.00	U	1.00	5	08/05/08	08/01/08	6020A	LCP
Thallium	1.00	U	1.00	5	08/04/08	08/01/08	6020A	LCP
Zinc	50.0	U	50.0	5	08/04/08	08/01/08	6020A	LCP

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

SDG: N/A

Client ID: SC27-S1 Matrix: Sediment

Matrix: Sedimer Percent Solid: 50.8

N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-01 Date Collected: 07/14/08 Date Received: 07/16/08

		Reporting			Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	7.3		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	7.3		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	0.47	U	0.47	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

125 of 20808/08 08:54



ARCADIS **Tidewater MGP**

N/A SDG: N/A

Client ID: SC44-S1 Matrix: Percent Solid: 76.0

Sediment

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-02 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.2		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	1.2		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	0.23	U	0.23	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

N/A



ARCADIS **Tidewater MGP**

N/A

Matrix:

Percent Solid: 73.4

Sediment

SDG:

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-03 Date Collected: 07/14/08 Date Received: 07/16/08

		Reporting			Date	Analytical			
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	1.2		0.01	1	08/05/08	%	9060	ES	
Total Organic Carbon (Run 2)	1.4		0.01	1	08/05/08	%	9060	ES	
Physiologically Available Cyanide	0.30	U	0.30	1	07/28/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

127 of 208/08/08 08:54

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ARCADIS **Tidewater MGP**

> N/A SDG:

Client ID: SC45-S1 Sediment

Percent Solid: 77.8

Matrix:

N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-04 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.76		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	0.56		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide		U	0.19	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

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ARCADIS : Tidewater MGP

SDG: N/A

Client ID: SC46-S1 Matrix: Sediment

N/A

Percent Solid: 78.4

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-05 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.34		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	0.32		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	0.27	U	0.27	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

129 of 208%/08 08:54



ARCADIS Tidewater MGP

SDG: N/A

Client ID: SC18-S1 Matrix: Sediment

N/A

Percent Solid: 24.3

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-06 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	5.8		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	5.6		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	0.73	U	0.73	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

130 of 208/08 08:54

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ARCADIS Tidewater MGP

N/A SDG; N/A

Client ID: SC7-S1 Matrix: Sediment Percent Solid: 62.5 Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-07 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	2.7		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	2.6		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	0.15	υ	0.15	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

131 of 208 8/08 08:54



ARCADIS **Tidewater MGP**

> SDG: N/A

Sediment Matrix: Percent Solid: 52.7

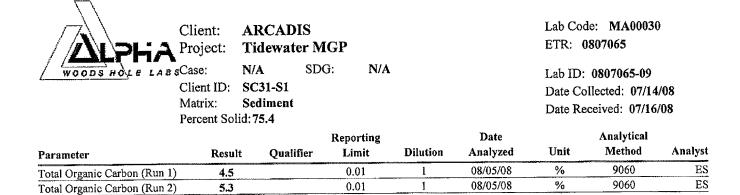
N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-08 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
	4.7		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 1)			0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	<u>4.7</u> 0.39	TI	0.39	1	07/28/08	mg/Kg	9010 PACN	ES
Physiologically Available Cyanide	0.39		0.57	• <u> </u>				

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



0.16

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N/A - Not Applicable

Physiologically Available Cyanide

1.5

133 of 208/08/08 08:54

9010 PACN

mg/Kg

07/28/08

ES



ARCADIS : Tidewater MGP

N/A SDG: N/A

Client ID: SC36-S1 Matrix: Sediment

Percent Solid: 63.3

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-10 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	5.9		0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	6.4		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	3.8		0.25	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable

134 of 20808/08 08:54



ARCADIS Tidewater MGP

SDG: N/A

SC37-S1 Sediment

Matrix: Sedime Percent Solid: 52.0

N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-11 Date Collected: 07/14/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.5	<u>Un dessentent Victories p</u> rocessories en	0.01	1	08/05/08	%	9060	ES
Total Organic Carbon (Run 2)	4.1		0.01	1	08/05/08	%	9060	ES
Physiologically Available Cyanide	0.36	U	0.36	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

135 of 2088 08:54

N/A



ARCADIS **Tidewater MGP**

N/A

Sediment

Matrix:

Percent Solid:82.7

SDG:

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-13 Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0.68		0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	0.75		0.01	1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.27	U	0.27	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

138 of 208 8/08 08:54



ARCADIS : Tidewater MGP

SDG: N/A

Client ID: SC2-S1 Matrix: Sediment

N/A

Percent Solid: 83.0

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-14 Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	2.2	5	0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	1.9	2	0.01	1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.30	U	0.30	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

141 of 208 B/08 DB:54

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ARCADIS : Tidewater MGP

N/A SDG: N/A

un-3

Matrix: Sediment Percent Solid:84.0 Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-15 Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	0,40	J	0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	0.22	J	0.01	1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.29	U	0.29	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

142 of 208 8/08 08:54



ARCADIS Tidewater MGP

SDG: N/A

Client ID: SC11-S1 Matrix: Sediment

N/A

Percent Solid: 74.7

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-16 Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	1.5	す	0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	1.0	J	0.01	· 1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.32	U	0.32	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

N/A SDG: N/A

Client ID: SC13-S1 Matrix: Sediment

Percent Solid: 80.9

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-17 Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Analytical			
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst	
Total Organic Carbon (Run 1)	2.2	J	0.01	1	08/06/08	%	9060	ES	
Total Organic Carbon (Run 2)	3.4	্য	0.01	1	08/06/08	%	9060	ES	
Physiologically Available Cyanide	0.26	Ų	0.26	1	07/28/08	mg/Kg	9010 PACN	ES	

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.



ARCADIS Tidewater MGP

N/A SDG: N/A

Client ID: SC14-S1 Matrix: Sediment Percent Solid: 30.5 Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-18 Date Collected: 07/15/08 Date Received: 07/16/08

			Reporting		Date	Analytical		
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.6		0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	4.6		0.01	1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.78	U	0.78	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

145 of 20808/08 08:54



ARCADIS Tidewater MGP

SDG: N/A

SC15-S1

Matrix: Sediment

N/A

Percent Solid: 30.6

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-19 Date Collected: 07/16/08 Date Received: 07/16/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	3.6		0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	3.8		0.01	1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.80	U	0.80	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

146 of 20808/08 08:54

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ARCADIS **Tidewater MGP**

N/A

Sediment

Matrix:

Percent Solid: 38.3

SDG: N/A

Lab Code: MA00030 ETR: 0807065

Lab ID: 0807065-20 Date Collected: 07/16/08 Date Received: 07/16/08

			Reporting		Date		Analytical	
Parameter	Result	Qualifier	Limit	Dilution	Analyzed	Unit	Method	Analyst
Total Organic Carbon (Run 1)	4.0	999.07	0.01	1	08/06/08	%	9060	ES
Total Organic Carbon (Run 2)	4.0		0.01	1	08/06/08	%	9060	ES
Physiologically Available Cyanide	0.63	υ	0.63	1	07/28/08	mg/Kg	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

147 of 208/08/08 08:54

	lient: oject:	ARCADIS Tidewater						ide: MA00030 0807065	•
	ase: ient ID: atrix:	N/A S Rinsate-1 Water	DG: 1	N/A			Date C	: 0807065-21 ollected: 07/16, eceived: 07/16/	
Parameter	Resul	t Qualifier	Reportiı Limit	÷	lution	Date Analyzed	Unit	Analytical Method	Analyst
Physiologically Available Cyanid	e 0.005	5 U	0.005		1	07/28/08	mg/L	9010 PACN	ES

N/A - Not Applicable U - The analyte was analyzed for but not detected at the sample specific level reported.

148 of 208 8/08 08:54

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810643-01 SC27-S1	Date Collected: 14-JUL-2008 08:50 Date Received : 17-JUL-2008
Sample Matrix:	SOIL	Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe:	rs: 1-Amber	

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
					PREP ANAL
Solids, Total	53	8	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	8040	mg/kg	75.5	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 3 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810643-02 SC44-S1	Date Collected: 14-JUL-2008 09:32 Date Received : 17-JUL-2008
Sample Matrix:	SOIL	Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe	rs: 1-Amber	

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
					PREP ANAL
48-1999	······································				
Solids, Total	73	ę	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	958	mg/kg	54,8	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 4 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810643-03 DUP-2	Date Collected: 14-JUL-2008 00:00 Date Received : 17-JUL-2008
Sample Matrix:	SOIL	Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe:	rs: 1-Amber	

PARAMETER	RESULT	UNITS	RDL I	REF METHOD	DATE ID
					PREP ANAL
					······································
Solids, Total	72	8	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	1260	mg/kg	55.6	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 5 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAC00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-04 SC45-S1 SOIL	Date Collected: Date Received : Date Reported :	
Condition of Sample:	Satisfactory	Field Prep: 1	None
Number & Type of Containe	rs: 1-Amber		

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
					PREP ANAL
Solids, Total	74	80 0	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	637	mg/kg	48.6	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 6 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-05 SC46-S1 SOIL	Date Collected: 14-JUL-2008 10:46 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe	ers: 1-Amber	

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
	-				PREP ANAL
					·····
Solids, Total	78	Q	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	451	mg/kg	46.2	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 7 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-06 SC18-S1 SOIL	Date Collected: 14-JUL-2008 11:25 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe	ers: 1-Amber	

PADAL/CORD	RESULT	UNITS	RDL	REF METHOD	DATE ID	
PARAMETER					PREP ANAL	
Solids, Total	23	5	0.10	30 2540G	0718 12:02 SD	
TPH, HEM-SGT	12200	mg/kg	156	1 9071B	0724 17:00 0725 11:30 AT	

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 8 of 26

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:	L0810643-07 SC7-S1	Date Collected: 14-JUL-2008 12:02 Date Received : 17-JUL-2008
Sample Matrix:	SOIL	Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None
Number & Type of Containe	ers: 1-Amber	

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
PARAMETER	1000				PREP ANAL
Solids, Total	59	ç	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	2750	mg/kg	67.8	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 9 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	SC8-51	Date Collected: 14-JUL-2008 12:38 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
- Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	49	ę	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	6890	mg/kg	81.6	1 9071B	0724 17:00 0725 11:30 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 10 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	SC31-S1	Date Collected: Date Received : Date Reported :	
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	73	ę	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	1290	mg/kg	54.8	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 11 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-10 SC36-S1 SOIL	Date Collected: 14-JUL-2008 13:18 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
	•				PREP ANAL
					0710 10:00 60
Solids, Total	61	8	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	4790	mg/kg	65.6	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 12 of 26

181 of 208

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-11 SC37-S1 SOIL	Date Collected: 14-JUL-2008 14:51 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 2-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	49	8	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	3960	mg/kg	81.6	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 13 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-12 SC1-S1 SOIL	Date Collected: 14-JUL-2008 11:0 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008	00
Condition of Sample:	Satisfactory	Field Prep: None	

Number & Type of Containers: 2-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	79	8	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	1290	mg/kg	50.6	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 14 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-13 SC2-S1 SOIL	Date Collected: 15-JUL-2008 12:04 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	87	સ્	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	520	mg/kg	41.4	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 15 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-14 DUP-3 SOIL	Date Collected: 15-JUL-2008 00:00 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	89	90 10	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	573	mg/kg	44.9	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 16 of 26

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LAO00065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-15 SC11-S1 SOIL	Date Collected: 15-JUL-2008 14:05 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
	nuo out	0.1.2.2.0			PREP ANAL
Solids, Total	70	8	0.10	30 2540G	0716 12:02 SD
TPH, HEM-SGT	5390	mg/kg	57.1	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 17 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-16 SC13-S1 SOIL	Date Collected: Date Received : Date Reported :	
Condition of Sample:	Satisfactory	Field Prep:	None

Number & Type of Containers: 1-Amber

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PRE P A NAL
Solids, Total	80	8	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	1730	mg/kg	45.0	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 18 of 26

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	SC14-S1	Date Collected: 15-JUL-2008 16:10 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
Solids, Total	29	90	0,10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	9320	mg/kg	138	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 19 of 26

188 of 208

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	SC15-S1	Date Collected: 16-JUL-2008 10:05 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

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PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID
					PREP ANAL
Solids, Total	30	ક્ષ	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	9630	mg/kg	120	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 20 of 26

189 of 208

MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number: Sample Matrix:	L0810643-19 SC16-S1 SOIL	Date Collected: 16-JUL-2008 12:22 Date Received : 17-JUL-2008 Date Reported : 30-JUL-2008
Condition of Sample:	Satisfactory	Field Prep: None

Number & Type of Containers: 1-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	date id Prep Anal
Solids, Total	38	95 15	0.10	30 2540G	0718 12:02 SD
TPH, HEM-SGT	5010	mg/kg	105	1 9071B	0728 10:15 0729 10:45 AT

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 21 of 26

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MA:M-MA086 NH:2003 CT:PH-0574 ME:MA0086 RI:LA000065 NY:11148 NJ:MA935 Army:USACE

Laboratory Sample Number:L0810643-20
RINSATE-1Date Collected:16-JUL-2008
09:10Sample Matrix:WATERDate Received:17-JUL-2008Condition of Sample:SatisfactoryField Prep:None

Number & Type of Containers: 2-Amber

PARAMETER	RESULT	UNITS	RDL	REF METHOD	DATE ID PREP ANAL
ТРН	ND	mg/l	4.00	74 1664A	0721 20:00 0722 18:00 JO

Comments: Complete list of References and Glossary of Terms found in Addendum I

07300812:11 Page 22 of 26

191 of 208

CHAIN OF CUSTODY

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WESTBORO, MA	RAYNHAM,MA	Project Information	ų			Report Information - Data Deliverables	ion - Data Dell	iverables	BillingIn	Billing Information	
TEL: 508-898-9220 FAX: 508-898-9193	300 288	Project Name: TIDE WOTER NGP	ervote	N NGF		CI FAX	MEMAIL		🗆 Same as	Client info PO #:	
Client Information		Project Location: POWHUCKET, RI	WHUC)	(et. RI	_	ADEX	C Add'l Deliverables	rables			
Client: Droodio	いた	Project #: 260	L C		Å	Regulatory Requirements/Report Limits	irements/Rep	ortLimits			; :
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