

**ATTACHMENT A  
DATA USABILITY SUMMARY REPORT  
FOURTH QUARTER 2010**

**HEMPSTEAD INTERSECTION STREET FORMER MGP SITE  
VILLAGES OF GARDEN CITY AND HEMPSTEAD  
LONG ISLAND, NEW YORK**

**Analyses Performed by:  
H2M LABORATORIES, INC.**

**Prepared For:**

**NATIONAL GRID  
175 EAST OLD COUNTRY RD.  
HICKSVILLE, NY 11801**

**Prepared by:**

**URS CORPORATION  
77 GOODELL STREET  
BUFFALO, NY 14203**

**FEBRUARY 2011**

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## I. INTRODUCTION

This Data Usability Summary Report (DUSR) has been prepared following the guidelines provided in New York State Department of Environmental Conservation (NYSDEC) Division of Environmental Remediation *DER-10, Technical Guidance for Site Investigation and Remediation, Appendix 2B - Guidance for Data Deliverables and Development of Data Usability Summary Reports*, May 2010.

Analytical data for twenty-one (21) groundwater samples, one (1) field duplicate, one (1) matrix spike/matrix spike duplicate (MS/MSD) pair, and three (3) trip blanks collected by URS personnel from October 21 to 28, 2010 are discussed in this DUSR. The samples were collected as part of the 2010 fourth quarter groundwater monitoring event at the Hempstead Intersection Street Former MGP Site.

## II. ANALYTICAL METHODOLOGIES AND DATA VALIDATION

The samples were analyzed by H2M Laboratories, Inc. (Melville, NY) for the following parameters:

- Benzene, toluene, ethylbenzene, and xylene (BTEX) – USEPA Method SW8260B, and
- Polynuclear aromatic hydrocarbons (PAHs) – USEPA Method SW8270C.

A limited data validation was performed on the samples in accordance with the guidelines presented in the following USEPA Region II documents:

- *Validating Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry SW-846 Method 8260B, SOP HW-24, Rev. 2, August 2008*; and
- *Validating Semivolatile Organic Compounds by Gas Chromatography/Mass Spectrometry SW-846 Method 8270D, SOP HW-22, Rev. 4, August 2008*.

The limited data validation included a review of completeness of all required deliverables; holding times; quality control (QC) results (instrument tunes, calibration standards, blanks, matrix spike recoveries, field duplicate analyses, laboratory control sample recoveries, and surrogate/internal standard recoveries) to determine if the data are within the protocol-required QC limits and specifications; a determination that all samples were analyzed using established and agreed upon analytical protocols; an evaluation of the raw data to confirm the results provided in the data summary sheets; and a review of laboratory data qualifiers.

Qualifications applied to the data during the data validation process include 'UJ' (estimated quantitation limit). The validated analytical results are presented in Tables A-1 and A-2. Copies of the validated laboratory results (i.e., Form 1's) are presented in Appendix A. Copies of the chain-of-custodies, case narratives, and documentation supporting the qualification of data are presented in Appendix B. Only problems affecting data usability are discussed in this report.

### **III. DATA DELIVERABLE COMPLETENESS**

Full deliverable data packages (i.e., NYSDEC ASP Category B or equivalent) were provided by the laboratory, and included all reporting forms and raw data necessary to fully evaluate and verify the reported analytical results.

### **IV. SAMPLE RECEIPT/HOLDING TIMES**

All samples were received by the laboratory intact, properly preserved, and under proper chain-of-custody (COC), except for the following instances.

- The cooler temperatures associated with samples collected from October 25 to 28, 2010 (i.e., 9-10°C) were above QC limits (i.e., 4°C ± 2°C). No qualification of the data was necessary, per USEPA Region II validation guidelines, since the cooler temperatures were less than or equal to 10°C.

All samples were analyzed within the required holding times.

## V. NON-CONFORMANCES

For PAH analyses, the percent differences (%Ds) between the initial calibration (ICAL) average relative response factors (RRFs) and the RRFs in the continuing calibration (CCAL) standards were greater than 20.0% for benzo(k)fluoranthene. The following PAH results were qualified 'UJ'.

Sample ID	Affected Compound
HIMW-05D, -05I, -08D, -12D, -12I, -12S, -13D, -13I, -13S, -14D, -14I, -15D, -15I, -20I, -20I-Dup, -20S,	Benzo(k)fluoranthene

Documentation supporting the qualification of data (i.e., Forms 5 and 7) is presented in Appendix B.

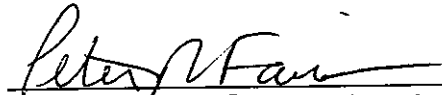
## VI. SAMPLE RESULTS AND REPORTING

All sample results were reported in accordance with method requirements and were adjusted for sample size and dilution factors. BTEX and PAH results detected below the quantitation limits were qualified 'J' by the laboratory. The results reported from secondary dilution analyses were qualified 'D' by the laboratory.

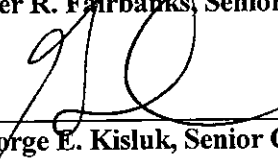
A field duplicate was collected from monitoring well location HIMW-20I, which exhibited good field and analytical precision.

## VIII. SUMMARY

All sample analyses were found to be compliant with the method and validation criteria, and the data are usable as reported, except for those results qualified 'UJ', which should be considered conditionally usable. URS does not recommend the re-collection of any samples at this time.

Prepared By:   
Peter R. Fairbanks, Senior Chemist

Date: 2/8/11

Reviewed By:   
George E. Kisluk, Senior Chemist

Date: 2/9/11

## DEFINITIONS OF USEPA REGION II DATA QUALIFIERS

- U – The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- J – The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ – The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R – The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.
- D – The sample results are reported from a separate secondary dilution analysis.
- NJ – The analysis indicates the presence of an analyte that has been “tentatively identified” and the associated numerical value represents its approximate concentration.

**TABLE A-1**  
**VALIDATED GROUNDWATER SAMPLE ANALYTICAL RESULTS**  
**NATIONAL GRID - HEMPSTEAD INTERSECTION STREET FORMER MGP SITE**

Location ID			HIMW-003D	HIMW-003I	HIMW-003S	HIMW-005D	HIMW-005I
Sample ID			HIMW-3D	HIMW-3I	HIMW-3S	HIMW-5D	HIMW-5I
Matrix			Groundwater	Groundwater	Groundwater	Groundwater	Groundwater
Depth Interval (ft)			-	-	-	-	-
Date Sampled			10/27/10	10/27/10	10/28/10	10/26/10	10/26/10
Parameter	Units	Criteria*					
<b>Volatile Organic Compounds</b>							
Benzene	UG/L	-	1 U	1 U	1 U	1 U	3
Ethylbenzene	UG/L	-	1 U	1 U	1 U	1 U	1
Toluene	UG/L	-	1 U	1 U	1 U	6	1 U
Xylene (total)	UG/L	-	1 U	1 U	1 U	210	150
Total BTEX	UG/L	100	ND	ND	ND	216	154
<b>Semivolatile Organic Compounds</b>							
2-Methylnaphthalene	UG/L	-	10 U	10 U	10 U	270 D	560 D
Acenaphthene	UG/L	-	10 U	10 U	10 U	4 J	16
Acenaphthylene	UG/L	-	10 U	10 U	10 U	51	230 DJ
Anthracene	UG/L	-	10 U	10 U	10 U	10 U	3 J
Benzo(a)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(a)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(b)fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(k)fluoranthene	UG/L	-	10 U	10 U	10 U	10 UJ	10 UJ
Chrysene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluorene	UG/L	-	10 U	10 U	10 U	3 J	23
Indeno(1,2,3-cd)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Naphthalene	UG/L	-	10 U	10 U	10 U	1,400 D	2,300 D
Phenanthrene	UG/L	-	10 U	10 U	10 U	10 U	20
Pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Total Polynuclear Aromatic Hydrocarbons	UG/L	100	ND	ND	ND	1,728	3,152

\*Criteria- Groundwater Plume Delineation/Design Criteria, Pre-Design Investigation Work Plan for In-Situ Solidification for the Hempstead Intersection Street Former MGP Site, Appendix E, Final, URS 2008.

Flags assigned during chemistry validation are shown.

 Concentration Exceeds Criteria

U - Not detected above the reported quantitation limit. J - The reported concentration is an estimated value.

UJ - Not detected. The reported quantitation limit is an estimated value.

D - Result reported from a secondary dilution analysis.

Made By\_PRF 01/27/11\_ Checked By\_AMK 01/27/11\_

Detection Limits shown are PQL



**TABLE A-1**  
**VALIDATED GROUNDWATER SAMPLE ANALYTICAL RESULTS**  
**NATIONAL GRID - HEMPSTEAD INTERSECTION STREET FORMER MGP SITE**

Location ID			HIMW-005S	HIMW-008D	HIMW-008I	HIMW-008S	HIMW-012D
Sample ID			HIMW-5S	HIMW-8D	HIMW-8I	HIMW-8S	HIMW-12D
Matrix			Groundwater	Groundwater	Groundwater	Groundwater	Groundwater
Depth Interval (ft)			-	-	-	-	-
Date Sampled			10/27/10	10/26/10	10/28/10	10/28/10	10/25/10
Parameter	Units	Criteria*					
<b>Volatile Organic Compounds</b>							
Benzene	UG/L	-	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	UG/L	-	1 U	1 U	1 U	1 U	1 U
Toluene	UG/L	-	1 U	1 U	1 U	1 U	1 U
Xylene (total)	UG/L	-	1 U	1 U	1 U	1 U	1 U
Total BTEX	UG/L	100	ND	ND	ND	ND	ND
<b>Semivolatile Organic Compounds</b>							
2-Methylnaphthalene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Acenaphthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Acenaphthylene	UG/L	-	10 U	10 U	10 U	1 J	10 U
Anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(a)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(a)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(b)fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(k)fluoranthene	UG/L	-	10 U	10 UJ	10 U	10 U	10 UJ
Chrysene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluorene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Indeno(1,2,3-cd)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Naphthalene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Phenanthrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Total Polynuclear Aromatic Hydrocarbons	UG/L	100	ND	ND	ND	1	ND

\*Criteria- Groundwater Plume Delineation/Design Criteria, Pre-Design Investigation Work Plan for In-Situ Solidification for the Hempstead Intersection Street Former MGP Site, Appendix E, Final, URS 2008.

Flags assigned during chemistry validation are shown.



Concentration Exceeds Criteria

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D - Result reported from a secondary dilution analysis.

Made By\_PRF 01/27/11\_; Checked By\_AMK 01/27/11\_

Detection Limits shown are PQL

**TABLE A-1**  
**VALIDATED GROUNDWATER SAMPLE ANALYTICAL RESULTS**  
**NATIONAL GRID - HEMPSTEAD INTERSECTION STREET FORMER MGP SITE**

Location ID			HIMW-012I	HIMW-012S	HIMW-013D	HIMW-013I	HIMW-013S
Sample ID			HIMW-12I	HIMW-12S	HIMW-13D	HIMW-13I	HIMW-13S
Matrix			Groundwater	Groundwater	Groundwater	Groundwater	Groundwater
Depth Interval (ft)			-	-	-	-	-
Date Sampled			10/25/10	10/25/10	10/21/10	10/21/10	10/22/10
Parameter	Units	Criteria*					
<b>Volatile Organic Compounds</b>							
Benzene	UG/L	-	48	1 U	4	96	1 U
Ethylbenzene	UG/L	-	1 U	1 U	1 U	1 U	1 U
Toluene	UG/L	-	1 U	1 U	1 U	1 U	1 U
Xylene (total)	UG/L	-	4	6	2	7	1 U
Total BTEX	UG/L	100	52	6	6	103	ND
<b>Semivolatile Organic Compounds</b>							
2-Methylnaphthalene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Acenaphthene	UG/L	-	40	10 U	6 J	11	10 U
Acenaphthylene	UG/L	-	43	10 U	13	90 D	10 U
Anthracene	UG/L	-	10 U	10 U	10 U	2 J	10 U
Benzo(a)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(a)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(b)fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(k)fluoranthene	UG/L	-	10 UJ	10 UJ	10 UJ	10 UJ	10 UJ
Chrysene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluorene	UG/L	-	25	10 U	10 U	15	10 U
Indeno(1,2,3-cd)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Naphthalene	UG/L	-	2 J	10 U	10 U	1 J	10 U
Phenanthrene	UG/L	-	8 J	10 U	10 U	14	10 U
Pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Total Polynuclear Aromatic Hydrocarbons	UG/L	100	118	ND	19	133	ND

\*Criteria- Groundwater Plume Delineation/Design Criteria, Pre-Design Investigation Work Plan for In-Situ Solidification for the Hempstead Intersection Street Former MGP Site, Appendix E, Final, URS 2008.

Flags assigned during chemistry validation are shown.

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D - Result reported from a secondary dilution analysis.

Made By\_PRF 01/27/11\_; Checked By\_AMK 01/27/11\_

Detection Limits shown are PQL

**TABLE A-1**  
**VALIDATED GROUNDWATER SAMPLE ANALYTICAL RESULTS**  
**NATIONAL GRID - HEMPSTEAD INTERSECTION STREET FORMER MGP SITE**

Location ID			HIMW-014D	HIMW-014I	HIMW-015D	HIMW-015I	HIMW-020I
Sample ID			HIMW-14D	HIMW-14I	HIMW-15D	HIMW-15I	DUP-102510
Matrix			Groundwater	Groundwater	Groundwater	Groundwater	Groundwater
Depth Interval (ft)			-	-	-	-	-
Date Sampled			10/21/10	10/21/10	10/22/10	10/22/10	10/25/10
Parameter	Units	Criteria*					Field Duplicate (1-1)
<b>Volatile Organic Compounds</b>							
Benzene	UG/L	-	1 U	19	1 U	22	47
Ethylbenzene	UG/L	-	1 U	2	1 U	1 U	3
Toluene	UG/L	-	1 U	1 U	1 U	1 U	2
Xylene (total)	UG/L	-	1 U	3	1 U	2	130
Total BTEX	UG/L	100	ND	24	ND	24	182
<b>Semivolatile Organic Compounds</b>							
2-Methylnaphthalene	UG/L	-	10 U	10 U	10 U	10 U	26
Acenaphthene	UG/L	-	10 U	15	10 U	6 J	16
Acenaphthylene	UG/L	-	10 U	21	10 U	24	210 D
Anthracene	UG/L	-	10 U	10 U	10 U	10 U	4 J
Benzo(a)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(a)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(b)fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Benzo(k)fluoranthene	UG/L	-	10 UJ	10 UJ	10 UJ	10 UJ	10 UJ
Chrysene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)anthracene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluoranthene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Fluorene	UG/L	-	10 U	8 J	10 U	10 U	34
Indeno(1,2,3-cd)pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Naphthalene	UG/L	-	10 U	1 J	10 U	10 U	120 D
Phenanthrene	UG/L	-	10 U	6 J	10 U	10 U	28
Pyrene	UG/L	-	10 U	10 U	10 U	10 U	10 U
Total Polynuclear Aromatic Hydrocarbons	UG/L	100	ND	51	ND	30	438

\*Criteria- Groundwater Plume Delineation/Design Criteria, Pre-Design Investigation Work Plan for In-Situ Solidification for the Hempstead Intersection Street Former MGP Site, Appendix E, Final, URS 2008.

Flags assigned during chemistry validation are shown.

 Concentration Exceeds Criteria

U - Not detected above the reported quantitation limit. J - The reported concentration is an estimated value.

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D - Result reported from a secondary dilution analysis.

Made By\_PRF 01/27/11\_; Checked By\_AMK 01/27/11\_

Detection Limits shown are PQL

**TABLE A-1**  
**VALIDATED GROUNDWATER SAMPLE ANALYTICAL RESULTS**  
**NATIONAL GRID - HEMPSTEAD INTERSECTION STREET FORMER MGP SITE**

Location ID			HIMW-020I	HIMW-020S
Sample ID			HIMW-20I	HIMW-20S
Matrix			Groundwater	Groundwater
Depth Interval (ft)			-	-
Date Sampled			10/25/10	10/26/10
Parameter	Units	Criteria*		
<b>Volatile Organic Compounds</b>				
Benzene	UG/L	-	46	1 U
Ethylbenzene	UG/L	-	3	1 U
Toluene	UG/L	-	2	1 U
Xylene (total)	UG/L	-	130	1 U
Total BTEX	UG/L	100	181	ND
<b>Semivolatile Organic Compounds</b>				
2-Methylnaphthalene	UG/L	-	26	10 U
Acenaphthene	UG/L	-	16	10 U
Acenaphthylene	UG/L	-	200 D	10 U
Anthracene	UG/L	-	4 J	10 U
Benzo(a)anthracene	UG/L	-	10 U	10 U
Benzo(a)pyrene	UG/L	-	10 U	10 U
Benzo(b)fluoranthene	UG/L	-	10 U	10 U
Benzo(g,h,i)perylene	UG/L	-	10 U	10 U
Benzo(k)fluoranthene	UG/L	-	10 UJ	10 UJ
Chrysene	UG/L	-	10 U	10 U
Dibenz(a,h)anthracene	UG/L	-	10 U	10 U
Fluoranthene	UG/L	-	10 U	10 U
Fluorene	UG/L	-	34	10 U
Indeno(1,2,3-cd)pyrene	UG/L	-	10 U	10 U
Naphthalene	UG/L	-	110 D	10 U
Phenanthrene	UG/L	-	29	10 U
Pyrene	UG/L	-	10 U	10 U
Total Polynuclear Aromatic Hydrocarbons	UG/L	100	419	ND

\*Criteria- Groundwater Plume Delineation/Design Criteria, Pre-Design Investigation Work Plan for In-Situ Solidification for the Hempstead Intersection Street Former MGP Site, Appendix E, Final, URS 2008.

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D - Result reported from a secondary dilution analysis.

Made By\_PRF 01/27/11\_; Checked By\_AMK 01/27/11\_

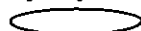
Detection Limits shown are PQL

**TABLE A-2**  
**VALIDATED FIELD QC SAMPLE ANALYTICAL RESULTS**  
**NATIONAL GRID - HEMPSTEAD INTERSECTION STREET FORMER MGP SITE**

Location ID			FIELDQC	FIELDQC	FIELDQC
Sample ID			TB 102110	TB-102510	TB-102710
Matrix			Water Quality	Water Quality	Water Quality
Depth Interval (ft)			-	-	-
Date Sampled			10/21/10	10/25/10	10/27/10
Parameter	Units	Criteria*	Trip Blank (1-1)	Trip Blank (1-1)	Trip Blank (1-1)
<b>Volatile Organic Compounds</b>					
Benzene	UG/L	-	1 U	1 U	1 U
Ethylbenzene	UG/L	-	1 U	1 U	1 U
Toluene	UG/L	-	1 U	1 U	1 U
Xylene (total)	UG/L	-	1 U	1 U	1 U
Total BTEX	UG/L	100	ND	ND	ND

\*Criteria- Groundwater Plume Delineation/Design Criteria, Pre-Design Investigation Work Plan for In-Situ Solidification for the Hempstead Intersection Street Former MGP Site, Appendix E, Final, URS 2008.

Flags assigned during chemistry validation are shown.



Concentration Exceeds Criteria

U - Not detected above the reported quantitation limit.

Made By\_PRF 01/27/11\_; Checked By\_AMK 01/27/11\_

Detection Limits shown are PQL

**APPENDIX A**

**VALIDATED FORM 1'S**

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-131

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010820-001A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0423.D  
 Level: (low/med) LOW Date Received: 10/22/10  
 % Moisture: not dec. Date Analyzed: 10/26/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	96	
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	7	

KEY-URS104 S31

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-13D

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010820-002A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0424.D  
 Level: (low/med) LOW Date Received: 10/22/10  
 % Moisture: not dec. Date Analyzed: 10/26/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	Q
71-43-2	Benzene	4	
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	2	



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-13S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water)

WATER

Lab Sample ID: 1010820-003A

Sample wt/vol: 5

(g/mL) ML

Lab File ID: 10\J0425.D

Level: (low/med)

LOW

Date Received: 10/22/10

% Moisture: not dec.

Date Analyzed: 10/26/10

GC Column: Rtx-624

ID: .18 (mm)

Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_

( $\mu$ L)

Soil Aliquot Volume \_\_\_\_\_ ( $\mu$ L)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(pg/L or $\mu$ g/Kg) <u>UG/L</u>	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-14D

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010820-004A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0426.D

Level: (low/med) LOW Date Received: 10/22/10

% Moisture: not dec. Date Analyzed: 10/26/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

KEY-URS104 S34

1A

EPA SAMPLE NO.

## VOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-141

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010820-005A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0427.D

Level: (low/med) LOW Date Received: 10/22/10

% Moisture: not dec. Date Analyzed: 10/26/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	19	
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	2	
1330-20-7	Xylene (total)	3	

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-15D

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010820-006A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0428.D  
 Level: (low/med) LOW Date Received: 10/22/10  
 % Moisture: not dec. Date Analyzed: 10/26/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-15I

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010820-007A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0429.D

Level: (low/med) LOW Date Received: 10/22/10

% Moisture: not dec. Date Analyzed: 10/26/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	22	
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	2	

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

TB 102110

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010820-008A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0418.D

Level: (low/med) LOW Date Received: 10/22/10

% Moisture: not dec. Date Analyzed: 10/25/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5D

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
Matrix: (soil/water) WATER Lab Sample ID: 1010958-001A  
Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0437.D  
Level: (low/med) LOW Date Received: 10/26/10  
% Moisture: not dec. Date Analyzed: 10/26/10  
GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	6	
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	210	

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5I

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010958-002A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0438.D

Level: (low/med) LOW Date Received: 10/26/10

% Moisture: not dec. Date Analyzed: 10/26/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	3	
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	
1330-20-7	Xylene (total)	150	



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-8D

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010958-003A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0439.D  
 Level: (low/med) LOW Date Received: 10/26/10  
 % Moisture: not dec. Date Analyzed: 10/26/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

KEY-URS104 S41

1A

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-12D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water)

WATERLab Sample ID: 1010958-004ASample wt/vol: 5(g/mL) MLLab File ID: 10\J0440.D

Level: (low/med)

LOWDate Received: 10/26/10

% Moisture: not dec.

Date Analyzed: 10/26/10GC Column: Rtx-624ID: .18 (mm)Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_

(µL)

Soil Aliquot Volume \_\_\_\_\_

(µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	<u>Q</u>
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-12I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS SAS No.: \_\_\_\_\_SDG No.: KEY-URS104

Matrix: (soil/water)

WATERLab Sample ID: 1010958-005ASample wt/vol: 5(g/mL) MLLab File ID: 10\J0441.D

Level: (low/med)

LOWDate Received: 10/26/10

% Moisture: not dec.

Date Analyzed: 10/26/10GC Column: Rtx-624ID: .18 (mm)Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_

(µL)

Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	Q
71-43-2	Benzene	48	
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	4	

KEY-URS104 S43

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-12S

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010958-006A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0442.D

Level: (low/med) LOW Date Received: 10/26/10

% Moisture: not dec. Date Analyzed: 10/26/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	6	

KEY-URS104 S44

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-20I

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010958-007A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0443.D  
 Level: (low/med) LOW Date Received: 10/26/10  
 % Moisture: not dec. Date Analyzed: 10/26/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	46	
108-88-3	Toluene	2	
100-41-4	Ethylbenzene	3	
1330-20-7	Xylene (total)	130	

KEY-URS104 S45

1A

EPA SAMPLE NO.

## VOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-20S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS SAS No.: \_\_\_\_\_SDG No.: KEY-URS104

Matrix: (soil/water)

WATERLab Sample ID: 1010958-008ASample wt/vol: 5(g/mL) MLLab File ID: 10\J0444.D

Level: (low/med)

LOWDate Received: 10/26/10

% Moisture: not dec.

Date Analyzed: 10/26/10GC Column: Rtx-624ID: .18 (mm)Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_

(µL)

Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

DUP-102510

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS SAS No.: \_\_\_\_\_SDG No.: KEY-URS104

Matrix: (soil/water)

WATERLab Sample ID: 1010958-009ASample wt/vol: 5(g/mL) MLLab File ID: 10\J0445.D

Level: (low/med)

LOWDate Received: 10/26/10

% Moisture: not dec.

Date Analyzed: 10/26/10GC Column: Rtx-624ID: .18 (mm)Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_

(µL)

Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	47	
108-88-3	Toluene	2	
100-41-4	Ethylbenzene	3	
1330-20-7	Xylene (total)	130	

## VOLATILE ORGANICS ANALYSIS DATA SHEET

TB-102510

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010958-010A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0447.D

Level: (low/med) LOW Date Received: 10/25/10

% Moisture: not dec. Date Analyzed: 10/27/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	<u>Q</u>
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U



## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-3D

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010A20-001A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0548.D

Level: (low/med) LOW Date Received: 10/28/10

% Moisture: not dec. Date Analyzed: 11/09/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-3I

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010A20-002A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0549.D  
 Level: (low/med) LOW Date Received: 10/28/10  
 % Moisture: not dec. Date Analyzed: 11/09/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-3S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water)

WATERLab Sample ID: 1010A20-003ASample wt/vol: 5(g/mL) MLLab File ID: 10\J0550.D

Level: (low/med)

LOWDate Received: 10/28/10

% Moisture: not dec.

Date Analyzed: 11/09/10GC Column: Rtx-624ID: .18 (mm)Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL)

Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	<u>Q</u>
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5S

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Matrix: (soil/water) WATER Lab Sample ID: 1010A20-004A  
 Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0551.D  
 Level: (low/med) LOW Date Received: 10/28/10  
 % Moisture: not dec. Date Analyzed: 11/09/10  
 GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00  
 Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-8I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water)

WATER

Lab Sample ID: 1010A20-005A

Sample wt/vol: 5

(g/mL) ML

Lab File ID: 10\J0552.D

Level: (low/med)

LOW

Date Received: 10/28/10

% Moisture: not dec.

Date Analyzed: 11/09/10

GC Column: Rtx-624

ID: .18 (mm)

Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL)

Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-8S

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010A20-006A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0553.D

Level: (low/med) LOW Date Received: 10/28/10

% Moisture: not dec. Date Analyzed: 11/09/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

TB-102710

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010A20-007A

Sample wt/vol: 5 (g/mL) ML Lab File ID: 10\J0547.D

Level: (low/med) LOW Date Received: 10/28/10

% Moisture: not dec. Date Analyzed: 11/09/10

GC Column: Rtx-624 ID: .18 (mm) Dilution Factor: 1.00

Soil Extract Volume: \_\_\_\_\_ (µL) Soil Aliquot Volume \_\_\_\_\_ (µL)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) <u>UG/L</u>	<u>Q</u>
71-43-2	Benzene	1	U
108-88-3	Toluene	1	U
100-41-4	Ethylbenzene	1	U
1330-20-7	Xylene (total)	1	U

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-13I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010820-001BSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1539.DLevel: (low/med) LOWDate Received: 10/22/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	1	J
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	<del>84</del> 90	<del>ED</del>
83-32-9	Acenaphthene	11	
86-73-7	Fluorene	15	
85-01-8	Phenanthrene	14	
120-12-7	Anthracene	2	J
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
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## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-13IDL

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010820-001BDLSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1550.DLevel: (low/med) LOWDate Received: 10/22/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 2.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	20	U
91-57-6	2-Methylnaphthalene	20	U
208-96-8	Acenaphthylene	90	D
83-32-9	Acenaphthene	11	DJ
86-73-7	Fluorene	16	DJ
85-01-8	Phenanthrene	14	DJ
120-12-7	Anthracene	20	U
206-44-0	Fluoranthene	20	U
129-00-0	Pyrene	20	U
56-55-3	Benzo(a)anthracene	20	U
218-01-9	Chrysene	20	U
205-99-2	Benzo(b)fluoranthene	20	U
207-08-9	Benzo(k)fluoranthene	20	U
50-32-8	Benzo(a)pyrene	20	U
193-39-5	Indeno(1,2,3-cd)pyrene	20	U
53-70-3	Dibenzo(a,h)anthracene	20	U
191-24-2	Benzo(g,h,i)perylene	20	U

(1) Cannot be separated from Diphenylamine

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-13D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010820-002B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1540.D

Level: (low/med) LOW

Date Received: 10/22/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 (µL)

Date Analyzed: 10/29/10

Injection Volume: 2 (µL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	13	
83-32-9	Acenaphthene	6	J
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11

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## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-13S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010820-003BSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1541.DLevel: (low/med) LOWDate Received: 10/22/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 (μL)Date Analyzed: 10/29/10Injection Volume: 2 (μL)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(μg/L or μg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U <sup>J</sup>
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-14D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010820-004BSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1542.DLevel: (low/med) LOWDate Received: 10/22/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U <sup>J</sup>
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
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EPA SAMPLE NO.

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-14I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010820-005BSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1543.DLevel: (low/med) LOWDate Received: 10/22/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	1	J
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	21	
83-32-9	Acenaphthene	15	
86-73-7	Fluorene	8	J
85-01-8	Phenanthrene	6	J
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
2

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-15D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS SAS No.: \_\_\_\_\_SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010820-006BSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1546.DLevel: (low/med) LOWDate Received: 10/22/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) <u>UG/L</u>	<u>Q</u>
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
2

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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-151

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010820-007B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1547.D

Level: (low/med) LOW

Date Received: 10/22/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 (µL)

Date Analyzed: 10/29/10

Injection Volume: 2 (µL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	24	
83-32-9	Acenaphthene	6	J
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
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SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010958-001B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1553.D

Level: (low/med) LOW

Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 (μL)

Date Analyzed: 10/29/10

Injection Volume: 2 (μL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(μg/L or μg/Kg) UG/L	Q
91-20-3	Naphthalene	<del>730</del> 1400	<del>ED</del>
91-57-6	2-Methylnaphthalene	<del>220</del> 270	<del>ED</del>
208-96-8	Acenaphthylene	51	
83-32-9	Acenaphthene	4	J
86-73-7	Fluorene	3	J
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
2



## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-5DDL

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010958-001BDLSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1572.DLevel: (low/med) LOWDate Received: 10/26/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/31/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 20.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) <u>UG/L</u>	Q
91-20-3	Naphthalene	1400	D
91-57-6	2-Methylnaphthalene	270	D
208-96-8	Acenaphthylene	64	DJ
83-32-9	Acenaphthene	200	U
86-73-7	Fluorene	200	U
85-01-8	Phenanthrene	200	U
120-12-7	Anthracene	200	U
206-44-0	Fluoranthene	200	U
129-00-0	Pyrene	200	U
56-55-3	Benzo (a) anthracene	200	U
218-01-9	Chrysene	200	U
205-99-2	Benzo (b) fluoranthene	200	U
207-08-9	Benzo (k) fluoranthene	200	U
50-32-8	Benzo (a) pyrene	200	U
193-39-5	Indeno (1, 2, 3-cd) pyrene	200	U
53-70-3	Dibenzo (a, h) anthracene	200	U
191-24-2	Benzo (g, h, i) perylene	200	U

(1) Cannot be separated from Diphenylamine

1/26/11

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010958-002B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1554.D

Level: (low/med) LOW

Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 (μL)

Date Analyzed: 10/29/10

Injection Volume: 2 (μL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(μg/L or μg/Kg)	UG/L	Q
91-20-3	Naphthalene	<del>1200</del>	2300	<del>F/D</del>
91-57-6	2-Methylnaphthalene	<del>400</del>	560	<del>F/D</del>
208-96-8	Acenaphthylene	<del>160</del>	230	<del>F/D</del>
83-32-9	Acenaphthene		16	
86-73-7	Fluorene		23	
85-01-8	Phenanthrene		20	
120-12-7	Anthracene		3	J
206-44-0	Fluoranthene		10	U
129-00-0	Pyrene		10	U
56-55-3	Benzo(a)anthracene		10	U
218-01-9	Chrysene		10	U
205-99-2	Benzo(b)fluoranthene		10	U
207-08-9	Benzo(k)fluoranthene		10	U J
50-32-8	Benzo(a)pyrene		10	U
193-39-5	Indeno(1,2,3-cd)pyrene		10	U
53-70-3	Dibenzo(a,h)anthracene		10	U
191-24-2	Benzo(g,h,i)perylene		10	U

(1) Cannot be separated from Diphenylamine

1/26/11

1C

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5IDL

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010958-002BDLSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1573.DLevel: (low/med) LOWDate Received: 10/26/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/31/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 50.00GPC Cleanup: (Y/N) N

pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/kg) UG/L	Q
91-20-3	Naphthalene	2300	D
91-57-6	2-Methylnaphthalene	560	D
208-96-8	Acenaphthylene	230	DJ
83-32-9	Acenaphthene	500	U
86-73-7	Fluorene	500	U
85-01-8	Phenanthrene	500	U
120-12-7	Anthracene	500	U
206-44-0	Fluoranthene	500	U
129-00-0	Pyrene	500	U
56-55-3	Benzo(a)anthracene	500	U
218-01-9	Chrysene	500	U
205-99-2	Benzo(b)fluoranthene	500	U
207-08-9	Benzo(k)fluoranthene	500	U
50-32-8	Benzo(a)pyrene	500	U
193-39-5	Indeno(1,2,3-cd)pyrene	500	U
53-70-3	Dibenzo(a,h)anthracene	500	U
191-24-2	Benzo(g,h,i)perylene	500	U

(1) Cannot be separated from Diphenylamine

1/26/11

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-8D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010958-003B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1555.D

Level: (low/med) LOW

Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 ( $\mu$ L)

Date Analyzed: 10/29/10

Injection Volume: 2 ( $\mu$ L)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11

1C

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-12D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010958-004BSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1556.DLevel: (low/med) LOWDate Received: 10/26/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) <u>UG/L</u>	<u>Q</u>
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-121

Lab Name: H2M LABS INC Contract: \_\_\_\_\_

Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104

Matrix: (soil/water) WATER Lab Sample ID: 1010958-005B

Sample wt/vol: 1000 (g/mL) ML Lab File ID: 10\R1557.D

Level: (low/med) LOW Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 (µL) Date Analyzed: 10/29/10

Injection Volume: 2 (µL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_ Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
91-20-3	Naphthalene	2	J
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	43	
83-32-9	Acenaphthene	40	
86-73-7	Fluorene	25	
85-01-8	Phenanthrene	8	J
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo (a) anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo (b) fluoranthene	10	U
207-08-9	Benzo (k) fluoranthene	10	U J
50-32-8	Benzo (a) pyrene	10	U
193-39-5	Indeno (1, 2, 3-cd) pyrene	10	U
53-70-3	Dibenzo (a, h) anthracene	10	U
191-24-2	Benzo (g, h, i) perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
2

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-12S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010958-006B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1558.D

Level: (low/med) LOW

Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 (μL)

Date Analyzed: 10/29/10

Injection Volume: 2 (μL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(μg/L or μg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
2

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-20I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010958-007B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1559.D

Level: (low/med) LOW

Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 ( $\mu$ L)

Date Analyzed: 10/29/10

Injection Volume: 2 ( $\mu$ L)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	<del>87</del> 110	<del>ED</del>
91-57-6	2-Methylnaphthalene	26	
208-96-8	Acenaphthylene	<del>150</del> 200	<del>ED</del>
83-32-9	Acenaphthene	16	
86-73-7	Fluorene	34	
85-01-8	Phenanthrene	29	
120-12-7	Anthracene	4	J
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/4  
2



1C

EPA SAMPLE NO.

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-20IDL

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010958-007BDLSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1574.DLevel: (low/med) LOWDate Received: 10/26/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/31/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 5.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

( $\mu$ g/L or  $\mu$ g/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: ( $\mu$ g/L or $\mu$ g/Kg) <u>UG/L</u>	Q
91-20-3	Naphthalene	110	D
91-57-6	2-Methylnaphthalene	30	DJ
208-96-8	Acenaphthylene	200	D
83-32-9	Acenaphthene	17	DJ
86-73-7	Fluorene	37	DJ
85-01-8	Phenanthrene	31	DJ
120-12-7	Anthracene	50	U
206-44-0	Fluoranthene	50	U
129-00-0	Pyrene	50	U
56-55-3	Benzo(a)anthracene	50	U
218-01-9	Chrysene	50	U
205-99-2	Benzo(b)fluoranthene	50	U
207-08-9	Benzo(k)fluoranthene	50	U
50-32-8	Benzo(a)pyrene	50	U
193-39-5	Indeno(1,2,3-cd)pyrene	50	U
53-70-3	Dibenzo(a,h)anthracene	50	U
191-24-2	Benzo(g,h,i)perylene	50	U

(1) Cannot be separated from Diphenylamine

1/26/14

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-20S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS SAS No.: \_\_\_\_\_SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010958-008ESample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1560.DLevel: (low/med) LOWDate Received: 10/26/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/29/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo (a) anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo (b) fluoranthene	10	U
207-08-9	Benzo (k) fluoranthene	10	U J
50-32-8	Benzo (a) pyrene	10	U
193-39-5	Indeno (1, 2, 3-cd) pyrene	10	U
53-70-3	Dibenzo (a, h) anthracene	10	U
191-24-2	Benzo (g, h, i) perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
2

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

DUP-102510

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010958-009B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 10\R1561.D

Level: (low/med) LOW

Date Received: 10/26/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 10/28/10

Concentrated Extract Volume: 1000 ( $\mu$ L)

Date Analyzed: 10/29/10

Injection Volume: 2 ( $\mu$ L)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	<del>85</del> 120	<del>ED</del>
91-57-6	2-Methylnaphthalene	26	
208-96-8	Acenaphthylene	<del>158</del> 210	<del>ED</del>
83-32-9	Acenaphthene	16	
86-73-7	Fluorene	34	
85-01-8	Phenanthrene	28	
120-12-7	Anthracene	4	J
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U J
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1/26/11  
02

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

DUP-102510DL

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010958-009BDLSample wt/vol: 1000 (g/mL) MLLab File ID: 10\R1575.DLevel: (low/med) LOWDate Received: 10/26/10% Moisture: Decanted: (Y/N) NDate Extracted: 10/28/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 10/31/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 5.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg)	UG/L	Q
91-20-3	Naphthalene	120		D
91-57-6	2-Methylnaphthalene	32		DJ
208-96-8	Acenaphthylene	210		D
83-32-9	Acenaphthene	19		DJ
86-73-7	Fluorene	41		DJ
85-01-8	Phenanthrene	33		DJ
120-12-7	Anthracene	50		U
206-44-0	Fluoranthene	50		U
129-00-0	Pyrene	50		U
56-55-3	Benzo(a)anthracene	50		U
218-01-9	Chrysene	50		U
205-99-2	Benzo(b)fluoranthene	50		U
207-08-9	Benzo(k)fluoranthene	50		U
50-32-8	Benzo(a)pyrene	50		U
193-39-5	Indeno(1,2,3-cd)pyrene	50		U
53-70-3	Dibenzo(a,h)anthracene	50		U
191-24-2	Benzo(g,h,i)perylene	50		U

(1) Cannot be separated from Diphenylamine

1/26/11

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-3D

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010A20-001B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: A\C56695.D

Level: (low/med) LOW

Date Received: 10/28/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 11/01/10

Concentrated Extract Volume: 1000 (µL)

Date Analyzed: 11/04/10

Injection Volume: 2 (µL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-3I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010A20-002B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: A\C56696.D

Level: (low/med) LOW

Date Received: 10/28/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 11/01/10

Concentrated Extract Volume: 1000 (µL)

Date Analyzed: 11/04/10

Injection Volume: 2 (µL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-3S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010A20-003B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: A\C56697.D

Level: (low/med) LOW

Date Received: 10/28/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 11/01/10

Concentrated Extract Volume: 1000 ( $\mu$ L)

Date Analyzed: 11/04/10

Injection Volume: 2 ( $\mu$ L)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-5S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010A20-004B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: A\C56698.D

Level: (low/med) LOW

Date Received: 10/28/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 11/01/10

Concentrated Extract Volume: 1000 (µL)

Date Analyzed: 11/04/10

Injection Volume: 2 (µL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(µg/L or µg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo (a) anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo (b) fluoranthene	10	U
207-08-9	Benzo (k) fluoranthene	10	U
50-32-8	Benzo (a) pyrene	10	U
193-39-5	Indeno (1,2,3-cd) pyrene	10	U
53-70-3	Dibenzo (a, h) anthracene	10	U
191-24-2	Benzo (g, h, i) perylene	10	U

(1) Cannot be separated from Diphenylamine



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

HIMW-8I

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2M

Case No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104

Matrix: (soil/water) WATER

Lab Sample ID: 1010A20-005B

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: A\C56699.D

Level: (low/med) LOW

Date Received: 10/28/10

% Moisture: Decanted: (Y/N) N

Date Extracted: 11/01/10

Concentrated Extract Volume: 1000 (μL)

Date Analyzed: 11/04/10

Injection Volume: 2 (μL)

Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: \_\_\_\_\_

Extraction: (Type) SEPF

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(μg/L or μg/Kg) UG/L	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	U
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a)anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b)fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U

(1) Cannot be separated from Diphenylamine

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

HIMW-8S

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Matrix: (soil/water) WATERLab Sample ID: 1010A20-006BSample wt/vol: 1000 (g/mL) MLLab File ID: A\C56700.DLevel: (low/med) LOWDate Received: 10/28/10% Moisture: Decanted: (Y/N) NDate Extracted: 11/01/10Concentrated Extract Volume: 1000 ( $\mu$ L)Date Analyzed: 11/04/10Injection Volume: 2 ( $\mu$ L)Dilution Factor: 1.00GPC Cleanup: (Y/N) N pH: \_\_\_\_\_Extraction: (Type) SEPF

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	( $\mu$ g/L or $\mu$ g/Kg) <u>UG/L</u>	Q
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	1	J
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo (a) anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo (b) fluoranthene	10	U
207-08-9	Benzo (k) fluoranthene	10	U
50-32-8	Benzo (a) pyrene	10	U
193-39-5	Indeno (1,2,3-cd) pyrene	10	U
53-70-3	Dibenzo (a,h) anthracene	10	U
191-24-2	Benzo (g,h,i) perylene	10	U

(1) Cannot be separated from Diphenylamine

**APPENDIX B**

**SUPPORT DOCUMENTATION**

# H2M LABS, INC.

575 Broad Hollow Rd, Melville, NY 11747-5076  
 Tel: (631) 694-3040 Fax: (631) 420-8436

PROJECT NAME/NUMBER

National Grid - MGP

SAMPLERS: (signature)/Client

*Cheryl / URS*

DELIVERABLES:

Full Cat B

TURNAROUND TIME:

Stand

00000

## EXTERNAL CHAIN OF CUSTODY

KEY-URS104

CLIENT: URS H2M SDG NO: URS125

Project Contact:  
 Peter Fairbanks  
 Phone Number:  
 716 856 5636  
 FIS/Quote #

NOTES:

Sample Container Description	ANALYSIS REQUESTED			
	ORGANIC	INORG.	PCB	Metal
2-40 ml Amber	X			
2-12 Amber glass	X			

Sample Container Description	Total No. of Containers	ANALYSIS REQUESTED				LAB I.D. NO.	REMARKS:
		ORGANIC	INORG.	PCB	Metal		
	4	X				1010820-003	
	5	X				-002-001	
	5	X				-001-002	
	7	X				-005	
	7	X				-004	
	7	X				-007	
	4	X				-006	
	8	X				-007	
	2	X				-008	

Retrieved by: (Signature)	Date	Time	Received by: (Signature)	Date	Time
<i>Cheryl</i>	10/22/10	14:05	<i>URS</i>	10/22/10	14:05
<i>Cheryl</i>	10/22/10	14:50	<i>URS</i>	10/22/10	14:50

### LABORATORY USE ONLY

Discrepancies Between Sample Labels and COC Record? Y or N  
 Explain:

Samples were:  
 1. Shipped  and Delivered  Ambient  or Chilled  Temp.  or  Ambient   
 2. Received in good condition  Y or N  
 3. Properly preserved  Y or N

COC Tada was:  
 1. Present on outer package: Y or N   
 2. Unbroken on outer package: Y or N   
 3. COC record present & complete upon sample receipt: Y or N

WHITE COPY (ORIGINAL)

YELLOW COPY - CLIENT

PINK COPY - LABORATORY

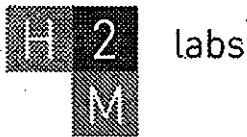
# H2M LABS, INC.

575 Broad Hollow Rd, Melville, NY 11747-5076  
Tel: (631) 694-3040 Fax: (631) 420-8436

50050

# EXTERNAL CHAIN OF CUSTODY

PROJECT NAME/NUMBER <b>National Blvd - Hempstead Map</b> 1176098		CLIENT: <b>URS</b>		H2M SDG NO: <b>KET-VPS104</b>	
SAMPLERS: (signature) Client <i>[Signature]</i>		Project Contact <b>Peter Fairbanks</b>		Phone Number <b>716 856 5630</b>	
DELIVERABLES: <b>Full Cat B</b>		TURNAROUND TIME: <b>Std</b>		PIS/Quote #	
DATE		TIME MATRIX		FIELD I.D.	
10/25/10	815	6W		HIMW-5I	
10/25/10	715			HIMW-5D	
10/25/10	1155			HIMW-12S	
10/25/10	1040			HIMW-12I	
10/25/10	0900			HIMW-12D	
10/25/10	1005			HIMW-20S	
10/25/10	1400	✓		HIMW-20I	
10/25/10	1200	6W		sup-102510	
10/26/10	1150	6W		HIMW-8D	
10/27/10		DI		TB	
Relinquished by: (Signature) <i>[Signature]</i>		Date <b>10/25/10</b>		Time <b>1352</b>	
Relinquished by: (Signature) <i>[Signature]</i>		Date <b>10/26/10</b>		Time <b>1115</b>	
Relinquished by: (Signature) <i>[Signature]</i>		Date <b>10/26/10</b>		Time <b>1415</b>	
Relinquished by: (Signature)		Date		Time	
Sample Container Description <b>40-ml Amber HCL</b>		Sample Container		Total No. of Containers	
ANALYSIS REQUESTED		ORGANIC		INORG.	
		BTEX		Metal	
		PAHs		Zn	
		PCBs			
		Pb			
		Cu			
		Cd			
		Mn			
		Fe			
		Ni			
		Co			
		Cr			
		Mg			
		Ca			
		K			
		Na			
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		Br			
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		Se			
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		Ag			
		Au			
		Pt			
		Ba			
		Sr			
		Zr			
		Nb			
		Ta			
		V			
		Cr			
		Mn			
		Fe			
		Co			
		Ni			
		Cu			
		Zn			
		Pb			
		Cd			
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		Ba			
		Sr			



H2M LABS INC  
 575 Broad Hollow Road  
 Melville, NY 11747  
 TEL: 631-694-3040 FAX: 631-420-8436  
 Website: www.h2mlabs.com

Key-URS104

Sample Receipt Checklist

Client Name: KEY-URS

Date and Time Received: 10/26/2010 4:45:00 PM

Work Order Number 1010958

RcptNo: 1

Received by: Melissa Watson

Completed by: *M. Watson*

Reviewed by: *JST*

Completed Date: 10-26-10

Reviewed Date: 10/28/10

Carrier name: H2M Pickup

- Chain of custody present? Yes  No
- Chain of custody signed when relinquished and received? Yes  No
- Chain of custody agrees with sample labels? Yes  No
- Are matrices correctly identified on Chain of custody? Yes  No
- Is it clear what analyses were requested? Yes  No
- Custody seals intact on sample bottles? Yes  No  Not Present
- Samples in proper container/bottle? Yes  No
- Were correct preservatives used and noted? Yes  No
- Sample containers intact? Yes  No
- Sufficient sample volume for indicated test? Yes  No
- Were container labels complete (ID, Pres, Date)? Yes  No
- All samples received within holding time? Yes  No
- Was an attempt made to cool the samples? Yes  No
- All samples received at a temp. of > 0° C to 6.0° C? Yes  No
- Response when temperature is outside of range:  
 Preservative added to bottles:
- Sample Temp. taken and recorded upon receipt? Yes  No  4 To 10°
- Water - Were bubbles absent in VOC vials? Yes  No  No Vials
- Water - Was there Chlorine Present? Yes  No  NA
- Water - pH acceptable upon receipt? Yes  No  No Water
- Are Samples considered acceptable? Yes  No
- Custody Seals present? Yes  No
- Traffic Report or Packing Lists present? Yes  No
- Airbill or Sticker? Air Bill  Sticker  Not Present
- Airbill No:
- Sample Tags Present? Yes  No
- Sample Tags Listed on COC? Yes  No
- Tag Numbers:
- Sample Condition? Intact  Broken  Leaking

Case Number: SDG: KEY-URS104

SAS:

Adjusted? \_\_\_\_\_ Checked by \_\_\_\_\_

Any No and/or NA (not applicable) response must be detailed in the comments section below.



labs

H2M LABS INC  
575 Broad Hollow Road  
Melville, NY 11747  
TEL: 631-694-3040 FAX: 631-420-8436  
Website: www.h2mlabs.com

## Sample Receipt Checklist

Client Contacted?  Yes  No

Person Contacted:

Comments:

Contact Mode:  Phone:  Fax:  Email:  In Person:

There was ice present in the coolers, but one cooler was greater than 6 degrees celcius.

Client Instructions:

Date Contacted:

Contacted By:

Regarding:

Corrective Action:

# H2M LABS, INC.

575 Broad Hollow Rd, Melville, NY 11747-5076  
 Tel: (631) 694-3040 Fax: (631) 420-8436

35655

## EXTERNAL CHAIN OF CUSTODY

PROJECT NAME/NUMBER National Grid Hempstead, NY 1176098.00004		CLIENT: URS Corporation		H2M SDG NO: KEY-VRS104	
SAMPLERS: (signature) Client Cory Friedman		NOTES:		Project Contact: Peter Fairbanks Phone Number: 716 856 5636 PIS/Quote #	
DELIVERABLES: Full Cat B		ANALYSIS REQUESTED		REMARKS:	
TURNAROUND TIME: Standard		ORGANIC		LAB I.D. NO.	
DATE	TIME	MATRIX	FIELD I.D.	LAB I.D. NO.	REMARKS:
10/28/10	1815	GW	H1MW-35	1010A7D-001	-003
10/27/10	1300	GW	H1MW-3I	-002	-002
10/27/10	1035	GW	H1MW-3D	-003	-001
10/27/10	0715	GW	H1MW-5S	-004	
10/28/10	0955	GW	H1MW-8I	-005	
10/28/10	1050	GW	H1MW-8S	-006	
10/27/10		DI	TB		
Relinquished by: (Signature)	Time	Date	Received by: (Signature)	Date	Time
<i>[Signature]</i>	1250	10/28/10	<i>[Signature]</i>	10/28/10	12:50
Relinquished by: (Signature)	Time	Date	Received by: (Signature)	Date	Time
<i>[Signature]</i>	1:35	10/28/10	<i>[Signature]</i>	10-28-10	13:35
Relinquished by: (Signature)	Time	Date	Received by: (Signature)	Date	Time
<i>[Signature]</i>			<i>[Signature]</i>		
Relinquished by: (Signature)	Time	Date	Received by: (Signature)	Date	Time
<i>[Signature]</i>			<i>[Signature]</i>		

### LABORATORY USE ONLY

Discrepancies Between Sample Labels and COC Record? Y or N  
 Explain:

Samples were:  
 1. Shipped or Hand Delivered ✓  
 2. Ambient or Chilled Temp 9°C on ice  
 3. Received in good condition: Y  
 4. Properly preserved: Y

COC Tags were:  
 1. Present on outer package: Y  
 2. Unbroken on outer package: Y  
 3. COC record present & complete upon sample receipt: Y

WHITE COPY ORIGINAL

YELLOW COPY - CLIENT

PINK COPY - LABORATORY





labs

H2M LABS INC
575 Broad Hollow Road
Melville, NY 11747
TEL: 631-694-3040 FAX: 631-420-8436
Website: www.h2mlabs.com

Key-URS 104

Sample Receipt Checklist

Client Name: KEY-URS

Date and Time Received: 10/28/2010 1:35:00 PM

Work Order Number 1010A20

RcptNo: 1

Received by: MelissaWatson

Completed by: M. Watson

Reviewed by: JSA

Completed Date: 10-28-10

Reviewed Date: 10/29/10

Carrier name: H2M Pickup

- Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Are matrices correctly identified on Chain of custody? Yes [checked] No [ ]
Is it clear what analyses were requested? Yes [checked] No [ ]
Custody seals intact on sample bottles? Yes [ ] No [ ] Not Present [checked]
Samples in proper container/bottle? Yes [checked] No [ ]
Were correct preservatives used and noted? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
Were container labels complete (ID, Pres, Date)? Yes [checked] No [ ]
All samples received within holding time? Yes [checked] No [ ]
Was an attempt made to cool the samples? Yes [checked] No [ ]
All samples received at a temp. of > 0° C to 6.0° C? Yes [ ] No [checked]
Response when temperature is outside of range:
Preservative added to bottles:
Sample Temp. taken and recorded upon receipt? Yes [checked] No [ ] To 9°
Water - Were bubbles absent in VOC vials? Yes [checked] No [ ] No Vials [ ]
Water - Was there Chlorine Present? Yes [ ] No [ ] NA [checked]
Water - pH acceptable upon receipt? Yes [checked] No [ ] No Water [ ]
Are Samples considered acceptable? Yes [checked] No [ ]
Custody Seals present? Yes [ ] No [checked]
Traffic Report or Packing Lists present? Yes [ ] No [checked]
Airbill or Sticker? Air Bill [ ] Sticker [ ] Not Present [checked]
Airbill No:
Sample Tags Present? Yes [ ] No [checked]
Sample Tags Listed on COC? Yes [ ] No [checked]
Tag Numbers:
Sample Condition? Intact [checked] Broken [ ] Leaking [ ]

Case Number: SDG: KEY-URS104

SAS:

Adjusted? \_\_\_\_\_ Checked by \_\_\_\_\_

Any No and/or NA (not applicable) response must be detailed in the comments section below.



labs

H2M LABS INC

575 Broad Hollow Road

Melville, NY 11747

TEL: 631-694-3040 FAX: 631-420-8436

Website: www.h2mlabs.com

# Sample Receipt Checklist

---

Client Contacted?	<input checked="" type="checkbox"/> Yes	<input type="checkbox"/> No	Person Contacted:		Comments:
Contact Mode:	<input type="checkbox"/> Phone:	<input type="checkbox"/> Fax:	<input type="checkbox"/> Email:	<input type="checkbox"/> In Person:	Ice was present in the cooler but the temperature was greater than 6 degrees celcius upon receipt.
Client Instructions:					
Date Contacted:	Contacted By:				
Regarding:					
Corrective Action:					

---

# H2M LABS, INC.

## SDG NARRATIVE FOR VOLATILE ORGANICS SAMPLES RECEIVED: 10/22/10, 10/26/10 & 10/28/10 SDG #: KEY-URS104

For Sample(s):

HIMW-13I	HIMW-5D	DUP-102510	
HIMW-13D	HIMW-5I	TB-102510	
HIMW-13S	HIMW-8D	<del>001A</del>	HIMW-3D
HIMW-14D	HIMW-12D	<del>002B</del>	- 3I
HIMW-14I	HIMW-12I	003A	- 3S
HIMW-15D	HIMW-12S	<del>004B</del>	- 5S
HIMW-15I	HIMW-20I	005A	- 8I
TB 102110	HIMW-20S	006B	- 8S
		007A	TB-102710

1/26/11  
CP

The above sample(s) was/were analyzed for a select list of volatile organic analytes (BTEX) by EPA method 8260B.

All QC data and calibrations met the requirements of the method, unless discussed below, and no problems were encountered with sample analysis. The following should be noted:


Sample HIMW-15I was analyzed as the matrix spike/matrix spike duplicate. All percent recoveries and RPD's were met.

A lab fortified blank was analyzed and indicates good method efficiency.

**I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or his designee, as verified by the following signature.**

Date Reported: November 15, 2010

\*\*\*\*\*  
\*  
\*  
\*\*\*\*\*

  
Joann M. Slavin  
Senior Vice President

KEY-URS104 S25

# H2M LABS, INC.

## SDG NARRATIVE FOR SEMIVOLATILE ORGANICS SAMPLES RECEIVED: 10/22/10, 10/26/10 & 10/28/10 SDG #: KEY-URS104

For Sample(s):

HIMW-13I	HIMW-15I	HIMW-12S	HIMW-5S
HIMW-13D	HIMW-5D	HIMW-20I	HIMW-8I
HIMW-13S	HIMW-5I	HIMW-20S	HIMW-8S
HIMW-14D	HIMW-8D	DUP-102510	HIMW-3S
HIMW-14I	HIMW-12D	HIMW-3D	
HIMW-15D	HIMW-12I	HIMW-3I	

1/26/11  
a

The above sample(s) was/were analyzed for a select list of semivolatile organic analytes (polynuclear aromatics) by EPA method 8270C.

All QC data and calibrations met the requirements of the method unless discussed below, and no problems were encountered with sample analysis. The following should be noted:

Sample HIMW-15I was analyzed as the matrix spike / matrix spike duplicate. All percent recoveries and RPD's were met. Lab fortified blanks were analyzed and indicate good method efficiency.

Sample HIMW-13I, HIMW-5D, HIMW-5I, HIMW-20I and DUP-102510 were reanalyzed at a dilution due to concentration levels of targeted analytes above the calibration range. Samples HIMW-5I, HIMW-12D, HIMW-3I, HIMW-5S and HIMW-8S had surrogate recoveries outside QC limits. All surrogate recoveries were diluted out in the dilution of sample HIMW-5I.

**I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or his designee, as verified by the following signature.**

Date Reported: November 17, 2010

\*\*\*\*\*  
\*  
\*  
\*\*\*\*\*

Joann M. Slavin  
Senior Vice President

KEY-URS104 S26

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Lab File ID: 10\R1519.D DFTPP Injection Date: 10/28/10  
 Instrument ID: HP5973R DFTPP Injection Time: 15:46

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	33.8
68	Less than 2% of mass 69	0.0 (0.0)1
69	Mass 69 relative abundance	45.6
70	Less than 2% of mass 69	0.3 (0.7)1
127	40.0 - 60.0% of mass 198	54.7
197	Less than 1% of mass 198	0.0
198	Base peak, 100% relative abundance	100.0
199	5.0 - 9.0% of mass 198	6.7
275	10.0 - 30.0% of mass 198	20.9
365	Greater than 1% of mass 198	2.1
441	Present, but less than mass 443	6.7
442	40.0 - 110.0% of mass 198	44.8
443	17.0 - 23.0% of mass 442	8.7 (19.5)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	SSTD025	SSTD025	10R1520.D	10/28/10	16:01
02	MB-26985	MB-26985	10R1537.D	10/29/10	0:25
03	LFB-26985	LFB-26985	10R1538.D	10/29/10	0:55
04	HIMW-13I	1010820-001B	10R1539.D	10/29/10	1:25
05	HIMW-13D	1010820-002B	10R1540.D	10/29/10	1:56
06	HIMW-13S	1010820-003B	10R1541.D	10/29/10	2:28
07	HIMW-14D	1010820-004B	10R1542.D	10/29/10	2:58
08	HIMW-14I	1010820-005B	10R1543.D	10/29/10	3:28

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Lab File ID: 10\R1544.D DFTPP Injection Date: 10/29/10  
 Instrument ID: HP5973R DFTPP Injection Time: 11:52

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	31.7
68	Less than 2% of mass 69	0.0 (0.0)1
69	Mass 69 relative abundance	44.0
70	Less than 2% of mass 69	0.2 (0.5)1
127	40.0 - 60.0% of mass 198	54.3
197	Less than 1% of mass 198	0.0
198	Base peak, 100% relative abundance	100.0
199	5.0 - 9.0% of mass 198	6.8
275	10.0 - 30.0% of mass 198	21.8
365	Greater than 1% of mass 198	2.3
441	Present, but less than mass 443	7.7
442	40.0 - 110.0% of mass 198	51.6
443	17.0 - 23.0% of mass 442	9.8 (18.9)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	SSTD025	SSTD025	10R1545.D	10/29/10	12:07
02	HIMW-15D	1010820-006B	10R1546.D	10/29/10	12:36
03	HIMW-15I	1010820-007B	10R1547.D	10/29/10	13:05
04	HIMW-15IMS	1010820-007BMS	10R1548.D	10/29/10	13:35
05	HIMW-15IMSD	1010820-007BMSD	10R1549.D	10/29/10	14:04
06	HIMW-13IDL	1010820-001BDL	10R1550.D	10/29/10	14:33
07	MB-26995	MB-26995	10R1551.D	10/29/10	15:02
08	LFB-26995	LFB-26995	10R1552.D	10/29/10	15:32
09	HIMW-5D	1010958-001B	10R1553.D	10/29/10	16:01
10	HIMW-5I	1010958-002B	10R1554.D	10/29/10	16:30
11	HIMW-8D	1010958-003B	10R1555.D	10/29/10	17:00
12	HIMW-12D	1010958-004B	10R1556.D	10/29/10	17:32
13	HIMW-12I	1010958-005B	10R1557.D	10/29/10	18:01
14	HIMW-12S	1010958-006B	10R1558.D	10/29/10	18:30
15	HIMW-20I	1010958-007B	10R1559.D	10/29/10	19:00
16	HIMW-20S	1010958-008B	10R1560.D	10/29/10	19:30
17	DUP-102510	1010958-009B	10R1561.D	10/29/10	20:00

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Lab File ID: 10\R1570.D DFTPP Injection Date: 10/31/10  
 Instrument ID: HP5973R DFTPP Injection Time: 14:59

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	31.7
68	Less than 2% of mass 69	0.0 (0.0)1
69	Mass 69 relative abundance	43.4
70	Less than 2% of mass 69	0.2 (0.5)1
127	40.0 - 60.0% of mass 198	53.1
197	Less than 1% of mass 198	0.0
198	Base peak, 100% relative abundance	100.0
199	5.0 - 9.0% of mass 198	6.7
275	10.0 - 30.0% of mass 198	21.7
365	Greater than 1% of mass 198	2.3
441	Present, but less than mass 443	7.4
442	40.0 - 110.0% of mass 198	52.7
443	17.0 - 23.0% of mass 442	10.1 (19.1)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	SSTD025	SSTD025	10R1571.D	10/31/10	15:13
02	HMW-5DDL	1010958-001BDL	10R1572.D	10/31/10	15:42
03	HMW-SIDL	1010958-002BDL	10R1573.D	10/31/10	16:12
04	HMW-20IDL	1010958-007BDL	10R1574.D	10/31/10	16:44
05	DUP-102510DL	1010958-009BDL	10R1575.D	10/31/10	17:14

## SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: H2M LABS INC

Contract: \_\_\_\_\_

Lab Code: H2MCase No.: KEY-URS

SAS No.: \_\_\_\_\_

SDG No.: KEY-URS104Instrument ID: HP5973RCalibration Date: 10/28/201 Time: 16:01Lab File ID: 10\R1520.DInit. Calib. Date(s): 10/20/10 10/20/10EPA Sample No. (SSTD050##): SSTD025Init. Calib. Times: 12:18 14:46GC Column: R-5SILMSID: .25 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Naphthalene	1.062	1.048		-1.3	
2-Methylnaphthalene	0.734	0.715		-2.5	
Acenaphthylene	1.916	1.885		-1.6	
Acenaphthene	1.206	1.194		-1.0	20.0
Fluorene	1.339	1.326		-1.0	
Phenanthrene	1.188	1.154		-2.9	
Anthracene	1.149	1.140		-0.8	
Fluoranthene	1.179	1.132		-4.0	20.0
Pyrene	1.346	1.344		-0.1	
Benzo(a)anthracene	1.120	1.093		-2.4	
Chrysene	1.058	1.056		-0.2	
Benzo(b)fluoranthene	1.351	1.202		-11.0	
Benzo(k)fluoranthene	0.917	1.140		24.3	
Benzo(a)pyrene	1.108	1.124		1.5	20.0
Indeno(1,2,3-cd)pyrene	1.165	1.187		1.9	
Dibenzo(a,h)anthracene	0.922	0.935		1.4	
Benzo(g,h,i)perylene	1.025	1.052		2.7	

All other compounds must meet a minimum RRF of 0.010.



## SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Instrument ID: HP5973R Calibration Date: 10/29/201 Time: 12:07  
 Lab File ID: 10\R1545.D Init. Calib. Date(s): 10/20/10 10/20/10  
 EPA Sample No. (SSTD050##): SSTD025 Init. Calib. Times: 12:18 14:46  
 GC Column: R-5SILMS ID: .25 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Naphthalene	1.062	1.057		-0.4	
2-Methylnaphthalene	0.734	0.715		-2.5	
Acenaphthylene	1.916	1.953		1.9	
Acenaphthene	1.206	1.211		0.4	20.0
Fluorene	1.339	1.359		1.5	
Phenanthrene	1.188	1.160		-2.4	
Anthracene	1.149	1.155		0.5	
Fluoranthene	1.179	1.144		-3.0	20.0
Pyrene	1.346	1.359		1.0	
Benzo(a)anthracene	1.120	1.090		-2.7	
Chrysene	1.058	1.083		2.3	
Benzo(b)fluoranthene	1.351	1.179		-12.7	
Benzo(k)fluoranthene	0.917	1.143		24.6	
Benzo(a)pyrene	1.108	1.129		1.9	20.0
Indeno(1,2,3-cd)pyrene	1.165	1.254		7.7	
Dibenzo(a,h)anthracene	0.922	0.971		5.3	
Benzo(g,h,i)perylene	1.025	1.104		7.8	

All other compounds must meet a minimum RRF of 0.010.

## SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: H2M LABS INC Contract: \_\_\_\_\_  
 Lab Code: H2M Case No.: KEY-URS SAS No.: \_\_\_\_\_ SDG No.: KEY-URS104  
 Instrument ID: HP5973R Calibration Date: 10/31/201 Time: 15:13  
 Lab File ID: 10\1571.D Init. Calib. Date(s): 10/20/10 10/20/10  
 EPA Sample No. (SSTD050##): SSTD025 Init. Calib. Times: 12:18 14:46  
 GC Column: R-5SILMS ID: .25 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Naphthalene	1.062	1.057		-0.4	
2-Methylnaphthalene	0.734	0.709		-3.4	
Acenaphthylene	1.916	1.900		-0.9	
Acenaphthene	1.206	1.197		-0.7	20.0
Fluorene	1.339	1.323		-1.2	
Phenanthrene	1.188	1.173		-1.3	
Anthracene	1.149	1.168		1.6	
Fluoranthene	1.179	1.135		-3.7	20.0
Pyrene	1.346	1.364		1.4	
Benzo(a)anthracene	1.120	1.102		-1.6	
Chrysene	1.058	1.075		1.6	
Benzo(b)fluoranthene	1.351	1.220		-9.7	
Benzo(k)fluoranthene	0.917	1.103		20.2	
Benzo(a)pyrene	1.108	1.122		1.3	20.0
Indeno(1,2,3-cd)pyrene	1.165	1.102		-5.4	
Dibenzo(a,h)anthracene	0.922	0.894		-3.0	
Benzo(g,h,i)perylene	1.025	0.865		-15.6	

All other compounds must meet a minimum RRF of 0.010.