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[www.alsglobal.com](http://www.alsglobal.com)

July 31, 2018

**Analytical Report for Service Request No: K1804985A**

Amy Dahl  
AECOM  
1111 Third Avenue, Suite 1600  
Seattle, WA 98101

**RE: Portland Harbor Pre-Remedial Design Investigation / 60566335**

Dear Amy,

Enclosed are the results of the sample(s) submitted to our laboratory May 25, 2018  
For your reference, these analyses have been assigned our service request number **K1804985**.

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. The test results meet requirements of the current NELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP-accredited analytes, refer to the certifications section at [www.alsglobal.com](http://www.alsglobal.com). All results are intended to be considered in their entirety, and ALS Group USA Corp. dba ALS Environmental (ALS) is not responsible for use of less than the complete report. Results apply only to the items submitted to the laboratory for analysis and individual items (samples) analyzed, as listed in the report.

Please contact me if you have any questions. My extension is 3364. You may also contact me via email at [howard.holmes@alsglobal.com](mailto:howard.holmes@alsglobal.com).

Respectfully submitted,

**ALS Group USA, Corp. dba ALS Environmental**

A handwritten signature in black ink, appearing to read "Howard Holmes".

Howard Holmes  
Project Manager



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## Acronyms

ASTM	American Society for Testing and Materials
A2LA	American Association for Laboratory Accreditation
CARB	California Air Resources Board
CAS Number	Chemical Abstract Service registry Number
CFC	Chlorofluorocarbon
CFU	Colony-Forming Unit
DEC	Department of Environmental Conservation
DEQ	Department of Environmental Quality
DHS	Department of Health Services
DOE	Department of Ecology
DOH	Department of Health
EPA	U. S. Environmental Protection Agency
ELAP	Environmental Laboratory Accreditation Program
GC	Gas Chromatography
GC/MS	Gas Chromatography/Mass Spectrometry
LOD	Limit of Detection
LOQ	Limit of Quantitation
LUFT	Leaking Underground Fuel Tank
M	Modified
MCL	Maximum Contaminant Level is the highest permissible concentration of a substance allowed in drinking water as established by the USEPA.
MDL	Method Detection Limit
MPN	Most Probable Number
MRL	Method Reporting Limit
NA	Not Applicable
NC	Not Calculated
NCASI	National Council of the Paper Industry for Air and Stream Improvement
ND	Not Detected
NIOSH	National Institute for Occupational Safety and Health
PQL	Practical Quantitation Limit
RCRA	Resource Conservation and Recovery Act
SIM	Selected Ion Monitoring
TPH	Total Petroleum Hydrocarbons
tr	Trace level is the concentration of an analyte that is less than the PQL but greater than or equal to the MDL.

## Inorganic Data Qualifiers

- \* The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- E The result is an estimate amount because the value exceeded the instrument calibration range.
- J The result is an estimated value.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.  
*DOD-QSM 4.2 definition* : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.
- H The holding time for this test is immediately following sample collection. The samples were analyzed as soon as possible after receipt by the laboratory.

## Metals Data Qualifiers

- # The control limit criteria is not applicable. See case narrative.
- J The result is an estimated value.
- E The percent difference for the serial dilution was greater than 10%, indicating a possible matrix interference in the sample.
- M The duplicate injection precision was not met.
- N The Matrix Spike sample recovery is not within control limits. See case narrative.
- S The reported value was determined by the Method of Standard Additions (MSA).
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.  
*DOD-QSM 4.2 definition* : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- W The post-digestion spike for furnace AA analysis is out of control limits, while sample absorbance is less than 50% of spike absorbance.
- i The MRL/MDL or LOQ/LOD is elevated due to a matrix interference.
- X See case narrative.
- + The correlation coefficient for the MSA is less than 0.995.
- Q See case narrative. One or more quality control criteria was outside the limits.

## Organic Data Qualifiers

- \* The result is an outlier. See case narrative.
- # The control limit criteria is not applicable. See case narrative.
- A A tentatively identified compound, a suspected aldol-condensation product.
- B The analyte was found in the associated method blank at a level that is significant relative to the sample result as defined by the DOD or NELAC standards.
- C The analyte was qualitatively confirmed using GC/MS techniques, pattern recognition, or by comparing to historical data.
- D The reported result is from a dilution.
- E The result is an estimated value.
- J The result is an estimated value.
- N The result is presumptive. The analyte was tentatively identified, but a confirmation analysis was not performed.
- P The GC or HPLC confirmation criteria was exceeded. The relative percent difference is greater than 40% between the two analytical results.
- U The analyte was analyzed for, but was not detected ("Non-detect") at or above the MRL/MDL.  
*DOD-QSM 4.2 definition* : Analyte was not detected and is reported as less than the LOD or as defined by the project. The detection limit is adjusted for dilution.
- i The MRL/MDL or LOQ/LOD is elevated due to a chromatographic interference.
- X See case narrative.
- Q See case narrative. One or more quality control criteria was outside the limits.

## Additional Petroleum Hydrocarbon Specific Qualifiers

- F The chromatographic fingerprint of the sample matches the elution pattern of the calibration standard.
- L The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of lighter molecular weight constituents than the calibration standard.
- H The chromatographic fingerprint of the sample resembles a petroleum product, but the elution pattern indicates the presence of a greater amount of heavier molecular weight constituents than the calibration standard.
- O The chromatographic fingerprint of the sample resembles an oil, but does not match the calibration standard.
- Y The chromatographic fingerprint of the sample resembles a petroleum product eluting in approximately the correct carbon range, but the elution pattern does not match the calibration standard.
- Z The chromatographic fingerprint does not resemble a petroleum product.

**ALS Group USA Corp. dba ALS Environmental (ALS) - Kelso**  
**State Certifications, Accreditations, and Licenses**

Agency	Web Site	Number
Alaska DEH	<a href="http://dec.alaska.gov/eh/lab/cs/csapproval.htm">http://dec.alaska.gov/eh/lab/cs/csapproval.htm</a>	UST-040
Arizona DHS	<a href="http://www.azdhs.gov/lab/license/env.htm">http://www.azdhs.gov/lab/license/env.htm</a>	AZ0339
Arkansas - DEQ	<a href="http://www.adeq.state.ar.us/techsvs/labcert.htm">http://www.adeq.state.ar.us/techsvs/labcert.htm</a>	88-0637
California DHS (ELAP)	<a href="http://www.cdpb.ca.gov/certlic/labs/Pages/ELAP.aspx">http://www.cdpb.ca.gov/certlic/labs/Pages/ELAP.aspx</a>	2795
DOD ELAP	<a href="http://www.denix.osd.mil/edqw/Accreditation/AccreditedLabs.cfm">http://www.denix.osd.mil/edqw/Accreditation/AccreditedLabs.cfm</a>	L16-58-R4
Florida DOH	<a href="http://www.doh.state.fl.us/lab/EnvLabCert/WaterCert.htm">http://www.doh.state.fl.us/lab/EnvLabCert/WaterCert.htm</a>	E87412
Hawaii DOH	<a href="http://health.hawaii.gov/">http://health.hawaii.gov/</a>	-
ISO 17025	<a href="http://www.pjlabs.com/">http://www.pjlabs.com/</a>	L16-57
Louisiana DEQ	<a href="http://www.deq.louisiana.gov/page/la-lab-accreditation">http://www.deq.louisiana.gov/page/la-lab-accreditation</a>	03016
Maine DHS	<a href="http://www.maine.gov/dhhs/">http://www.maine.gov/dhhs/</a>	WA01276
Minnesota DOH	<a href="http://www.health.state.mn.us/accreditation">http://www.health.state.mn.us/accreditation</a>	053-999-457
Nevada DEP	<a href="http://ndep.nv.gov/bsdw/labservice.htm">http://ndep.nv.gov/bsdw/labservice.htm</a>	WA01276
New Jersey DEP	<a href="http://www.nj.gov/dep/enforcement/oqa.html">http://www.nj.gov/dep/enforcement/oqa.html</a>	WA005
New York - DOH	<a href="https://www.wadsworth.org/regulatory/elap">https://www.wadsworth.org/regulatory/elap</a>	12060
North Carolina DEQ	<a href="https://deq.nc.gov/about/divisions/water-resources/water-resources-data/water-sciences-home-page/laboratory-certification-branch/non-field-lab-certification">https://deq.nc.gov/about/divisions/water-resources/water-resources-data/water-sciences-home-page/laboratory-certification-branch/non-field-lab-certification</a>	605
Oklahoma DEQ	<a href="http://www.deq.state.ok.us/CSDnew/labcert.htm">http://www.deq.state.ok.us/CSDnew/labcert.htm</a>	9801
Oregon – DEQ (NELAP)	<a href="http://public.health.oregon.gov/LaboratoryServices/EnvironmentalLaboratoryAccreditation/Pages/index.aspx">http://public.health.oregon.gov/LaboratoryServices/EnvironmentalLaboratoryAccreditation/Pages/index.aspx</a>	WA100010
South Carolina DHEC	<a href="http://www.scdhec.gov/environment/EnvironmentalLabCertification/">http://www.scdhec.gov/environment/EnvironmentalLabCertification/</a>	61002
Texas CEQ	<a href="http://www.tceq.texas.gov/field/qa/env_lab_accreditation.html">http://www.tceq.texas.gov/field/qa/env_lab_accreditation.html</a>	T104704427
Washington DOE	<a href="http://www.ecy.wa.gov/programs/eap/labs/lab-accreditation.html">http://www.ecy.wa.gov/programs/eap/labs/lab-accreditation.html</a>	C544
Wyoming (EPA Region 8)	<a href="https://www.epa.gov/region8-waterops/epa-region-8-certified-drinking-water">https://www.epa.gov/region8-waterops/epa-region-8-certified-drinking-water</a>	-
Kelso Laboratory Website	<a href="http://www.alsglobal.com">www.alsglobal.com</a>	NA

Analyses were performed according to our laboratory's NELAP-approved quality assurance program. A complete listing of specific NELAP-certified analytes, can be found in the certification section at [www.alsglobal.com](http://www.alsglobal.com) or at the accreditation bodies web site.

Please refer to the certification and/or accreditation body's web site if samples are submitted for compliance purposes. The states highlighted above, require the analysis be listed on the state certification if used for compliance purposes and if the method/analyte is offered by that state.



## Case Narrative

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**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Received:** 05/25/2018

#### CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples designated for Tier IV validation deliverables including summary forms and all of the associated raw data for each of the analyses. When appropriate to the method, method blank results have been reported with each analytical test.

#### Sample Receipt:

Eleven sediment samples were received for analysis at ALS Environmental on 05/25/2018. The samples were received in good condition and consistent with the accompanying chain of custody form. The samples were stored in a refrigerator at 4°C upon receipt at the laboratory.

#### Semivolatiles by GC/MS:

Method 8270D SIM-PAH, 07/26/2018: The matrix spike recovery of Phenanthrene, Fluoranthene, Pyrene, Chrysene and Benzo(b)fluoranthene for sample PDI-SG-S204 was outside control criteria. Recovery in the Laboratory Control Sample (LCS) was acceptable, which indicated the analytical batch was in control. The matrix spike outlier suggested a potential high bias in this matrix. No further corrective action was appropriate.

Method 8270D SIM-PAH, 07/26/2018: The recovery of Phenanthrene, Fluoranthene, Chrysene, Benzo(b)fluoranthene and Benzo(a)pyrene in the Duplicate Matrix Spike (DMS) KWG1803256-2 was outside the recovery control limits listed in the results summary. The DMSS is used to evaluate batch precision. The relative percent difference (RPD) was within control limits indicating the quality of the sample data was not significantly affected. No further corrective action was taken.

A handwritten signature in black ink is placed over a horizontal line intended for approval. The signature appears to be a stylized version of the name "Howard Johnson".

Approved by \_\_\_\_\_

Date \_\_\_\_\_ 07/31/2018



## Chain of Custody

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K1804985

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Kelso, WA 98626
Ph: 360-577-7222 Fax: 360-636-1068
Client Contact

## SURFACE SEDIMENT CHAIN OF CUSTODY

Client Contact		Project Contact: Amy Dahl / Chelsey Cook Tel: (206) 438-2261 / (206) 438-2010			Site Contact: Jennifer Ray / Michaela McCool Laboratory Contact: Howard-Holmes			Carrier: courier			
AECOM 1111 3rd Ave Suite 1600 Seattle, WA 98101 Phone: (206) 438-2700 Fax: 1-(866) 495-5288 Project Name: Portland Harbor Pre-Remedial Design Investigation and Baseline Sampling		Analysis Turnaround Time Calendar (C) or Work Days (W)						5/25/2018 COC No: 3 1 of 1 pages			
Portland, OR Project #: 60566335 Study: Surface Sediment - SMA		<input type="checkbox"/> 21 days <input type="checkbox"/> Other _____									
Sample Identification		Sample Date	Sample Time	Matrix	QC Sample	Sampler's Initials	Total No. of Cont.	Preparation	Hold at 20°C	Pesticides, PAHs, Total Solids 1669M, 8270-SIM, 160.3	Sample Specific Notes:
PDI-SG-S204		5/3/2018	16:25	SS			1		x	H	Frozen 5/3/18 08:00
PDI-SG-S147		5/4/2018	17:19	SS			1		x	H	Frozen 5/4/18 18:30
PDI-SG-S084		5/8/2018	13:40	SS			1		x	H	Frozen 5/8/18 18:20
PDI-SG-S090		5/9/2018	14:34	SS			1		x	H	Frozen 5/9/18 18:50
PDI-SG-S010		5/9/2018	17:30	SS			1		x	H	Frozen 5/9/18 18:50
PDI-SG-S255		5/11/2018	12:40	SS			1		x	H	Frozen 5/14/18 10:45
PDI-SG-S097		5/13/2018	11:45	SS			1		x	H	Frozen 5/13/18 18:45
PDI-SG-S115		5/12/2018	12:21	SS			1		x	H	Frozen 5/12/18 17:15
PDI-SG-S078		5/12/2018	15:50	SS			1		x	H	Frozen 5/12/18 17:15
PDI-SG-S155		5/14/2018	11:15	SS			1		x	H	Frozen 5/15/18 11:15
PDI-SG-S135		5/14/2018	10:15	SS			1		x	H	Frozen 5/15/18 11:15
PDI-SG-S157		5/14/2018	15:45	SS			1		x	H	Frozen 5/15/18 11:15
Container Type: WMG=Wide Mouth Glass Jar, P=HDPE, PP=Polypropylene, AG=amber glass, G=glass, RC=Resin Column Preservative: HCl = Hydrochloric Acid, H <sub>3</sub> PO <sub>4</sub> = Phosphoric Acid, HNO <sub>3</sub> = Nitric Acid Fraction: D = Dissolved, PRT = Particulate, T = Total (unfiltered)											
Sample Disposal <input type="checkbox"/> Return To Client <input checked="" type="checkbox"/> Disposal By Lab <input checked="" type="checkbox"/> Archive For 12 Months											
Special Instructions/QC Requirements & Comments: Please hold in freezer pending approval of analyses marked H.											
Relinquished by: 	Company: AECOM	Date/Time: 5-25-18	Received by:	Company: ARV	Date/Time: 5/25/18 12:35						
Relinquished by: 	Company: ARV	Date/Time: 5/25/18 1405	Received by:	Company: ARV	Date/Time: 5/25/18 1405						
Relinquished by: 	Company:	Date/Time:	Received by:	Company:	Date/Time:						



PC 44

## Cooler Receipt and Preservation Form

Client AECOM  
 Received: S125118 Opened: S125118 By: BR Unloaded: S125118 By: BR

Service Request K1804985

1. Samples were received via?  **USPS**  **FedEx**  **UPS**  **DHL**  **PDX**  **Courier**  **Hand Delivered**
2. Samples were received in: (circle)  **Cooler**  **Box**  **Envelope**  **Other**  **NA**
3. Were custody seals on coolers?  **Y**  **N** If yes, how many and where? IF FRONT  
 If present, were custody seals intact?  **Y**  **N** If present, were they signed and dated?  **Y**  **N**

Raw Cooler Temp	Corrected Cooler Temp	Raw Temp Blank	Corrected Temp Blank	Corr. Factor	Thermometer ID	Cooler/COC ID	Tracking Number	NA Filed
33	31	49	49	-0.2	351	NA		
38	38	24	39	0.8	374			
38	38	21	21	0.8	325			

4. Packing material:  **Inserts**  **Buggies**  **Bubble Wrap**  **Gel Packs**  **Wet Ice**  **Dry Ice**  **Sleeves** \_\_\_\_\_
5. Were custody papers properly filled out (ink, signed, etc.)?  **NA**  **Y**  **N**
6. Were samples received in good condition (temperature, unbroken)? *Indicate in the table below.*  
 If applicable, tissue samples were received:  **Frozen**  **Partially Thawed**  **Thawed**
7. Were all sample labels complete (i.e analysis, preservation, etc.)?  **NA**  **Y**  **N**
8. Did all sample labels and tags agree with custody papers? *Indicate major discrepancies in the table on page 2.*  **NA**  **Y**  **N**
9. Were appropriate bottles/containers and volumes received for the tests indicated?  **NA**  **Y**  **N**
10. Were the pH-preserved bottles (*see SMO GEN SOP*) received at the appropriate pH? *Indicate in the table below.*  **NA**  **Y**  **N**
11. Were VOA vials received without headspace? *Indicate in the table below.*  **NA**  **Y**  **N**
12. Was C12/Res negative?  **NA**  **Y**  **N**

Sample ID on Bottle	Sample ID on COC	Identified by:

Sample ID	Bottle Count Bottle Type	Out of Temp	Head- space	Broke	pH	Reagent	Volume added	Reagent Lot Number	Initials	Time

Notes, Discrepancies, & Resolutions:

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## Total Solids

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**COLUMBIA ANALYTICAL SERVICES, INC.**

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## Analytical Report

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment  
**Analysis Method:** 160.3 Modified  
**Prep Method:** None

**Service Request:** K1804985  
**Date Collected:** 05/03/18 - 05/14/18  
**Date Received:** 05/25/18  
**Units:** Percent  
**Basis:** As Received

**Solids, Total**

Sample Name	Lab Code	Result	MRL	MDL	Dil.	Date Analyzed	Q
PDI-SG-S204	K1804985-001	<b>70.3</b>	-	-	1	06/28/18 15:55	
PDI-SG-S147	K1804985-002	<b>54.8</b>	-	-	1	06/28/18 15:55	
PDI-SG-S084	K1804985-003	<b>82.3</b>	-	-	1	06/28/18 15:55	
PDI-SG-S090	K1804985-004	<b>75.2</b>	-	-	1	06/28/18 15:55	
PDI-SG-S010	K1804985-005	<b>75.9</b>	-	-	1	06/28/18 15:55	
PDI-SG-S255	K1804985-006	<b>62.4</b>	-	-	1	06/28/18 15:55	
PDI-SG-S097	K1804985-007	<b>62.5</b>	-	-	1	06/28/18 15:55	
PDI-SG-S115	K1804985-008	<b>74.0</b>	-	-	1	06/28/18 15:55	
PDI-SG-S078	K1804985-009	<b>55.1</b>	-	-	1	06/28/18 15:55	
PDI-SG-S135	K1804985-010	<b>70.5</b>	-	-	1	06/28/18 15:55	
PDI-SG-S157	K1804985-011	<b>60.6</b>	-	-	1	06/28/18 15:55	

**COLUMBIA ANALYTICAL SERVICES, INC.**

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## QA/QC Report

<b>Client:</b>	AECOM	<b>Service Request:</b>	K1804985
<b>Project</b>	Portland Harbor Pre-Remedial Design Investigation/60566335	<b>Date Collected:</b>	05/03/18 - 05/14/18
<b>Sample Matrix:</b>	Sediment	<b>Date Received:</b>	05/25/18
<b>Analysis Method:</b>	160.3 Modified	<b>Units:</b>	Percent
<b>Prep Method:</b>	None	<b>Basis:</b>	As Received

**Duplicate Sample Summary****Inorganic Parameters**

<b>Sample Name:</b>	<b>Lab Code:</b>	<b>MRL</b>	<b>Sample Result</b>	<b>Duplicate Result</b>	<b>Average</b>	<b>RPD</b>	<b>RPD Limit</b>	<b>Date Analyzed</b>
PDI-SG-S204	K1804985-001DUP	-	70.3	69.4	69.9	1	20	06/28/18
PDI-SG-S157	K1804985-011DUP	-	60.6	61.2	60.9	<1	20	06/28/18

Results flagged with an asterisk (\*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.



## Polynuclear Aromatic Hydrocarbons

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[www.alsglobal.com](http://www.alsglobal.com)

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985

**Cover Page - Organic Analysis Data Package**  
**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name</b>	<b>Lab Code</b>	<b>Date Collected</b>	<b>Date Received</b>
PDI-SG-S204	K1804985-001	05/03/2018	05/25/2018
PDI-SG-S147	K1804985-002	05/04/2018	05/25/2018
PDI-SG-S084	K1804985-003	05/08/2018	05/25/2018
PDI-SG-S090	K1804985-004	05/09/2018	05/25/2018
PDI-SG-S010	K1804985-005	05/09/2018	05/25/2018
PDI-SG-S255	K1804985-006	05/11/2018	05/25/2018
PDI-SG-S097	K1804985-007	05/13/2018	05/25/2018
PDI-SG-S115	K1804985-008	05/12/2018	05/25/2018
PDI-SG-S078	K1804985-009	05/12/2018	05/25/2018
PDI-SG-S135	K1804985-010	05/14/2018	05/25/2018
PDI-SG-S157	K1804985-011	05/14/2018	05/25/2018
PDI-SG-S204MS	KWG1803256-1	05/03/2018	05/25/2018
PDI-SG-S204DMS	KWG1803256-2	05/03/2018	05/25/2018

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/03/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

**Sample Name:** PDI-SG-S204      **Units:** ug/Kg  
**Lab Code:** K1804985-001      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	7.9	JD	18	7.5	50	07/03/18	07/10/18	KWG1803256	
2-Methylnaphthalene	ND	U	18	6.0	50	07/03/18	07/10/18	KWG1803256	
Acenaphthylene	6.5	JD	8.9	2.3	50	07/03/18	07/10/18	KWG1803256	
Acenaphthene	6.7	JD	8.9	2.4	50	07/03/18	07/10/18	KWG1803256	
Fluorene	9.2	D	8.9	2.6	50	07/03/18	07/10/18	KWG1803256	
Phenanthrene	69	D	8.9	3.3	50	07/03/18	07/10/18	KWG1803256	
Anthracene	12	D	8.9	1.9	50	07/03/18	07/10/18	KWG1803256	
Fluoranthene	170	D	8.9	2.5	50	07/03/18	07/10/18	KWG1803256	
Pyrene	250	D	8.9	2.5	50	07/03/18	07/10/18	KWG1803256	
Benz(a)anthracene	55	D	8.9	1.9	50	07/03/18	07/10/18	KWG1803256	
Chrysene	86	D	8.9	2.8	50	07/03/18	07/10/18	KWG1803256	
Benzo(b)fluoranthene†	85	D	8.9	3.3	50	07/03/18	07/10/18	KWG1803256	
Benzo(k)fluoranthene	26	D	8.9	2.9	50	07/03/18	07/10/18	KWG1803256	
Benzo(a)pyrene	54	D	8.9	3.7	50	07/03/18	07/10/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	54	D	8.9	4.8	50	07/03/18	07/10/18	KWG1803256	
Dibenz(a,h)anthracene	16	D	8.9	4.3	50	07/03/18	07/10/18	KWG1803256	
Benzo(g,h,i)perylene	73	D	8.9	4.8	50	07/03/18	07/10/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	64	42-106	07/10/18	Acceptable
Fluoranthene-d10	69	45-109	07/10/18	Acceptable
Terphenyl-d14	77	41-102	07/10/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/04/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	PDI-SG-S147	<b>Units:</b>	ug/Kg
<b>Lab Code:</b>	K1804985-002	<b>Basis:</b>	Dry
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM		

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	370	D	2.3	0.75	5	07/03/18	07/10/18	KWG1803256	
2-Methylnaphthalene	100	D	2.3	0.60	5	07/03/18	07/10/18	KWG1803256	
Acenaphthylene	32	D	1.2	0.23	5	07/03/18	07/10/18	KWG1803256	
Acenaphthene	59	D	1.2	0.24	5	07/03/18	07/10/18	KWG1803256	
Fluorene	61	D	1.2	0.26	5	07/03/18	07/10/18	KWG1803256	
Phenanthrene	260	D	1.2	0.33	5	07/03/18	07/10/18	KWG1803256	
Anthracene	96	D	1.2	0.19	5	07/03/18	07/10/18	KWG1803256	
Fluoranthene	360	D	1.2	0.25	5	07/03/18	07/10/18	KWG1803256	
Pyrene	410	D	1.2	0.25	5	07/03/18	07/10/18	KWG1803256	
Benz(a)anthracene	140	D	1.2	0.19	5	07/03/18	07/10/18	KWG1803256	
Chrysene	220	D	1.2	0.28	5	07/03/18	07/10/18	KWG1803256	
Benzo(b)fluoranthene†	170	D	1.2	0.33	5	07/03/18	07/10/18	KWG1803256	
Benzo(k)fluoranthene	60	D	1.2	0.29	5	07/03/18	07/10/18	KWG1803256	
Benzo(a)pyrene	160	D	1.2	0.37	5	07/03/18	07/10/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	120	D	1.2	0.48	5	07/03/18	07/10/18	KWG1803256	
Dibenz(a,h)anthracene	22	D	1.2	0.43	5	07/03/18	07/10/18	KWG1803256	
Benzo(g,h,i)perylene	130	D	1.2	0.48	5	07/03/18	07/10/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	74	42-106	07/10/18	Acceptable
Fluoranthene-d10	82	45-109	07/10/18	Acceptable
Terphenyl-d14	92	41-102	07/10/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/08/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	PDI-SG-S084	<b>Units:</b>	ug/Kg
<b>Lab Code:</b>	K1804985-003	<b>Basis:</b>	Dry
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM		

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	22	D	1.6	0.75	5	07/03/18	07/11/18	KWG1803256	
2-Methylnaphthalene	5.1	D	1.6	0.60	5	07/03/18	07/11/18	KWG1803256	
Acenaphthylene	61	D	0.76	0.23	5	07/03/18	07/11/18	KWG1803256	
Acenaphthene	4.4	D	0.76	0.24	5	07/03/18	07/11/18	KWG1803256	
Fluorene	6.1	D	0.76	0.26	5	07/03/18	07/11/18	KWG1803256	
Phenanthrene	14	D	0.76	0.33	5	07/03/18	07/11/18	KWG1803256	
Anthracene	29	D	0.76	0.19	5	07/03/18	07/11/18	KWG1803256	
Fluoranthene	840	D	7.6	2.5	50	07/03/18	07/12/18	KWG1803256	
Pyrene	1500	D	7.6	2.5	50	07/03/18	07/12/18	KWG1803256	
Benz(a)anthracene	480	D	7.6	1.9	50	07/03/18	07/12/18	KWG1803256	
Chrysene	580	D	7.6	2.8	50	07/03/18	07/12/18	KWG1803256	
Benzo(b)fluoranthene†	630	D	7.6	3.3	50	07/03/18	07/12/18	KWG1803256	
Benzo(k)fluoranthene	180	D	0.76	0.29	5	07/03/18	07/11/18	KWG1803256	
Benzo(a)pyrene	740	D	7.6	3.7	50	07/03/18	07/12/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	510	D	7.6	4.8	50	07/03/18	07/12/18	KWG1803256	
Dibenz(a,h)anthracene	60	D	0.76	0.43	5	07/03/18	07/11/18	KWG1803256	
Benzo(g,h,i)perylene	540	D	7.6	4.8	50	07/03/18	07/12/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	66	42-106	07/11/18	Acceptable
Fluoranthene-d10	75	45-109	07/11/18	Acceptable
Terphenyl-d14	90	41-102	07/11/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/09/2018  
**Date Received:** 05/25/2018

## Polynuclear Aromatic Hydrocarbons

**Sample Name:** PDI-SG-S090      **Units:** ug/Kg  
**Lab Code:** K1804985-004      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	54	D	17	7.5	50	07/03/18	07/11/18	KWG1803256	
2-Methylnaphthalene	13	JD	17	6.0	50	07/03/18	07/11/18	KWG1803256	
Acenaphthylene	130	D	8.3	2.3	50	07/03/18	07/11/18	KWG1803256	
Acenaphthene	7.2	JD	8.3	2.4	50	07/03/18	07/11/18	KWG1803256	
Fluorene	11	D	8.3	2.6	50	07/03/18	07/11/18	KWG1803256	
Phenanthrene	34	D	8.3	3.3	50	07/03/18	07/11/18	KWG1803256	
Anthracene	67	D	8.3	1.9	50	07/03/18	07/11/18	KWG1803256	
Fluoranthene	1400	D	8.3	2.5	50	07/03/18	07/11/18	KWG1803256	
Pyrene	2500	D	8.3	2.5	50	07/03/18	07/11/18	KWG1803256	
Benz(a)anthracene	1200	D	8.3	1.9	50	07/03/18	07/11/18	KWG1803256	
Chrysene	1500	D	8.3	2.8	50	07/03/18	07/11/18	KWG1803256	
Benzo(b)fluoranthene†	1700	D	8.3	3.3	50	07/03/18	07/11/18	KWG1803256	
Benzo(k)fluoranthene	550	D	8.3	2.9	50	07/03/18	07/11/18	KWG1803256	
Benzo(a)pyrene	2100	D	8.3	3.7	50	07/03/18	07/11/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	1300	D	8.3	4.8	50	07/03/18	07/11/18	KWG1803256	
Dibenz(a,h)anthracene	190	D	8.3	4.3	50	07/03/18	07/11/18	KWG1803256	
Benzo(g,h,i)perylene	1300	D	8.3	4.8	50	07/03/18	07/11/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	63	42-106	07/11/18	Acceptable
Fluoranthene-d10	68	45-109	07/11/18	Acceptable
Terphenyl-d14	82	41-102	07/11/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/09/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

**Sample Name:** PDI-SG-S010      **Units:** ug/Kg  
**Lab Code:** K1804985-005      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	8.6		0.33	0.15	1	07/03/18	07/10/18	KWG1803256	
2-Methylnaphthalene	3.2		0.33	0.12	1	07/03/18	07/10/18	KWG1803256	
Acenaphthylene	3.3		0.17	0.046	1	07/03/18	07/10/18	KWG1803256	
Acenaphthene	5.6		0.17	0.047	1	07/03/18	07/10/18	KWG1803256	
Fluorene	7.0		0.17	0.052	1	07/03/18	07/10/18	KWG1803256	
Phenanthrene	100	D	1.7	0.66	10	07/03/18	07/12/18	KWG1803256	
Anthracene	5.6		0.17	0.038	1	07/03/18	07/10/18	KWG1803256	
Fluoranthene	140	D	1.7	0.49	10	07/03/18	07/12/18	KWG1803256	
Pyrene	160	D	1.7	0.50	10	07/03/18	07/12/18	KWG1803256	
Benz(a)anthracene	35		0.17	0.038	1	07/03/18	07/10/18	KWG1803256	
Chrysene	60		0.17	0.055	1	07/03/18	07/10/18	KWG1803256	
Benzo(b)fluoranthene†	61		0.17	0.066	1	07/03/18	07/10/18	KWG1803256	
Benzo(k)fluoranthene	21		0.17	0.057	1	07/03/18	07/10/18	KWG1803256	
Benzo(a)pyrene	56		0.17	0.073	1	07/03/18	07/10/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	48		0.17	0.096	1	07/03/18	07/10/18	KWG1803256	
Dibenz(a,h)anthracene	6.5		0.17	0.086	1	07/03/18	07/10/18	KWG1803256	
Benzo(g,h,i)perylene	52		0.17	0.095	1	07/03/18	07/10/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	64	42-106	07/10/18	Acceptable
Fluoranthene-d10	74	45-109	07/10/18	Acceptable
Terphenyl-d14	86	41-102	07/10/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/11/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

**Sample Name:** PDI-SG-S255      **Units:** ug/Kg  
**Lab Code:** K1804985-006      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	2.2		0.40	0.15	1	07/03/18	07/10/18	KWG1803256	
2-Methylnaphthalene	0.96		0.40	0.12	1	07/03/18	07/10/18	KWG1803256	
Acenaphthylene	11		0.20	0.046	1	07/03/18	07/10/18	KWG1803256	
Acenaphthene	1.3		0.20	0.047	1	07/03/18	07/10/18	KWG1803256	
Fluorene	2.4		0.20	0.052	1	07/03/18	07/10/18	KWG1803256	
Phenanthrene	58		0.20	0.066	1	07/03/18	07/10/18	KWG1803256	
Anthracene	10		0.20	0.038	1	07/03/18	07/10/18	KWG1803256	
Fluoranthene	180 D		2.0	0.49	10	07/03/18	07/12/18	KWG1803256	
Pyrene	240 D		2.0	0.50	10	07/03/18	07/12/18	KWG1803256	
Benz(a)anthracene	65		0.20	0.038	1	07/03/18	07/10/18	KWG1803256	
Chrysene	100 D		2.0	0.55	10	07/03/18	07/12/18	KWG1803256	
Benzo(b)fluoranthene†	110 D		2.0	0.66	10	07/03/18	07/12/18	KWG1803256	
Benzo(k)fluoranthene	33		0.20	0.057	1	07/03/18	07/10/18	KWG1803256	
Benzo(a)pyrene	110 D		2.0	0.73	10	07/03/18	07/12/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	88 D		2.0	0.96	10	07/03/18	07/12/18	KWG1803256	
Dibenz(a,h)anthracene	11		0.20	0.086	1	07/03/18	07/10/18	KWG1803256	
Benzo(g,h,i)perylene	96 D		2.0	0.95	10	07/03/18	07/12/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	60	42-106	07/10/18	Acceptable
Fluoranthene-d10	67	45-109	07/10/18	Acceptable
Terphenyl-d14	80	41-102	07/10/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/13/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	PDI-SG-S097	<b>Units:</b>	ug/Kg
<b>Lab Code:</b>	K1804985-007	<b>Basis:</b>	Dry
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM		

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	44	D	2.0	0.75	5	07/03/18	07/11/18	KWG1803256	
2-Methylnaphthalene	17	D	2.0	0.60	5	07/03/18	07/11/18	KWG1803256	
Acenaphthylene	13	D	1.0	0.23	5	07/03/18	07/11/18	KWG1803256	
Acenaphthene	33	D	1.0	0.24	5	07/03/18	07/11/18	KWG1803256	
Fluorene	27	D	1.0	0.26	5	07/03/18	07/11/18	KWG1803256	
Phenanthrene	190	D	1.0	0.33	5	07/03/18	07/11/18	KWG1803256	
Anthracene	36	D	1.0	0.19	5	07/03/18	07/11/18	KWG1803256	
Fluoranthene	360	D	1.0	0.25	5	07/03/18	07/11/18	KWG1803256	
Pyrene	470	D	2.0	0.50	10	07/03/18	07/12/18	KWG1803256	
Benz(a)anthracene	170	D	1.0	0.19	5	07/03/18	07/11/18	KWG1803256	
Chrysene	230	D	1.0	0.28	5	07/03/18	07/11/18	KWG1803256	
Benzo(b)fluoranthene†	230	D	1.0	0.33	5	07/03/18	07/11/18	KWG1803256	
Benzo(k)fluoranthene	81	D	1.0	0.29	5	07/03/18	07/11/18	KWG1803256	
Benzo(a)pyrene	200	D	1.0	0.37	5	07/03/18	07/11/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	140	D	1.0	0.48	5	07/03/18	07/11/18	KWG1803256	
Dibenz(a,h)anthracene	28	D	1.0	0.43	5	07/03/18	07/11/18	KWG1803256	
Benzo(g,h,i)perylene	140	D	1.0	0.48	5	07/03/18	07/11/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	72	42-106	07/11/18	Acceptable
Fluoranthene-d10	81	45-109	07/11/18	Acceptable
Terphenyl-d14	96	41-102	07/11/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/12/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	PDI-SG-S115	<b>Units:</b>	ug/Kg
<b>Lab Code:</b>	K1804985-008	<b>Basis:</b>	Dry
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM		

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	62	D	1.7	0.75	5	07/03/18	07/11/18	KWG1803256	
2-Methylnaphthalene	40	D	1.7	0.60	5	07/03/18	07/11/18	KWG1803256	
Acenaphthylene	38	D	0.84	0.23	5	07/03/18	07/11/18	KWG1803256	
Acenaphthene	12	D	0.84	0.24	5	07/03/18	07/11/18	KWG1803256	
Fluorene	17	D	0.84	0.26	5	07/03/18	07/11/18	KWG1803256	
Phenanthrene	210	D	0.84	0.33	5	07/03/18	07/11/18	KWG1803256	
Anthracene	56	D	0.84	0.19	5	07/03/18	07/11/18	KWG1803256	
Fluoranthene	400	D	4.2	1.3	25	07/03/18	07/12/18	KWG1803256	
Pyrene	540	D	4.2	1.3	25	07/03/18	07/12/18	KWG1803256	
Benz(a)anthracene	230	D	0.84	0.19	5	07/03/18	07/11/18	KWG1803256	
Chrysene	310	D	0.84	0.28	5	07/03/18	07/11/18	KWG1803256	
Benzo(b)fluoranthene†	290	D	0.84	0.33	5	07/03/18	07/11/18	KWG1803256	
Benzo(k)fluoranthene	98	D	0.84	0.29	5	07/03/18	07/11/18	KWG1803256	
Benzo(a)pyrene	300	D	0.84	0.37	5	07/03/18	07/11/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	220	D	0.84	0.48	5	07/03/18	07/11/18	KWG1803256	
Dibenz(a,h)anthracene	44	D	0.84	0.43	5	07/03/18	07/11/18	KWG1803256	
Benzo(g,h,i)perylene	230	D	0.84	0.48	5	07/03/18	07/11/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	60	42-106	07/11/18	Acceptable
Fluoranthene-d10	69	45-109	07/11/18	Acceptable
Terphenyl-d14	82	41-102	07/11/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/12/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	PDI-SG-S078	<b>Units:</b>	ug/Kg
<b>Lab Code:</b>	K1804985-009	<b>Basis:</b>	Dry
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM		

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	21	D	2.3	0.75	5	07/03/18	07/12/18	KWG1803256	
2-Methylnaphthalene	9.4	D	2.3	0.60	5	07/03/18	07/12/18	KWG1803256	
Acenaphthylene	7.9	D	1.2	0.23	5	07/03/18	07/12/18	KWG1803256	
Acenaphthene	17	D	1.2	0.24	5	07/03/18	07/12/18	KWG1803256	
Fluorene	20	D	1.2	0.26	5	07/03/18	07/12/18	KWG1803256	
Phenanthrene	110	D	1.2	0.33	5	07/03/18	07/12/18	KWG1803256	
Anthracene	28	D	1.2	0.19	5	07/03/18	07/12/18	KWG1803256	
Fluoranthene	190	D	1.2	0.25	5	07/03/18	07/12/18	KWG1803256	
Pyrene	240	D	1.2	0.25	5	07/03/18	07/12/18	KWG1803256	
Benz(a)anthracene	94	D	1.2	0.19	5	07/03/18	07/12/18	KWG1803256	
Chrysene	120	D	1.2	0.28	5	07/03/18	07/12/18	KWG1803256	
Benzo(b)fluoranthene†	130	D	1.2	0.33	5	07/03/18	07/12/18	KWG1803256	
Benzo(k)fluoranthene	45	D	1.2	0.29	5	07/03/18	07/12/18	KWG1803256	
Benzo(a)pyrene	130	D	1.2	0.37	5	07/03/18	07/12/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	99	D	1.2	0.48	5	07/03/18	07/12/18	KWG1803256	
Dibenz(a,h)anthracene	16	D	1.2	0.43	5	07/03/18	07/12/18	KWG1803256	
Benzo(g,h,i)perylene	100	D	1.2	0.48	5	07/03/18	07/12/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	67	42-106	07/12/18	Acceptable
Fluoranthene-d10	78	45-109	07/12/18	Acceptable
Terphenyl-d14	88	41-102	07/12/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/14/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

**Sample Name:** PDI-SG-S135      **Units:** ug/Kg  
**Lab Code:** K1804985-010      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	46	D	18	7.5	50	07/03/18	07/11/18	KWG1803256	
2-Methylnaphthalene	6.8	JD	18	6.0	50	07/03/18	07/11/18	KWG1803256	
Acenaphthylene	83	D	8.9	2.3	50	07/03/18	07/11/18	KWG1803256	
Acenaphthene	12	D	8.9	2.4	50	07/03/18	07/11/18	KWG1803256	
Fluorene	13	D	8.9	2.6	50	07/03/18	07/11/18	KWG1803256	
Phenanthrene	180	D	8.9	3.3	50	07/03/18	07/11/18	KWG1803256	
Anthracene	60	D	8.9	1.9	50	07/03/18	07/11/18	KWG1803256	
Fluoranthene	1200	D	8.9	2.5	50	07/03/18	07/11/18	KWG1803256	
Pyrene	1600	D	8.9	2.5	50	07/03/18	07/11/18	KWG1803256	
Benz(a)anthracene	550	D	8.9	1.9	50	07/03/18	07/11/18	KWG1803256	
Chrysene	680	D	8.9	2.8	50	07/03/18	07/11/18	KWG1803256	
Benzo(b)fluoranthene†	750	D	8.9	3.3	50	07/03/18	07/11/18	KWG1803256	
Benzo(k)fluoranthene	270	D	8.9	2.9	50	07/03/18	07/11/18	KWG1803256	
Benzo(a)pyrene	800	D	8.9	3.7	50	07/03/18	07/11/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	610	D	8.9	4.8	50	07/03/18	07/11/18	KWG1803256	
Dibenz(a,h)anthracene	82	D	8.9	4.3	50	07/03/18	07/11/18	KWG1803256	
Benzo(g,h,i)perylene	600	D	8.9	4.8	50	07/03/18	07/11/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	67	42-106	07/11/18	Acceptable
Fluoranthene-d10	81	45-109	07/11/18	Acceptable
Terphenyl-d14	88	41-102	07/11/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** 05/14/2018  
**Date Received:** 05/25/2018

**Polynuclear Aromatic Hydrocarbons**

**Sample Name:** PDI-SG-S157      **Units:** ug/Kg  
**Lab Code:** K1804985-011      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	8.8	0.41	0.15	1	07/03/18	07/10/18	KWG1803256	
2-Methylnaphthalene	5.8	0.41	0.12	1	07/03/18	07/10/18	KWG1803256	
Acenaphthylene	2.6	0.21	0.046	1	07/03/18	07/10/18	KWG1803256	
Acenaphthene	5.3	0.21	0.047	1	07/03/18	07/10/18	KWG1803256	
Fluorene	7.9	0.21	0.052	1	07/03/18	07/10/18	KWG1803256	
Phenanthrene	47	0.21	0.066	1	07/03/18	07/10/18	KWG1803256	
Anthracene	9.1	0.21	0.038	1	07/03/18	07/10/18	KWG1803256	
Fluoranthene	77	0.21	0.049	1	07/03/18	07/10/18	KWG1803256	
Pyrene	80	0.21	0.050	1	07/03/18	07/10/18	KWG1803256	
Benz(a)anthracene	23	0.21	0.038	1	07/03/18	07/10/18	KWG1803256	
Chrysene	39	0.21	0.055	1	07/03/18	07/10/18	KWG1803256	
Benzo(b)fluoranthene†	27	0.21	0.066	1	07/03/18	07/10/18	KWG1803256	
Benzo(k)fluoranthene	9.7	0.21	0.057	1	07/03/18	07/10/18	KWG1803256	
Benzo(a)pyrene	19	0.21	0.073	1	07/03/18	07/10/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	14	0.21	0.096	1	07/03/18	07/10/18	KWG1803256	
Dibenz(a,h)anthracene	2.7	0.21	0.086	1	07/03/18	07/10/18	KWG1803256	
Benzo(g,h,i)perylene	15	0.21	0.095	1	07/03/18	07/10/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	61	42-106	07/10/18	Acceptable
Fluoranthene-d10	74	45-109	07/10/18	Acceptable
Terphenyl-d14	82	41-102	07/10/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

## Analytical Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Collected:** NA  
**Date Received:** NA

**Polynuclear Aromatic Hydrocarbons**

**Sample Name:** Method Blank      **Units:** ug/Kg  
**Lab Code:** KWG1803256-4      **Basis:** Dry  
**Extraction Method:** EPA 3541      **Level:** Low  
**Analysis Method:** 8270D SIM

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Naphthalene	ND	U	0.25	0.15	1	07/03/18	07/10/18	KWG1803256	
2-Methylnaphthalene	ND	U	0.25	0.12	1	07/03/18	07/10/18	KWG1803256	
Acenaphthylene	ND	U	0.13	0.046	1	07/03/18	07/10/18	KWG1803256	
Acenaphthene	ND	U	0.13	0.047	1	07/03/18	07/10/18	KWG1803256	
Fluorene	ND	U	0.13	0.052	1	07/03/18	07/10/18	KWG1803256	
Phenanthrene	ND	U	0.13	0.066	1	07/03/18	07/10/18	KWG1803256	
Anthracene	ND	U	0.13	0.038	1	07/03/18	07/10/18	KWG1803256	
Fluoranthene	ND	U	0.13	0.049	1	07/03/18	07/10/18	KWG1803256	
Pyrene	ND	U	0.13	0.050	1	07/03/18	07/10/18	KWG1803256	
Benz(a)anthracene	ND	U	0.13	0.038	1	07/03/18	07/10/18	KWG1803256	
Chrysene	ND	U	0.13	0.055	1	07/03/18	07/10/18	KWG1803256	
Benzo(b)fluoranthene†	ND	U	0.13	0.066	1	07/03/18	07/10/18	KWG1803256	
Benzo(k)fluoranthene	ND	U	0.13	0.057	1	07/03/18	07/10/18	KWG1803256	
Benzo(a)pyrene	ND	U	0.13	0.073	1	07/03/18	07/10/18	KWG1803256	
Indeno(1,2,3-cd)pyrene	ND	U	0.13	0.096	1	07/03/18	07/10/18	KWG1803256	
Dibenz(a,h)anthracene	ND	U	0.13	0.086	1	07/03/18	07/10/18	KWG1803256	
Benzo(g,h,i)perylene	ND	U	0.13	0.095	1	07/03/18	07/10/18	KWG1803256	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Fluorene-d10	61	42-106	07/10/18	Acceptable
Fluoranthene-d10	66	45-109	07/10/18	Acceptable
Terphenyl-d14	75	41-102	07/10/18	Acceptable

## † Analyte Comments

Benzo(b)fluoranthene This analyte cannot be separated from Benzo(j)fluoranthene.

Comments: \_\_\_\_\_

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985

**Surrogate Recovery Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Extraction Method:** EPA 3541      **Units:** Percent  
**Analysis Method:** 8270D SIM      **Level:** Low

<b>Sample Name</b>	<b>Lab Code</b>	<b>Sur1</b>	<b>Sur2</b>	<b>Sur3</b>
PDI-SG-S204	K1804985-001	64 D	69 D	77 D
PDI-SG-S147	K1804985-002	74 D	82 D	92 D
PDI-SG-S084	K1804985-003	66 D	75 D	90 D
PDI-SG-S090	K1804985-004	63 D	68 D	82 D
PDI-SG-S010	K1804985-005	64	74	86
PDI-SG-S255	K1804985-006	60	67	80
PDI-SG-S097	K1804985-007	72 D	81 D	96 D
PDI-SG-S115	K1804985-008	60 D	69 D	82 D
PDI-SG-S078	K1804985-009	67 D	78 D	88 D
PDI-SG-S135	K1804985-010	67 D	81 D	88 D
PDI-SG-S157	K1804985-011	61	74	82
Method Blank	KWG1803256-4	61	66	75
PDI-SG-S204MS	KWG1803256-1	75 D	68 D	80 D
PDI-SG-S204DMS	KWG1803256-2	73 D	67 D	78 D
Lab Control Sample	KWG1803256-3	56	62	72

**Surrogate Recovery Control Limits (%)**

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Sur1 = Fluorene-d10	42-106
Sur2 = Fluoranthene-d10	45-109
Sur3 = Terphenyl-d14	41-102

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Results flagged with an asterisk (\*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/10/2018  
**Time Analyzed:** 15:36

**Internal Standard Area and RT Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS20\DATA\071018\0710F017.D  
**Instrument ID:** MS20  
**Analysis Method:** 8270D SIM

**Lab Code:** KWG1803364-2  
**Analysis Lot:** KWG1803364

	Naphthalene-d8		Acenaphthene-d10		Phenanthrene-d10	
	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>
<b>Results ==&gt;</b>	89,619	5.93	47,018	8.24	100,099	11.42
<b>Upper Limit ==&gt;</b>	179,238	6.43	94,036	8.74	200,198	11.92
<b>Lower Limit ==&gt;</b>	44,810	5.43	23,509	7.74	50,050	10.92
<b>ICAL Result ==&gt;</b>	90,101	6.06	44,197	8.42	87,517	11.64

**Associated Analyses**

Method Blank	KWG1803256-4	78,166	5.94	43,527	8.25	87,936	11.43
Lab Control Sample	KWG1803256-3	76,732	5.93	40,988	8.24	84,096	11.42
PDI-SG-S010	K1804985-005	69,421	5.94	39,089	8.25	75,395	11.43
PDI-SG-S255	K1804985-006	70,351	5.94	39,166	8.25	78,045	11.43
PDI-SG-S157	K1804985-011	68,836	5.94	38,733	8.25	71,998	11.43
PDI-SG-S204	K1804985-001	71,423	5.94	41,045	8.25	82,385	11.43
PDI-SG-S147	K1804985-002	75,771	5.93	41,527	8.25	82,291	11.43
PDI-SG-S084	K1804985-003	77,730	5.93	42,090	8.25	85,825	11.43
PDI-SG-S090	K1804985-004	76,666	5.94	44,290	8.25	88,964	11.43
PDI-SG-S135	K1804985-010	78,031	5.94	44,853	8.25	91,701	11.43
PDI-SG-S097	K1804985-007	82,281	5.93	44,652	8.24	89,720	11.42
PDI-SG-S115	K1804985-008	80,286	5.93	42,575	8.25	86,490	11.42

Results flagged with an asterisk (\*) indicate values outside control criteria.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/10/2018  
**Time Analyzed:** 15:36

**Internal Standard Area and RT Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS20\DATA\071018\0710F017.D  
**Instrument ID:** MS20  
**Analysis Method:** 8270D SIM

**Lab Code:** KWG1803364-2  
**Analysis Lot:** KWG1803364

	Chrysene-d12		Perylene-d12	
	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>
<b>Results ==&gt;</b>	110,826	18.78	122,835	23.09
<b>Upper Limit ==&gt;</b>	221,652	19.28	245,670	23.59
<b>Lower Limit ==&gt;</b>	55,413	18.28	61,418	22.59
<b>ICAL Result ==&gt;</b>	105,110	19.00	102,151	23.35

*Associated Analyses*

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Method Blank	KWG1803256-4	95,568	18.79	108,069	23.09
Lab Control Sample	KWG1803256-3	88,463	18.78	101,452	23.08
PDI-SG-S010	K1804985-005	78,327	18.81	89,758	23.13
PDI-SG-S255	K1804985-006	79,457	18.80	92,556	23.13
PDI-SG-S157	K1804985-011	76,943	18.82	89,793	23.14
PDI-SG-S204	K1804985-001	86,335	18.80	97,504	23.11
PDI-SG-S147	K1804985-002	85,995	18.80	99,481	23.12
PDI-SG-S084	K1804985-003	84,276	18.80	102,887	23.10
PDI-SG-S090	K1804985-004	91,310	18.79	107,749	23.09
PDI-SG-S135	K1804985-010	96,836	18.79	109,351	23.09
PDI-SG-S097	K1804985-007	88,491	18.79	102,725	23.10
PDI-SG-S115	K1804985-008	85,290	18.79	101,456	23.10

Results flagged with an asterisk (\*) indicate values outside control criteria.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/12/2018  
**Time Analyzed:** 06:21

**Internal Standard Area and RT Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS20\DATA\071218\0712F002.D  
**Instrument ID:** MS20  
**Analysis Method:** 8270D SIM

**Lab Code:** KWG1803435-2  
**Analysis Lot:** KWG1803435

	Naphthalene-d8		Acenaphthene-d10		Phenanthrene-d10	
	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>
<b>Results ==&gt;</b>	88,527	5.93	47,488	8.25	99,112	11.42
<b>Upper Limit ==&gt;</b>	177,054	6.43	94,976	8.75	198,224	11.92
<b>Lower Limit ==&gt;</b>	44,264	5.43	23,744	7.75	49,556	10.92
<b>ICAL Result ==&gt;</b>	90,101	6.06	44,197	8.42	87,517	11.64

**Associated Analyses**

PDI-SG-S078	K1804985-009	83,640	5.93	45,475	8.24	88,725	11.42
PDI-SG-S010DL	K1804985-005	76,768	5.93	43,852	8.25	88,314	11.42
PDI-SG-S255DL	K1804985-006	80,247	5.93	44,926	8.25	90,900	11.42
PDI-SG-S084DL	K1804985-003	78,738	5.93	44,617	8.25	90,124	11.43
PDI-SG-S097DL	K1804985-007	83,295	5.93	45,749	8.25	91,783	11.42
PDI-SG-S115DL	K1804985-008	80,083	5.93	44,661	8.25	92,248	11.42

Results flagged with an asterisk (\*) indicate values outside control criteria.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/12/2018  
**Time Analyzed:** 06:21

**Internal Standard Area and RT Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS20\DATA\071218\0712F002.D  
**Instrument ID:** MS20  
**Analysis Method:** 8270D SIM

**Lab Code:** KWG1803435-2  
**Analysis Lot:** KWG1803435

	Chrysene-d12		Perylene-d12	
	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>
<b>Results ==&gt;</b>	109,372	18.78	120,089	23.09
<b>Upper Limit ==&gt;</b>	218,744	19.28	240,178	23.59
<b>Lower Limit ==&gt;</b>	54,686	18.28	60,045	22.59
<b>ICAL Result ==&gt;</b>	105,110	19.00	102,151	23.35

*Associated Analyses*

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PDI-SG-S078	K1804985-009	92,097	18.79	106,793	23.09
PDI-SG-S010DL	K1804985-005	93,198	18.79	105,736	23.09
PDI-SG-S255DL	K1804985-006	96,322	18.79	108,770	23.09
PDI-SG-S084DL	K1804985-003	95,946	18.79	109,520	23.09
PDI-SG-S097DL	K1804985-007	92,958	18.79	108,239	23.09
PDI-SG-S115DL	K1804985-008	96,123	18.79	110,888	23.09

Results flagged with an asterisk (\*) indicate values outside control criteria.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/25/2018  
**Time Analyzed:** 23:20

**Internal Standard Area and RT Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS14\DATA\072518\0725F028.D  
**Instrument ID:** MS14  
**Analysis Method:** 8270D SIM

**Lab Code:** KWG1803644-2  
**Analysis Lot:** KWG1803644

	Naphthalene-d8		Acenaphthene-d10		Phenanthrene-d10	
	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>
<b>Results ==&gt;</b>	47,711	4.80	22,281	6.36	49,366	7.59
<b>Upper Limit ==&gt;</b>	95,422	5.30	44,562	6.86	98,732	8.09
<b>Lower Limit ==&gt;</b>	23,856	4.30	11,141	5.86	24,683	7.09
<b>ICAL Result ==&gt;</b>	60,855	4.80	25,959	6.35	53,916	7.59

*Associated Analyses*

PDI-SG-S204MS	KWG1803256-1	47,011	4.80	24,742	6.36	50,688	7.59
PDI-SG-S204DMS	KWG1803256-2	49,241	4.81	25,955	6.36	53,157	7.60

Results flagged with an asterisk (\*) indicate values outside control criteria.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/25/2018  
**Time Analyzed:** 23:20

**Internal Standard Area and RT Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS14\DATA\072518\0725F028.D  
**Instrument ID:** MS14  
**Analysis Method:** 8270D SIM

**Lab Code:** KWG1803644-2  
**Analysis Lot:** KWG1803644

	Chrysene-d12		Perylene-d12	
	<u>Area</u>	<u>RT</u>	<u>Area</u>	<u>RT</u>
<b>Results ==&gt;</b>	52,395	10.16	54,650	13.32
<b>Upper Limit ==&gt;</b>	104,790	10.66	109,300	13.82
<b>Lower Limit ==&gt;</b>	26,198	9.66	27,325	12.82
<b>ICAL Result ==&gt;</b>	68,964	10.15	73,742	13.30

*Associated Analyses*

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PDI-SG-S204MS	KWG1803256-1	49,546	10.16	58,017	13.34
PDI-SG-S204DMS	KWG1803256-2	53,359	10.17	61,997	13.35

Results flagged with an asterisk (\*) indicate values outside control criteria.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Extracted:** 07/03/2018  
**Date Analyzed:** 07/26/2018

**Matrix Spike/Duplicate Matrix Spike Summary**  
**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	PDI-SG-S204	<b>Units:</b>	ug/Kg
<b>Lab Code:</b>	K1804985-001	<b>Basis:</b>	Dry
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM	<b>Extraction Lot:</b>	KWG1803256

<b>Analyte Name</b>	<b>Sample Result</b>	PDI-SG-S204MS			PDI-SG-S204DMS			<b>%Rec Limits</b>	<b>RPD</b>	<b>RPD Limit</b>			
		KWG1803256-1			KWG1803256-2								
		Matrix Spike			Duplicate Matrix Spike								
		<b>Result</b>	<b>Spike Amount</b>	<b>%Rec</b>	<b>Result</b>	<b>Spike Amount</b>	<b>%Rec</b>						
Naphthalene	7.9	47.0	88.7	44	55.9	88.4	54	37-104	17	40			
2-Methylnaphthalene	ND	55.1	88.7	62	60.2	88.4	68	39-115	9	40			
Acenaphthylene	6.5	54.3	88.7	54	59.2	88.4	60	39-115	9	40			
Acenaphthene	6.7	62.5	88.7	63	62.1	88.4	63	41-116	1	40			
Fluorene	9.2	75.2	88.7	74	67.0	88.4	65	43-117	12	40			
Phenanthrene	69	255	88.7	209 *	222	88.4	173 *	42-119	14	40			
Anthracene	12	75.7	88.7	71	70.8	88.4	66	42-124	7	40			
Fluoranthene	170	381	88.7	241 *	332	88.4	187 *	42-130	14	40			
Pyrene	250	373	88.7	143 *	338	88.4	103	33-125	10	40			
Benz(a)anthracene	55	145	88.7	101	154	88.4	112	42-123	6	40			
Chrysene	86	232	88.7	165 *	266	88.4	204 *	40-134	14	40			
Benzo(b)fluoranthene	85	229	88.7	161 *	239	88.4	174 *	27-139	5	40			
Benzo(k)fluoranthene	26	120	88.7	106	121	88.4	107	40-125	0	40			
Benzo(a)pyrene	54	151	88.7	109	182	88.4	144 *	39-130	18	40			
Indeno(1,2,3-cd)pyrene	54	114	88.7	67	142	88.4	100	37-143	22	40			
Dibenz(a,h)anthracene	16	51.1	88.7	40	60.3	88.4	50	39-141	16	40			
Benzo(g,h,i)perylene	73	140	88.7	76	164	88.4	103	35-140	16	40			

Results flagged with an asterisk (\*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Extracted:** 07/03/2018  
**Date Analyzed:** 07/10/2018

**Lab Control Spike Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Extraction Method:** EPA 3541  
**Analysis Method:** 8270D SIM

**Units:** ug/Kg  
**Basis:** Dry  
**Level:** Low  
**Extraction Lot:** KWG1803256

Lab Control Sample

KWG1803256-3

**Lab Control Spike**

<b>Analyte Name</b>	<b>Result</b>	<b>Spike</b>	<b>%Rec</b>	<b>%Rec</b> Limits
		<b>Amount</b>		
Naphthalene	55.0	100	55	42-107
2-Methylnaphthalene	54.2	100	54	40-116
Acenaphthylene	56.9	100	57	41-112
Acenaphthene	53.9	100	54	43-113
Fluorene	54.4	100	54	44-114
Phenanthrene	53.4	100	53	44-115
Anthracene	56.6	100	57	45-121
Fluoranthene	57.8	100	58	47-123
Pyrene	67.2	100	67	41-121
Benz(a)anthracene	71.9	100	72	42-123
Chrysene	66.8	100	67	46-130
Benzo(b)fluoranthene	65.6	100	66	46-125
Benzo(k)fluoranthene	62.3	100	62	47-125
Benzo(a)pyrene	65.3	100	65	45-128
Indeno(1,2,3-cd)pyrene	66.4	100	66	45-128
Dibenz(a,h)anthracene	65.6	100	66	44-128
Benzo(g,h,i)perylene	58.7	100	59	43-125

Results flagged with an asterisk (\*) indicate values outside control criteria.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Extracted:** 07/03/2018  
**Date Analyzed:** 07/10/2018  
**Time Analyzed:** 16:54

**Method Blank Summary**  
**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	Method Blank	<b>Instrument ID:</b>	MS20
<b>Lab Code:</b>	KWG1803256-4	<b>File ID:</b>	J:\MS20\DATA\071018\0710F019.D
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM	<b>Extraction Lot:</b>	KWG1803256

This Method Blank applies to the following analyses:

<b>Sample Name</b>	<b>Lab Code</b>	<b>File ID</b>	<b>Date Analyzed</b>	<b>Time Analyzed</b>
Lab Control Sample	KWG1803256-3	J:\MS20\DATA\071018\0710F020.D	07/10/18	17:34
PDI-SG-S010	K1804985-005	J:\MS20\DATA\071018\0710F023.D	07/10/18	19:32
PDI-SG-S255	K1804985-006	J:\MS20\DATA\071018\0710F024.D	07/10/18	20:11
PDI-SG-S157	K1804985-011	J:\MS20\DATA\071018\0710F025.D	07/10/18	20:50
PDI-SG-S204	K1804985-001	J:\MS20\DATA\071018\0710F026.D	07/10/18	21:29
PDI-SG-S147	K1804985-002	J:\MS20\DATA\071018\0710F029.D	07/10/18	23:27
PDI-SG-S084	K1804985-003	J:\MS20\DATA\071018\0710F030.D	07/11/18	00:06
PDI-SG-S090	K1804985-004	J:\MS20\DATA\071018\0710F031.D	07/11/18	00:45
PDI-SG-S135	K1804985-010	J:\MS20\DATA\071018\0710F032.D	07/11/18	01:25
PDI-SG-S097	K1804985-007	J:\MS20\DATA\071018\0710F033.D	07/11/18	02:04
PDI-SG-S115	K1804985-008	J:\MS20\DATA\071018\0710F034.D	07/11/18	02:43
PDI-SG-S078	K1804985-009	J:\MS20\DATA\071218\0712F007.D	07/12/18	09:37
PDI-SG-S010	K1804985-005	J:\MS20\DATA\071218\0712F008.D	07/12/18	10:16
PDI-SG-S255	K1804985-006	J:\MS20\DATA\071218\0712F009.D	07/12/18	10:55
PDI-SG-S084	K1804985-003	J:\MS20\DATA\071218\0712F010.D	07/12/18	11:35
PDI-SG-S097	K1804985-007	J:\MS20\DATA\071218\0712F011.D	07/12/18	12:14
PDI-SG-S115	K1804985-008	J:\MS20\DATA\071218\0712F012.D	07/12/18	12:54
PDI-SG-S204MS	KWG1803256-1	J:\MS14\DATA\072518\0725F044.D	07/26/18	06:13
PDI-SG-S204DMS	KWG1803256-2	J:\MS14\DATA\072518\0725F045.D	07/26/18	06:39

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Extracted:** 07/03/2018  
**Date Analyzed:** 07/10/2018  
**Time Analyzed:** 17:34

**Lab Control Sample Summary**  
**Polynuclear Aromatic Hydrocarbons**

<b>Sample Name:</b>	Lab Control Sample	<b>Instrument ID:</b>	MS20
<b>Lab Code:</b>	KWG1803256-3	<b>File ID:</b>	J:\MS20\DATA\071018\0710F020.D
<b>Extraction Method:</b>	EPA 3541	<b>Level:</b>	Low
<b>Analysis Method:</b>	8270D SIM	<b>Extraction Lot:</b>	KWG1803256

This Lab Control Sample applies to the following analyses:

<b>Sample Name</b>	<b>Lab Code</b>	<b>File ID</b>	<b>Date Analyzed</b>	<b>Time Analyzed</b>
Method Blank	KWG1803256-4	J:\MS20\DATA\071018\0710F019.D	07/10/18	16:54
PDI-SG-S010	K1804985-005	J:\MS20\DATA\071018\0710F023.D	07/10/18	19:32
PDI-SG-S255	K1804985-006	J:\MS20\DATA\071018\0710F024.D	07/10/18	20:11
PDI-SG-S157	K1804985-011	J:\MS20\DATA\071018\0710F025.D	07/10/18	20:50
PDI-SG-S204	K1804985-001	J:\MS20\DATA\071018\0710F026.D	07/10/18	21:29
PDI-SG-S147	K1804985-002	J:\MS20\DATA\071018\0710F029.D	07/10/18	23:27
PDI-SG-S084	K1804985-003	J:\MS20\DATA\071018\0710F030.D	07/11/18	00:06
PDI-SG-S090	K1804985-004	J:\MS20\DATA\071018\0710F031.D	07/11/18	00:45
PDI-SG-S135	K1804985-010	J:\MS20\DATA\071018\0710F032.D	07/11/18	01:25
PDI-SG-S097	K1804985-007	J:\MS20\DATA\071018\0710F033.D	07/11/18	02:04
PDI-SG-S115	K1804985-008	J:\MS20\DATA\071018\0710F034.D	07/11/18	02:43
PDI-SG-S078	K1804985-009	J:\MS20\DATA\071218\0712F007.D	07/12/18	09:37
PDI-SG-S010	K1804985-005	J:\MS20\DATA\071218\0712F008.D	07/12/18	10:16
PDI-SG-S255	K1804985-006	J:\MS20\DATA\071218\0712F009.D	07/12/18	10:55
PDI-SG-S084	K1804985-003	J:\MS20\DATA\071218\0712F010.D	07/12/18	11:35
PDI-SG-S097	K1804985-007	J:\MS20\DATA\071218\0712F011.D	07/12/18	12:14
PDI-SG-S115	K1804985-008	J:\MS20\DATA\071218\0712F012.D	07/12/18	12:54
PDI-SG-S204MS	KWG1803256-1	J:\MS14\DATA\072518\0725F044.D	07/26/18	06:13
PDI-SG-S204DMS	KWG1803256-2	J:\MS14\DATA\072518\0725F045.D	07/26/18	06:39

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/10/2018  
**Time Analyzed:** 14:56

**Tune Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS20\DATA\071018\0710F016.D

**Instrument ID:** MS20

**Column:**

**Analysis Method:** 8270D SIM  
**Analysis Lot:** KWG1803364

Target Mass	Relative to Mass	Lower Limit%	Upper Limit%	Relative Abundance %	Raw Abundance	Result Pass/Fail
51	198	10	80	26.5	159168	PASS
68	69	0	2	0.0	0	PASS
69	198	0	100	27.6	165866	PASS
70	69	0	2	0.5	901	PASS
127	198	10	80	39.4	236352	PASS
197	198	0	2	0.0	0	PASS
198	442	30	100	41.3	600106	PASS
199	198	5	9	6.7	40312	PASS
275	198	10	60	37.4	224725	PASS
365	442	1	50	2.5	36480	PASS
441	443	0	100	82.7	243157	PASS
442	442	100	100	100.0	1454250	PASS
443	442	15	24	20.2	294037	PASS

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed	Q
Continuing Calibration Verification	KWG1803364-2	J:\MS20\DATA\071018\0710F017.D	07/10/2018	15:36	
Method Blank	KWG1803256-4	J:\MS20\DATA\071018\0710F019.D	07/10/2018	16:54	
Lab Control Sample	KWG1803256-3	J:\MS20\DATA\071018\0710F020.D	07/10/2018	17:34	
PDI-SG-S010	K1804985-005	J:\MS20\DATA\071018\0710F023.D	07/10/2018	19:32	
PDI-SG-S255	K1804985-006	J:\MS20\DATA\071018\0710F024.D	07/10/2018	20:11	
PDI-SG-S157	K1804985-011	J:\MS20\DATA\071018\0710F025.D	07/10/2018	20:50	
PDI-SG-S204	K1804985-001	J:\MS20\DATA\071018\0710F026.D	07/10/2018	21:29	
PDI-SG-S147	K1804985-002	J:\MS20\DATA\071018\0710F029.D	07/10/2018	23:27	
PDI-SG-S084	K1804985-003	J:\MS20\DATA\071018\0710F030.D	07/11/2018	00:06	
PDI-SG-S090	K1804985-004	J:\MS20\DATA\071018\0710F031.D	07/11/2018	00:45	
PDI-SG-S135	K1804985-010	J:\MS20\DATA\071018\0710F032.D	07/11/2018	01:25	
PDI-SG-S097	K1804985-007	J:\MS20\DATA\071018\0710F033.D	07/11/2018	02:04	
PDI-SG-S115	K1804985-008	J:\MS20\DATA\071018\0710F034.D	07/11/2018	02:43	

Results flagged with an asterisk (\*) indicate the analysis performed outside specified tune window

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/12/2018  
**Time Analyzed:** 05:41

**Tune Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS20\DATA\071218\0712F001.D

**Instrument ID:** MS20

**Column:**

**Analysis Method:** 8270D SIM  
**Analysis Lot:** KWG1803435

Target Mass	Relative to Mass	Lower Limit%	Upper Limit%	Relative Abundance %	Raw Abundance	Result Pass/Fail
51	198	10	80	24.7	141898	PASS
68	69	0	2	0.0	0	PASS
69	198	0	100	26.5	152260	PASS
70	69	0	2	0.2	333	PASS
127	198	10	80	37.3	214101	PASS
197	198	0	2	0.0	0	PASS
198	442	30	100	38.7	573760	PASS
199	198	5	9	6.8	39146	PASS
275	198	10	60	37.5	215018	PASS
365	442	1	50	2.4	35437	PASS
441	443	0	100	84.8	249045	PASS
442	442	100	100	100.0	1480704	PASS
443	442	15	24	19.8	293546	PASS

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed	Q
Continuing Calibration Verification	KWG1803435-2	J:\MS20\DATA\071218\0712F002.D	07/12/2018	06:21	
PDI-SG-S078	K1804985-009	J:\MS20\DATA\071218\0712F007.D	07/12/2018	09:37	
PDI-SG-S010	K1804985-005	J:\MS20\DATA\071218\0712F008.D	07/12/2018	10:16	
PDI-SG-S255	K1804985-006	J:\MS20\DATA\071218\0712F009.D	07/12/2018	10:55	
PDI-SG-S084	K1804985-003	J:\MS20\DATA\071218\0712F010.D	07/12/2018	11:35	
PDI-SG-S097	K1804985-007	J:\MS20\DATA\071218\0712F011.D	07/12/2018	12:14	
PDI-SG-S115	K1804985-008	J:\MS20\DATA\071218\0712F012.D	07/12/2018	12:54	

Results flagged with an asterisk (\*) indicate the analysis performed outside specified tune window

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/25/2018  
**Time Analyzed:** 22:54

**Tune Summary**  
**Polynuclear Aromatic Hydrocarbons**

**File ID:** J:\MS14\DATA\072518\0725F027.D

**Instrument ID:** MS14

**Column:**

**Analysis Method:** 8270D SIM  
**Analysis Lot:** KWG1803644

Target Mass	Relative to Mass	Lower Limit%	Upper Limit%	Relative Abundance %	Raw Abundance	Result Pass/Fail
443	442	15	24	18.6	46781	PASS
51	198	10	80	33.6	42605	PASS
68	69	0	2	0.0	0	PASS
69	198	0	100	36.2	45874	PASS
70	69	0	2	0.6	287	PASS
127	198	10	80	43.1	54618	PASS
197	198	0	2	0.0	0	PASS
198	442	30	100	50.5	126613	PASS
199	198	5	9	6.5	8255	PASS
275	198	10	60	36.6	46320	PASS
365	442	1	50	2.8	6992	PASS
441	443	0	100	78.1	36552	PASS
442	442	100	100	100.0	250837	PASS

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed	Q
Continuing Calibration Verification	KWG1803644-2	J:\MS14\DATA\072518\0725F028.D	07/25/2018	23:20	
PDI-SG-S204MS	KWG1803256-1	J:\MS14\DATA\072518\0725F044.D	07/26/2018	06:13	
PDI-SG-S204DMS	KWG1803256-2	J:\MS14\DATA\072518\0725F045.D	07/26/2018	06:39	

Results flagged with an asterisk (\*) indicate the analysis performed outside specified tune window

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 07/11/2018

**Initial Calibration Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration ID:** CAL15779  
**Instrument ID:** MS14

**Column:** MS

Level ID	File ID	Level ID	File ID
A	J:\MS14\DATA\071118\0711F003.D	G	J:\MS14\DATA\071118\0711F009.D
B	J:\MS14\DATA\071118\0711F004.D	H	J:\MS14\DATA\071118\0711F010.D
C	J:\MS14\DATA\071118\0711F005.D	I	J:\MS14\DATA\071118\0711F011.D
D	J:\MS14\DATA\071118\0711F006.D	J	J:\MS14\DATA\071118\0711F012.D
E	J:\MS14\DATA\071118\0711F007.D		
F	J:\MS14\DATA\071118\0711F008.D		

Analyte Name	Level			Level			Level			Level			Level		
	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF
Naphthalene	A	2.0	1.42	B	4.0	1.31	C	8.0	1.19	D	20	1.15	E	100	1.13
	F	200	1.11	G	400	1.09	H	1000	1.06	I	1600	1.03	J	2000	1.07
2-Methylnaphthalene	A	2.0	0.789	B	4.0	0.741	C	8.0	0.734	D	20	0.715	E	100	0.718
	F	200	0.692	G	400	0.662	H	1000	0.650	I	1600	0.627	J	2000	0.635
Acenaphthylene	A	2.0	2.52	B	4.0	2.35	C	8.0	2.29	D	20	2.33	E	100	2.37
	F	200	2.41	G	400	2.42	H	1000	2.41	I	1600	2.43	J	2000	2.34
Acenaphthene	A	2.0	1.44	B	4.0	1.38	C	8.0	1.33	D	20	1.38	E	100	1.38
	F	200	1.39	G	400	1.38	H	1000	1.35	I	1600	1.38	J	2000	1.32
Fluorene	A	2.0	1.81	B	4.0	1.68	C	8.0	1.62	D	20	1.64	E	100	1.70
	F	200	1.68	G	400	1.67	H	1000	1.60	I	1600	1.63	J	2000	1.56
Phenanthrene	A	2.0	1.33	B	4.0	1.33	C	8.0	1.25	D	20	1.23	E	100	1.25
	F	200	1.25	G	400	1.22	H	1000	1.23	I	1600	1.22	J	2000	1.16
Anthracene	A	2.0	1.20	B	4.0	1.17	C	8.0	1.17	D	20	1.14	E	100	1.21
	F	200	1.22	G	400	1.22	H	1000	1.22	I	1600	1.21	J	2000	1.17
Fluoranthene	A	2.0	1.50	B	4.0	1.44	C	8.0	1.44	D	20	1.41	E	100	1.54
	F	200	1.60	G	400	1.58	H	1000	1.63	I	1600	1.65	J	2000	1.62
Pyrene	A	2.0	1.49	B	4.0	1.46	C	8.0	1.38	D	20	1.37	E	100	1.34
	F	200	1.29	G	400	1.32	H	1000	1.33	I	1600	1.39	J	2000	1.36
Benz(a)anthracene	A	2.0	1.50	B	4.0	1.32	C	8.0	1.26	D	20	1.19	E	100	1.20
	F	200	1.23	G	400	1.27	H	1000	1.32	I	1600	1.34	J	2000	1.31
Chrysene	A	2.0	1.23	B	4.0	1.21	C	8.0	1.23	D	20	1.20	E	100	1.24
	F	200	1.25	G	400	1.25	H	1000	1.26	I	1600	1.28	J	2000	1.24
Benzo(b)fluoranthene	A	2.0	1.19	B	4.0	1.16	C	8.0	1.15	D	20	1.16	E	100	1.20
	F	200	1.26	G	400	1.30	H	1000	1.34	I	1600	1.32	J	2000	1.27
Benzo(k)fluoranthene	A	2.0	1.16	B	4.0	1.17	C	8.0	1.13	D	20	1.19	E	100	1.23
	F	200	1.26	G	400	1.29	H	1000	1.29	I	1600	1.29	J	2000	1.25
Benzo(a)pyrene	A	2.0	1.09	B	4.0	1.01	C	8.0	1.03	D	20	1.04	E	100	1.05
	F	200	1.07	G	400	1.11	H	1000	1.15	I	1600	1.16	J	2000	1.12

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 07/11/2018

**Initial Calibration Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration ID:** CAL15779  
**Instrument ID:** MS14

**Column:** MS

Analyte Name	Level	ID	Amt	RRF	Level	ID	Amt	RRF	Level	ID	Amt	RRF	Level	ID	Amt	RRF	
	A	2.0	1.23	B	4.0	1.10	C	8.0	1.06	D	20	1.06	E	100	1.07		
Indeno(1,2,3-cd)pyrene	F	200	1.08		G	400	1.05		H	1000	1.01		I	1600	0.996	J	2000 0.959
Dibenz(a,h)anthracene	A	2.0	1.12		B	4.0	1.10		C	8.0	1.09		D	20	1.17	E	100 1.12
	F	200	1.10		G	400	1.07		H	1000	1.02		I	1600	1.00	J	2000 0.971
Benzo(g,h,i)perylene	A	2.0	1.44		B	4.0	1.37		C	8.0	1.31		D	20	1.34	E	100 1.28
	F	200	1.27		G	400	1.21		H	1000	1.12		I	1600	1.07	J	2000 1.04
Fluorene-d10					B	4.0	1.49		C	8.0	1.33		D	20	1.28	E	100 1.25
					F	200	1.24		G	400	1.25		H	1000	1.21	I	1600 1.24
Fluoranthene-d10	A	2.0	1.18		B	4.0	1.19		C	8.0	1.15		D	20	1.11	E	100 1.20
	F	200	1.26		G	400	1.31		H	1000	1.40		I	1600	1.42	J	2000 1.39
Terphenyl-d14					B	4.0	1.05		C	8.0	0.934		D	20	0.867	E	100 0.832
					F	200	0.823		G	400	0.835		H	1000	0.837	I	1600 0.830
															J	2000 0.801	

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 07/11/2018

**Initial Calibration Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration ID:** CAL15779  
**Instrument ID:** MS14

**Column:** MS

<b>Analyte Name</b>	<b>Compound Type</b>	<b>Calibration Evaluation</b>				<b>RRF Evaluation</b>		
		<b>Fit Type</b>	<b>Eval.</b>	<b>Result</b>	<b>Q</b>	<b>Control Criteria</b>	<b>Average RRF</b>	<b>Q</b>
Naphthalene	MS	AverageRF	% RSD	10.7		≤ 20	1.15	0.70
2-Methylnaphthalene	MS	AverageRF	% RSD	7.5		≤ 20	0.696	0.40
Acenaphthylene	MS	AverageRF	% RSD	2.7		≤ 20	2.39	0.90
Acenaphthene	MS	AverageRF	% RSD	2.3		≤ 20	1.37	0.90
Fluorene	MS	AverageRF	% RSD	4.1		≤ 20	1.66	0.90
Phenanthrene	MS	AverageRF	% RSD	4.1		≤ 20	1.25	0.70
Anthracene	MS	AverageRF	% RSD	2.4		≤ 20	1.19	0.70
Fluoranthene	MS	AverageRF	% RSD	5.8		≤ 20	1.54	0.60
Pyrene	MS	AverageRF	% RSD	4.5		≤ 20	1.37	0.60
Benz(a)anthracene	MS	AverageRF	% RSD	6.8		≤ 20	1.29	0.80
Chrysene	MS	AverageRF	% RSD	1.9		≤ 20	1.24	0.70
Benzo(b)fluoranthene	MS	AverageRF	% RSD	5.7		≤ 20	1.24	0.70
Benzo(k)fluoranthene	MS	AverageRF	% RSD	4.7		≤ 20	1.23	0.70
Benzo(a)pyrene	MS	AverageRF	% RSD	4.7		≤ 20	1.08	0.70
Indeno(1,2,3-cd)pyrene	MS	AverageRF	% RSD	6.9		≤ 20	1.06	0.50
Dibenz(a,h)anthracene	MS	AverageRF	% RSD	5.7		≤ 20	1.08	0.40
Benzo(g,h,i)perylene	MS	AverageRF	% RSD	10.7		≤ 20	1.24	0.50
Fluorene-d10	SURR	AverageRF	% RSD	6.9		≤ 20	1.28	0.01
Fluoranthene-d10	SURR	AverageRF	% RSD	9.0		≤ 20	1.26	0.01
Terphenyl-d14	SURR	AverageRF	% RSD	8.9		≤ 20	0.867	0.01

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 07/11/2018  
**Date Analyzed:** 07/11/2018

**Second Source Calibration Verification**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration Type:** Internal Standard  
**Analysis Method:** 8270D SIM

**Calibration ID:** CAL15779  
**Units:** ng/ml

**File ID:** J:\MS14\DATA\071118\0711F013.D

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Naphthalene	400	370	1.15	1.06	-8	NA	± 30 %	AverageRF
2-Methylnaphthalene	400	390	0.696	0.683	-2	NA	± 30 %	AverageRF
Acenaphthylene	400	390	2.39	2.33	-2	NA	± 30 %	AverageRF
Acenaphthene	400	380	1.37	1.32	-4	NA	± 30 %	AverageRF
Fluorene	400	390	1.66	1.60	-4	NA	± 30 %	AverageRF
Phenanthrene	400	380	1.25	1.19	-4	NA	± 30 %	AverageRF
Anthracene	400	400	1.19	1.18	-1	NA	± 30 %	AverageRF
Fluoranthene	400	410	1.54	1.60	4	NA	± 30 %	AverageRF
Pyrene	400	350	1.37	1.21	-12	NA	± 30 %	AverageRF
Benz(a)anthracene	400	380	1.29	1.22	-6	NA	± 30 %	AverageRF
Chrysene	400	390	1.24	1.20	-3	NA	± 30 %	AverageRF
Benzo(b)fluoranthene	400	410	1.24	1.27	2	NA	± 30 %	AverageRF
Benzo(k)fluoranthene	400	420	1.23	1.28	4	NA	± 30 %	AverageRF
Benzo(a)pyrene	400	400	1.08	1.08	0	NA	± 30 %	AverageRF
Indeno(1,2,3-cd)pyrene	400	370	1.06	0.993	-7	NA	± 30 %	AverageRF
Dibenz(a,h)anthracene	400	380	1.08	1.03	-4	NA	± 30 %	AverageRF
Benzo(g,h,i)perylene	400	370	1.24	1.14	-9	NA	± 30 %	AverageRF

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 11/02/2017

**Initial Calibration Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration ID:** CAL15594  
**Instrument ID:** MS20

**Column:** MS

Level ID	File ID	Level ID	File ID
A	J:\MS20\DATA\110217\1102F003.D	G	J:\MS20\DATA\110217\1102F009.D
B	J:\MS20\DATA\110217\1102F004.D	H	J:\MS20\DATA\110217\1102F010.D
C	J:\MS20\DATA\110217\1102F005.D	I	J:\MS20\DATA\110217\1102F011.D
D	J:\MS20\DATA\110217\1102F006.D	J	J:\MS20\DATA\110217\1102F012.D
E	J:\MS20\DATA\110217\1102F007.D		
F	J:\MS20\DATA\110217\1102F008.D		

Analyte Name	Level			Level			Level			Level			Level		
	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF
Naphthalene	A	2.0	1.21	B	4.0	1.04	C	8.0	1.03	D	20	0.996	E	100	0.990
	F	200	0.989	G	400	1.02	H	1000	1.01	I	1600	0.972	J	2000	0.967
2-Methylnaphthalene	A	2.0	0.781	B	4.0	0.671	C	8.0	0.661	D	20	0.679	E	100	0.678
	F	200	0.694	G	400	0.692	H	1000	0.678	I	1600	0.645	J	2000	0.644
Acenaphthylene	A	2.0	2.23	B	4.0	1.95	C	8.0	1.92	D	20	1.97	E	100	1.97
	F	200	2.00	G	400	2.07	H	1000	2.08	I	1600	2.05	J	2000	2.02
Acenaphthene	A	2.0	1.45	B	4.0	1.25	C	8.0	1.25	D	20	1.24	E	100	1.21
	F	200	1.22	G	400	1.25	H	1000	1.23	I	1600	1.21	J	2000	1.19
Fluorene	A	2.0	1.81	B	4.0	1.46	C	8.0	1.44	D	20	1.44	E	100	1.46
	F	200	1.47	G	400	1.51	H	1000	1.48	I	1600	1.45	J	2000	1.44
Phenanthrene	A	2.0	1.50	B	4.0	1.16	C	8.0	1.16	D	20	1.16	E	100	1.12
	F	200	1.13	G	400	1.15	H	1000	1.14	I	1600	1.10	J	2000	1.09
Anthracene	A	2.0	1.32	B	4.0	1.05	C	8.0	1.04	D	20	1.03	E	100	1.03
	F	200	1.06	G	400	1.10	H	1000	1.11	I	1600	1.08	J	2000	1.06
Fluoranthene	A	2.0	1.51	B	4.0	1.20	C	8.0	1.19	D	20	1.24	E	100	1.23
	F	200	1.27	G	400	1.32	H	1000	1.30	I	1600	1.25	J	2000	1.24
Pyrene	A	2.0	1.34	B	4.0	1.10	C	8.0	1.09	D	20	1.10	E	100	1.08
	F	200	1.09	G	400	1.12	H	1000	1.13	I	1600	1.12	J	2000	1.11
Benz(a)anthracene	A	2.0	1.36	B	4.0	1.13	C	8.0	1.03	D	20	1.02	E	100	0.986
	F	200	1.01	G	400	1.05	H	1000	1.09	I	1600	1.09	J	2000	1.10
Chrysene	A	2.0	1.26	B	4.0	1.06	C	8.0	1.08	D	20	1.07	E	100	1.06
	F	200	1.06	G	400	1.07	H	1000	1.06	I	1600	1.06	J	2000	1.04
Benzo(b)fluoranthene	A	2.0	1.22	B	4.0	1.06	C	8.0	1.06	D	20	1.10	E	100	1.10
	F	200	1.12	G	400	1.17	H	1000	1.19	I	1600	1.20	J	2000	1.18
Benzo(k)fluoranthene	A	2.0	1.19	B	4.0	1.07	C	8.0	1.09	D	20	1.09	E	100	1.15
	F	200	1.18	G	400	1.20	H	1000	1.19	I	1600	1.19	J	2000	1.18
Benzo(a)pyrene	A	2.0	0.956	B	4.0	0.883	C	8.0	0.866	D	20	0.903	E	100	0.955
	F	200	1.00	G	400	1.06	H	1000	1.06	I	1600	1.08	J	2000	1.08

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 11/02/2017

**Initial Calibration Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration ID:** CAL15594  
**Instrument ID:** MS20

**Column:** MS

<b>Analyte Name</b>	Level A			Level B			Level C			Level D			Level E		
	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF	ID	Amt	RRF
Indeno(1,2,3-cd)pyrene	A	2.0	1.05	B	4.0	0.927	C	8.0	0.901	D	20	0.943	E	100	0.987
	F	200	1.03	G	400	1.10	H	1000	1.08	I	1600	1.07	J	2000	1.07
Dibenz(a,h)anthracene	A	2.0	1.02	B	4.0	0.967	C	8.0	0.953	D	20	1.01	E	100	1.05
	F	200	1.08	G	400	1.13	H	1000	1.11	I	1600	1.10	J	2000	1.10
Benzo(g,h,i)perylene	A	2.0	1.36	B	4.0	1.19	C	8.0	1.17	D	20	1.25	E	100	1.21
	F	200	1.23	G	400	1.27	H	1000	1.18	I	1600	1.17	J	2000	1.16
Fluorene-d10				B	4.0	1.51	C	8.0	1.31	D	20	1.26	E	100	1.23
	F	200	1.23	G	400	1.27	H	1000	1.25	I	1600	1.23	J	2000	1.21
Fluoranthene-d10	A	2.0	1.49	B	4.0	1.12	C	8.0	1.07	D	20	1.08	E	100	1.06
	F	200	1.10	G	400	1.15	H	1000	1.17	I	1600	1.16	J	2000	1.15
Terphenyl-d14	A	2.0	1.04	B	4.0	0.848	C	8.0	0.825	D	20	0.837	E	100	0.817
	F	200	0.826	G	400	0.831	H	1000	0.840	I	1600	0.837	J	2000	0.829

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 11/02/2017

**Initial Calibration Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration ID:** CAL15594  
**Instrument ID:** MS20

**Column:** MS

<b>Analyte Name</b>	<b>Compound Type</b>	<b>Calibration Evaluation</b>				<b>RRF Evaluation</b>		
		<b>Fit Type</b>	<b>Eval.</b>	<b>Result</b>	<b>Q</b>	<b>Control Criteria</b>	<b>Average RRF</b>	<b>Q</b>
Naphthalene	MS	AverageRF	% RSD	6.8		≤ 20	1.02	0.70
2-Methylnaphthalene	MS	AverageRF	% RSD	5.7		≤ 20	0.682	0.40
Acenaphthylene	MS	AverageRF	% RSD	4.4		≤ 20	2.03	0.90
Acenaphthene	MS	AverageRF	% RSD	5.8		≤ 20	1.25	0.90
Fluorene	MS	AverageRF	% RSD	7.5		≤ 20	1.50	0.90
Phenanthrene	MS	AverageRF	% RSD	10.0		≤ 20	1.17	0.70
Anthracene	MS	AverageRF	% RSD	7.9		≤ 20	1.09	0.70
Fluoranthene	MS	AverageRF	% RSD	7.2		≤ 20	1.27	0.60
Pyrene	MS	AverageRF	% RSD	6.7		≤ 20	1.13	0.60
Benz(a)anthracene	MS	AverageRF	% RSD	9.8		≤ 20	1.09	0.80
Chrysene	MS	AverageRF	% RSD	5.7		≤ 20	1.08	0.70
Benzo(b)fluoranthene	MS	AverageRF	% RSD	5.0		≤ 20	1.14	0.70
Benzo(k)fluoranthene	MS	AverageRF	% RSD	4.4		≤ 20	1.15	0.70
Benzo(a)pyrene	MS	AverageRF	% RSD	8.4		≤ 20	0.984	0.70
Indeno(1,2,3-cd)pyrene	MS	AverageRF	% RSD	7.1		≤ 20	1.02	0.50
Dibenz(a,h)anthracene	MS	AverageRF	% RSD	5.9		≤ 20	1.05	0.40
Benzo(g,h,i)perylene	MS	AverageRF	% RSD	5.1		≤ 20	1.22	0.50
Fluorene-d10	SURR	AverageRF	% RSD	7.1		≤ 20	1.28	0.01
Fluoranthene-d10	SURR	AverageRF	% RSD	10.7		≤ 20	1.15	0.01
Terphenyl-d14	SURR	AverageRF	% RSD	7.7		≤ 20	0.853	0.01

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Calibration Date:** 11/02/2017  
**Date Analyzed:** 11/02/2017

**Second Source Calibration Verification**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration Type:** Internal Standard  
**Analysis Method:** 8270D SIM

**Calibration ID:** CAL15594  
**Units:** ng/ml

**File ID:** J:\MS20\DATA\110217\1102F013.D

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Naphthalene	400	360	1.02	0.924	-10	NA	± 30 %	AverageRF
2-Methylnaphthalene	400	380	0.682	0.644	-6	NA	± 30 %	AverageRF
Acenaphthylene	400	380	2.03	1.93	-5	NA	± 30 %	AverageRF
Acenaphthene	400	380	1.25	1.17	-6	NA	± 30 %	AverageRF
Fluorene	400	370	1.50	1.39	-7	NA	± 30 %	AverageRF
Phenanthrene	400	370	1.17	1.08	-8	NA	± 30 %	AverageRF
Anthracene	400	380	1.09	1.03	-6	NA	± 30 %	AverageRF
Fluoranthene	400	400	1.27	1.28	0	NA	± 30 %	AverageRF
Pyrene	400	380	1.13	1.08	-5	NA	± 30 %	AverageRF
Benz(a)anthracene	400	370	1.09	0.996	-8	NA	± 30 %	AverageRF
Chrysene	400	380	1.08	1.03	-5	NA	± 30 %	AverageRF
Benzo(b)fluoranthene	400	380	1.14	1.10	-4	NA	± 30 %	AverageRF
Benzo(k)fluoranthene	400	390	1.15	1.13	-2	NA	± 30 %	AverageRF
Benzo(a)pyrene	400	400	0.984	0.973	-1	NA	± 30 %	AverageRF
Indeno(1,2,3-cd)pyrene	400	370	1.02	0.946	-7	NA	± 30 %	AverageRF
Dibenz(a,h)anthracene	400	380	1.05	1.00	-5	NA	± 30 %	AverageRF
Benzo(g,h,i)perylene	400	360	1.22	1.11	-9	NA	± 30 %	AverageRF

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/10/2018

**Continuing Calibration Verification Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration Type:** Internal Standard  
**Analysis Method:** 8270D SIM

**Calibration Date:** 11/02/2017  
**Calibration ID:** CAL15594  
**Analysis Lot:** KWG1803364  
**Units:** ng/ml

**File ID:** J:\MS20\DATA\071018\0710F017.D

Analyte Name	Expected	Result	Min RF	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Naphthalene	400	410	0.70	1.02	1.03	1	NA	± 20	AverageRF
2-Methylnaphthalene	400	400	0.40	0.682	0.684	0	NA	± 20	AverageRF
Acenaphthylene	400	420	0.90	2.03	2.12	4	NA	± 20	AverageRF
Acenaphthene	400	390	0.90	1.25	1.22	-2	NA	± 20	AverageRF
Fluorene	400	400	0.90	1.50	1.49	0	NA	± 20	AverageRF
Phenanthrene	400	380	0.70	1.17	1.10	-6	NA	± 20	AverageRF
Anthracene	400	350	0.70	1.09	0.956	-12	NA	± 20	AverageRF
Fluoranthene	400	410	0.60	1.27	1.31	3	NA	± 20	AverageRF
Pyrene	400	440	0.60	1.13	1.25	11	NA	± 20	AverageRF
Benz(a)anthracene	400	450	0.80	1.09	1.23	13	NA	± 20	AverageRF
Chrysene	400	420	0.70	1.08	1.13	4	NA	± 20	AverageRF
Benzo(b)fluoranthene	400	420	0.70	1.14	1.20	5	NA	± 20	AverageRF
Benzo(k)fluoranthene	400	410	0.70	1.15	1.19	4	NA	± 20	AverageRF
Benzo(a)pyrene	400	430	0.70	0.984	1.07	8	NA	± 20	AverageRF
Indeno(1,2,3-cd)pyrene	400	440	0.50	1.02	1.11	9	NA	± 20	AverageRF
Dibenz(a,h)anthracene	400	430	0.40	1.05	1.12	7	NA	± 20	AverageRF
Benzo(g,h,i)perylene	400	390	0.50	1.22	1.19	-3	NA	± 20	AverageRF
Fluorene-d10	400	390	0.01	1.28	1.26	-2	NA	± 20	AverageRF
Fluoranthene-d10	400	410	0.01	1.15	1.18	2	NA	± 20	AverageRF
Terphenyl-d14	400	410	0.01	0.853	0.880	3	NA	± 20	AverageRF

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/12/2018

**Continuing Calibration Verification Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration Type:** Internal Standard  
**Analysis Method:** 8270D SIM

**Calibration Date:** 11/02/2017  
**Calibration ID:** CAL15594  
**Analysis Lot:** KWG1803435  
**Units:** ng/ml

**File ID:** J:\MS20\DATA\071218\0712F002.D

Analyte Name	Expected	Result	Min RF	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Naphthalene	400	410	0.70	1.02	1.04	2	NA	± 20	AverageRF
2-Methylnaphthalene	400	410	0.40	0.682	0.691	1	NA	± 20	AverageRF
Acenaphthylene	400	410	0.90	2.03	2.10	4	NA	± 20	AverageRF
Acenaphthene	400	390	0.90	1.25	1.23	-2	NA	± 20	AverageRF
Fluorene	400	400	0.90	1.50	1.49	0	NA	± 20	AverageRF
Phenanthrene	400	380	0.70	1.17	1.12	-4	NA	± 20	AverageRF
Anthracene	400	360	0.70	1.09	0.979	-10	NA	± 20	AverageRF
Fluoranthene	400	410	0.60	1.27	1.31	3	NA	± 20	AverageRF
Pyrene	400	440	0.60	1.13	1.25	11	NA	± 20	AverageRF
Benz(a)anthracene	400	450	0.80	1.09	1.22	13	NA	± 20	AverageRF
Chrysene	400	420	0.70	1.08	1.14	5	NA	± 20	AverageRF
Benzo(b)fluoranthene	400	420	0.70	1.14	1.20	5	NA	± 20	AverageRF
Benzo(k)fluoranthene	400	420	0.70	1.15	1.21	5	NA	± 20	AverageRF
Benzo(a)pyrene	400	440	0.70	0.984	1.08	9	NA	± 20	AverageRF
Indeno(1,2,3-cd)pyrene	400	440	0.50	1.02	1.12	11	NA	± 20	AverageRF
Dibenz(a,h)anthracene	400	430	0.40	1.05	1.14	8	NA	± 20	AverageRF
Benzo(g,h,i)perylene	400	390	0.50	1.22	1.18	-3	NA	± 20	AverageRF
Fluorene-d10	400	390	0.01	1.28	1.25	-2	NA	± 20	AverageRF
Fluoranthene-d10	400	410	0.01	1.15	1.19	3	NA	± 20	AverageRF
Terphenyl-d14	400	420	0.01	0.853	0.901	6	NA	± 20	AverageRF

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985  
**Date Analyzed:** 07/25/2018

**Continuing Calibration Verification Summary**  
**Polynuclear Aromatic Hydrocarbons**

**Calibration Type:** Internal Standard  
**Analysis Method:** 8270D SIM

**Calibration Date:** 07/11/2018  
**Calibration ID:** CAL15779  
**Analysis Lot:** KWG1803644  
**Units:** ng/ml

**File ID:** J:\MS14\DATA\072518\0725F028.D

Analyte Name	Expected	Result	Min RF	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Naphthalene	400	380	0.70	1.15	1.10	-4	NA	± 20	AverageRF
2-Methylnaphthalene	400	400	0.40	0.696	0.698	0	NA	± 20	AverageRF
Acenaphthylene	400	410	0.90	2.39	2.42	1	NA	± 20	AverageRF
Acenaphthene	400	410	0.90	1.37	1.40	2	NA	± 20	AverageRF
Fluorene	400	430	0.90	1.66	1.76	6	NA	± 20	AverageRF
Phenanthrene	400	400	0.70	1.25	1.25	0	NA	± 20	AverageRF
Anthracene	400	390	0.70	1.19	1.17	-2	NA	± 20	AverageRF
Fluoranthene	400	370	0.60	1.54	1.44	-7	NA	± 20	AverageRF
Pyrene	400	430	0.60	1.37	1.46	7	NA	± 20	AverageRF
Benz(a)anthracene	400	410	0.80	1.29	1.34	4	NA	± 20	AverageRF
Chrysene	400	410	0.70	1.24	1.28	4	NA	± 20	AverageRF
Benzo(b)fluoranthene	400	430	0.70	1.24	1.32	7	NA	± 20	AverageRF
Benzo(k)fluoranthene	400	440	0.70	1.23	1.34	9	NA	± 20	AverageRF
Benzo(a)pyrene	400	410	0.70	1.08	1.12	3	NA	± 20	AverageRF
Indeno(1,2,3-cd)pyrene	400	400	0.50	1.06	1.06	0	NA	± 20	AverageRF
Dibenz(a,h)anthracene	400	400	0.40	1.08	1.08	0	NA	± 20	AverageRF
Benzo(g,h,i)perylene	400	410	0.50	1.24	1.27	2	NA	± 20	AverageRF
Fluorene-d10	400	450	0.01	1.28	1.44	13	NA	± 20	AverageRF
Fluoranthene-d10	400	420	0.01	1.26	1.32	5	NA	± 20	AverageRF
Terphenyl-d14	400	420	0.01	0.867	0.922	6	NA	± 20	AverageRF

Results flagged with an asterisk (\*) indicate values outside control criteria.

† SPCC Compound

‡ CCC Compound

**Client:**  
**Project:**

AECOM  
Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985

**Analysis Run Log**  
**Polynuclear Aromatic Hydrocarbons**

**Analysis Method:** 8270D SIM

**Analysis Lot:** KWG1803644  
**Instrument ID:** MS14

<b>File ID</b>	<b>Sample Name</b>	<b>Lab Code</b>	<b>Date Analysis Started</b>	<b>Start Time</b>	<b>Q</b>	<b>Date Analysis Finished</b>	<b>Finish Time</b>
0725F027.D	GC/MS Tuning - Decafluorotriphenylphosph	KWG1803644-1	7/25/2018	22:54		7/25/2018	23:14
0725F028.D	Continuing Calibration Verification	KWG1803644-2	7/25/2018	23:20		7/25/2018	23:39
0725F029.D	ZZZZZZ	ZZZZZZ	7/25/2018	23:45		7/26/2018	00:04
0725F030.D	ZZZZZZ	ZZZZZZ	7/26/2018	00:11		7/26/2018	00:30
0725F031.D	ZZZZZZ	ZZZZZZ	7/26/2018	00:37		7/26/2018	00:56
0725F032.D	ZZZZZZ	ZZZZZZ	7/26/2018	01:03		7/26/2018	01:22
0725F033.D	ZZZZZZ	ZZZZZZ	7/26/2018	01:29		7/26/2018	01:48
0725F034.D	ZZZZZZ	ZZZZZZ	7/26/2018	01:55		7/26/2018	02:14
0725F035.D	ZZZZZZ	ZZZZZZ	7/26/2018	02:20		7/26/2018	02:39
0725F036.D	ZZZZZZ	ZZZZZZ	7/26/2018	02:46		7/26/2018	03:05
0725F037.D	ZZZZZZ	ZZZZZZ	7/26/2018	03:12		7/26/2018	03:31
0725F038.D	ZZZZZZ	ZZZZZZ	7/26/2018	03:38		7/26/2018	03:57
0725F039.D	ZZZZZZ	ZZZZZZ	7/26/2018	04:04		7/26/2018	04:23
0725F040.D	ZZZZZZ	ZZZZZZ	7/26/2018	04:30		7/26/2018	04:49
0725F041.D	ZZZZZZ	ZZZZZZ	7/26/2018	04:56		7/26/2018	05:15
0725F042.D	ZZZZZZ	ZZZZZZ	7/26/2018	05:22		7/26/2018	05:41
0725F043.D	ZZZZZZ	ZZZZZZ	7/26/2018	05:47		7/26/2018	06:06
0725F044.D	PDI-SG-S204MS	KWG1803256-1	7/26/2018	06:13		7/26/2018	06:32
0725F045.D	PDI-SG-S204DMS	KWG1803256-2	7/26/2018	06:39		7/26/2018	06:58

Results flagged with an asterisk (\*) indicate the holding time was exceeded for the analysis

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985

**Analysis Run Log**  
**Polynuclear Aromatic Hydrocarbons**

**Analysis Method:** 8270D SIM

**Analysis Lot:** KWG1803364

**Instrument ID:** MS20

<b>File ID</b>	<b>Sample Name</b>	<b>Lab Code</b>	<b>Date Analysis Started</b>	<b>Start Time</b>	<b>Q</b>	<b>Date Analysis Finished</b>	<b>Finish Time</b>
0710F016.D	GC/MS Tuning - Decafluorotriphenylphosph	KWG1803364-1	7/10/2018	14:56		7/10/2018	15:26
0710F017.D	Continuing Calibration Verification	KWG1803364-2	7/10/2018	15:36		7/10/2018	16:05
0710F019.D	Method Blank	KWG1803256-4	7/10/2018	16:54		7/10/2018	17:23
0710F020.D	Lab Control Sample	KWG1803256-3	7/10/2018	17:34		7/10/2018	18:03
0710F021.D	ZZZZZZ	ZZZZZZ	7/10/2018	18:13		7/10/2018	18:42
0710F022.D	ZZZZZZ	ZZZZZZ	7/10/2018	18:52		7/10/2018	19:21
0710F023.D	PDI-SG-S010	K1804985-005	7/10/2018	19:32		7/10/2018	20:01
0710F024.D	PDI-SG-S255	K1804985-006	7/10/2018	20:11		7/10/2018	20:40
0710F025.D	PDI-SG-S157	K1804985-011	7/10/2018	20:50		7/10/2018	21:19
0710F026.D	PDI-SG-S204	K1804985-001	7/10/2018	21:29		7/10/2018	21:58
0710F027.D	ZZZZZZ	ZZZZZZ	7/10/2018	22:08		7/10/2018	22:37
0710F028.D	ZZZZZZ	ZZZZZZ	7/10/2018	22:48		7/10/2018	23:17
0710F029.D	PDI-SG-S147	K1804985-002	7/10/2018	23:27		7/10/2018	23:56
0710F030.D	PDI-SG-S084	K1804985-003	7/11/2018	00:06		7/11/2018	00:35
0710F031.D	PDI-SG-S090	K1804985-004	7/11/2018	00:45		7/11/2018	01:14
0710F032.D	PDI-SG-S135	K1804985-010	7/11/2018	01:25		7/11/2018	01:54
0710F033.D	PDI-SG-S097	K1804985-007	7/11/2018	02:04		7/11/2018	02:33
0710F034.D	PDI-SG-S115	K1804985-008	7/11/2018	02:43		7/11/2018	03:12

Results flagged with an asterisk (\*) indicate the holding time was exceeded for the analysis

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335

**Service Request:** K1804985

**Analysis Run Log**  
**Polynuclear Aromatic Hydrocarbons**

**Analysis Method:** 8270D SIM

**Analysis Lot:** KWG1803435

**Instrument ID:** MS20

<b>File ID</b>	<b>Sample Name</b>	<b>Lab Code</b>	<b>Date Analysis Started</b>	<b>Start Time</b>	<b>Q</b>	<b>Date Analysis Finished</b>	<b>Finish Time</b>
0712F001.D	GC/MS Tuning - Decafluorotriphenylphosph	KWG1803435-1	7/12/2018	05:41		7/12/2018	06:11
0712F002.D	Continuing Calibration Verification	KWG1803435-2	7/12/2018	06:21		7/12/2018	06:50
0712F003.D	ZZZZZZ	ZZZZZZ	7/12/2018	07:00		7/12/2018	07:29
0712F004.D	ZZZZZZ	ZZZZZZ	7/12/2018	07:39		7/12/2018	08:08
0712F005.D	ZZZZZZ	ZZZZZZ	7/12/2018	08:18		7/12/2018	08:47
0712F006.D	ZZZZZZ	ZZZZZZ	7/12/2018	08:57		7/12/2018	09:26
0712F007.D	PDI-SG-S078	K1804985-009	7/12/2018	09:37		7/12/2018	10:06
0712F008.D	PDI-SG-S010	K1804985-005	7/12/2018	10:16		7/12/2018	10:45
0712F009.D	PDI-SG-S255	K1804985-006	7/12/2018	10:55		7/12/2018	11:24
0712F010.D	PDI-SG-S084	K1804985-003	7/12/2018	11:35		7/12/2018	12:04
0712F011.D	PDI-SG-S097	K1804985-007	7/12/2018	12:14		7/12/2018	12:43
0712F012.D	PDI-SG-S115	K1804985-008	7/12/2018	12:54		7/12/2018	13:23
0712F013.D	ZZZZZZ	ZZZZZZ	7/12/2018	13:33		7/12/2018	14:02
0712F014.D	ZZZZZZ	ZZZZZZ	7/12/2018	14:13		7/12/2018	14:42
0712F015.D	ZZZZZZ	ZZZZZZ	7/12/2018	14:52		7/12/2018	15:21
0712F016.D	ZZZZZZ	ZZZZZZ	7/12/2018	15:32		7/12/2018	16:01
0712F017.D	ZZZZZZ	ZZZZZZ	7/12/2018	16:12		7/12/2018	16:41
0712F018.D	ZZZZZZ	ZZZZZZ	7/12/2018	16:51		7/12/2018	17:20
0712F019.D	ZZZZZZ	ZZZZZZ	7/12/2018	17:31		7/12/2018	18:00

Results flagged with an asterisk (\*) indicate the holding time was exceeded for the analysis

## QA/QC Results

**Client:** AECOM  
**Project:** Portland Harbor Pre-Remedial Design Investigation/60566335  
**Sample Matrix:** Sediment

**Service Request:** K1804985  
**Date Extracted:** 07/03/2018

**Extraction Prep Log**  
**Polynuclear Aromatic Hydrocarbons**

**Extraction Method:** EPA 3541  
**Analysis Method:** 8270D SIM

**Extraction Lot:** KWG1803256  
**Level:** Low

Sample Name	Lab Code	Date Collected	Date Received	Sample Amount	Final Volume	% Solids	Note
PDI-SG-S204	K1804985-001	05/03/18	05/25/18	40.268g	1mL	70.3	
PDI-SG-S147	K1804985-002	05/04/18	05/25/18	40.342g	1mL	54.8	
PDI-SG-S084DL	K1804985-003	05/08/18	05/25/18	40.336g	1mL	82.3	
PDI-SG-S084	K1804985-003	05/08/18	05/25/18	40.336g	1mL	82.3	
PDI-SG-S090	K1804985-004	05/09/18	05/25/18	40.141g	1mL	75.2	
PDI-SG-S010DL	K1804985-005	05/09/18	05/25/18	40.113g	1mL	75.9	
PDI-SG-S010	K1804985-005	05/09/18	05/25/18	40.113g	1mL	75.9	
PDI-SG-S255	K1804985-006	05/11/18	05/25/18	40.393g	1mL	62.4	
PDI-SG-S255DL	K1804985-006	05/11/18	05/25/18	40.393g	1mL	62.4	
PDI-SG-S097	K1804985-007	05/13/18	05/25/18	40.037g	1mL	62.5	
PDI-SG-S097DL	K1804985-007	05/13/18	05/25/18	40.037g	1mL	62.5	
PDI-SG-S115DL	K1804985-008	05/12/18	05/25/18	40.425g	1mL	74	
PDI-SG-S115	K1804985-008	05/12/18	05/25/18	40.425g	1mL	74	
PDI-SG-S078	K1804985-009	05/12/18	05/25/18	40.393g	1mL	55.1	
PDI-SG-S135	K1804985-010	05/14/18	05/25/18	40.282g	1mL	70.5	
PDI-SG-S157	K1804985-011	05/14/18	05/25/18	40.484g	1mL	60.6	
Method Blank	KWG1803256-4	NA	NA	40.484g	1mL	NA	
PDI-SG-S204MS	KWG1803256-1	05/03/18	05/25/18	40.098g	1mL	70.3	
PDI-SG-S204DMS	KWG1803256-2	05/03/18	05/25/18	40.221g	1mL	70.3	
Lab Control Sample	KWG1803256-3	NA	NA	20.000g	1mL	NA	

Results flagged with an asterisk (\*) indicate the holding time was exceeded for the analysis