



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY
REGION 5 CHICAGO REGIONAL LABORATORY
536 SOUTH CLARK STREET
CHICAGO, ILLINOIS 60605

Date: 8/21/2014
Subject: Review of Region 5 Data for SE Chicago Petcoke
To: Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago, IL 60605
From: Troy Stroock, Analyst
US EPA Region 5 Chicago Regional Laboratory

The data transmitted under this cover memo successfully passed CRL's data review procedures as documented in the current Quality Management Plan and applicable Standard Operating Procedures. In accordance with EPA's *Guidance on Environmental Data Verification and Data Validation* (Document EPA QA/G-8), CRL verified and validated the data but does not perform data quality assessment based on project plans.

This report was reviewed and the information provided herein accurately represents the analysis performed.

X _____

Please contact the analyst with any technical report issues, Amanda Wroble at (312)-353-0375 for sample project concerns, and Sylvia Griffin at (312)-353-9073 with data transmittal questions. Thank you.

Attached are Results for: SE Chicago Petcoke

/ /

Data Management Coordinator and Date Transmitted

Analyses included in this report:

SVOA PAHs in solids by press. fluid extr. (SIM)



Environmental Protection Agency Region 5
Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605
Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

Reported:
Aug-21-14 14:53

ANALYSIS CASE NARRATIVE

Analyst Phone number: 312.353.8362

General Information

2 filter samples were received by the Chicago Regional Laboratory (CRL) on 7/1/2014, and two new filters were purchased by the CRL on 7/8/2014, all for extraction and analysis by Standard Operating Procedure (SOP) MS026 rev. 8 (based on SW-846 Method 8270D). These samples were analyzed for polycyclic aromatic hydrocarbons (PAHs), including the target analytes listed in Table A.3.7 in the appendix of the SOP, along with perylene, an additional PAH of interest to the data user. All holding times were met for sample preparation and analysis.

Pen and ink change MS026 R8.0-PI02 applies to this data set, which addresses analysis and reporting of PAHs in the filter material. That pen and ink change states that PAHs will be reported in units of $\mu\text{g}/\text{kg}$, but the client has since requested results to be reported in the same units (ng/wipe) as used for wipe samples for this project. Clean matrix QC samples (method blank, RL verification, and blank spike QC samples) were prepared with Ottawa sand and hydromatrix, as described in SOP MS026 Rev 8 for solids, not with clean wipes as described in Pen and ink change MS026 R8.0-PI01 for preparation and analysis of dust wipe samples.

The filter samples received for work order 1407001 were cut approximately in half with metal shears before sample preparation in order to reserve the other portion for possible split analysis off-site. For sample 1407001-15, the filter dimensions tested were approximately 3.75"x6", and the mass of the filamentous material pulled off the metal backing that was tested was measured at 1.415 g. For sample 1407001-16, the filter dimensions tested were approximately 4.75"x 4.5", and the mass of the filamentous material pulled off the metal backing that was tested was measured at 1.49 g.

Filter sample 1407004-01 and associated matrix spike QC samples were cut from a single filter purchased new from a store, as was filter sample 1407004-02 and associated matrix spike QC samples. Subsamples were cut in dimensions of approximately 3.5" x 7", and masses of the filamentous material pulled from the metal backing for testing ranged from 0.825 g to 1.119 g.

Extracts were filtered through 0.45 μm particle filters twice during concentration of extracts to remove particulates observed in the extracts after extraction or observed to form during nitrogen blow-down of extracts.

The mass spectrometer was operated in selected ion monitoring (SIM) mode, which is more sensitive than full scan. Mass spectra acquired in SIM mode are limited to a couple of ions per analyte, resulting in poorer confirmation of peak identification by mass spectrum in unknown samples, as mass spectra acquired in SIM cannot be searched against a mass spectral library.



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Sample Analysis and Results

Many of the target PAHs were measured above the RL in the filter samples in work order 1407001, and a few PAHs were also observed in the filter samples in work order 1407004. Perylene was qualified as potentially high bias (“K”) in sample ID 1407001-16 due to poor peak shape and a qualifier to quantification ion ratio that was much higher than in the calibration standard. Full scan analysis of the extract showed that a large hydrocarbon peak elutes at the expected perylene retention time, so this target analyte measurement is suspected to be high bias or false positive.

The data reported herein meet the Data Quality Objectives referenced in the SE Chicago Petcoke QAPP dated 12/16/2013 and addendum named SE Chicago Petcoke QAPP HVAC filter update 063014.docx, with the exception that PAH concentrations were reported for sample ID 1407001-15 from a 10-fold dilution, resulting in reporting of some results as non-detects at 10-fold higher reporting limits (RLs) than listed in the QAPP.

Quality Controls

Initial Calibration:

Most of the target analytes were contained in the primary and second source standards used for calibration of the instrument and initial calibration verification, respectively. However, no certified second source standard was analyzed or evaluated for 1-methylnaphthalene, benzo(e)pyrene, or perylene. Response factors for these analytes were similar to structural isomers that met criteria for initial calibration verification.

Continuing Calibration Verification (CCV):

One of the CCVs bracketing analysis of the method blank, RL verification, and blank spike QC samples associated with the extraction batch slightly exceeded the % difference criteria of $\leq \pm 25\%$ relative to the expected concentration for pyrene. The concentration or RL of this target analyte in these QC samples is qualified as estimated (“J”) for measurements above and below the RL, respectively. One or both CCVs bracketing analysis of the 10-fold dilution of sample ID 1407001-15 exceeded the % difference criteria of $\leq \pm 25\%$ relative to the expected concentration for nitrobenzene-d5, 2-methylnaphthalene, 1-methylnaphthalene, phenanthrene, and fluorene. Of these target analytes, only phenanthrene was measured in this sample above the RL. The concentration of phenanthrene is qualified as estimated in this field sample, and the RLs for the other non-detected target analytes that did not meet the CCV criteria are qualified as estimated (“J”).

Internal Standards:

Internal standard responses in all reported field samples and associated batch QC samples were within SOP criteria of 50-200% of the associated peak responses in the continuing calibration verification (CCV) injections. However,



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responses of some internal standards in one of the two CCV injections bracketing analysis of all of the field samples and matrix spike QC samples reported for this work order except for the 10-fold dilution of sample 1407001-15 were slightly low relative to their responses in the midpoint initial calibration standard. Internal standard responses in one of the CCVs bracketing reported analytical runs of sample ID 1407004-01, 1407004-02, and all matrix spike QC samples ranged between 37.2 and 50.4% of the response in the midpoint initial calibration standard, and internal standard responses in one of the CCVs bracketing the reported analytical run of the 1 mL extract volume of sample ID 1407001-16 ranged between 47.8 and 72.3% of the midpoint initial calibration standard. One of the two CCVs bracketing analysis of these samples met the internal standard performance criteria, and the CCVs with low internal standard responses met the % difference criteria for all target analytes and surrogates, so recalibration of the instrument was deemed unnecessary. No field sample data is qualified as a result of low response observed in 1 of the 2 CCVs bracketing the samples.

Surrogates Spike Recovery:

The non-ionizable surrogates listed in section 8.2 of the SOP (nitrobenzene-d5, 2-fluorobiphenyl, terphenyl-d14) were used for evaluation of successful preparation and analysis of each field sample and QC sample listed in this report. Nitrobenzene-d5 was not recovered in filter sample 1407001-15, and nitrobenzene-d5 recovery was low in one of the matrix spike QC samples (in B407046-MS2, nitrobenzene recovery was 47.5%,; lower control limit is 50%), but the other two surrogates were recovered acceptably in both of these samples, so no data is qualified due to surrogate recovery.

Target Compounds Spike Recovery:

Recovery of 1-methylnaphthalene is evaluated against control limits provided in the SOP for 2-methylnaphthalene, and recovery of perylene and benzo(e)pyrene is evaluated against control limits for benzo(a)pyrene. Once sufficient data are acquired, historical spike recovery statistics for these analytes will be used to generate more appropriate recovery criteria. An RL study has been completed for benzo(e)pyrene and 1-methylnaphthalene and is summarized in the SOP in Attachment 3 table A.3.7. A similar RL study has not been completed for perylene, so the concentration of this analyte is qualified as estimated ("J") in all field samples until it is included in the SOP.

One of the blank spike (LCS) QC samples, B407047-BSD1, was lost during pressurized fluid extraction due to an equipment malfunction. The other LCS QC sample, B407047-BS1, was analyzed successfully, so no further actions were taken due to loss of one of the LCS QC samples in the batch.

4 replicate matrix spike QC samples were prepared and analyzed for each of field samples 1407004-01 and -02 and are reported as B407047-MS1 through -MS8. In spite of background observed in the unspiked samples for a few target analytes, all target analytes were recovered acceptably in all matrix spikes except for phenanthrene in B407047-MS3 (63.4% recovery; lower control limit of 68.1%), fluorene in B407047-MS4 (45.4% recovery; lower control limit 57.1%), and a recovery of a number of target analytes was low in B407047-MS8. Part of the extract of



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B407047-MS8 was lost during the second particle filtration (approximately 2 mL of 10 mL), resulting in slightly low bias recovery for most of the target analytes. The target analytes in the other 3 replicate matrix spike QC samples were recovered acceptably, so no data is qualified as a result of matrix spike recovery.

RL Verification:

For this analysis, RL verification QC samples were prepared at two (nominal) spike levels: 20 ng/wipe (B407047-MRL1) and 100 ng/wipe (B407047-MRL2), and the same acceptance criteria are used to evaluate these QC samples as used for the blank spikes. All target analytes were recovered at their respective RLs within the acceptance criteria. Results are only reported for the individual RL verification QC sample corresponding to that analyte's RL.

All other quality controls not mentioned here meet the criteria in the SOP.

Signature _____, Date _____



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ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
3M Filtrete Allergy Defense Filter	1407004-01	Wipe	Jul-09-14 08:15	Jul-09-14 08:15
True Blue BASIC protection MERV 7 Filter	1407004-02	Wipe	Jul-09-14 08:15	Jul-09-14 08:15



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Semivolatiles by GC/MS, EPA 8270D (modified) US EPA Region 5 Chicago Regional Laboratory

3M Filtrete Allergy Defense Filter (1407004-01) Wipe **Sampled: Jul-09-14 08:15** **Received: Jul-09-14 08:15**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Naphthalene	U			100	ng/Wipe	1	B407047	Jul-10-14	Jul-24-14
2-Methylnaphthalene	U			100	"	"	"	"	"
1-Methylnaphthalene	U			100	"	"	"	"	"
Acenaphthylene	U			20.0	"	"	"	"	"
Acenaphthene	169			100	"	"	"	"	"
Fluorene	270			20.0	"	"	"	"	"
Phenanthrene	U			100	"	"	"	"	"
Anthracene	U			20.0	"	"	"	"	"
Fluoranthene	U			100	"	"	"	"	"
Pyrene	U			100	"	"	"	"	"
Benzo (a) anthracene	U			20.0	"	"	"	"	"
Chrysene	U			20.0	"	"	"	"	"
Benzo(b)fluoranthene	U			20.0	"	"	"	"	"
Benzo(k)fluoranthene	U			20.0	"	"	"	"	"
Benzo (e) pyrene	U			20.0	"	"	"	"	"
Benzo(a)pyrene	U			20.0	"	"	"	"	"
Perylene	U	J		20.0	"	"	"	"	"
Indeno(1,2,3-cd)pyrene	U			20.0	"	"	"	"	"
Dibenz(a,h)anthracene	U			20.0	"	"	"	"	"
Benzo(g,h,i)perylene	U			20.0	"	"	"	"	"

Surogate	Result	%REC Limits	Batch	Prepared	Analyzed
<i>Surogate: Nitrobenzene-d5</i>	180	72.0%	50-94.9	"	"
<i>Surogate: 2-Fluorobiphenyl</i>	194	77.7%	48.1-108	"	"
<i>Surogate: Terphenyl-d14</i>	217	86.7%	59.4-127	"	"

True Blue BASIC protection MERV 7 Filter (1407004-02) Wipe **Sampled: Jul-09-14 08:15** **Received: Jul-09-14 08:15**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Naphthalene	210			100	ng/Wipe	1	B407047	Jul-10-14	Jul-25-14
2-Methylnaphthalene	126			100	"	"	"	"	"
1-Methylnaphthalene	U			100	"	"	"	"	"
Acenaphthylene	U			20.0	"	"	"	"	"
Acenaphthene	U			100	"	"	"	"	"
Fluorene	73.6			20.0	"	"	"	"	"



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Semivolatiles by GC/MS, EPA 8270D (modified) US EPA Region 5 Chicago Regional Laboratory

True Blue BASIC protection MERV 7 Filter (1407004-02) Wipe Sampled: Jul-09-14 08:15 Received: Jul-09-14 08:15

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Phenanthrene	U			100	ng/Wipe	1	B407047	Jul-10-14	Jul-25-14
Anthracene	U			20.0	"	"	"	"	"
Fluoranthene	U			100	"	"	"	"	"
Pyrene	U			100	"	"	"	"	"
Benzo (a) anthracene	U			20.0	"	"	"	"	"
Chrysene	U			20.0	"	"	"	"	"
Benzo(b)fluoranthene	U			20.0	"	"	"	"	"
Benzo(k)fluoranthene	U			20.0	"	"	"	"	"
Benzo (e) pyrene	U			20.0	"	"	"	"	"
Benzo(a)pyrene	U			20.0	"	"	"	"	"
Perylene	U	J		20.0	"	"	"	"	"
Indeno(1,2,3-cd)pyrene	U			20.0	"	"	"	"	"
Dibenz(a,h)anthracene	U			20.0	"	"	"	"	"
Benzo(g,h,i)perylene	U			20.0	"	"	"	"	"

Surogate	Result	%REC	%REC Limits	Batch	Prepared	Analyzed
<i>Surogate: Nitrobenzene-d5</i>	193	77.0%	50-94.9	"	"	"
<i>Surogate: 2-Fluorobiphenyl</i>	222	88.7%	48.1-108	"	"	"
<i>Surogate: Terphenyl-d14</i>	219	87.7%	59.4-127	"	"	"



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Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

Blank (B407047-BLK1)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	U			100	ng/Wipe						
2-Methylnaphthalene	U			100	"						
1-Methylnaphthalene	U			100	"						
Acenaphthylene	U			20.0	"						
Acenaphthene	U			100	"						
Fluorene	U			20.0	"						
Phenanthrene	U			100	"						
Anthracene	U			20.0	"						
Fluoranthene	U			100	"						
Pyrene	U	J		100	"						
Benzo (a) anthracene	U			20.0	"						
Chrysene	U			20.0	"						
Benzo(b)fluoranthene	U			20.0	"						
Benzo(k)fluoranthene	U			20.0	"						
Benzo (e) pyrene	U			20.0	"						
Benzo(a)pyrene	U			20.0	"						
Perylene	U			20.0	"						
Indeno(1,2,3-cd)pyrene	U			20.0	"						
Dibenz(a,h)anthracene	U			20.0	"						
Benzo(g,h,i)perylene	U			20.0	"						
Surrogate: Nitrobenzene-d5	170				"	250.0		68.1%	50-94.9		
Surrogate: 2-Fluorobiphenyl	167				"	250.0		66.6%	48.1-108		
Surrogate: Terphenyl-d14	214				"	250.0		85.4%	59.4-127		

Blank (B407047-BLK2)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	U			100	ng/Wipe						
2-Methylnaphthalene	U			100	"						
1-Methylnaphthalene	U			100	"						
Acenaphthylene	U			20.0	"						
Acenaphthene	U			100	"						



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Batch B407047 - Solvent Extraction

Blank (B407047-BLK2)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Fluorene	U			20.0	ng/Wipe					
Phenanthrene	U			100	"					
Anthracene	U			20.0	"					
Fluoranthene	U			100	"					
Pyrene	U	J		100	"					
Benzo (a) anthracene	U			20.0	"					
Chrysene	U			20.0	"					
Benzo(b)fluoranthene	U			20.0	"					
Benzo(k)fluoranthene	U			20.0	"					
Benzo (e) pyrene	U			20.0	"					
Benzo(a)pyrene	U			20.0	"					
Perylene	U			20.0	"					
Indeno(1,2,3-cd)pyrene	U			20.0	"					
Dibenz(a,h)anthracene	U			20.0	"					
Benzo(g,h,i)perylene	U			20.0	"					
Surrogate: Nitrobenzene-d5	178				"	250.0		71.2% 50-94.9		
Surrogate: 2-Fluorobiphenyl	171				"	250.0		68.4% 48.1-108		
Surrogate: Terphenyl-d14	215				"	250.0		86.0% 59.4-127		

LCS (B407047-BS1)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Naphthalene	123			100	ng/Wipe	250.0		49.3% 43.1-108		
2-Methylnaphthalene	139			100	"	250.0		55.6% 45.5-117		
1-Methylnaphthalene	135			100	"	250.0		53.9% 45.5-117		
Acenaphthylene	149			20.0	"	250.0		59.7% 52.7-117		
Acenaphthene	152			100	"	250.0		60.8% 51.7-116		
Fluorene	176			20.0	"	250.0		70.5% 57.9-120		
Phenanthrene	200			100	"	250.0		80.2% 68.1-114		
Anthracene	175			20.0	"	250.0		70.0% 69.7-116		
Fluoranthene	212			100	"	250.0		84.8% 70.8-122		
Pyrene	213	J		100	"	250.0		85.3% 71.8-117		
Benzo (a) anthracene	210			20.0	"	250.0		84.1% 67.6-115		



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Batch B407047 - Solvent Extraction

LCS (B407047-BS1)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Chrysene	213			20.0	ng/Wipe	250.0		85.1%	68.5-117	
Benzo(b)fluoranthene	220			20.0	"	250.0		88.0%	68.9-128	
Benzo(k)fluoranthene	217			20.0	"	250.0		86.9%	62.8-135	
Benzo (e) pyrene	212			20.0	"	250.0		85.0%	68.9-133	
Benzo(a)pyrene	179			20.0	"	250.0		71.8%	68.9-133	
Perylene	190			20.0	"	250.0		76.2%	68.9-133	
Indeno(1,2,3-cd)pyrene	208			20.0	"	250.0		83.3%	70-129	
Dibenz(a,h)anthracene	211			20.0	"	250.0		84.5%	69.1-131	
Benzo(g,h,i)perylene	192			20.0	"	250.0		77.0%	53.9-139	
<i>Surrogate: Nitrobenzene-d5</i>	<i>129</i>				"	<i>250.0</i>		<i>51.7%</i>	<i>50-94.9</i>	
<i>Surrogate: 2-Fluorobiphenyl</i>	<i>144</i>				"	<i>250.0</i>		<i>57.6%</i>	<i>48.1-108</i>	
<i>Surrogate: Terphenyl-d14</i>	<i>210</i>				"	<i>250.0</i>		<i>84.2%</i>	<i>59.4-127</i>	

MRL Check (B407047-MRL1)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Acenaphthylene	15.0			20.0	ng/Wipe	20.00		75.1%	52.7-117	
Fluorene	18.4			20.0	"	20.00		91.9%	57.9-120	
Anthracene	17.2			20.0	"	20.00		86.0%	69.7-116	
Benzo (a) anthracene	18.9			20.0	"	20.00		94.6%	67.6-115	
Chrysene	17.7			20.0	"	20.00		88.5%	68.5-117	
Benzo(b)fluoranthene	18.8			20.0	"	20.00		94.2%	68.9-128	
Benzo(k)fluoranthene	18.3			20.0	"	20.00		91.4%	62.8-135	
Benzo (e) pyrene	18.2			20.0	"	20.00		91.0%	68.9-133	
Benzo(a)pyrene	18.3			20.0	"	20.00		91.4%	68.9-133	
Perylene	16.6			20.0	"	20.00		82.9%	68.9-133	
Indeno(1,2,3-cd)pyrene	16.5			20.0	"	20.00		82.4%	70-129	
Dibenz(a,h)anthracene	16.9			20.0	"	20.00		84.6%	69.1-131	
Benzo(g,h,i)perylene	16.4			20.0	"	20.00		82.0%	53.9-139	
<i>Surrogate: Nitrobenzene-d5</i>	<i>180</i>				"	<i>250.0</i>		<i>72.1%</i>	<i>50-94.9</i>	
<i>Surrogate: 2-Fluorobiphenyl</i>	<i>176</i>				"	<i>250.0</i>		<i>70.2%</i>	<i>48.1-108</i>	
<i>Surrogate: Terphenyl-d14</i>	<i>214</i>				"	<i>250.0</i>		<i>85.6%</i>	<i>59.4-127</i>	



Environmental Protection Agency Region 5 Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605
Phone: (312)353-8370 Fax: (312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

Reported:
Aug-21-14 14:53

Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

MRL Check (B407047-MRL2)

Prepared: Jul-10-14 Analyzed: Jul-23-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Naphthalene	68.7			100	ng/Wipe	100.0		68.7%	43.1-108	
2-Methylnaphthalene	73.7			100	"	100.0		73.7%	45.5-117	
1-Methylnaphthalene	67.5			100	"	100.0		67.5%	45.5-117	
Acenaphthene	71.1			100	"	100.0		71.1%	51.7-116	
Phenanthrene	87.7			100	"	100.0		87.7%	68.1-114	
Fluoranthene	88.6			100	"	100.0		88.6%	70.8-122	
Pyrene	88.4	J		100	"	100.0		88.4%	71.8-117	
Surrogate: Nitrobenzene-d5	171				"	250.0		68.5%	50-94.9	
Surrogate: 2-Fluorobiphenyl	168				"	250.0		67.3%	48.1-108	
Surrogate: Terphenyl-d14	213				"	250.0		85.2%	59.4-127	

Matrix Spike (B407047-MS1)

Source: 1407004-01

Prepared: Jul-10-14 Analyzed: Jul-24-14

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Naphthalene	188			100	ng/Wipe	250.0	19.9	67.2%	43.1-108	
2-Methylnaphthalene	216			100	"	250.0	23.4	77.1%	45.5-117	
1-Methylnaphthalene	201			100	"	250.0	16.0	74.1%	45.5-117	
Acenaphthylene	201			20.0	"	250.0	U	80.2%	52.7-117	
Acenaphthene	371			100	"	250.0	169	80.9%	51.7-116	
Fluorene	441			20.0	"	250.0	270	68.4%	57.9-120	
Phenanthrene	255			100	"	250.0	77.1	71.2%	68.1-114	
Anthracene	222			20.0	"	250.0	U	88.8%	69.7-116	
Fluoranthene	214			100	"	250.0	17.5	78.5%	70.8-122	
Pyrene	197			100	"	250.0	U	78.9%	71.8-117	
Benzo (a) anthracene	218			20.0	"	250.0	U	87.1%	67.6-115	
Chrysene	192			20.0	"	250.0	U	76.6%	68.5-117	
Benzo(b)fluoranthene	232			20.0	"	250.0	15.1	86.8%	68.9-128	
Benzo(k)fluoranthene	225			20.0	"	250.0	14.9	83.9%	62.8-135	
Benzo (e) pyrene	201			20.0	"	250.0	U	80.4%	68.9-133	
Benzo(a)pyrene	205			20.0	"	250.0	U	82.2%	68.9-133	
Perylene	179			20.0	"	250.0	U	71.6%	68.9-133	
Indeno(1,2,3-cd)pyrene	224			20.0	"	250.0	U	89.6%	70-129	
Dibenz(a,h)anthracene	229			20.0	"	250.0	U	91.7%	69.1-131	



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536 South Clark Street, Chicago, IL 60605
Phone: (312)353-8370 Fax: (312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

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Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

Matrix Spike (B407047-MS1)		Source: 1407004-01			Prepared: Jul-10-14 Analyzed: Jul-24-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	219			20.0	ng/Wipe	250.0	U	87.5%	53.9-139		
<i>Surrogate: Nitrobenzene-d5</i>	<i>183</i>				"	<i>250.0</i>		<i>73.1%</i>	<i>50-94.9</i>		
<i>Surrogate: 2-Fluorobiphenyl</i>	<i>201</i>				"	<i>250.0</i>		<i>80.5%</i>	<i>48.1-108</i>		
<i>Surrogate: Terphenyl-d14</i>	<i>189</i>				"	<i>250.0</i>		<i>75.5%</i>	<i>59.4-127</i>		

Matrix Spike (B407047-MS2)		Source: 1407004-01			Prepared: Jul-10-14 Analyzed: Jul-24-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	135			100	ng/Wipe	250.0	19.9	46.0%	43.1-108		
2-Methylnaphthalene	173			100	"	250.0	23.4	59.9%	45.5-117		
1-Methylnaphthalene	161			100	"	250.0	16.0	58.1%	45.5-117		
Acenaphthylene	183			20.0	"	250.0	U	73.2%	52.7-117		
Acenaphthene	384			100	"	250.0	169	86.0%	51.7-116		
Fluorene	475			20.0	"	250.0	270	81.7%	57.9-120		
Phenanthrene	252			100	"	250.0	77.1	70.1%	68.1-114		
Anthracene	225			20.0	"	250.0	U	89.8%	69.7-116		
Fluoranthene	216			100	"	250.0	17.5	79.2%	70.8-122		
Pyrene	232			100	"	250.0	U	92.9%	71.8-117		
Benzo (a) anthracene	230			20.0	"	250.0	U	92.0%	67.6-115		
Chrysene	212			20.0	"	250.0	U	85.0%	68.5-117		
Benzo(b)fluoranthene	250			20.0	"	250.0	15.1	93.9%	68.9-128		
Benzo(k)fluoranthene	218			20.0	"	250.0	14.9	81.3%	62.8-135		
Benzo (e) pyrene	191			20.0	"	250.0	U	76.3%	68.9-133		
Benzo(a)pyrene	205			20.0	"	250.0	U	82.0%	68.9-133		
Perylene	179			20.0	"	250.0	U	71.6%	68.9-133		
Indeno(1,2,3-cd)pyrene	219			20.0	"	250.0	U	87.6%	70-129		
Dibenz(a,h)anthracene	232			20.0	"	250.0	U	92.6%	69.1-131		
Benzo(g,h,i)perylene	225			20.0	"	250.0	U	89.9%	53.9-139		
<i>Surrogate: Nitrobenzene-d5</i>	<i>119</i>				"	<i>250.0</i>		<i>47.5%</i>	<i>50-94.9</i>		
<i>Surrogate: 2-Fluorobiphenyl</i>	<i>173</i>				"	<i>250.0</i>		<i>69.2%</i>	<i>48.1-108</i>		
<i>Surrogate: Terphenyl-d14</i>	<i>225</i>				"	<i>250.0</i>		<i>90.0%</i>	<i>59.4-127</i>		



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536 South Clark Street, Chicago, IL 60605
Phone: (312)353-8370 Fax: (312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

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Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

Matrix Spike (B407047-MS3)		Source: 1407004-01			Prepared: Jul-10-14 Analyzed: Jul-24-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	177			100	ng/Wipe	250.0	19.9	63.0%	43.1-108		
2-Methylnaphthalene	208			100	"	250.0	23.4	73.7%	45.5-117		
1-Methylnaphthalene	197			100	"	250.0	16.0	72.3%	45.5-117		
Acenaphthylene	194			20.0	"	250.0	U	77.7%	52.7-117		
Acenaphthene	401			100	"	250.0	169	92.7%	51.7-116		
Fluorene	474			20.0	"	250.0	270	81.5%	57.9-120		
Phenanthrene	236			100	"	250.0	77.1	63.4%	68.1-114		
Anthracene	229			20.0	"	250.0	U	91.6%	69.7-116		
Fluoranthene	210			100	"	250.0	17.5	77.0%	70.8-122		
Pyrene	193			100	"	250.0	U	77.3%	71.8-117		
Benzo (a) anthracene	213			20.0	"	250.0	U	85.1%	67.6-115		
Chrysene	232			20.0	"	250.0	U	92.6%	68.5-117		
Benzo(b)fluoranthene	225			20.0	"	250.0	15.1	84.0%	68.9-128		
Benzo(k)fluoranthene	208			20.0	"	250.0	14.9	77.1%	62.8-135		
Benzo (e) pyrene	215			20.0	"	250.0	U	86.2%	68.9-133		
Benzo(a)pyrene	212			20.0	"	250.0	U	85.0%	68.9-133		
Perylene	190			20.0	"	250.0	U	75.9%	68.9-133		
Indeno(1,2,3-cd)pyrene	222			20.0	"	250.0	U	88.6%	70-129		
Dibenz(a,h)anthracene	224			20.0	"	250.0	U	89.8%	69.1-131		
Benzo(g,h,i)perylene	221			20.0	"	250.0	U	88.5%	53.9-139		
Surrogate: Nitrobenzene-d5	159				"	250.0		63.8%	50-94.9		
Surrogate: 2-Fluorobiphenyl	199				"	250.0		79.6%	48.1-108		
Surrogate: Terphenyl-d14	199				"	250.0		79.6%	59.4-127		

Matrix Spike (B407047-MS4)		Source: 1407004-01			Prepared: Jul-10-14 Analyzed: Jul-25-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	188			100	ng/Wipe	250.0	19.9	67.2%	43.1-108		
2-Methylnaphthalene	217			100	"	250.0	23.4	77.5%	45.5-117		
1-Methylnaphthalene	201			100	"	250.0	16.0	74.1%	45.5-117		
Acenaphthylene	211			20.0	"	250.0	U	84.3%	52.7-117		
Acenaphthene	322			100	"	250.0	169	61.0%	51.7-116		
Fluorene	384			20.0	"	250.0	270	45.4%	57.9-120		



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536 South Clark Street, Chicago, IL 60605
Phone: (312) 353-8370 Fax: (312) 886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

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Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

Matrix Spike (B407047-MS4)		Source: 1407004-01			Prepared: Jul-10-14 Analyzed: Jul-25-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Phenanthrene	254			100	ng/Wipe	250.0	77.1	70.7%	68.1-114		
Anthracene	226			20.0	"	250.0	U	90.4%	69.7-116		
Fluoranthene	222			100	"	250.0	17.5	81.9%	70.8-122		
Pyrene	217			100	"	250.0	U	86.7%	71.8-117		
Benzo (a) anthracene	234			20.0	"	250.0	U	93.7%	67.6-115		
Chrysene	203			20.0	"	250.0	U	81.1%	68.5-117		
Benzo(b)fluoranthene	236			20.0	"	250.0	15.1	88.4%	68.9-128		
Benzo(k)fluoranthene	221			20.0	"	250.0	14.9	82.4%	62.8-135		
Benzo (e) pyrene	227			20.0	"	250.0	U	90.8%	68.9-133		
Benzo(a)pyrene	210			20.0	"	250.0	U	84.1%	68.9-133		
Perylene	185			20.0	"	250.0	U	74.1%	68.9-133		
Indeno(1,2,3-cd)pyrene	222			20.0	"	250.0	U	88.6%	70-129		
Dibenz(a,h)anthracene	227			20.0	"	250.0	U	90.8%	69.1-131		
Benzo(g,h,i)perylene	220			20.0	"	250.0	U	88.1%	53.9-139		
Surrogate: Nitrobenzene-d5	180				"	250.0		71.9%	50-94.9		
Surrogate: 2-Fluorobiphenyl	206				"	250.0		82.5%	48.1-108		
Surrogate: Terphenyl-d14	208				"	250.0		83.2%	59.4-127		

Matrix Spike (B407047-MS5)		Source: 1407004-02			Prepared: Jul-10-14 Analyzed: Jul-25-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	378			100	ng/Wipe	250.0	210	67.2%	43.1-108		
2-Methylnaphthalene	318			100	"	250.0	126	76.8%	45.5-117		
1-Methylnaphthalene	242			100	"	250.0	42.9	79.8%	45.5-117		
Acenaphthylene	206			20.0	"	250.0	U	82.5%	52.7-117		
Acenaphthene	296			100	"	250.0	98.8	78.9%	51.7-116		
Fluorene	284			20.0	"	250.0	73.6	84.2%	57.9-120		
Phenanthrene	256			100	"	250.0	58.2	79.1%	68.1-114		
Anthracene	216			20.0	"	250.0	U	86.3%	69.7-116		
Fluoranthene	225			100	"	250.0	U	90.0%	70.8-122		
Pyrene	232			100	"	250.0	25.6	82.7%	71.8-117		
Benzo (a) anthracene	206			20.0	"	250.0	10.4	78.2%	67.6-115		
Chrysene	242			20.0	"	250.0	10.6	92.6%	68.5-117		



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536 South Clark Street, Chicago, IL 60605
Phone: (312)353-8370 Fax: (312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

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Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

Matrix Spike (B407047-MS5)		Source: 1407004-02		Prepared: Jul-10-14 Analyzed: Jul-25-14							
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Benzo(b)fluoranthene	226			20.0	ng/Wipe	250.0	U	90.2%	68.9-128		
Benzo(k)fluoranthene	226			20.0	"	250.0	U	90.2%	62.8-135		
Benzo (e) pyrene	219			20.0	"	250.0	U	87.8%	68.9-133		
Benzo(a)pyrene	243			20.0	"	250.0	14.2	91.5%	68.9-133		
Perylene	202			20.0	"	250.0	U	80.9%	68.9-133		
Indeno(1,2,3-cd)pyrene	237			20.0	"	250.0	U	94.8%	70-129		
Dibenz(a,h)anthracene	244			20.0	"	250.0	U	97.8%	69.1-131		
Benzo(g,h,i)perylene	228			20.0	"	250.0	U	91.3%	53.9-139		
Surrogate: Nitrobenzene-d5	189				"	250.0		75.7%	50-94.9		
Surrogate: 2-Fluorobiphenyl	212				"	250.0		84.7%	48.1-108		
Surrogate: Terphenyl-d14	215				"	250.0		85.9%	59.4-127		

Matrix Spike (B407047-MS6)		Source: 1407004-02		Prepared: Jul-10-14 Analyzed: Jul-25-14							
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	377			100	ng/Wipe	250.0	210	66.7%	43.1-108		
2-Methylnaphthalene	318			100	"	250.0	126	76.9%	45.5-117		
1-Methylnaphthalene	235			100	"	250.0	42.9	76.8%	45.5-117		
Acenaphthylene	203			20.0	"	250.0	U	81.1%	52.7-117		
Acenaphthene	321			100	"	250.0	98.8	88.8%	51.7-116		
Fluorene	294			20.0	"	250.0	73.6	88.0%	57.9-120		
Phenanthrene	251			100	"	250.0	58.2	77.0%	68.1-114		
Anthracene	218			20.0	"	250.0	U	87.3%	69.7-116		
Fluoranthene	212			100	"	250.0	U	84.9%	70.8-122		
Pyrene	239			100	"	250.0	25.6	85.5%	71.8-117		
Benzo (a) anthracene	214			20.0	"	250.0	10.4	81.3%	67.6-115		
Chrysene	244			20.0	"	250.0	10.6	93.5%	68.5-117		
Benzo(b)fluoranthene	216			20.0	"	250.0	U	86.6%	68.9-128		
Benzo(k)fluoranthene	216			20.0	"	250.0	U	86.4%	62.8-135		
Benzo (e) pyrene	208			20.0	"	250.0	U	83.1%	68.9-133		
Benzo(a)pyrene	234			20.0	"	250.0	14.2	88.0%	68.9-133		
Perylene	196			20.0	"	250.0	U	78.6%	68.9-133		
Indeno(1,2,3-cd)pyrene	216			20.0	"	250.0	U	86.4%	70-129		



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536 South Clark Street, Chicago, IL 60605
Phone: (312)353-8370 Fax: (312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
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Project Manager: Motria Caudill

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Batch B407047 - Solvent Extraction

Matrix Spike (B407047-MS6)		Source: 1407004-02			Prepared: Jul-10-14 Analyzed: Jul-25-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Dibenz(a,h)anthracene	219			20.0	ng/Wipe	250.0	U	87.8%	69.1-131		
Benzo(g,h,i)perylene	202			20.0	"	250.0	U	80.8%	53.9-139		
<i>Surrogate: Nitrobenzene-d5</i>	<i>192</i>				"	<i>250.0</i>		<i>76.7%</i>	<i>50-94.9</i>		
<i>Surrogate: 2-Fluorobiphenyl</i>	<i>214</i>				"	<i>250.0</i>		<i>85.5%</i>	<i>48.1-108</i>		
<i>Surrogate: Terphenyl-d14</i>	<i>213</i>				"	<i>250.0</i>		<i>85.3%</i>	<i>59.4-127</i>		

Matrix Spike (B407047-MS7)		Source: 1407004-02			Prepared: Jul-10-14 Analyzed: Jul-25-14						
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	420			100	ng/Wipe	250.0	210	84.0%	43.1-108		
2-Methylnaphthalene	359			100	"	250.0	126	93.2%	45.5-117		
1-Methylnaphthalene	260			100	"	250.0	42.9	86.8%	45.5-117		
Acenaphthylene	212			20.0	"	250.0	U	84.9%	52.7-117		
Acenaphthene	340			100	"	250.0	98.8	96.5%	51.7-116		
Fluorene	317			20.0	"	250.0	73.6	97.3%	57.9-120		
Phenanthrene	279			100	"	250.0	58.2	88.2%	68.1-114		
Anthracene	220			20.0	"	250.0	U	87.9%	69.7-116		
Fluoranthene	222			100	"	250.0	U	88.6%	70.8-122		
Pyrene	239			100	"	250.0	25.6	85.4%	71.8-117		
Benzo (a) anthracene	213			20.0	"	250.0	10.4	81.0%	67.6-115		
Chrysene	250			20.0	"	250.0	10.6	95.6%	68.5-117		
Benzo(b)fluoranthene	222			20.0	"	250.0	U	88.9%	68.9-128		
Benzo(k)fluoranthene	210			20.0	"	250.0	U	83.9%	62.8-135		
Benzo (e) pyrene	212			20.0	"	250.0	U	84.8%	68.9-133		
Benzo(a)pyrene	250			20.0	"	250.0	14.2	94.4%	68.9-133		
Perylene	198			20.0	"	250.0	U	79.2%	68.9-133		
Indeno(1,2,3-cd)pyrene	228			20.0	"	250.0	U	91.3%	70-129		
Dibenz(a,h)anthracene	234			20.0	"	250.0	U	93.4%	69.1-131		
Benzo(g,h,i)perylene	218			20.0	"	250.0	U	87.0%	53.9-139		
<i>Surrogate: Nitrobenzene-d5</i>	<i>186</i>				"	<i>250.0</i>		<i>74.5%</i>	<i>50-94.9</i>		
<i>Surrogate: 2-Fluorobiphenyl</i>	<i>232</i>				"	<i>250.0</i>		<i>92.8%</i>	<i>48.1-108</i>		
<i>Surrogate: Terphenyl-d14</i>	<i>215</i>				"	<i>250.0</i>		<i>86.0%</i>	<i>59.4-127</i>		



Environmental Protection Agency Region 5 Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605
Phone: (312)353-8370 Fax: (312)886-2591

Air Division, US EPA Region 5
77 West Jackson Boulevard
Chicago IL, 60605

Project: SE Chicago Petcoke
Project Number: [none]
Project Manager: Motria Caudill

Reported:
Aug-21-14 14:53

Semivolatiles by GC/MS, EPA 8270D (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B407047 - Solvent Extraction

Matrix Spike (B407047-MS8)		Source: 1407004-02		Prepared: Jul-10-14 Analyzed: Jul-25-14							
Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	301			100	ng/Wipe	250.0	210	36.6%	43.1-108		
2-Methylnaphthalene	228			100	"	250.0	126	40.7%	45.5-117		
1-Methylnaphthalene	172			100	"	250.0	42.9	51.8%	45.5-117		
Acenaphthylene	157			20.0	"	250.0	U	62.7%	52.7-117		
Acenaphthene	238			100	"	250.0	98.8	55.6%	51.7-116		
Fluorene	222			20.0	"	250.0	73.6	59.3%	57.9-120		
Phenanthrene	183			100	"	250.0	58.2	49.8%	68.1-114		
Anthracene	154			20.0	"	250.0	U	61.7%	69.7-116		
Fluoranthene	158			100	"	250.0	U	63.1%	70.8-122		
Pyrene	166			100	"	250.0	25.6	56.1%	71.8-117		
Benzo (a) anthracene	145			20.0	"	250.0	10.4	54.0%	67.6-115		
Chrysene	168			20.0	"	250.0	10.6	62.9%	68.5-117		
Benzo(b)fluoranthene	153			20.0	"	250.0	U	61.3%	68.9-128		
Benzo(k)fluoranthene	154			20.0	"	250.0	U	61.5%	62.8-135		
Benzo (e) pyrene	150			20.0	"	250.0	U	59.9%	68.9-133		
Benzo(a)pyrene	171			20.0	"	250.0	14.2	62.6%	68.9-133		
Perylene	142			20.0	"	250.0	U	56.8%	68.9-133		
Indeno(1,2,3-cd)pyrene	155			20.0	"	250.0	U	61.8%	70-129		
Dibenz(a,h)anthracene	158			20.0	"	250.0	U	63.4%	69.1-131		
Benzo(g,h,i)perylene	146			20.0	"	250.0	U	58.5%	53.9-139		
Surrogate: Nitrobenzene-d5	149				"	250.0		59.6%	50-94.9		
Surrogate: 2-Fluorobiphenyl	159				"	250.0		63.5%	48.1-108		
Surrogate: Terphenyl-d14	150				"	250.0		60.0%	59.4-127		



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Notes and Definitions

- J The identification of the analyte is acceptable; the reported value is an estimate.
- U Not Detected
- NR Not Reported