



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

**Date:** 3/22/2016  
**Subject:** Review of Region 5 Data for L'Anse, Michigan Biomass Utility  
**To:** Air Division, US EPA Region 5  
77 West Jackson Boulevard  
Chicago, IL 60605  
**From:** Troy Stroock, Chemist  
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, Robert Thompson at (312)-353-9078 for sample project concerns, and Sylvia Griffin at (312)-353-9073 with data transmittal questions. Thank you.

**Attached are Results for: L'Anse, Michigan Biomass Utility**

\_\_\_\_\_  
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

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

Project: L'Anse, Michigan Biomass Utility  
Project Number: [none]  
Project Manager: Molly Smith

**Reported:**  
Mar-22-16 13:23

## ANALYSIS CASE NARRATIVE

Analyst Phone number: 312.353.8362

### General Information

5 surface wipe samples were received by the Chicago Regional Laboratory (CRL) on 2/4/2016 for preparation and analysis of polycyclic aromatic hydrocarbons (PAHs) by CRL Standard Operating Procedure (SOP) MS026 V1. The procedures are based on SW-846 methods 3545 A (pressurized fluid extraction) and 8270D (Semivolatile Organic Compounds by GC/MS). Pen and Ink change 5670 documents modifications to the SOP to accommodate testing of surface wipes for the analytes of interest, and pen and ink change MS026 Rev 1 PI07 was also relevant to this data set regarding reporting of perylene as a target analyte. Only measurements of target PAHs and associated neutral surrogates are presented in this report.

Holding times were met for preparation and analysis of all field samples.

### Sample Analysis and Results

Higher levels of PAHs were measured in laboratory sample IDs 1602022-01 and -02, while sample ID 1602022-03 had lower levels of PAHs, even though many of them were measured above the reporting limit (RL). These extracts required 10 to 100-fold dilutions to interpolate all target analyte concentration measurements. All measured PAH concentrations were below the RL in samples 1602022-04 and -05 except for anthracene in 1602022-04, which was measured just above the RL.

### Quality Controls

Please refer to the report for qualifiers added by analyte; the key at the end of the report contains descriptions of each data qualifier added and the expected impact on the data. All other quality controls not mentioned below met SOP criteria.

3 neutral surrogates were used for data evaluation: nitrobenzene-d5, 2-fluorobiphenyl, and terphenyl-d14. 2-methylnaphthalene-d10 and fluoranthene-d10 were also added to all field samples and QC samples, and this data is included in the report, but it is intended for the client's information purposes only. Of the surrogates used for data evaluation, nitrobenzene-d5 did not meet continuing calibration verification (CCV) acceptance criteria in one or both CCVs bracketing analysis of all field samples and QC sample B16C023-MRL1 (20 ng/wipe reporting limit check). The affected data is qualified appropriately. Recovery of nitrobenzene-d5 was also above



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the upper acceptance limit in all field samples and batch QC samples, but sample data is only qualified if 2 or more surrogates are outside of control limits, so no qualifiers were added.

Also, dibenz(ah)anthracene did not meet the CCV acceptance criteria in one of the CCVs bracketing analysis of field samples 1602022-01, -02, -04, and -05. The affected data is qualified appropriately.

Standard additions were performed at 20 ng/wipe (B16C023-MRL1), 100 ng/wipe (B16C023-MRL2), and 500 ng/wipe (B16C023-BS1 and -BSD1) in order to evaluate performance by target analyte. Recovery of benzo(a)anthracene and benzo(b)fluoranthene was slightly above the upper acceptance limit in B16C023-MRL1, but recovery of these analytes was acceptable in the other QC samples. Concentrations of these chemicals were measured in the field samples at concentrations >100 ng/wipe where recovery was demonstrated to be acceptable (laboratory sample IDs 1602022-01, -02, and -03), or they were not found above the RL (laboratory sample IDs 1602022-04 and -05). Therefore, no target analytes are qualified in any field samples for high bias recovery in B16C023-MRL1.

Benzo(b)fluoranthene and benzo(k)fluoranthene met the SOP resolution criteria (>50%) in the CCV standards where they are at the same concentration, but in the field samples where they were at different relative concentration the resolution criteria was not met. Therefore, the concentrations of these two target analytes are qualified as estimated in the field samples where they were measured above the RL (in laboratory sample IDs 1602022-01, -02, and -03).

Internal standard responses were above the upper acceptance limit in the ending CCV standards analyzed with both sequences included in this data package. All internal standards met the SOP criteria in the opening CCVs in both sequences, and all target analytes and surrogates met the CCV criteria in these injections as well (except dibenz(ah)anthracene as noted above). The same problem was not observed for any corresponding field samples, so no additional data qualifiers were added.

In field samples 1602022-01, -02, and -03, indeno(123-cd)pyrene did not meet the ion ratio criteria specified in the SOP for the qualifier ion to quantitation ion ( $\pm 30\%$  of the ratio in the midpoint initial calibration standard). Dibenz(ah)anthracene coelutes close to indeno(123-cd)pyrene and also produces a response for the indeno(123-cd)pyrene qualifier ion, thereby enhancing its response in the standard used to estimate what the qualifier ion ratio should be. In the field samples, the concentration of dibenz(ah)anthracene is low relative to that of indeno(123-cd)pyrene, so the qualifier ion response is not the same as in a calibration standard with both indeno(123-cd)pyrene and dibenz(ah)anthracene in it. In the analyst's opinion, qualitative identification of indeno(123-cd)pyrene is reasonable in the field samples (i.e., similar in the calibration standards and field samples) in spite of not meeting the ion ratio criteria in the SOP which is complicated slightly by a coelution, so no additional data qualifiers were added.



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**ANALYTICAL REPORT FOR SAMPLES**

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
L-1P	1602022-01	Wipe	Feb-01-16 14:00	Feb-04-16 09:45
L-2P	1602022-02	Wipe	Feb-01-16 14:15	Feb-04-16 09:45
L-3P	1602022-03	Wipe	Feb-01-16 15:35	Feb-04-16 09:45
L-4P	1602022-04	Wipe	Feb-01-16 15:55	Feb-04-16 09:45
L-5P	1602022-05	Wipe	Feb-01-16 16:20	Feb-04-16 09:45



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

L-1P (1602022-01) Wipe Sampled: Feb-01-16 14:00 Received: Feb-04-16 09:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Naphthalene</b>	<b>385</b>			100	ng/Wipe	1	B16C023	Feb-12-16	Mar-03-16
<b>2-Methylnaphthalene</b>	<b>430</b>			100	"	"	"	"	"
<b>1-Methylnaphthalene</b>	<b>115</b>			100	"	"	"	"	"
<b>Acenaphthylene</b>	<b>497</b>			20.0	"	"	"	"	"
<b>Acenaphthene</b>	<b>475</b>			100	"	"	"	"	"
<b>Fluorene</b>	<b>3310</b>			1000	"	10	"	"	Mar-10-16
<b>Phenanthrene</b>	<b>23100</b>			10000	"	100	"	"	Mar-10-16
<b>Anthracene</b>	<b>8420</b>			200	"	10	"	"	Mar-10-16
<b>Fluoranthene</b>	<b>44600</b>			10000	"	100	"	"	Mar-10-16
<b>Pyrene</b>	<b>23000</b>			10000	"	"	"	"	"
<b>Benzo (a) anthracene</b>	<b>4290</b>			200	"	10	"	"	Mar-10-16
<b>Chrysene</b>	<b>13000</b>			2000	"	100	"	"	Mar-10-16
<b>Benzo(b)fluoranthene</b>	<b>9690</b>	J		200	"	10	"	"	Mar-10-16
<b>Benzo(k)fluoranthene</b>	<b>4090</b>	J		200	"	"	"	"	"
<b>Benzo (e) pyrene</b>	<b>4180</b>			200	"	"	"	"	"
<b>Benzo(a)pyrene</b>	<b>2690</b>			200	"	"	"	"	"
<b>Perylene</b>	<b>349</b>			20.0	"	1	"	"	Mar-03-16
<b>Indeno(1,2,3-cd)pyrene</b>	<b>1990</b>			200	"	10	"	"	Mar-10-16
<b>Dibenz(a,h)anthracene</b>	<b>476</b>	(CCV), J		20.0	"	1	"	"	Mar-03-16
<b>Benzo(g,h,i)perylene</b>	<b>1890</b>			200	"	10	"	"	Mar-10-16

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
<i>Surogate: Nitrobenzene-d5</i>	270	(CCV), J	108%	50-94.9	"	"	Mar-03-16
<i>Surogate: 2-Fluorobiphenyl</i>	185		74.2%	48.1-108	"	"	"
<i>Surogate: Terphenyl-d14</i>	231		92.3%	59.4-127	"	"	"
<i>Surogate: 2-Methylnaphthalene-d10</i>	209		83.5%	45.5-117	"	"	"
<i>Surogate: Fluoranthene-d10</i>	280		112%	70.8-122	"	"	"



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

L-2P (1602022-02) Wipe Sampled: Feb-01-16 14:15 Received: Feb-04-16 09:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Naphthalene	582			100	ng/Wipe	1	B16C023	Feb-12-16	Mar-03-16
2-Methylnaphthalene	840			100	"	"	"	"	"
1-Methylnaphthalene	183			100	"	"	"	"	"
Acenaphthylene	871			20.0	"	"	"	"	"
Acenaphthene	824			100	"	"	"	"	"
Fluorene	7870			1000	"	10	"	"	Mar-11-16
Phenanthrene	57400			10000	"	100	"	"	Mar-11-16
Anthracene	19900			2000	"	"	"	"	"
Fluoranthene	82400			10000	"	"	"	"	"
Pyrene	38800			10000	"	"	"	"	"
Benzo (a) anthracene	6990			200	"	10	"	"	Mar-11-16
Chrysene	21000			2000	"	100	"	"	Mar-11-16
Benzo(b)fluoranthene	13900	J		2000	"	"	"	"	"
Benzo(k)fluoranthene	6500	J		200	"	10	"	"	Mar-11-16
Benzo (e) pyrene	6520			200	"	"	"	"	"
Benzo(a)pyrene	4480			200	"	"	"	"	"
Perylene	555			20.0	"	1	"	"	Mar-03-16
Indeno(1,2,3-cd)pyrene	3250			200	"	10	"	"	Mar-11-16
Dibenz(a,h)anthracene	763	(CCV), J		20.0	"	1	"	"	Mar-03-16
Benzo(g,h,i)perylene	3090			200	"	10	"	"	Mar-11-16

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Nitrobenzene-d5	292	(CCV), J	117%	50-94.9	"	"	Mar-03-16
Surrogate: 2-Fluorobiphenyl	183		73.2%	48.1-108	"	"	"
Surrogate: Terphenyl-d14	243		97.0%	59.4-127	"	"	"
Surrogate: 2-Methylnaphthalene-d10	211		84.2%	45.5-117	"	"	"
Surrogate: Fluoranthene-d10	300		120%	70.8-122	"	"	"



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

L-3P (1602022-03) Wipe    Sampled: Feb-01-16 15:35    Received: Feb-04-16 09:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Naphthalene	U			100	ng/Wipe	1	B16C023	Feb-12-16	Mar-10-16
2-Methylnaphthalene	U			100	"	"	"	"	"
1-Methylnaphthalene	U			100	"	"	"	"	"
Acenaphthylene	48.8			20.0	"	"	"	"	"
Acenaphthene	U			100	"	"	"	"	"
Fluorene	352			100	"	"	"	"	"
Phenanthrene	1860			1000	"	10	"	"	Mar-10-16
Anthracene	1050			200	"	"	"	"	"
Fluoranthene	3350			1000	"	"	"	"	"
Pyrene	1880			1000	"	"	"	"	"
Benzo (a) anthracene	387			20.0	"	1	"	"	Mar-10-16
Chrysene	976			20.0	"	"	"	"	"
Benzo(b)fluoranthene	807	J		20.0	"	"	"	"	"
Benzo(k)fluoranthene	309	J		20.0	"	"	"	"	"
Benzo (e) pyrene	357			20.0	"	"	"	"	"
Benzo(a)pyrene	290			20.0	"	"	"	"	"
Perylene	46.5			20.0	"	"	"	"	"
Indeno(1,2,3-cd)pyrene	198			20.0	"	"	"	"	"
Dibenz(a,h)anthracene	49.0			20.0	"	"	"	"	"
Benzo(g,h,i)perylene	194			20.0	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
<i>Surrogate: Nitrobenzene-d5</i>	292	(CCV), J	117%	50-94.9	"	"	"
<i>Surrogate: 2-Fluorobiphenyl</i>	179		71.8%	48.1-108	"	"	"
<i>Surrogate: Terphenyl-d14</i>	245		98.0%	59.4-127	"	"	"
<i>Surrogate: 2-Methylnaphthalene-d10</i>	194		77.6%	45.5-117	"	"	"
<i>Surrogate: Fluoranthene-d10</i>	239		95.4%	70.8-122	"	"	"



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

L-4P (1602022-04) Wipe Sampled: Feb-01-16 15:55 Received: Feb-04-16 09:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Naphthalene	U			100	ng/Wipe	1	B16C023	Feb-12-16	Mar-03-16
2-Methylnaphthalene	U			100	"	"	"	"	"
1-Methylnaphthalene	U			100	"	"	"	"	"
Acenaphthylene	U			20.0	"	"	"	"	"
Acenaphthene	U			100	"	"	"	"	"
Fluorene	U			100	"	"	"	"	"
Phenanthrene	U			100	"	"	"	"	"
Anthracene	20.6			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			20.0	"	"	"	"	"
Indeno(1,2,3-cd)pyrene	U			20.0	"	"	"	"	"
Dibenz(a,h)anthracene	U	(CCV), J		20.0	"	"	"	"	"
Benzo(g,h,i)perylene	U			20.0	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Nitrobenzene-d5	286	(CCV), J	114%	50-94.9	"	"	"
Surrogate: 2-Fluorobiphenyl	177		70.9%	48.1-108	"	"	"
Surrogate: Terphenyl-d14	234		93.8%	59.4-127	"	"	"
Surrogate: 2-Methylnaphthalene-d10	197		78.9%	45.5-117	"	"	"
Surrogate: Fluoranthene-d10	258		103%	70.8-122	"	"	"



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

L-5P (1602022-05) Wipe    Sampled: Feb-01-16 16:20    Received: Feb-04-16 09:45

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Naphthalene	U			100	ng/Wipe	1	B16C023	Feb-12-16	Mar-03-16
2-Methylnaphthalene	U			100	"	"	"	"	"
1-Methylnaphthalene	U			100	"	"	"	"	"
Acenaphthylene	U			20.0	"	"	"	"	"
Acenaphthene	U			100	"	"	"	"	"
Fluorene	U			100	"	"	"	"	"
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			20.0	"	"	"	"	"
Indeno(1,2,3-cd)pyrene	U			20.0	"	"	"	"	"
Dibenz(a,h)anthracene	U	(CCV), J		20.0	"	"	"	"	"
Benzo(g,h,i)perylene	U			20.0	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Nitrobenzene-d5	296	(CCV), J	118%	50-94.9	"	"	"
Surrogate: 2-Fluorobiphenyl	186		74.5%	48.1-108	"	"	"
Surrogate: Terphenyl-d14	240		95.8%	59.4-127	"	"	"
Surrogate: 2-Methylnaphthalene-d10	213		85.1%	45.5-117	"	"	"
Surrogate: Fluoranthene-d10	265		106%	70.8-122	"	"	"



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

### Batch B16C023 - Solvent Extraction

Blank (B16C023-BLK1)

Prepared: Feb-12-16 Analyzed: Mar-02-16

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			100	"						
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			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	254				"	250.0		102%	50-94.9		
Surrogate: 2-Fluorobiphenyl	187				"	250.0		74.7%	48.1-108		
Surrogate: Terphenyl-d14	230				"	250.0		92.1%	59.4-127		
Surrogate: 2-Methylnaphthalene-d10	199				"	250.0		79.5%	45.5-117		
Surrogate: Fluoranthene-d10	246				"	250.0		98.2%	70.8-122		

Blank (B16C023-BLK2)

Prepared: Feb-12-16 Analyzed: Mar-02-16

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	"						



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**Reported:**  
Mar-22-16 13:23

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

### Batch B16C023 - Solvent Extraction

Blank (B16C023-BLK2)

Prepared: Feb-12-16 Analyzed: Mar-02-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Acenaphthene	U			100	ng/Wipe						
Fluorene	U			100	"						
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			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	276				"	250.0		110%	50-94.9		
Surrogate: 2-Fluorobiphenyl	186				"	250.0		74.2%	48.1-108		
Surrogate: Terphenyl-d14	244				"	250.0		97.6%	59.4-127		
Surrogate: 2-Methylnaphthalene-d10	199				"	250.0		79.4%	45.5-117		
Surrogate: Fluoranthene-d10	252				"	250.0		101%	70.8-122		

LCS (B16C023-BS1)

Prepared: Feb-12-16 Analyzed: Mar-02-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Naphthalene	356			100	ng/Wipe	500.0		71.2%	43.1-108		
2-Methylnaphthalene	402			100	"	500.0		80.4%	45.5-117		
1-Methylnaphthalene	406			100	"	500.0		81.2%	45.5-117		
Acenaphthylene	443			20.0	"	500.0		88.6%	52.7-117		
Acenaphthene	397			100	"	500.0		79.4%	51.7-116		
Fluorene	453			100	"	500.0		90.6%	57.9-120		
Phenanthrene	395			100	"	500.0		79.0%	68.1-114		
Anthracene	468			20.0	"	500.0		93.7%	69.7-116		



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Air Division, US EPA Region 5  
77 West Jackson Boulevard  
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Project: L'Anse, Michigan Biomass Utility  
Project Number: [none]  
Project Manager: Molly Smith

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

### Batch B16C023 - Solvent Extraction

LCS (B16C023-BS1)

Prepared: Feb-12-16 Analyzed: Mar-02-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Fluoranthene	480			100	ng/Wipe	500.0		96.1% 70.8-122		
Pyrene	443			100	"	500.0		88.6% 71.8-117		
Benzo (a) anthracene	553			20.0	"	500.0		111% 67.6-115		
Chrysene	412			20.0	"	500.0		82.4% 68.5-117		
Benzo(b)fluoranthene	526			20.0	"	500.0		105% 68.9-128		
Benzo(k)fluoranthene	474			20.0	"	500.0		94.8% 62.8-135		
Benzo (e) pyrene	466			20.0	"	500.0		93.2% 68.9-133		
Benzo(a)pyrene	511			20.0	"	500.0		102% 68.9-133		
Perylene	400			20.0	"	500.0		80.0% 68.9-133		
Indeno(1,2,3-cd)pyrene	508			20.0	"	500.0		102% 70-129		
Dibenz(a,h)anthracene	537			20.0	"	500.0		107% 69.1-131		
Benzo(g,h,i)perylene	455			20.0	"	500.0		91.0% 53.9-139		
Surrogate: Nitrobenzene-d5	263				"	250.0		105% 50-94.9		
Surrogate: 2-Fluorobiphenyl	186				"	250.0		74.4% 48.1-108		
Surrogate: Terphenyl-d14	254				"	250.0		102% 59.4-127		
Surrogate: 2-Methylnaphthalene-d10	199				"	250.0		79.7% 45.5-117		
Surrogate: Fluoranthene-d10	251				"	250.0		100% 70.8-122		

LCS Dup (B16C023-BSD1)

Prepared: Feb-12-16 Analyzed: Mar-02-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Naphthalene	358			100	ng/Wipe	500.0		71.6% 43.1-108	0.580	30.2
2-Methylnaphthalene	401			100	"	500.0		80.3% 45.5-117	0.182	36.4
1-Methylnaphthalene	393			100	"	500.0		78.5% 45.5-117	3.33	36.4
Acenaphthylene	436			20.0	"	500.0		87.2% 52.7-117	1.53	30
Acenaphthene	395			100	"	500.0		79.0% 51.7-116	0.508	30
Fluorene	446			100	"	500.0		89.1% 57.9-120	1.69	30
Phenanthrene	384			100	"	500.0		76.8% 68.1-114	2.78	30
Anthracene	458			20.0	"	500.0		91.7% 69.7-116	2.12	30
Fluoranthene	463			100	"	500.0		92.7% 70.8-122	3.65	30
Pyrene	414			100	"	500.0		82.9% 71.8-117	6.73	30
Benzo (a) anthracene	527			20.0	"	500.0		105% 67.6-115	4.89	30
Chrysene	386			20.0	"	500.0		77.1% 68.5-117	6.61	30



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

### Batch B16C023 - Solvent Extraction

#### LCS Dup (B16C023-BSD1)

Prepared: Feb-12-16 Analyzed: Mar-02-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	Limit	RPD	RPD Limit
Benzo(b)fluoranthene	511			20.0	ng/Wipe	500.0		102%	68.9-128	2.87	30
Benzo(k)fluoranthene	456			20.0	"	500.0		91.2%	62.8-135	3.90	30
Benzo (e) pyrene	451			20.0	"	500.0		90.2%	68.9-133	3.31	30
Benzo(a)pyrene	505			20.0	"	500.0		101%	68.9-133	1.16	30
Perylene	383			20.0	"	500.0		76.6%	68.9-133	4.37	30
Indeno(1,2,3-cd)pyrene	498			20.0	"	500.0		99.7%	70-129	1.92	30
Dibenz(a,h)anthracene	530			20.0	"	500.0		106%	69.1-131	1.45	30
Benzo(g,h,i)perylene	448			20.0	"	500.0		89.5%	53.9-139	1.65	30
Surrogate: Nitrobenzene-d5	279				"	250.0		112%	50-94.9		
Surrogate: 2-Fluorobiphenyl	183				"	250.0		73.1%	48.1-108		
Surrogate: Terphenyl-d14	226				"	250.0		90.5%	59.4-127		
Surrogate: 2-Methylnaphthalene-d10	200				"	250.0		79.9%	45.5-117		
Surrogate: Fluoranthene-d10	244				"	250.0		97.6%	70.8-122		

#### MRL Check (B16C023-MRL1)

Prepared: Feb-12-16 Analyzed: Mar-10-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	Limit	RPD	RPD Limit
Acenaphthylene	22.5			20.0	ng/Wipe	20.00		112%	52.7-117		
Anthracene	18.9			20.0	"	20.00		94.4%	69.7-116		
Benzo (a) anthracene	24.2			20.0	"	20.00		121%	67.6-115		
Chrysene	16.7			20.0	"	20.00		83.7%	68.5-117		
Benzo(b)fluoranthene	26.1			20.0	"	20.00		130%	68.9-128		
Benzo(k)fluoranthene	21.7			20.0	"	20.00		108%	62.8-135		
Benzo (e) pyrene	19.3			20.0	"	20.00		96.6%	68.9-133		
Benzo(a)pyrene	19.4			20.0	"	20.00		97.1%	68.9-133		
Perylene	15.9			20.0	"	20.00		79.4%	68.9-133		
Indeno(1,2,3-cd)pyrene	23.4			20.0	"	20.00		117%	70-129		
Dibenz(a,h)anthracene	22.5			20.0	"	20.00		113%	69.1-131		
Benzo(g,h,i)perylene	18.8			20.0	"	20.00		94.2%	53.9-139		
Surrogate: Nitrobenzene-d5	281	(CCV), J			"	250.0		112%	50-94.9		
Surrogate: 2-Fluorobiphenyl	188				"	250.0		75.2%	48.1-108		
Surrogate: Terphenyl-d14	249				"	250.0		99.4%	59.4-127		
Surrogate: 2-Methylnaphthalene-d10	199				"	250.0		79.4%	45.5-117		



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### Batch B16C023 - Solvent Extraction

#### MRL Check (B16C023-MRL1)

Prepared: Feb-12-16 Analyzed: Mar-10-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<i>Surrogate: Fluoranthene-d10</i>	244				ng/Wipe	250.0		97.7%	70.8-122		

#### MRL Check (B16C023-MRL2)

Prepared: Feb-12-16 Analyzed: Mar-02-16

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<b>Naphthalene</b>	<b>84.4</b>			100	ng/Wipe	100.0		84.4%	43.1-108		
<b>2-Methylnaphthalene</b>	<b>90.0</b>			100	"	100.0		90.0%	45.5-117		
<b>1-Methylnaphthalene</b>	<b>86.6</b>			100	"	100.0		86.6%	45.5-117		
<b>Acenaphthene</b>	<b>81.6</b>			100	"	100.0		81.6%	51.7-116		
<b>Fluorene</b>	<b>94.1</b>			100	"	100.0		94.1%	57.9-120		
<b>Phenanthrene</b>	<b>83.4</b>			100	"	100.0		83.4%	68.1-114		
<b>Fluoranthene</b>	<b>97.7</b>			100	"	100.0		97.7%	70.8-122		
<b>Pyrene</b>	<b>91.2</b>			100	"	100.0		91.2%	71.8-117		
<i>Surrogate: Nitrobenzene-d5</i>	284				"	250.0		114%	50-94.9		
<i>Surrogate: 2-Fluorobiphenyl</i>	192				"	250.0		77.0%	48.1-108		
<i>Surrogate: Terphenyl-d14</i>	245				"	250.0		97.9%	59.4-127		
<i>Surrogate: 2-Methylnaphthalene-d10</i>	210				"	250.0		83.8%	45.5-117		
<i>Surrogate: Fluoranthene-d10</i>	251				"	250.0		100%	70.8-122		



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**Notes and Definitions**

- J The identification of the analyte is acceptable; the reported value is an estimate.
- (CCV) Continuing calibration verification criteria not met for this analyte
- U Not Detected
- NR Not Reported