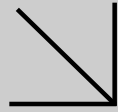




Supplemental Report 1

The original report has been revised to include the Level III deliverables package.

**WORK ORDER NUMBER: 17-03-1557**

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For**Client:** Andersen Environmental**Client Project Name:** Burbank Airport / 9836002041

Attention: Brian Martasin
5261 West Imperial Highway
Los Angeles, CA 90045-6231

A handwritten signature in black ink, appearing to read "S. Nowak".

Approved for release on 04/28/2017 by:
Stephen Nowak
Project Manager

ResultLink ▶

Email your PM ▶

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



Calscience

Contents

Client Project Name: Burbank Airport / 9836002041
 Work Order Number: 17-03-1557

1	Work Order Narrative.	3
2	Sample Summary.	4
3	Detections Summary.	5
4	Client Sample Data.	6
	4.1 ASTM D-2216 (M) Moisture Content (Solid).	6
	4.2 EPA 8015B (M) Diesel and Motor Oil Ranges (Solid).	7
	4.3 EPA 6010B/7471A CAC Title 22 Metals (Solid).	9
	4.4 EPA 7471A Mercury (Solid).	13
	4.5 EPA 8082 PCB Aroclors (Solid).	14
	4.6 EPA 8270C SIM PAHs (Solid).	16
5	Quality Control Sample Data.	20
	5.1 MS/MSD.	20
	5.2 Sample Duplicate.	25
	5.3 LCS/LCSD.	26
6	Sample Analysis Summary.	31
7	Glossary of Terms and Qualifiers.	32
8	Chain-of-Custody/Sample Receipt Form.	33
9	Level III Data Package - Case Narrative.	45
	9.1 ASTM D-2216 (M) Moisture Content (Solid).	50
	9.2 EPA 8015 (M) Diesel and Motor Oil Ranges (Solid).	60
	9.3 EPA 6010B Title 22 Metals (Solid).	131
	9.4 EPA 7471A Mercury (Solid).	190
	9.5 EPA 8082 PCB Aroclors (Solid).	221
	9.6 EPA 8270C SIM PAHs (Solid).	299

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 03/21/17. They were assigned to Work Order 17-03-1557.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Sample Summary

Client: Andersen Environmental	Work Order:	17-03-1557
5261 West Imperial Highway	Project Name:	Burbank Airport / 9836002041
Los Angeles, CA 90045-6231	PO Number:	
	Date/Time Received:	03/21/17 15:37
	Number of Containers:	30

Attn: Brian Martasin

Sample Identification	Lab Number	Collection Date and Time	Number of Containers	Matrix
B-DU1-S-06-8	17-03-1557-1	03/16/17 12:25	1	Solid
B-DU1-S-SG-02-8S	17-03-1557-2	03/16/17 07:25	1	Solid
B-DU1-S-SG-07-8S	17-03-1557-3	03/16/17 12:55	1	Solid
B-DU1-S-SG-08-8S	17-03-1557-4	03/16/17 15:10	1	Solid
B-DU1-S-SG-10-8S	17-03-1557-5	03/16/17 09:22	1	Solid
B-DU1-S-SG-11-8S	17-03-1557-6	03/16/17 10:47	1	Solid
B-DU1-S-04-8	17-03-1557-7	03/16/17 14:25	1	Solid
B-DU1-S-SG-01-8S	17-03-1557-8	03/17/17 13:00	1	Solid
B-DU1-S-SG-03-8S	17-03-1557-9	03/17/17 12:20	1	Solid
B-DU1-S-SG-04-8S	17-03-1557-10	03/17/17 10:30	1	Solid
B-DU1-S-03-8	17-03-1557-11	03/17/17 13:35	1	Solid
B-DU1-S-01-8	17-03-1557-12	03/20/17 14:20	1	Solid
B-DU1-S-02-8	17-03-1557-13	03/20/17 11:00	1	Solid
B-DU1-S-05-8	17-03-1557-14	03/20/17 10:20	1	Solid
B-DU1-S-07-8	17-03-1557-15	03/20/17 09:45	1	Solid
B-DU1-S-08-8	17-03-1557-16	03/20/17 13:05	1	Solid
B-DU1-S-09-8	17-03-1557-17	03/20/17 09:21	1	Solid
B-DU1-S-SG-05-8S	17-03-1557-18	03/20/17 08:20	1	Solid
B-DU1-S-SG-06-8S	17-03-1557-19	03/20/17 12:42	1	Solid
B-DU1-S-SG-09-8S	17-03-1557-20	03/20/17 07:40	1	Solid
COMPOSITE	17-03-1557-21	03/16/17 00:00	1	Solid
B-DU1-ISM1-8	17-03-1557-22	03/16/17 00:00	1	Solid
B-DU1-ISM2-8	17-03-1557-23	03/16/17 00:00	1	Solid
B-DU1-ISM3-8	17-03-1557-24	03/16/17 00:00	1	Solid
B-DU1-ISM1-8	17-03-1557-25	03/16/17 00:00	1	Solid
B-DU1-ISM2-8	17-03-1557-26	03/16/17 00:00	1	Solid
B-DU1-ISM3-8	17-03-1557-27	03/16/17 00:00	1	Solid
B-DU1-ISM1-8	17-03-1557-28	03/16/17 00:00	1	Solid
B-DU1-ISM2-8	17-03-1557-29	03/16/17 00:00	1	Solid
B-DU1-ISM3-8	17-03-1557-30	03/16/17 00:00	1	Solid

Detections Summary

Client: Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Work Order: 17-03-1557
 Project Name: Burbank Airport / 9836002041
 Received: 03/21/17

Attn: Brian Martasin

Page 1 of 1

Client SampleID

Analyte	Result	Qualifiers	RL	Units	Method	Extraction
B-DU1-ISM1-8 (17-03-1557-25)						
Moisture	3.0		0.10	%	ASTM D-2216 (M)	N/A
B-DU1-ISM2-8 (17-03-1557-26)						
Moisture	2.8		0.10	%	ASTM D-2216 (M)	N/A
B-DU1-ISM3-8 (17-03-1557-27)						
Moisture	3.0		0.10	%	ASTM D-2216 (M)	N/A
B-DU1-ISM1-8 (17-03-1557-28)						
Arsenic	1.37		0.739	mg/kg	EPA 6010B	EPA 3050B
Barium	61.1		0.493	mg/kg	EPA 6010B	EPA 3050B
Chromium	4.85		0.246	mg/kg	EPA 6010B	EPA 3050B
Cobalt	4.92		0.246	mg/kg	EPA 6010B	EPA 3050B
Copper	6.93		0.493	mg/kg	EPA 6010B	EPA 3050B
Lead	4.39		0.493	mg/kg	EPA 6010B	EPA 3050B
Nickel	4.31		0.246	mg/kg	EPA 6010B	EPA 3050B
Vanadium	16.5		0.246	mg/kg	EPA 6010B	EPA 3050B
Zinc	24.3		0.985	mg/kg	EPA 6010B	EPA 3050B
B-DU1-ISM2-8 (17-03-1557-29)						
Barium	66.9		0.490	mg/kg	EPA 6010B	EPA 3050B
Chromium	4.74		0.245	mg/kg	EPA 6010B	EPA 3050B
Cobalt	4.75		0.245	mg/kg	EPA 6010B	EPA 3050B
Copper	6.71		0.490	mg/kg	EPA 6010B	EPA 3050B
Lead	3.46		0.490	mg/kg	EPA 6010B	EPA 3050B
Nickel	3.81		0.245	mg/kg	EPA 6010B	EPA 3050B
Vanadium	15.8		0.245	mg/kg	EPA 6010B	EPA 3050B
Zinc	24.3		0.980	mg/kg	EPA 6010B	EPA 3050B
B-DU1-ISM3-8 (17-03-1557-30)						
Arsenic	0.950		0.735	mg/kg	EPA 6010B	EPA 3050B
Barium	56.9		0.490	mg/kg	EPA 6010B	EPA 3050B
Chromium	4.12		0.245	mg/kg	EPA 6010B	EPA 3050B
Cobalt	4.22		0.245	mg/kg	EPA 6010B	EPA 3050B
Copper	6.12		0.490	mg/kg	EPA 6010B	EPA 3050B
Lead	3.74		0.490	mg/kg	EPA 6010B	EPA 3050B
Nickel	3.48		0.245	mg/kg	EPA 6010B	EPA 3050B
Vanadium	13.4		0.245	mg/kg	EPA 6010B	EPA 3050B
Zinc	23.2		0.980	mg/kg	EPA 6010B	EPA 3050B

Subcontracted analyses, if any, are not included in this summary.

* MDL is shown

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: N/A
 Method: ASTM D-2216 (M)
 Units: %

Project: Burbank Airport / 9836002041

Page 1 of 1

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM1-8	17-03-1557-25-A	03/16/17 00:00	Solid	N/A	03/23/17	03/23/17 21:00	H0323MOIB3
<u>Parameter</u>		<u>Result</u>	<u>RL</u>		<u>DF</u>		<u>Qualifiers</u>
Moisture		3.0	0.10		1.00		
B-DU1-ISM2-8	17-03-1557-26-A	03/16/17 00:00	Solid	N/A	03/23/17	03/23/17 21:00	H0323MOIB3
<u>Parameter</u>		<u>Result</u>	<u>RL</u>		<u>DF</u>		<u>Qualifiers</u>
Moisture		2.8	0.10		1.00		
B-DU1-ISM3-8	17-03-1557-27-A	03/16/17 00:00	Solid	N/A	03/23/17	03/23/17 21:00	H0323MOIB3
<u>Parameter</u>		<u>Result</u>	<u>RL</u>		<u>DF</u>		<u>Qualifiers</u>
Moisture		3.0	0.10		1.00		
Method Blank	099-05-014-6770	N/A	Solid	N/A	03/23/17	03/23/17 21:00	H0323MOIB3
<u>Parameter</u>		<u>Result</u>	<u>RL</u>		<u>DF</u>		<u>Qualifiers</u>
Moisture		ND	0.10		1.00		

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3550B
 Method: EPA 8015B (M)
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 1 of 2

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM1-8	17-03-1557-25-C	03/16/17 00:00	Solid	GC 45	03/24/17	03/28/17 14:08	170324B12

Comment(s): - Results are reported on a dry weight basis.
 - Motor Oil Range Organics (C17-C44) uses a Diesel Range Organics (C10-C28) standard for quantitation and quality control.

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel	ND	5.1	1.00	
TPH as Motor Oil	ND	5.1	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
n-Octacosane	102	61-145	

B-DU1-ISM2-8	17-03-1557-26-C	03/16/17 00:00	Solid	GC 45	03/24/17	03/28/17 14:29	170324B12
---------------------	------------------------	---------------------------	--------------	--------------	-----------------	---------------------------	------------------

Comment(s): - Results are reported on a dry weight basis.
 - Motor Oil Range Organics (C17-C44) uses a Diesel Range Organics (C10-C28) standard for quantitation and quality control.

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel	ND	5.0	1.00	
TPH as Motor Oil	ND	5.0	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
n-Octacosane	105	61-145	

B-DU1-ISM3-8	17-03-1557-27-C	03/16/17 00:00	Solid	GC 45	03/24/17	03/28/17 14:51	170324B12
---------------------	------------------------	---------------------------	--------------	--------------	-----------------	---------------------------	------------------

Comment(s): - Results are reported on a dry weight basis.
 - Motor Oil Range Organics (C17-C44) uses a Diesel Range Organics (C10-C28) standard for quantitation and quality control.

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel	ND	5.1	1.00	
TPH as Motor Oil	ND	5.1	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
n-Octacosane	101	61-145	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3550B
Method: EPA 8015B (M)
Units: mg/kg

Project: Burbank Airport / 9836002041

Page 2 of 2

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-14-353-13	N/A	Solid	GC 45	03/24/17	03/28/17 11:16	170324B12

Comment(s): - Motor Oil Range Organics (C17-C44) uses a Diesel Range Organics (C10-C28) standard for quantitation and quality control.

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel	ND	5.0	1.00	
TPH as Motor Oil	ND	5.0	1.00	
<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>	
n-Octacosane	103	61-145		

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3050B
 Method: EPA 6010B
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 1 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM1-8	17-03-1557-28-A	03/16/17 00:00	Solid	ICP 7300	03/29/17	03/29/17 18:56	170329L05

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Antimony	ND	0.739	0.985	
Arsenic	1.37	0.739	0.985	
Barium	61.1	0.493	0.985	
Beryllium	ND	0.246	0.985	
Cadmium	ND	0.493	0.985	
Chromium	4.85	0.246	0.985	
Cobalt	4.92	0.246	0.985	
Copper	6.93	0.493	0.985	
Lead	4.39	0.493	0.985	
Molybdenum	ND	0.246	0.985	
Nickel	4.31	0.246	0.985	
Selenium	ND	0.739	0.985	
Silver	ND	0.246	0.985	
Thallium	ND	0.739	0.985	
Vanadium	16.5	0.246	0.985	
Zinc	24.3	0.985	0.985	

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3050B
 Method: EPA 6010B
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 2 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM2-8	17-03-1557-29-A	03/16/17 00:00	Solid	ICP 7300	03/29/17	03/29/17 18:57	170329L05

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Antimony	ND	0.735	0.980	
Arsenic	ND	0.735	0.980	
Barium	66.9	0.490	0.980	
Beryllium	ND	0.245	0.980	
Cadmium	ND	0.490	0.980	
Chromium	4.74	0.245	0.980	
Cobalt	4.75	0.245	0.980	
Copper	6.71	0.490	0.980	
Lead	3.46	0.490	0.980	
Molybdenum	ND	0.245	0.980	
Nickel	3.81	0.245	0.980	
Selenium	ND	0.735	0.980	
Silver	ND	0.245	0.980	
Thallium	ND	0.735	0.980	
Vanadium	15.8	0.245	0.980	
Zinc	24.3	0.980	0.980	

Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3050B
 Method: EPA 6010B
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 3 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM3-8	17-03-1557-30-A	03/16/17 00:00	Solid	ICP 7300	03/29/17	03/29/17 18:58	170329L05

Parameter	Result	RL	DF	Qualifiers
Antimony	ND	0.735	0.980	
Arsenic	0.950	0.735	0.980	
Barium	56.9	0.490	0.980	
Beryllium	ND	0.245	0.980	
Cadmium	ND	0.490	0.980	
Chromium	4.12	0.245	0.980	
Cobalt	4.22	0.245	0.980	
Copper	6.12	0.490	0.980	
Lead	3.74	0.490	0.980	
Molybdenum	ND	0.245	0.980	
Nickel	3.48	0.245	0.980	
Selenium	ND	0.735	0.980	
Silver	ND	0.245	0.980	
Thallium	ND	0.735	0.980	
Vanadium	13.4	0.245	0.980	
Zinc	23.2	0.980	0.980	



 Return to Contents

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3050B
 Method: EPA 6010B
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 4 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	097-01-002-24537	N/A	Solid	ICP 7300	03/29/17	03/29/17 17:07	170329L05

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Antimony	ND	0.732	0.976	
Arsenic	ND	0.732	0.976	
Barium	ND	0.488	0.976	
Beryllium	ND	0.244	0.976	
Cadmium	ND	0.488	0.976	
Chromium	ND	0.244	0.976	
Cobalt	ND	0.244	0.976	
Copper	ND	0.488	0.976	
Lead	ND	0.488	0.976	
Molybdenum	ND	0.244	0.976	
Nickel	ND	0.244	0.976	
Selenium	ND	0.732	0.976	
Silver	ND	0.244	0.976	
Thallium	ND	0.732	0.976	
Vanadium	ND	0.244	0.976	
Zinc	ND	0.976	0.976	

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 7471A Total
 Method: EPA 7471A
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 1 of 1

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM1-8	17-03-1557-25-B	03/16/17 00:00	Solid	Mercury 08	03/29/17	03/29/17 15:54	170329L01

Comment(s): - Results are reported on a dry weight basis.

Parameter	Result	RL	DF	Qualifiers
Mercury	ND	0.0818	1.00	

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM2-8	17-03-1557-26-B	03/16/17 00:00	Solid	Mercury 08	03/29/17	03/29/17 15:56	170329L01

Comment(s): - Results are reported on a dry weight basis.

Parameter	Result	RL	DF	Qualifiers
Mercury	ND	0.0830	1.00	

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM3-8	17-03-1557-27-A	03/16/17 00:00	Solid	Mercury 08	03/29/17	03/29/17 15:59	170329L01

Comment(s): - Results are reported on a dry weight basis.

Parameter	Result	RL	DF	Qualifiers
Mercury	ND	0.0859	1.00	

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-16-272-2910	N/A	Solid	Mercury 08	03/29/17	03/29/17 15:08	170329L01

Parameter	Result	RL	DF	Qualifiers
Mercury	ND	0.0833	1.00	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3545
 Method: EPA 8082
 Units: ug/kg

Project: Burbank Airport / 9836002041

Page 1 of 2

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM1-8	17-03-1557-28-B	03/16/17 00:00	Solid	GC 66	03/27/17	03/28/17 03:41	170327L05

Parameter	Result	RL	DF	Qualifiers
Aroclor-1016	ND	50	1.00	
Aroclor-1221	ND	50	1.00	
Aroclor-1232	ND	50	1.00	
Aroclor-1242	ND	50	1.00	
Aroclor-1248	ND	50	1.00	
Aroclor-1254	ND	50	1.00	
Aroclor-1260	ND	50	1.00	
Aroclor-1262	ND	50	1.00	
Aroclor-1268	ND	50	1.00	

Surrogate	Rec. (%)	Control Limits	Qualifiers
Decachlorobiphenyl	77	24-168	
2,4,5,6-Tetrachloro-m-Xylene	65	25-145	

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM2-8	17-03-1557-29-B	03/16/17 00:00	Solid	GC 66	03/27/17	03/28/17 03:59	170327L05

Parameter	Result	RL	DF	Qualifiers
Aroclor-1016	ND	50	1.00	
Aroclor-1221	ND	50	1.00	
Aroclor-1232	ND	50	1.00	
Aroclor-1242	ND	50	1.00	
Aroclor-1248	ND	50	1.00	
Aroclor-1254	ND	50	1.00	
Aroclor-1260	ND	50	1.00	
Aroclor-1262	ND	50	1.00	
Aroclor-1268	ND	50	1.00	

Surrogate	Rec. (%)	Control Limits	Qualifiers
Decachlorobiphenyl	73	24-168	
2,4,5,6-Tetrachloro-m-Xylene	61	25-145	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3545
 Method: EPA 8082
 Units: ug/kg

Project: Burbank Airport / 9836002041

Page 2 of 2

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM3-8	17-03-1557-30-B	03/16/17 00:00	Solid	GC 66	03/27/17	03/28/17 04:35	170327L05

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Aroclor-1016	ND	50	1.00	
Aroclor-1221	ND	50	1.00	
Aroclor-1232	ND	50	1.00	
Aroclor-1242	ND	50	1.00	
Aroclor-1248	ND	50	1.00	
Aroclor-1254	ND	50	1.00	
Aroclor-1260	ND	50	1.00	
Aroclor-1262	ND	50	1.00	
Aroclor-1268	ND	50	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
Decachlorobiphenyl	76	24-168	
2,4,5,6-Tetrachloro-m-Xylene	70	25-145	

Method Blank	099-12-535-4123	N/A	Solid	GC 66	03/27/17	03/28/17 02:30	170327L05
---------------------	------------------------	------------	--------------	--------------	-----------------	---------------------------	------------------

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Aroclor-1016	ND	50	1.00	
Aroclor-1221	ND	50	1.00	
Aroclor-1232	ND	50	1.00	
Aroclor-1242	ND	50	1.00	
Aroclor-1248	ND	50	1.00	
Aroclor-1254	ND	50	1.00	
Aroclor-1260	ND	50	1.00	
Aroclor-1262	ND	50	1.00	
Aroclor-1268	ND	50	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
Decachlorobiphenyl	70	24-168	
2,4,5,6-Tetrachloro-m-Xylene	69	25-145	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3545
 Method: EPA 8270C SIM PAHs
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 1 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM1-8	17-03-1557-25-D	03/16/17 00:00	Solid	GC/MS EEE	03/24/17	03/27/17 13:57	170324L17

Comment(s): - Results are reported on a dry weight basis.

Parameter	Result	RL	DF	Qualifiers
Naphthalene	ND	0.010	1.00	
2-Methylnaphthalene	ND	0.010	1.00	
1-Methylnaphthalene	ND	0.010	1.00	
Acenaphthylene	ND	0.010	1.00	
Acenaphthene	ND	0.010	1.00	
Fluorene	ND	0.010	1.00	
Phenanthrene	ND	0.010	1.00	
Anthracene	ND	0.010	1.00	
Fluoranthene	ND	0.010	1.00	
Pyrene	ND	0.010	1.00	
Benzo (a) Anthracene	ND	0.010	1.00	
Chrysene	ND	0.010	1.00	
Benzo (k) Fluoranthene	ND	0.010	1.00	
Benzo (b) Fluoranthene	ND	0.010	1.00	
Benzo (a) Pyrene	ND	0.010	1.00	
Indeno (1,2,3-c,d) Pyrene	ND	0.010	1.00	
Dibenz (a,h) Anthracene	ND	0.010	1.00	
Benzo (g,h,i) Perylene	ND	0.010	1.00	

Surrogate	Rec. (%)	Control Limits	Qualifiers
2-Fluorobiphenyl	83	13-127	
Nitrobenzene-d5	79	17-137	
p-Terphenyl-d14	94	4-160	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3545
 Method: EPA 8270C SIM PAHs
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 2 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM2-8	17-03-1557-26-D	03/16/17 00:00	Solid	GC/MS EEE	03/24/17	03/27/17 14:17	170324L17

Comment(s): - Results are reported on a dry weight basis.

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	ND	0.010	1.00	
2-Methylnaphthalene	ND	0.010	1.00	
1-Methylnaphthalene	ND	0.010	1.00	
Acenaphthylene	ND	0.010	1.00	
Acenaphthene	ND	0.010	1.00	
Fluorene	ND	0.010	1.00	
Phenanthrene	ND	0.010	1.00	
Anthracene	ND	0.010	1.00	
Fluoranthene	ND	0.010	1.00	
Pyrene	ND	0.010	1.00	
Benzo (a) Anthracene	ND	0.010	1.00	
Chrysene	ND	0.010	1.00	
Benzo (k) Fluoranthene	ND	0.010	1.00	
Benzo (b) Fluoranthene	ND	0.010	1.00	
Benzo (a) Pyrene	ND	0.010	1.00	
Indeno (1,2,3-c,d) Pyrene	ND	0.010	1.00	
Dibenz (a,h) Anthracene	ND	0.010	1.00	
Benzo (g,h,i) Perylene	ND	0.010	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
2-Fluorobiphenyl	91	13-127	
Nitrobenzene-d5	86	17-137	
p-Terphenyl-d14	98	4-160	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 3545
 Method: EPA 8270C SIM PAHs
 Units: mg/kg

Project: Burbank Airport / 9836002041

Page 3 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
B-DU1-ISM3-8	17-03-1557-27-D	03/16/17 00:00	Solid	GC/MS EEE	03/24/17	03/27/17 14:38	170324L17

Comment(s): - Results are reported on a dry weight basis.

Parameter	Result	RL	DF	Qualifiers
Naphthalene	ND	0.010	1.00	
2-Methylnaphthalene	ND	0.010	1.00	
1-Methylnaphthalene	ND	0.010	1.00	
Acenaphthylene	ND	0.010	1.00	
Acenaphthene	ND	0.010	1.00	
Fluorene	ND	0.010	1.00	
Phenanthrene	ND	0.010	1.00	
Anthracene	ND	0.010	1.00	
Fluoranthene	ND	0.010	1.00	
Pyrene	ND	0.010	1.00	
Benzo (a) Anthracene	ND	0.010	1.00	
Chrysene	ND	0.010	1.00	
Benzo (k) Fluoranthene	ND	0.010	1.00	
Benzo (b) Fluoranthene	ND	0.010	1.00	
Benzo (a) Pyrene	ND	0.010	1.00	
Indeno (1,2,3-c,d) Pyrene	ND	0.010	1.00	
Dibenz (a,h) Anthracene	ND	0.010	1.00	
Benzo (g,h,i) Perylene	ND	0.010	1.00	

Surrogate	Rec. (%)	Control Limits	Qualifiers
2-Fluorobiphenyl	99	13-127	
Nitrobenzene-d5	101	17-137	
p-Terphenyl-d14	90	4-160	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.

Analytical Report

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3545
Method: EPA 8270C SIM PAHs
Units: mg/kg

Project: Burbank Airport / 9836002041

Page 4 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-14-035-386	N/A	Solid	GC/MS EEE	03/24/17	03/27/17 11:15	170324L17

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	ND	0.010	1.00	
2-Methylnaphthalene	ND	0.010	1.00	
1-Methylnaphthalene	ND	0.010	1.00	
Acenaphthylene	ND	0.010	1.00	
Acenaphthene	ND	0.010	1.00	
Fluorene	ND	0.010	1.00	
Phenanthrene	ND	0.010	1.00	
Anthracene	ND	0.010	1.00	
Fluoranthene	ND	0.010	1.00	
Pyrene	ND	0.010	1.00	
Benzo (a) Anthracene	ND	0.010	1.00	
Chrysene	ND	0.010	1.00	
Benzo (k) Fluoranthene	ND	0.010	1.00	
Benzo (b) Fluoranthene	ND	0.010	1.00	
Benzo (a) Pyrene	ND	0.010	1.00	
Indeno (1,2,3-c,d) Pyrene	ND	0.010	1.00	
Dibenz (a,h) Anthracene	ND	0.010	1.00	
Benzo (g,h,i) Perylene	ND	0.010	1.00	

<u>Surrogate</u>	<u>Rec. (%)</u>	<u>Control Limits</u>	<u>Qualifiers</u>
2-Fluorobiphenyl	107	13-127	
Nitrobenzene-d5	109	17-137	
p-Terphenyl-d14	104	4-160	

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Calscience

Quality Control - Spike/Spike Duplicate

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3550B
Method: EPA 8015B (M)

Project: Burbank Airport / 9836002041

Page 1 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	MS/MSD Batch Number				
B-DU1-ISM3-8	Sample	Solid	GC 45	03/24/17	03/28/17 14:51	170324S12				
B-DU1-ISM3-8	Matrix Spike	Solid	GC 45	03/24/17	03/28/17 12:42	170324S12				
B-DU1-ISM3-8	Matrix Spike Duplicate	Solid	GC 45	03/24/17	03/28/17 13:03	170324S12				
<u>Parameter</u>	<u>Sample Conc.</u>	<u>Spike Added</u>	<u>MS Conc.</u>	<u>MS %Rec.</u>	<u>MSD Conc.</u>	<u>MSD %Rec.</u>	<u>%Rec. CL</u>	<u>RPD</u>	<u>RPD CL</u>	<u>Qualifiers</u>
TPH as Diesel	ND	400.0	379.0	95	393.3	98	61-145	4	0-25	



Calscience

Quality Control - Spike/Spike Duplicate

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3050B
Method: EPA 6010B

Project: Burbank Airport / 9836002041

Page 2 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	MS/MSD Batch Number
17-03-1813-1	Sample	Solid	ICP 7300	03/29/17	03/29/17 17:12	170329S05
17-03-1813-1	Matrix Spike	Solid	ICP 7300	03/29/17	03/29/17 17:13	170329S05
17-03-1813-1	Matrix Spike Duplicate	Solid	ICP 7300	03/29/17	03/29/17 17:16	170329S05

Parameter	Sample Conc.	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Antimony	ND	25.00	11.32	45	11.05	44	50-115	2	0-20	3
Arsenic	2.141	25.00	28.57	106	26.28	97	75-125	8	0-20	
Barium	65.88	25.00	95.97	120	78.14	49	75-125	20	0-20	3
Beryllium	0.3817	25.00	26.81	106	25.14	99	75-125	6	0-20	
Cadmium	ND	25.00	28.11	112	26.45	106	75-125	6	0-20	
Chromium	12.69	25.00	39.67	108	35.64	92	75-125	11	0-20	
Cobalt	7.098	25.00	35.02	112	32.08	100	75-125	9	0-20	
Copper	10.42	25.00	37.01	106	33.11	91	75-125	11	0-20	
Lead	3.987	25.00	32.73	115	30.39	106	75-125	7	0-20	
Molybdenum	ND	25.00	24.76	99	23.57	94	75-125	5	0-20	
Nickel	8.515	25.00	34.90	106	31.89	93	75-125	9	0-20	
Selenium	ND	25.00	27.32	109	25.65	103	75-125	6	0-20	
Silver	ND	12.50	13.18	105	12.25	98	75-125	7	0-20	
Thallium	ND	25.00	25.36	101	24.63	99	75-125	3	0-20	
Vanadium	24.52	25.00	50.22	103	43.83	77	75-125	14	0-20	
Zinc	29.81	25.00	57.58	111	49.16	77	75-125	16	0-20	

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 7471A Total
Method: EPA 7471A

Project: Burbank Airport / 9836002041

Page 3 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	MS/MSD Batch Number				
17-03-1569-1	Sample	Sediment	Mercury 08	03/29/17	03/29/17 15:13	170329S01				
17-03-1569-1	Matrix Spike	Sediment	Mercury 08	03/29/17	03/29/17 15:20	170329S01				
17-03-1569-1	Matrix Spike Duplicate	Sediment	Mercury 08	03/29/17	03/29/17 15:22	170329S01				
Parameter	Sample Conc.	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Mercury	0.06019	0.8350	0.7189	79	0.8023	89	76-136	11	0-16	

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3545
Method: EPA 8082

Project: Burbank Airport / 9836002041

Page 4 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	MS/MSD Batch Number
B-DU1-ISM1-8	Sample	Solid	GC 66	03/27/17	03/28/17 03:41	170327S05
B-DU1-ISM1-8	Matrix Spike	Solid	GC 66	03/27/17	03/28/17 03:05	170327S05
B-DU1-ISM1-8	Matrix Spike Duplicate	Solid	GC 66	03/27/17	03/28/17 03:23	170327S05

<u>Parameter</u>	<u>Sample Conc.</u>	<u>Spike Added</u>	<u>MS Conc.</u>	<u>MS %Rec.</u>	<u>MSD Conc.</u>	<u>MSD %Rec.</u>	<u>%Rec. CL</u>	<u>RPD</u>	<u>RPD CL</u>	<u>Qualifiers</u>
Aroclor-1016	ND	100.0	82.50	82	81.00	81	50-135	2	0-20	
Aroclor-1260	ND	100.0	85.00	85	86.00	86	50-135	1	0-20	

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3545
Method: EPA 8270C SIM PAHs

Project: Burbank Airport / 9836002041

Page 5 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	MS/MSD Batch Number
17-03-1753-6	Sample	Solid	GC/MS EEE	03/24/17	03/27/17 12:36	170324S17
17-03-1753-6	Matrix Spike	Solid	GC/MS EEE	03/24/17	03/27/17 11:56	170324S17
17-03-1753-6	Matrix Spike Duplicate	Solid	GC/MS EEE	03/24/17	03/27/17 12:16	170324S17

Parameter	Sample Conc.	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Naphthalene	ND	0.1000	0.1123	112	0.1093	109	20-150	3	0-33	
2-Methylnaphthalene	ND	0.1000	0.1381	138	0.1341	134	29-137	3	0-31	3
1-Methylnaphthalene	0.01270	0.1000	0.1196	107	0.1189	106	34-136	1	0-29	
Acenaphthylene	ND	0.1000	0.1041	104	0.1049	105	29-131	1	0-32	
Acenaphthene	ND	0.1000	0.1076	108	0.1064	106	29-137	1	0-28	
Fluorene	ND	0.1000	0.1148	115	0.1155	115	36-132	1	0-27	
Phenanthrene	ND	0.1000	0.1057	106	0.1076	108	20-144	2	0-27	
Anthracene	ND	0.1000	0.1124	112	0.1092	109	26-134	3	0-27	
Fluoranthene	ND	0.1000	0.1066	107	0.1059	106	20-151	1	0-26	
Pyrene	ND	0.1000	0.1054	105	0.1084	108	20-150	3	0-32	
Benzo (a) Anthracene	ND	0.1000	0.1010	101	0.1040	104	24-150	3	0-24	
Chrysene	ND	0.1000	0.1012	101	0.09957	100	25-145	2	0-28	
Benzo (k) Fluoranthene	ND	0.1000	0.08966	90	0.08590	86	28-148	4	0-26	
Benzo (b) Fluoranthene	ND	0.1000	0.1132	113	0.1146	115	21-153	1	0-26	
Benzo (a) Pyrene	ND	0.1000	0.09968	100	0.09760	98	29-149	2	0-22	
Indeno (1,2,3-c,d) Pyrene	ND	0.1000	0.1004	100	0.09966	100	20-154	1	0-25	
Dibenz (a,h) Anthracene	ND	0.1000	0.1039	104	0.1048	105	20-132	1	0-26	
Benzo (g,h,i) Perylene	ND	0.1000	0.1071	107	0.1078	108	20-148	1	0-27	

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Sample Duplicate

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: N/A
Method: ASTM D-2216 (M)

Project: Burbank Airport / 9836002041

Page 1 of 1

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	Duplicate Batch Number
17-03-1556-25	Sample	Solid	N/A	03/23/17 00:00	03/23/17 21:00	H0323MOID3
17-03-1556-25	Sample Duplicate	Solid	N/A	03/23/17 00:00	03/23/17 21:00	H0323MOID3

<u>Parameter</u>	<u>Sample Conc.</u>	<u>DUP Conc.</u>	<u>RPD</u>	<u>RPD CL</u>	<u>Qualifiers</u>
Moisture	4.600	4.700	2	0-10	

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Andersen Environmental	Date Received:	03/21/17
5261 West Imperial Highway	Work Order:	17-03-1557
Los Angeles, CA 90045-6231	Preparation:	EPA 3550B
Project: Burbank Airport / 9836002041	Method:	EPA 8015B (M)

Page 1 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-14-353-13	LCS	Solid	GC 45	03/24/17	03/28/17 11:37	170324B12

<u>Parameter</u>	<u>Spike Added</u>	<u>Conc. Recovered</u>	<u>LCS %Rec.</u>	<u>%Rec. CL</u>	<u>Qualifiers</u>
TPH as Diesel	400.0	373.8	93	61-145	



RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3050B
Method: EPA 6010B

Project: Burbank Airport / 9836002041

Page 2 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
097-01-002-24537	LCS	Solid	ICP 7300	03/29/17	03/29/17 17:08	170329L05
<u>Parameter</u>	<u>Spike Added</u>	<u>Conc. Recovered</u>	<u>LCS %Rec.</u>	<u>%Rec. CL</u>	<u>ME CL</u>	<u>Qualifiers</u>
Antimony	25.00	20.99	84	80-120	73-127	
Arsenic	25.00	22.47	90	80-120	73-127	
Barium	25.00	25.38	102	80-120	73-127	
Beryllium	25.00	22.33	89	80-120	73-127	
Cadmium	25.00	25.54	102	80-120	73-127	
Chromium	25.00	24.93	100	80-120	73-127	
Cobalt	25.00	26.08	104	80-120	73-127	
Copper	25.00	24.17	97	80-120	73-127	
Lead	25.00	25.61	102	80-120	73-127	
Molybdenum	25.00	23.03	92	80-120	73-127	
Nickel	25.00	24.87	99	80-120	73-127	
Selenium	25.00	23.31	93	80-120	73-127	
Silver	12.50	11.95	96	80-120	73-127	
Thallium	25.00	24.20	97	80-120	73-127	
Vanadium	25.00	23.74	95	80-120	73-127	
Zinc	25.00	25.45	102	80-120	73-127	

Total number of LCS compounds: 16

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass

Return to Contents

Quality Control - LCS

Andersen Environmental
 5261 West Imperial Highway
 Los Angeles, CA 90045-6231

Date Received: 03/21/17
 Work Order: 17-03-1557
 Preparation: EPA 7471A Total
 Method: EPA 7471A

Project: Burbank Airport / 9836002041

Page 3 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-16-272-2910	LCS	Solid	Mercury 08	03/29/17	03/29/17 15:11	170329L01

<u>Parameter</u>	<u>Spike Added</u>	<u>Conc. Recovered</u>	<u>LCS %Rec.</u>	<u>%Rec. CL</u>	<u>Qualifiers</u>
Mercury	0.8350	0.7369	88	85-121	

Return to Contents 

RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Andersen Environmental	Date Received:	03/21/17
5261 West Imperial Highway	Work Order:	17-03-1557
Los Angeles, CA 90045-6231	Preparation:	EPA 3545
	Method:	EPA 8082
Project: Burbank Airport / 9836002041		Page 4 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-12-535-4123	LCS	Solid	GC 66	03/27/17	03/28/17 02:47	170327L05
<u>Parameter</u>		<u>Spike Added</u>	<u>Conc. Recovered</u>	<u>LCS %Rec.</u>	<u>%Rec. CL</u>	<u>Qualifiers</u>
Aroclor-1016		100.0	88.50	88	50-135	
Aroclor-1260		100.0	88.00	88	50-135	



Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Andersen Environmental
5261 West Imperial Highway
Los Angeles, CA 90045-6231

Date Received: 03/21/17
Work Order: 17-03-1557
Preparation: EPA 3545
Method: EPA 8270C SIM PAHs

Project: Burbank Airport / 9836002041

Page 5 of 5

Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number	
099-14-035-386	LCS	Solid	GC/MS EEE	03/24/17	03/27/17 11:35	170324L17	
<u>Parameter</u>		<u>Spike Added</u>	<u>Conc. Recovered</u>	<u>LCS %Rec.</u>	<u>%Rec. CL</u>	<u>ME CL</u>	<u>Qualifiers</u>
Naphthalene		0.1000	0.09899	99	51-129	38-142	
2-Methylnaphthalene		0.1000	0.1132	113	50-127	37-140	
1-Methylnaphthalene		0.1000	0.1008	101	54-132	41-145	
Acenaphthylene		0.1000	0.09919	99	50-123	38-135	
Acenaphthene		0.1000	0.1036	104	53-125	41-137	
Fluorene		0.1000	0.1041	104	55-127	43-139	
Phenanthrene		0.1000	0.1044	104	50-122	38-134	
Anthracene		0.1000	0.1059	106	50-132	36-146	
Fluoranthene		0.1000	0.1039	104	55-127	43-139	
Pyrene		0.1000	0.1021	102	50-134	36-148	
Benzo (a) Anthracene		0.1000	0.09779	98	50-133	36-147	
Chrysene		0.1000	0.1023	102	51-129	38-142	
Benzo (k) Fluoranthene		0.1000	0.09607	96	49-150	32-167	
Benzo (b) Fluoranthene		0.1000	0.1068	107	50-142	35-157	
Benzo (a) Pyrene		0.1000	0.09631	96	50-134	36-148	
Indeno (1,2,3-c,d) Pyrene		0.1000	0.09683	97	50-148	34-164	
Dibenz (a,h) Anthracene		0.1000	0.1004	100	50-133	36-147	
Benzo (g,h,i) Perylene		0.1000	0.1028	103	50-130	37-143	

Total number of LCS compounds: 18

Total number of ME compounds: 0

Total number of ME compounds allowed: 1

LCS ME CL validation result: Pass



Return to Contents

Sample Analysis Summary Report

Work Order: 17-03-1557

Page 1 of 1

<u>Method</u>	<u>Extraction</u>	<u>Chemist ID</u>	<u>Instrument</u>	<u>Analytical Location</u>
ASTM D-2216 (M)	N/A	1050	N/A	1
EPA 6010B	EPA 3050B	935	ICP 7300	1
EPA 7471A	EPA 7471A Total	868	Mercury 08	1
EPA 8015B (M)	EPA 3550B	972	GC 45	1
EPA 8082	EPA 3545	1028	GC 66	1
EPA 8270C SIM PAHs	EPA 3545	907	GC/MS EEE	1

Glossary of Terms and Qualifiers

Work Order: 17-03-1557

Page 1 of 1

<u>Qualifiers</u>	<u>Definition</u>
*	See applicable analysis comment.
<	Less than the indicated value.
>	Greater than the indicated value.
1	Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control.
4	The MS/MSD RPD was out of control due to suspected matrix interference.
5	The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference.
6	Surrogate recovery below the acceptance limit.
7	Surrogate recovery above the acceptance limit.
B	Analyte was present in the associated method blank.
BU	Sample analyzed after holding time expired.
BV	Sample received after holding time expired.
CI	See case narrative.
E	Concentration exceeds the calibration range.
ET	Sample was extracted past end of recommended max. holding time.
HD	The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
HDH	The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected).
HDL	The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected).
J	Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.
JA	Analyte positively identified but quantitation is an estimate.
ME	LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean).
ND	Parameter not detected at the indicated reporting limit.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.
SG	The sample extract was subjected to Silica Gel treatment prior to analysis.
X	% Recovery and/or RPD out-of-range.
Z	Analyte presence was not confirmed by second column or GC/MS analysis.
	Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.
	Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.
	A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

17-03-1557

LAB_SDG	LAB_SAMPLE_NUMBER	FIELD_SAMPLE	SAMPLING_DATE	SAMPLING_TIME	MATRIX	DEPTH	PROJECT_ID
17-03-1279	17-03-1279-2	B-DU1-S-06-8	20170316	12:25:00	Soil	8	Burbank Airport / 9836002041
17-03-1279	17-03-1279-5	B-DU1-S-SG-02-8S	20170316	07:25:00	Soil	8	Burbank Airport / 9836002041
17-03-1279	17-03-1279-8	B-DU1-S-SG-07-8S	20170316	12:55:00	Soil	8	Burbank Airport / 9836002041
17-03-1279	17-03-1279-11	B-DU1-S-SG-08-8S	20170316	15:10:00	Soil	8	Burbank Airport / 9836002041
17-03-1279	17-03-1279-14	B-DU1-S-SG-10-8S	20170316	09:22:00	Soil	8	Burbank Airport / 9836002041
17-03-1279	17-03-1279-17	B-DU1-S-SG-11-8S	20170316	10:47:00	Soil	8	Burbank Airport / 9836002041
17-03-1279	17-03-1279-20	B-DU1-S-04-8	20170316	14:25:00	Soil	8	Burbank Airport / 9836002041
17-03-1382	17-03-1382-2	B-DU1-S-SG-01-8S	20170317	13:00:00	Soil	8	Burbank Airport / 9836002041
17-03-1382	17-03-1382-5	B-DU1-S-SG-03-8S	20170317	12:20:00	Soil	8	Burbank Airport / 9836002041
17-03-1382	17-03-1382-8	B-DU1-S-SG-04-8S	20170317	10:30:00	Soil	8	Burbank Airport / 9836002041
17-03-1382	17-03-1382-11	B-DU1-S-03-8	20170317	13:35:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-2	B-DU1-S-01-8	20170320	14:20:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-5	B-DU1-S-02-8	20170320	11:00:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-8	B-DU1-S-05-8	20170320	10:20:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-11	B-DU1-S-07-8	20170320	09:45:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-14	B-DU1-S-08-8	20170320	13:05:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-17	B-DU1-S-09-8	20170320	09:21:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-20	B-DU1-S-SG-05-8S	20170320	08:20:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-23	B-DU1-S-SG-06-8S	20170320	12:42:00	Soil	8	Burbank Airport / 9836002041
17-03-1524	17-03-1524-26	B-DU1-S-SG-09-8S	20170320	07:40:00	Soil	8	Burbank Airport / 9836002041

B-DU1-ISM1-8
 B-DU1-ISM2-8
 B-DU1-ISM3-8

Number	Sample ID	Lab ID	Type		Matrix			Preservative				Sampling Information		
			Grab	Discrete	Water	Soil	Vapor	Other	Cold (4° C)	HNO3	NaHSO4	HCl	Date	Time
1	B-D41-S-06-3	1	X		X								3/19/17	02:21
2	-8	2											3/19/17	52:21
3	-15	3											3/19/17	12:30
4	B-D41-S-SG-02-35	4											3/19/17	07:02
5	-85	5											3/19/17	07:02
6	-155	6											3/19/17	07:02
7	B-D41-S-SG-07-35	7											3/19/17	12:50
8	-85	8											3/19/17	12:55
9	755	9											3/19/17	13:05
10	B-D41-S-SG-08-35	10											3/19/17	01:50
11	-85	11											3/19/17	01:51
12	-155	12											3/19/17	01:51
13	B-D41-S-SG-10-35	13											3/19/17	01:51
14	-85	14											3/19/17	01:51
15	-155	15											3/19/17	01:51
16	B-D41-S-SG-11-35	16											3/19/17	01:51
17	-85	17											3/19/17	01:51
18	-155	18											3/19/17	01:51
19	B-D41-S-04-3	19											3/19/17	01:51
20	-8	20											3/19/17	01:51

Method	Container	Turnaround Time																	
			VOCs, EPA Method 826B	PAHs, EPA Method 8270C SIM	Metals, EPA Method 6010B/7471A	Lead, EPA Method 6010B	Arsenic, EPA Method 6010B	STLC Lead EPA Method	TCLP Lead EPA Method	PCBs, EPA Method 8082	TPH full chain, EPA Method 8015M	TPHd, TPHmo, EPA Method 8015M	Composite	Hold	4 or 8-ounce Glass	250-ml Poly Bottle	EZ Draw (EPA 5035)	Acetate Liner	1 Amber Bottle

Reinquished by	Date	Time	Received by	Date	Time	Remarks
<i>[Signature]</i>	3/16/17	1605	<i>[Signature]</i>	3/16/17	1705	Sample condition (circle): Chilled Intact
<i>[Signature]</i>	3/16/17	1835	<i>[Signature]</i>	3/16/17	1835	Take sample from center of liner.
						Hold all samples until call for DUA
						Received for 15H composites (Separate case)
						Hold 15S samples as needed after 3+8' analysis



Return to Contents

Analytical Laboratory: Calscience
 Project Name: Burbank Airport
 Project Address: 2627 N. Hollywood Way, Burbank
 Project Manager: B. Martasin
 Sampled by: G. Baader
 Phone/Email: 310-854-6300, brian_martasin@efiglobal.com
 Phone/Email: Robert Cheung, 510-529-5948, RCheung@Geosyntec.com

Method	Container	Turnaround Time	Sampling Information	
			Date	Time
OCPS, EPA Method 8081A			3/16/17	4:35
PAHs, EPA Method 8270C SIM				
Metals, EPA Method 6010B/7471A				
Lead, EPA Method 6010B				
Arsenic, EPA Method 6010B				
STLC Lead EPA Method				
TCLP Lead EPA Method				
PCBs, EPA Method 8082				
TPH full chain, EPA Method 8015M				
TPH, TPHm, EPA Method 8015M				
Composite				
Hold				
4 or 8-ounce Glass				
250-ml Poly Bottle				
EZ Draw (EPA 5035)				
Acetate Liner				
1-l Amber Bottle				
24 hours				
48 hours				
Normal				

Number	Sample ID	Lab ID	Type		Matrix		Preservative		Sampling Information											
			Grab	Discrete	Water	Soil	Vapor	Other	Cold (4° C)	HNO3	NaHSO4	HCl	Date	Time						
21	B-DU-5-04-15	21	X		X															
22																				
23																				

Relinquished by: *[Signature]*
 Date: 3/16/17 18:35
 Received by: *[Signature]*
 Date: 3/16/17 18:35
 Remarks: Sample condition (circle): Chilled Intact
 Take sample from center of liner unless otherwise indicated.
 Hold all samples until all for DU received, analysis on separate COC.



SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 1

CLIENT: Andersen

DATE: 03 / 16 / 2017

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)
 Thermometer ID: SC3B (CF: 0.0°C); Temperature (w/o CF): 2.0 °C (w/ CF): 2.0 °C; Blank Sample
 Sample(s) outside temperature criteria (PM/APM contacted by: _____)
 Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling
 Sample(s) received at ambient temperature; placed on ice for transport by courier
 Ambient Temperature: Air Filter

Checked by: 1091

CUSTODY SEAL:

Cooler Present and Intact Present but Not Intact Not Present N/A Checked by: 1091
 Sample(s) Present and Intact Present but Not Intact Not Present N/A Checked by: 778

SAMPLE CONDITION:	Yes	No	N/A
Chain-of-Custody (COC) document(s) received with samples	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
COC document(s) received complete	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers			
<input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time			
Sampler's name indicated on COC	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Sample container label(s) consistent with COC	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Sample container(s) intact and in good condition	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Proper containers for analyses requested	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Sufficient volume/mass for analyses requested	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Samples received within holding time	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Aqueous samples for certain analyses received within 15-minute holding time			
<input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Proper preservation chemical(s) noted on COC and/or sample container	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Unpreserved aqueous sample(s) received for certain analyses			
<input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals			
Container(s) for certain analysis free of headspace	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500)			
<input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)			
Tedlar™ bag(s) free of condensation	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

CONTAINER TYPE: (Trip Blank Lot Number: _____)

Aqueous: VOA VOA_h VOA_{na2} 100PJ 100PJ_{na2} 125AGB 125AGB_h 125AGB_p 125PB
 125PB_{znna} 250AGB 250CGB 250CGB_s 250PB 250PB_n 500AGB 500AGJ 500AGJ_s
 500PB 1AGB 1AGB_{na2} 1AGB_s 1PB 1PB_{na} _____ _____ _____ _____

Solid: 4ozCGJ 8ozCGJ 16ozCGJ Sleeve (P) EnCores® (____) TerraCores® (____) _____

Air: Tedlar™ Canister Sorbent Tube PUF _____ **Other Matrix** (____): _____ _____

Container: **A** = Amber, **B** = Bottle, **C** = Clear, **E** = Envelope, **G** = Glass, **J** = Jar, **P** = Plastic, and **Z** = Ziploc/Resealable Bag
 Preservative: **b** = buffered, **f** = filtered, **h** = HCl, **n** = HNO₃, **na** = NaOH, **na₂** = Na₂S₂O₃, **p** = H₃PO₄, **s** = H₂SO₄, **u** = ultra-pure, **x** = Na₂SO₃+NaHSO₄.H₂O, **znna** = Zn (CH₃CO₂)₂ + NaOH

Labeled/Checked by: 778
 Reviewed by: 679

Return to Contents

Analytical Laboratory: Calscience Job # 9836002041

Project Name: Burbank Airport

Project Address: 2627 N. Hollywood Way, Burbank

Project Manager: B. Martasin

Sampled by: G. Baader

Phone/Email: 310-854-6300, brian_martasin@efiglobal.com

Phone/Email: Robert Cheung, 510-529-5948, RCheung@Geosyntec.com

Number	Sample ID	Lab ID	Type		Matrix			Preservative			Sampling Information		
			Grab	Discrete	Water	Soil	Vapor	Other	Cold (4° C)	HNO3	NaHSO4	HCl	Date
1	B-DUI-556-01-35		X	X	X	X	X	X	X	X	X	3/17/17	12:55
2	-85		X	X	X	X	X	X	X	X	X	3/17/17	13:00
3	-155		X	X	X	X	X	X	X	X	X	3/17/17	13:10
4	B-DUI-556-03-35		X	X	X	X	X	X	X	X	X	3/17/17	12:15
5	-85		X	X	X	X	X	X	X	X	X	3/17/17	12:20
6	-155		X	X	X	X	X	X	X	X	X	3/17/17	13:35
7	B-DUI-556-04-35		X	X	X	X	X	X	X	X	X	3/17/17	10:25
8	-85		X	X	X	X	X	X	X	X	X	3/17/17	10:30
9	-155		X	X	X	X	X	X	X	X	X	3/17/17	10:50
10	B-DUI-5-03-33		X	X	X	X	X	X	X	X	X	3/17/17	13:30
11	-8		X	X	X	X	X	X	X	X	X	3/17/17	13:35
12	-15		X	X	X	X	X	X	X	X	X	3/17/17	13:50
13													
14													
15													
16													
17													
18													
19													
20													

Relinquished by: [Signature]

Date: 3/17/17 Time: 1520 Received by: [Signature] Date: 3/17/17 Time: 1825

Date: 3/17/17 Time: 1825 Received by: [Signature] Date: 03/17/17 Time: 1825

Method		Container		Turnaround Time
Method	Container	Method	Container	
VOCs, EPA Method 8260B	PAHs, EPA Method 8270C SIM	Metals, EPA Method 6010B/7471A	Lead, EPA Method 6010B	17-03-1382 4- or 8-ounce Glass 250-ml Poly Bottle EZ Draw (EPA 5035) 1- Amber Bottle 48 hours Normal
Lead, EPA Method 6010B	Arsenic, EPA Method 6010B	STLC Lead EPA Method	TCLP Lead EPA Method	
PCBs, EPA Method 8082	TPH full chain, EPA Method 8015M	TPHD, TPHmo, EPA Method 8015M	Composite	
Hold	4- or 8-ounce Glass	250-ml Poly Bottle	EZ Draw (EPA 5035)	
1- Amber Bottle	48 hours	Normal		

Sample condition (circle): Chilled Intact
 bold in bold all B-DUI received for SFH
 Composites. Take sample from one of B-DUI
 unless otherwise indicated. Add 15' samples
 into C B-DUI & Sample away 8' 15' 0's needed.

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 1

CLIENT: Andersen

DATE: 03 / 17 / 2017

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)
 Thermometer ID: SC3B (CF: 0.0°C); Temperature (w/o CF): 1.8 °C (w/ CF): 1.8 °C; Blank Sample
 Sample(s) outside temperature criteria (PM/APM contacted by: _____)
 Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling
 Sample(s) received at ambient temperature; placed on ice for transport by courier
 Ambient Temperature: Air Filter

Checked by: 1091

CUSTODY SEAL:

Cooler Present and Intact Present but Not Intact Not Present N/A
 Sample(s) Present and Intact Present but Not Intact Not Present N/A

Checked by: 1091

Checked by: 1053

SAMPLE CONDITION:

	Yes	No	N/A
Chain-of-Custody (COC) document(s) received with samples	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
COC document(s) received complete	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/> Sampling date <input type="checkbox"/> Sampling time <input type="checkbox"/> Matrix <input type="checkbox"/> Number of containers			
<input type="checkbox"/> No analysis requested <input type="checkbox"/> Not relinquished <input type="checkbox"/> No relinquished date <input type="checkbox"/> No relinquished time			
Sampler's name indicated on COC	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Sample container label(s) consistent with COC	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Sample container(s) intact and in good condition	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Proper containers for analyses requested	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Sufficient volume/mass for analyses requested	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Samples received within holding time	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Aqueous samples for certain analyses received within 15-minute holding time			
<input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfide <input type="checkbox"/> Dissolved Oxygen	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Proper preservation chemical(s) noted on COC and/or sample container	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Unpreserved aqueous sample(s) received for certain analyses			
<input type="checkbox"/> Volatile Organics <input type="checkbox"/> Total Metals <input type="checkbox"/> Dissolved Metals			
Container(s) for certain analysis free of headspace	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/> Volatile Organics <input type="checkbox"/> Dissolved Gases (RSK-175) <input type="checkbox"/> Dissolved Oxygen (SM 4500)			
<input type="checkbox"/> Carbon Dioxide (SM 4500) <input type="checkbox"/> Ferrous Iron (SM 3500) <input type="checkbox"/> Hydrogen Sulfide (Hach)			
Tedlar™ bag(s) free of condensation	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

CONTAINER TYPE:

(Trip Blank Lot Number: _____)

Aqueous: VOA VOA_h VOA_{na2} 100PJ 100PJ_{na2} 125AGB 125AGB_h 125AGB_p 125PB
 125PB_{z_{na}} 250AGB 250CGB 250CGB_s 250PB 250PB_n 500AGB 500AGJ 500AGJ_s
 500PB 1AGB 1AGB_{na2} 1AGB_s 1PB 1PB_{na} _____ _____ _____
Solid: 4ozCGJ 8ozCGJ 16ozCGJ Sleeve (1) EnCores® (_____) TerraCores® (_____) _____
Air: Tedlar™ Canister Sorbent Tube PUF _____ **Other Matrix** (_____): _____ _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag

Preservative: b = buffered, f = filtered, h = HCl, n = HNO₃, na = NaOH, na₂ = Na₂S₂O₃, p = H₃PO₄, Labeled/Checked by: 1053

s = H₂SO₄, u = ultra-pure, x = Na₂SO₃+NaHSO₄.H₂O, z_{na} = Zn (CH₃CO₂)₂ + NaOH Reviewed by: 681

Return to Contents

SAMPLE RECEIPT CHECKLIST

COOLER 1 OF 1

CLIENT: EFI

DATE: 03 / 31 / 2017

TEMPERATURE: (Criteria: 0.0°C – 6.0°C, not frozen except sediment/tissue)
Thermometer ID: SC3B (CF: 0.0°C); Temperature (w/o CF): 3.8 °C (w/ CF): 3.8 °C; [X] Blank [] Sample
[] Sample(s) outside temperature criteria (PM/APM contacted by: _____)
[] Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling
[] Sample(s) received at ambient temperature; placed on ice for transport by courier
Ambient Temperature: [] Air [] Filter
Checked by: 678

CUSTODY SEAL:
Cooler [] Present and Intact [] Present but Not Intact [X] Not Present [] N/A Checked by: 678
Sample(s) [] Present and Intact [] Present but Not Intact [X] Not Present [] N/A Checked by: 1110

Table with columns: SAMPLE CONDITION, Yes, No, N/A. Rows include Chain-of-Custody (COC) document(s) received with samples, COC document(s) received complete, Sampler's name indicated on COC, Sample container label(s) consistent with COC, Sample container(s) intact and in good condition, Proper containers for analyses requested, Sufficient volume/mass for analyses requested, Samples received within holding time, Aqueous samples for certain analyses received within 15-minute holding time, Proper preservation chemical(s) noted on COC and/or sample container, Unpreserved aqueous sample(s) received for certain analyses, Container(s) for certain analysis free of headspace, Tedlar™ bag(s) free of condensation.

CONTAINER TYPE: (Trip Blank Lot Number: _____)

Aqueous: [] VOA [] VOA h [] VOAna2 [] 100PJ [] 100PJna2 [] 125AGB [] 125AGBh [] 125AGBp [] 125PB
[] 125PBz nna [] 250AGB [] 250CGB [] 250CGBs [] 250PB [] 250PBn [] 500AGB [] 500AGJ [] 500AGJs
[] 500PB [] 1AGB [] 1AGBna2 [] 1AGBs [] 1PB [] 1PBna [] _____ [] _____ [] _____ [] _____
Solid: [] 4ozCGJ [] 8ozCGJ [] 16ozCGJ [X] Sleeve (P) [] EnCores® () [] TerraCores® () [] _____
Air: [] Tedlar™ [] Canister [] Sorbent Tube [] PUF [] _____ Other Matrix (): [] _____ [] _____

Container: A = Amber, B = Bottle, C = Clear, E = Envelope, G = Glass, J = Jar, P = Plastic, and Z = Ziploc/Resealable Bag
Preservative: b = buffered, f = filtered, h = HCl, n = HNO3, na = NaOH, na2 = Na2S2O3, p = H3PO4, Labeled/Checked by: 1110
s = H2SO4, u = ultra-pure, x = Na2SO3+NaHSO4.H2O, z nna = Zn (CH3CO2)2 + NaOH Reviewed by: 679

Case Narrative

Client Project Name: Burbank Airport / 9836002041

Work Order Number: 17-03-1557

CONDITION UPON RECEIPT:

Eurofins Calscience, Inc. received (20) solid samples on March 16, 17, and 21, 2017. A total of (20) containers were received in good condition and at temperatures of 1.8°C, 2.0°C, and 3.8°C, which were within the recommended temperature of 0°C – 6°C.

Client Sample ID	Lab Sample ID	Date & Time Sampled	Date & Time Received
B-DU1-S-06-8	17-03-1557-1	03/16/17 12:25	03/21/17 15:37
B-DU1-S-SG-02-8S	17-03-1557-2	03/16/17 07:25	03/21/17 15:37
B-DU1-S-SG-07-8S	17-03-1557-3	03/16/17 12:55	03/21/17 15:37
B-DU1-S-SG-08-8S	17-03-1557-4	03/16/17 15:10	03/21/17 15:37
B-DU1-S-SG-10-8S	17-03-1557-5	03/16/17 09:22	03/21/17 15:37
B-DU1-S-SG-11-8S	17-03-1557-6	03/16/17 10:47	03/21/17 15:37
B-DU1-S-04-8	17-03-1557-7	03/16/17 14:25	03/21/17 15:37
B-DU1-S-SG-01-8S	17-03-1557-8	03/17/17 13:00	03/21/17 15:37
B-DU1-S-SG-03-8S	17-03-1557-9	03/17/17 12:20	03/21/17 15:37
B-DU1-S-SG-04-8S	17-03-1557-10	03/17/17 10:30	03/21/17 15:37
B-DU1-S-03-8	17-03-1557-11	03/17/17 13:35	03/21/17 15:37
B-DU1-S-01-8	17-03-1557-12	03/20/17 14:20	03/21/17 15:37
B-DU1-S-02-8	17-03-1557-13	03/20/17 11:00	03/21/17 15:37
B-DU1-S-05-8	17-03-1557-14	03/20/17 10:20	03/21/17 15:37
B-DU1-S-07-8	17-03-1557-15	03/20/17 09:45	03/21/17 15:37
B-DU1-S-08-8	17-03-1557-16	03/20/17 13:05	03/21/17 15:37
B-DU1-S-09-8	17-03-1557-17	03/20/17 09:21	03/21/17 15:37
B-DU1-S-SG-05-8S	17-03-1557-18	03/20/17 08:20	03/21/17 15:37
B-DU1-S-SG-06-8S	17-03-1557-19	03/20/17 12:42	03/21/17 15:37
B-DU1-S-SG-09-8S	17-03-1557-20	03/20/17 07:40	03/21/17 15:37
COMPOSITE	17-03-1557-21	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM1-8	17-03-1557-22	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM2-8	17-03-1557-23	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM3-8	17-03-1557-24	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM1-8	17-03-1557-25	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM2-8	17-03-1557-26	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM3-8	17-03-1557-27	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM1-8	17-03-1557-28	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM2-8	17-03-1557-29	03/16/17 00:00	03/21/17 15:37
B-DU1-ISM3-8	17-03-1557-30	03/16/17 00:00	03/21/17 15:37

Case Narrative

Client Project Name: Burbank Airport / 9836002041

Work Order Number: 17-03-1557

DATA SUMMARY:

As per the chain of custody (COC), the samples were analyzed using one or more of the following methodologies:

- ASTM D-2216 (M) Moisture Content (Solid)
- EPA 6010B Title 22 Metals (Solid)
- EPA 7471A Mercury (Solid)
- EPA 8015B (M) Diesel and Motor Oil Ranges (Solid)
- EPA 8082 PCB Aroclors (Solid)
- EPA 8270C SIM PAHs (Solid)

The samples were analyzed within the suggested EPA holding time for the requested methods unless otherwise noted.

Sample results were reported in the RL format.

The sample data is reported in dry weight. The instrument printouts do not reflect the correction for dry weight.

Any dilutions made to the sample(s) and/or QC will be noted in the following narrative. Reporting limits have been adjusted accordingly.

Manual integrations made to the data will be noted in the following narrative. The before and amended chromatograms have been included in the data package.

All sample and analytical QC are within acceptance criteria unless otherwise noted.

ASTM D-2216 (M) Moisture Content (Solid):

Samples -25 through -27 were analyzed for % Moisture by ASTM D-2216 (M). The samples were prepared and analyzed on 03/23/17 in batch #s H0323MOIB3 / H0323MOID3.

Balance Calibration/Verification:

All values were within acceptance criteria.

Sample and QC:

A sample from a different work order was used as the sample duplicate for quality control. The method blank was non-detect and the duplicate analysis was within acceptance criteria.

Case Narrative

Client Project Name: Burbank Airport / 9836002041

Work Order Number: 17-03-1557

EPA 6010B Title 22 Metals (Solid):

Samples -28 through -30 were analyzed for Metals by EPA 6010B. The samples were prepared and analyzed on 03/29/17 in batch #s 170329L05 / 170329S05 on ICP 7300.

Initial Calibration, Initial Calibration Verification, and Initial Calibration Blank:

All values were within acceptance criteria.

Continuing Calibration Verification and Continuing Calibration Blank:

All values were within acceptance criteria.

ICS A/AB:

All values were within acceptance criteria.

Sample and QC:

The method blank was non-detect and the LCS was within acceptance criteria.

A non-client sample was used for the MS/MSD; refer to the MS/MSD summary form for further information.

EPA 7471A Mercury (Solid):

Samples -25 through -27 were analyzed for Mercury by EPA 7471A. The samples were prepared and analyzed on 03/29/17 in batch #s 170329L01 / 170329S01 on Mercury 08.

Initial Calibration, Initial Calibration Verification, and Initial Calibration Blank:

All values were within acceptance criteria.

Continuing Calibration Verification and Continuing Calibration Blank:

All values were within acceptance criteria.

Sample and QC:

The method blank was non-detect and the LCS was within acceptance criteria.

A non-client sample was used for the MS/MSD; refer to the MS/MSD summary form for further information.

Case Narrative

Client Project Name: Burbank Airport / 9836002041

Work Order Number: 17-03-1557

EPA 8015B (M) Diesel and Motor Oil Ranges (Solid):

Samples -25 through -27 were analyzed for Diesel and Motor Oil Ranges by EPA 8015B (M). The samples were prepared on 03/24/17 and analyzed on 03/28/17 in batch #s 170324B12 / 170324S12 on GC 45.

Initial Calibration and Initial Calibration Verification:

The initial calibration was performed on 12/02/16 on GC 45. The ICAL was within the 20% RSD acceptance criteria and the ICV was within the 30% D acceptance criteria.

For the Diesel and Motor Oil Ranges, Diesel is used for the initial calibration and spiking standards. The surrogate recoveries for the samples and QC were calculated from the 5-point surrogate curve analyzed with the Diesel ICAL.

Continuing Calibration Verification:

All values were within the 20% D acceptance criteria.

Sample and QC:

Sample -27 was used for the MS/MSD. The method blank was non-detect; the LCS, MS/MSD and all surrogate recoveries were within acceptance criteria.

EPA 8082 PCB Aroclors (Solid):

Samples -28 through -30 were analyzed for Polychlorinated Biphenyls Aroclors by EPA 8082. The samples were prepared on 03/27/17 and analyzed on 03/28/17 in batch #s 170327L05 / 170327S05 on GC 66.

Initial Calibration and Initial Calibration Verification:

The initial calibration was performed on 02/22/17 on GC 66. The ICAL was within the 20% RSD acceptance criteria and the ICV was within the 15% D acceptance criteria for Aroclors 1016 and 1260. Single point response factors were generated for all other Aroclors.

Continuing Calibration Verification:

All values were within the 15% D acceptance criteria.

Case Narrative

Client Project Name: Burbank Airport / 9836002041

Work Order Number: 17-03-1557

Sample and QC:

Sample -28 was used for the MS/MSD. The method blank was non-detect; the LCS, MS/MSD and all surrogate recoveries were within acceptance criteria.

EPA 8270C SIM PAHs (Solid):

Samples -25 through -27 were analyzed for Polynuclear Aromatic Hydrocarbons by EPA 8270C SIM. The samples were prepared on 03/24/17 and analyzed on 03/27/17 in batch #s 170324L17 / 170324S17 on GC/MS EEE.

Initial Calibration and Initial Calibration Verification:

The initial calibration was performed on 03/13/17 on GC/MS EEE. The ICAL was within the 15% RSD acceptance criteria and the ICV was within the 20% D acceptance criteria.

Continuing Calibration Verification:

All values were within the 20% D acceptance criteria.

Tuning Standards:

All instrument tuning standards (DFTPP) were within acceptance criteria.

Sample and QC:

The method blank was non-detect; the LCS, all surrogate and internal standard recoveries were within acceptance criteria.

A non-client sample was used for the MS/MSD; refer to the MS/MSD summary form for further information.

Manual integration was performed on one or more of the samples to correct the peak and/or baseline integration.

ASTM D-2216 (M)
Moisture Content
(Solid)

RAW DATA

RAW DATA SHEET
FOR METHOD: ASTM D-2216 (M)

WORK ORDER: 17-03-1557
INSTRUMENT: N/A
EXTRACTION: N/A
D/T EXTRACTED: 2017-03-23 00:00
DATA FILE: NONE

ANALYZED BY: 1,050
D/T ANALYZED: 2017-03-23 21:00
REVIEWED BY: 1,050
D/T REVIEWED: 2017-03-27 09:57

25 **CLIENT SAMPLE NUMBER: B-DU1-ISM1-8**

LCS/MB BATCH: H0323MOIB3 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 1.00 g
MS/MSD BATCH: H0323MOID3 **FINAL VOLUME / WEIGHT:** DEFAULT: 1.00 ml
UNITS: % **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Moisture	3.00	1.00	3.00	0.10	

**RAW DATA SHEET
FOR METHOD: ASTM D-2216 (M)**

WORK ORDER: 17-03-1557
INSTRUMENT: N/A
EXTRACTION: N/A
D/T EXTRACTED: 2017-03-23 00:00
DATA FILE: NONE

ANALYZED BY: 1,050
D/T ANALYZED: 2017-03-23 21:00
REVIEWED BY: 1,050
D/T REVIEWED: 2017-03-27 09:57

26 **CLIENT SAMPLE NUMBER:** B-DU1-ISM2-8

LCS/MB BATCH: H0323MOIB3 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 1.00 g
MS/MSD BATCH: H0323MOID3 **FINAL VOLUME / WEIGHT:** DEFAULT: 1.00 ml
UNITS: % **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Moisture	2.80	1.00	2.80	0.10	



**RAW DATA SHEET
FOR METHOD: ASTM D-2216 (M)**

WORK ORDER: 17-03-1557
INSTRUMENT: N/A
EXTRACTION: N/A
D/T EXTRACTED: 2017-03-23 00:00
DATA FILE: NONE

ANALYZED BY: 1,050
D/T ANALYZED: 2017-03-23 21:00
REVIEWED BY: 1,050
D/T REVIEWED: 2017-03-27 09:57

27 **CLIENT SAMPLE NUMBER:** B-DU1-ISM3-8

<u>LCS/MB BATCH:</u> H0323MOIB3	<u>SAMPLE VOLUME / WEIGHT:</u> DEFAULT: 1.00 g
<u>MS/MSD BATCH:</u> H0323MOID3	<u>FINAL VOLUME / WEIGHT:</u> DEFAULT: 1.00 ml
<u>UNITS:</u> %	<u>ADJUSTMENT RATIO TO PF:</u> 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Moisture	3.00	1.00	3.00	0.10	

**METHOD BLANK ASSOCIATION SUMMARY
FOR METHOD: ASTM D-2216 (M)**

MB SAMPLE ID: 099-05-014-6770
MB BATCH ID: H0323MOIB3
INSTRUMENT: N/A
EXTRACTION: N/A
D/T EXTRACTED: 2017-03-23 00:00

ANALYZED BY: 1,050
D/T ANALYZED: 2017-03-23 21:00
REVIEWED BY: 1,050
D/T REVIEWED: 2017-03-27 09:57
MATRIX: Soil

DATA FILE: NONE

CLIENT WORK ORDER: 17-03-1557

<u>S#</u>	<u>RUN TYPE</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
25	B-DU1-ISM1-8		2017-03-23 21:00	NONE
26	B-DU1-ISM2-8		2017-03-23 21:00	NONE
27	B-DU1-ISM3-8		2017-03-23 21:00	NONE

**RAW DATA SHEET
FOR METHOD: ASTM D-2216 (M)**

WORK ORDER: 099-05-014
INSTRUMENT: N/A
EXTRACTION: N/A
D/T EXTRACTED: 2017-03-23 00:00
DATA FILE: NONE

ANALYZED BY: 1,050
D/T ANALYZED: 2017-03-23 21:00
REVIEWED BY: 1,050
D/T REVIEWED: 2017-03-27 09:57

MB **CLIENT SAMPLE NUMBER:** Method Blank

LCS/MB BATCH: H0323MOIB3 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 1.00 g
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 1.00 ml
UNITS: % **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Moisture	0.000	1.00	ND	0.10	



DUPLICATE REPORT FOR METHOD: ASTM D-2216 (M)

DUP SAMPLE ID: 17-03-1556-25
DUP BATCH: H0323MOID3
INSTRUMENTS:
SAMPLE: N/A
DUP SAMPLE: N/A

EXTRACTION: N/A
D/T EXTRACTED:
SAMPLE: 2017-03-23 00:00
DUP SAMPLE: 2017-03-23 00:00

ANALYZED BY: 1,050
D/T ANALYZED:
SAMPLE: 2017-03-23 21:00
DUP SAMPLE: 2017-03-23 21:00
REVIEWED BY: 1,050
D/T REVIEWED: 2017-03-27 09:57

<u>COMPOUND</u>	<u>SAMPLE CONC</u>	<u>DUP CONC</u>	<u>% RPD</u>	<u>CONTROL LIMIT</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Moisture	4.600	4.700	2	0-10	PASS	

Data Files:

<u>TYPE</u>	<u>DATA FILE</u>	<u>DATA FILE PATH</u>
SDP	NONE	

Moisture Content Raw Data Logbook

METHOD ASTM D2216(M)	MATRIX <input checked="" type="checkbox"/> Solid <input type="checkbox"/> Other	DATE Preparation: 3/23/17 Analysis: ↓	ANALYST(S) 1050/1009 ↓	BATCH NUMBER MB: H0323 MOI B3 Sample Duplicate: ↓ D3
-------------------------	---	---	------------------------------	--

CEL ID #	DISH	DRY DISH	MOISTURE	% MOISTURE	RL (%)	COMMENTS	ALUMINUM DISH SUPPLY ID
MB	1.88	1.88	0.00	0.00	0.16		608-66-15

CEL ID #	% MOISTURE	RPD	CONTROL LIMIT	OVEN ID	THERMOMETER ID / CF (°C)	OVEN TEMP (°C)	BATCH PREP TIME (h:m)
Sample 17-03-1556-25-A	4.6	2	0 - 10	<input checked="" type="checkbox"/> IO 07	<input checked="" type="checkbox"/> TSS 107A CF: -2	Start: 103	Start: 1300
Duplicate ↓	4.7			<input type="checkbox"/> IO 08	<input type="checkbox"/> TSS 108A CF: _____	End: 103	End: 2100

Instructions: CEL ID consists of Work Order Number and Container ID

CEL ID #	DISH	DISH + SAMPLE (as-received)	SAMPLE (as-received)	DISH + SAMPLE (oven-dried)	MOISTURE	% MOISTURE		BALANCE ID #	RL (%)	QUAL	COMMENTS
						WET-BASED	DRY-BASED				
17-03-1556-25A	1.91	12.14	10.23	11.66	0.48	4.7		63	0.10		
↓	1.86	12.05	10.19	11.58	0.47	4.6					
↓	1.87	12.28	10.41	11.82	0.46	4.4					
↓	1.87	12.38	10.51	11.89	0.49	4.7					
17-03-1557-25A	1.89	12.48	10.59	12.16	0.32	3.0					
↓	1.88	12.05	10.17	11.77	0.28	2.8					
↓	1.89	12.32	10.43	12.01	0.31	3.0					
17-03-1558-25A	1.86	12.35	10.49	12.04	0.31	3.0					
↓	1.88	12.44	10.56	12.12	0.32	3.0					
↓	1.88	12.06	10.18	11.78	0.28	2.8					
17-03-1618-1A	1.89	12.47	10.58	9.88	2.59	24.9					

BATCH TIME
 Time (24 Hrs): 2100
 InMets: 1009/1050

BALANCE CALIBRATION CHECK LOG

Eurofins Calscience

Date performed: 03/23/17 Initials: III

ID	Class 2 Weight (g)	Reading (g)	Acceptance Range	Pass? (circle one)	Comment (If not passed, note removal or corrective action)
25	1	1.00	0.98 - 1.02	⊗ N	IO Lab
	100	100.01	98.00 - 102.00	⊗ N	
	500	500.03	498.00 - 502.00	⊗ N	
62	0.002	.0018	0.00180 - 0.00220	⊗ N	IO Lab
	1	.9993	0.99900 - 1.00100	⊗ N	
	100	99.9964	99.90000 - 100.10000	⊗ N	
26	1	1.01	0.98 - 1.02	⊗ N	IO Lab
	100	99.99	98.00 - 102.00	⊗ N	
55	1	1.00	0.98 - 1.02	⊗ N	IO Lab
	100	99.97	98.00 - 102.00	⊗ N	
	500	499.89	498.00 - 502.00	⊗ N	
11	1	1.00	0.98 - 1.02	⊗ N	IO Lab
	100	99.99	98.00 - 102.00	⊗ N	
66	0.002	.0018	0.00180 - 0.00220	⊗ N	Metals
	1	.9995	0.99900 - 1.00100	⊗ N	
	100	100.0001	99.90000 - 100.10000	⊗ N	
53	0.1	.10	0.09 - 0.11	⊗ N	Extractions
	1	1.07	0.98 - 1.02	⊗ N	
	100	100.06	98.00 - 102.00	⊗ N	
	500	499.99	498 - 502	⊗ N	
70	1	.99	0.98 - 1.02	⊗ N	Extractions
	100	99.83	98.00 - 102.00	⊗ N	
	500	499.17	498.00 - 502.00	⊗ N	
57	100	100.0	98.0-102.0	⊗ N	Extractions
	1000	1000.0	998.0-1002.0	⊗ N	
	2000	2000.0	1998.0-2002.0	⊗ N	
52	0.002	.0018	0.0018 - 0.0022	⊗ N	Extractions
	1	9993	0.9990 - 1.0010	⊗ N	
	100	99.9961	99.9000 - 100.1000	⊗ N	
71	0.002	.0021	0.0018 - 0.0022	⊗ N	BOD Room
	1	.9994	0.9990 - 1.0010	⊗ N	
	100	99.9966	99.9000 - 100.1000	⊗ N	
63	0.1	.10	0.09 - 0.11	⊗ N	BOD Room
	100	100.00	98.00 - 102.00	⊗ N	
64	1	1.00	0.98 - 1.02	⊗ N	Metals Clean Room
	10	10.00	9.8 - 10.2	⊗ N	
	100	100.01	98.00 - 102.00	⊗ N	
72	0.002	.0020	0.0018 - 0.0022	⊗ N	Oil & Grease Room
	1	.9994	0.9990 - 1.0010	⊗ N	
	100	100.0003	99.9000 - 100.1000	⊗ N	
30	1	1.00	0.98 - 1.02	⊗ N	Oil & Grease Room
	100	100.00	98.00 - 102.00	⊗ N	

EPA 8015B (M)
Diesel + Motor Oil

RAW DATA

EPA 8015B (M) Diesel + Motor Oil

INITIAL CALIBRATION

INITIAL CALIBRATION QUALITY CONTROL SUMMARY FOR METHOD: EPA 8015B (M)

ICAL WORK ORDER: 099-14-354-18-5901
ICAL BATCH ID: 1612021007
INSTRUMENT: GC 45

ANALYZED BY: 972
ICAL D/T ANALYZED: 2016-12-02 19:43
REVIEWED BY: 27
D/T REVIEWED: 2017-02-16 16:05

COMPOUND	COMP. CALIB. TYPE	MODEL	1	2	3	4	5	6	7	8	9	Avg. RF	Min. RF	%RSD CL	%RSD CL	R or R ²	R or R ² CL	STATUS
TPH as Diesel	C	Avg RF	69.187	81.182	81.639	84.113	83.988					80.02	0.00	8	0-20			PASS
												2						

Data Files:

Level #	D/T Analyzed	Data File
1	2016-12-02 19:43	T:\GC_45\GC_45_data\2016\161202\16120207.d\Report.txt16120207
2	2016-12-02 20:05	T:\GC_45\GC_45_data\2016\161202\16120208.d\Report.txt16120208
3	2016-12-02 20:27	T:\GC_45\GC_45_data\2016\161202\16120209.d\Report.txt16120209
4	2016-12-02 20:48	T:\GC_45\GC_45_data\2016\161202\16120210.d\Report.txt16120210
5	2016-12-02 21:09	T:\GC_45\GC_45_data\2016\161202\16120211.d\Report.txt16120211

LR - E: Linear Regression (Equal Weight) LR - IC: Linear Regression (Inverse Concentration Weight) LR - ISC: Linear Regression (Inverse Square Concentration Weight)
QR - E: Quadratic Regression (Equal Weight)



INITIAL CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8015B (M)

ICV WORK ORDER: 099-14-354-18-5901
INITIAL BATCH: 1612021007
INSTRUMENT: GC 45

ANALYZED BY: 972
D/T ANALYZED: 2016-12-02 19:43
INITIAL: 2016-12-02 21:31
ICV: 27
REVIEWED BY: 2017-02-16 16:05
D/T REVIEWED:

DATA FILE: T:\GC_45\data\201616120216120212.d\Report.txt16120212

<u>COMPOUND NAME</u>	<u>COMP TYPE</u>	<u>CALIB MODEL</u>	<u>MIN RF</u>	<u>AVG RF</u>	<u>ICV RF</u>	<u>AMOUNT</u>	<u>ICV CONC</u>	<u>ICV %D</u>	<u>ICV %D CL</u>	<u>STATUS</u>
TPH as Diesel	C	Avg Resp	0.00	80.022	82.011			-2	0-30	PASS

MIN RF: Method Specified Minimum Response Factor

Report Date : 05-Dec-2016 12:22

Page 1

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 18-OCT-2016 12:03
 End Cal Date : 03-DEC-2016 01:29
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Cal Date : 05-Dec-2016 12:21 uj3k
 Curve Type : Average

Calibration File Names:

Level 1: /chem1/SVOA/GC_45.i/161202.b/16120219.d
 Level 2: /chem1/SVOA/GC_45.i/161202.b/16120220.d
 Level 3: /chem1/SVOA/GC_45.i/161202.b/16120221.d
 Level 4: /chem1/SVOA/GC_45.i/161202.b/16120222.d
 Level 5: /chem1/SVOA/GC_45.i/161202.b/16120223.d

Compound	5.000 Level 1	200.000 Level 2	400.000 Level 3	800.000 Level 4	1600.000 Level 5	RRF	% RSD
S 1 TPH as Jet A rf	86545	105289	90515	85391	87961	91140	9
S 5 TPH as JP5 rf	69215	101437	91291	87968	106079	91198	16
S 8 TPH Gas/Diesel rf	37970	53614	55393	55706	53469	51230	15
S 15 TPH as Diesel rf	69187	81182	81639	84113	83988	80022	8

Report Date : 05-Dec-2016 12:22

Page 2

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 18-OCT-2016 12:03
 End Cal Date : 03-DEC-2016 01:29
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Cal Date : 05-Dec-2016 12:21 uj3k
 Curve Type : Average

Compound	5.000 Level 1	200.000 Level 2	400.000 Level 3	800.000 Level 4	1600.000 Level 5	RRF	% RSD
S 27 Diesel Range Organics rf	66446	76346	76541	77565	77319	74844	6
S 32 Oil Range Organics rf	61212	79825	76603	79618	78958	75243	11
S 36 TPH as Motor Oil rf	61342	80625	77575	80771	80149	76092	11

Report Date : 05-Dec-2016 12:22

Page 3

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 18-OCT-2016 12:03
 End Cal Date : 03-DEC-2016 01:29
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Cal Date : 05-Dec-2016 12:21 uj3k
 Curve Type : Average

	5.000	200.000	400.000	800.000	1600.000		
Compound	Level 1	Level 2	Level 3	Level 4	Level 5	RRF	% RSD
S 53 Crude Oil rf	46307	41013	42600	48242	48611	45355	8
S 65 Hydraulic Oil rf	70102	84275	85720	85748	86559	82481	8

Report Date : 05-Dec-2016 12:22

Page 8

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 18-OCT-2016 12:03
 End Cal Date : 03-DEC-2016 01:29
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Cal Date : 05-Dec-2016 12:21 uj3k
 Curve Type : Average

Compound	5.000 Level 1	200.000 Level 2	400.000 Level 3	800.000 Level 4	1600.000 Level 5	RRF	% RSD
S 206 NWTPH_Diesel rf	54683	80240	89681	88534	88358	80299	18
S 207 NWTPH_Diesel range rf	54683	80240	89681	88534	88358	80299	18
S 209 NWTPH_Motor Oil rf	81215	73070	86473	80435	85496	81338	7
\$ 93 n-Octacosane	86757	86415	85727	86538	85868	86261	1
\$ 94 C28 n-Octacosane	92246	92350	96946	93190	93357	93618	2

Data File: /chem1/SVOA/GC_45.i/161202.b/16120207.d
 Report Date: 05-Dec-2016 10:25

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/161202.b/16120207.d
 Lab Smp Id:
 Inj Date : 02-DEC-2016 19:43
 Operator : 682 Inst ID: GC_45.i
 Smp Info : ICAL D5 C28 0.625 L102516B
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Meth Date : 05-Dec-2016 10:25 uj3k Quant Type: ESTD
 Cal Date : 03-DEC-2016 00:02 Cal File: 16120219.d
 Als bottle: 7 Calibration Sample, Level: 1
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: ICAL_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.666-8.944			345934	5.00000	4.323 (a)
S 27 Diesel Range Organics	2.374-8.944			332229	5.00000	4.438 (a)
\$ 93 n-Octacosane	8.730	8.730	0.000	54223	0.62500	0.628

QC Flag Legend

a - Target compound detected but, quantitated amount
 Below Limit Of Quantitation(BLOQ).

Data File: /chem1/SV00A/GC_45.i/161202.b/16120207.d

Date : 02-DEC-2016 19:43

Client ID:

Sample Info: ICAL D5 C28 0.625 L102516B

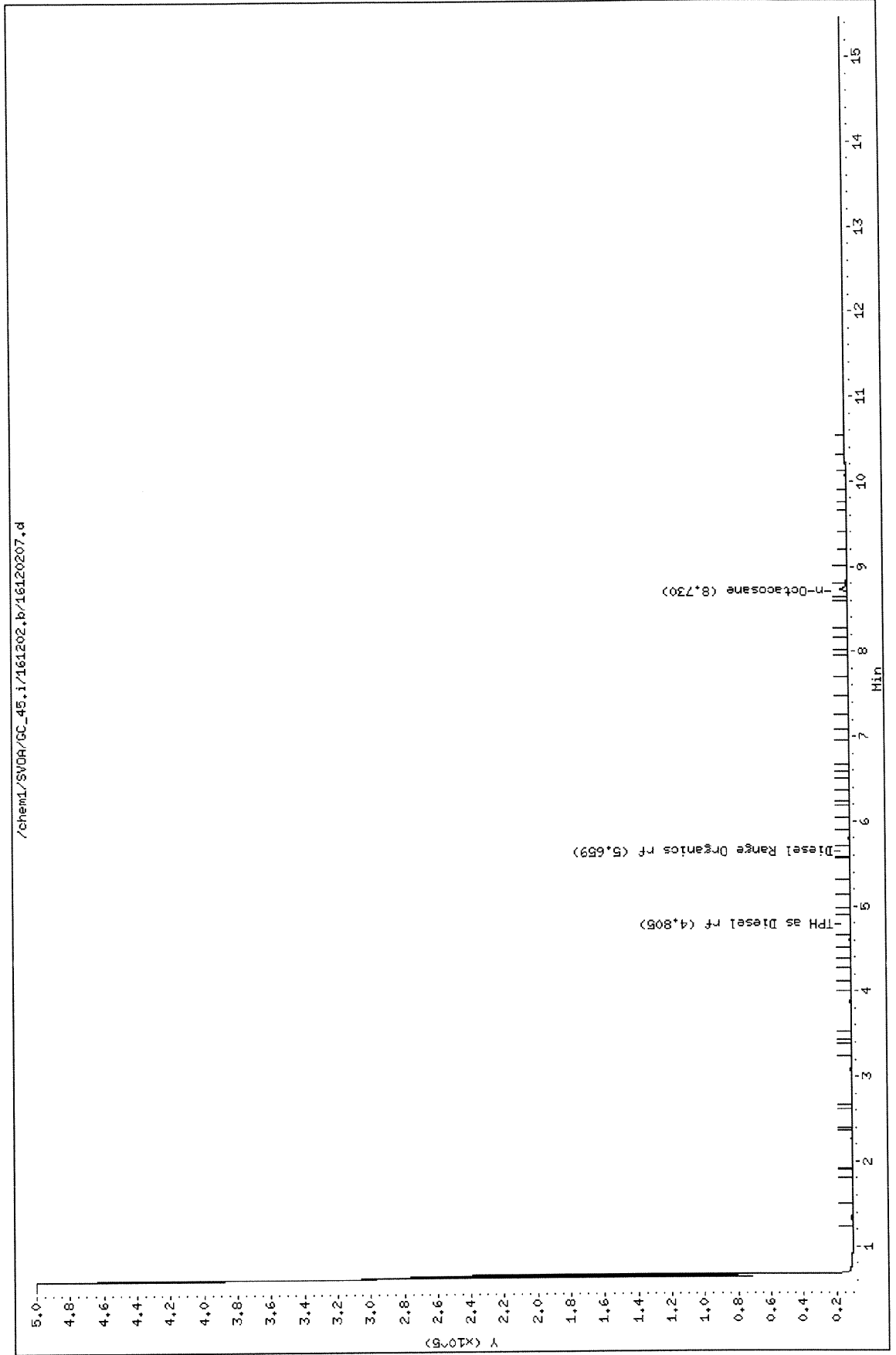
Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:

/chem1/SV00A/GC_45.i/161202.b/16120207.d



Data File: /chem1/SVOA/GC_45.i/161202.b/16120208.d
 Report Date: 05-Dec-2016 10:25

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/161202.b/16120208.d
 Lab Smp Id:
 Inj Date : 02-DEC-2016 20:05
 Operator : 682 Inst ID: GC_45.i
 Smp Info : ICAL D200 C28 25 L102516C
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Meth Date : 05-Dec-2016 10:25 uj3k Quant Type: ESTD
 Cal Date : 03-DEC-2016 00:23 Cal File: 16120220.d
 Als bottle: 8 Calibration Sample, Level: 2
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: ICAL_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.666-8.944			16236450	200.000	202.900
S 27 Diesel Range Organics	2.374-8.944			15269183	200.000	204.014
\$ 93 n-Octacosane	8.713	8.713	0.000	2160363	25.0000	25.044



Data File: /chem1/SVDA/GC_45.i/161202.b/16120208.d

Date : 02-DEC-2016 20:05

Client ID:

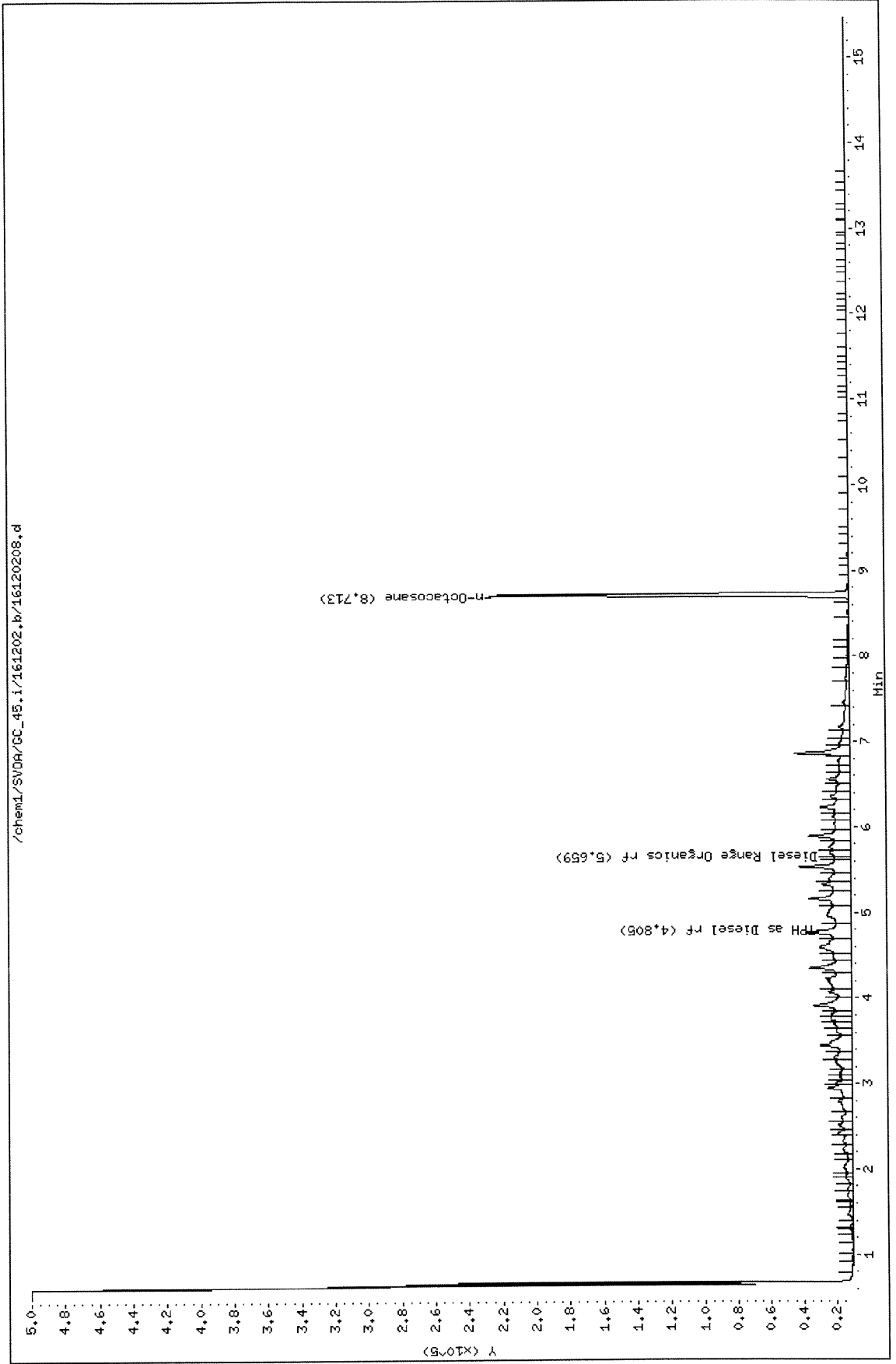
Sample Info: ICAL D200 C28 25 L102516C

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_45.i/161202.b/16120209.d
 Report Date: 05-Dec-2016 10:25

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/161202.b/16120209.d
 Lab Smp Id:
 Inj Date : 02-DEC-2016 20:27
 Operator : 682 Inst ID: GC_45.i
 Smp Info : ICAL D400 C28 50 L102516D
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Meth Date : 05-Dec-2016 10:25 uj3k Quant Type: ESTD
 Cal Date : 03-DEC-2016 00:46 Cal File: 16120221.d
 Als bottle: 9 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: ICAL_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.666-8.944			32655500	400.000	408.083
S 27 Diesel Range Organics	2.374-8.944			30616597	400.000	409.074
\$ 93 n-Octacosane	8.717	8.717	0.000	4286357	50.0000	49.690



Page 1

Data File: /chem1/SV09/GC_45.i/161202.k/16120209.d

Date : 02-DEC-2016 20:27

Client ID:

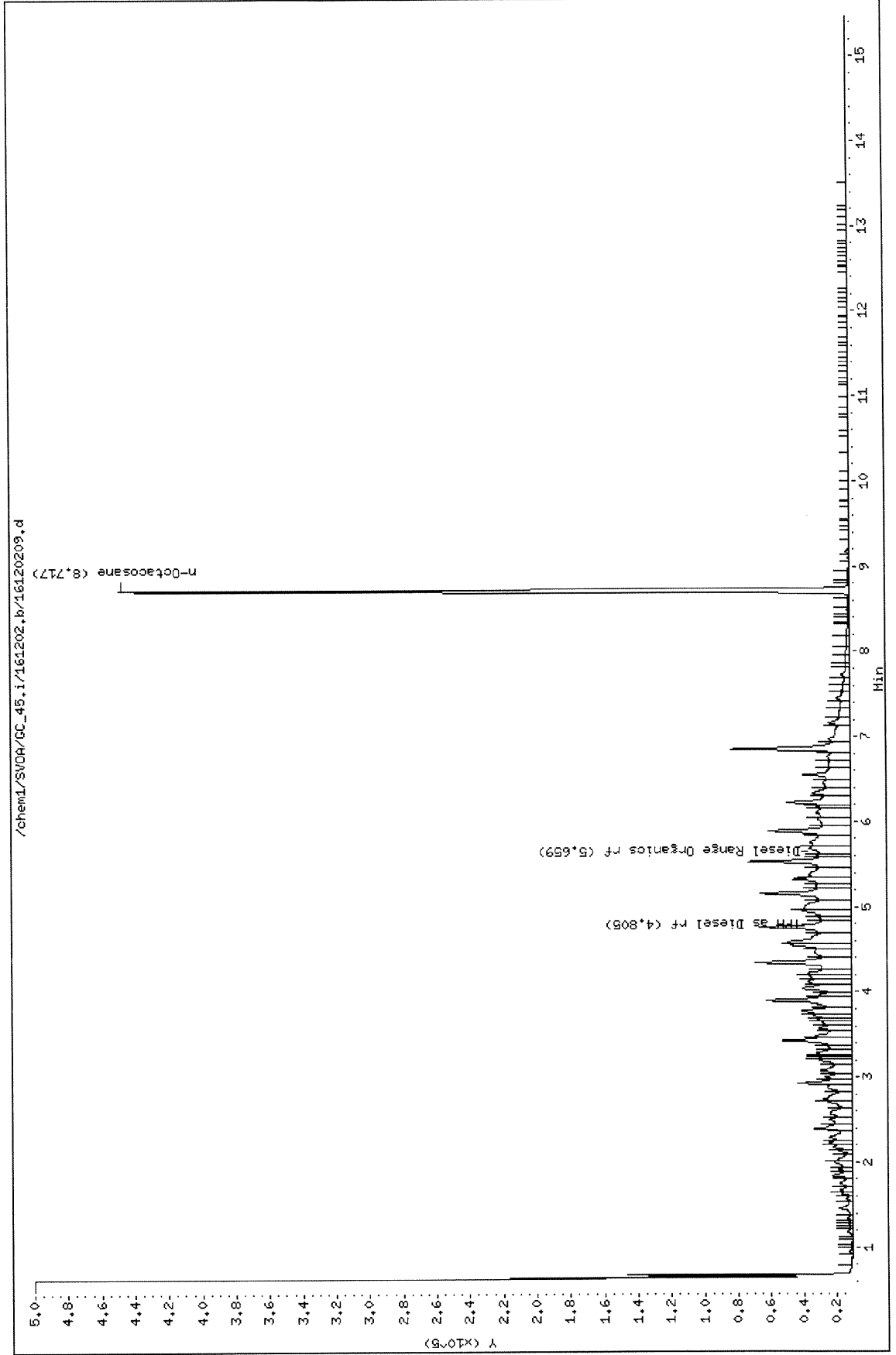
Sample Info: ICAL D400 C28 50 L102516D

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_45.i/161202.b/16120210.d
 Report Date: 05-Dec-2016 10:25

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/161202.b/16120210.d
 Lab Smp Id:
 Inj Date : 02-DEC-2016 20:48
 Operator : 682 Inst ID: GC_45.i
 Smp Info : ICAL D800 C28 100 L102516E
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Meth Date : 05-Dec-2016 10:25 uj3k Quant Type: ESTD
 Cal Date : 03-DEC-2016 01:08 Cal File: 16120222.d
 Als bottle: 10 Calibration Sample, Level: 4
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: ICAL_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.666-8.944			67290511	800.000	840.903
S 27 Diesel Range Organics	2.374-8.944			62052115	800.000	829.091
\$ 93 n-Octacosane	8.724	8.724	0.000	8653751	100.000	100.320



Data File: /chem1/SVDA/GC_45.1/161202.b/16120210.d

Date : 02-DEC-2016 20:48

Client ID:

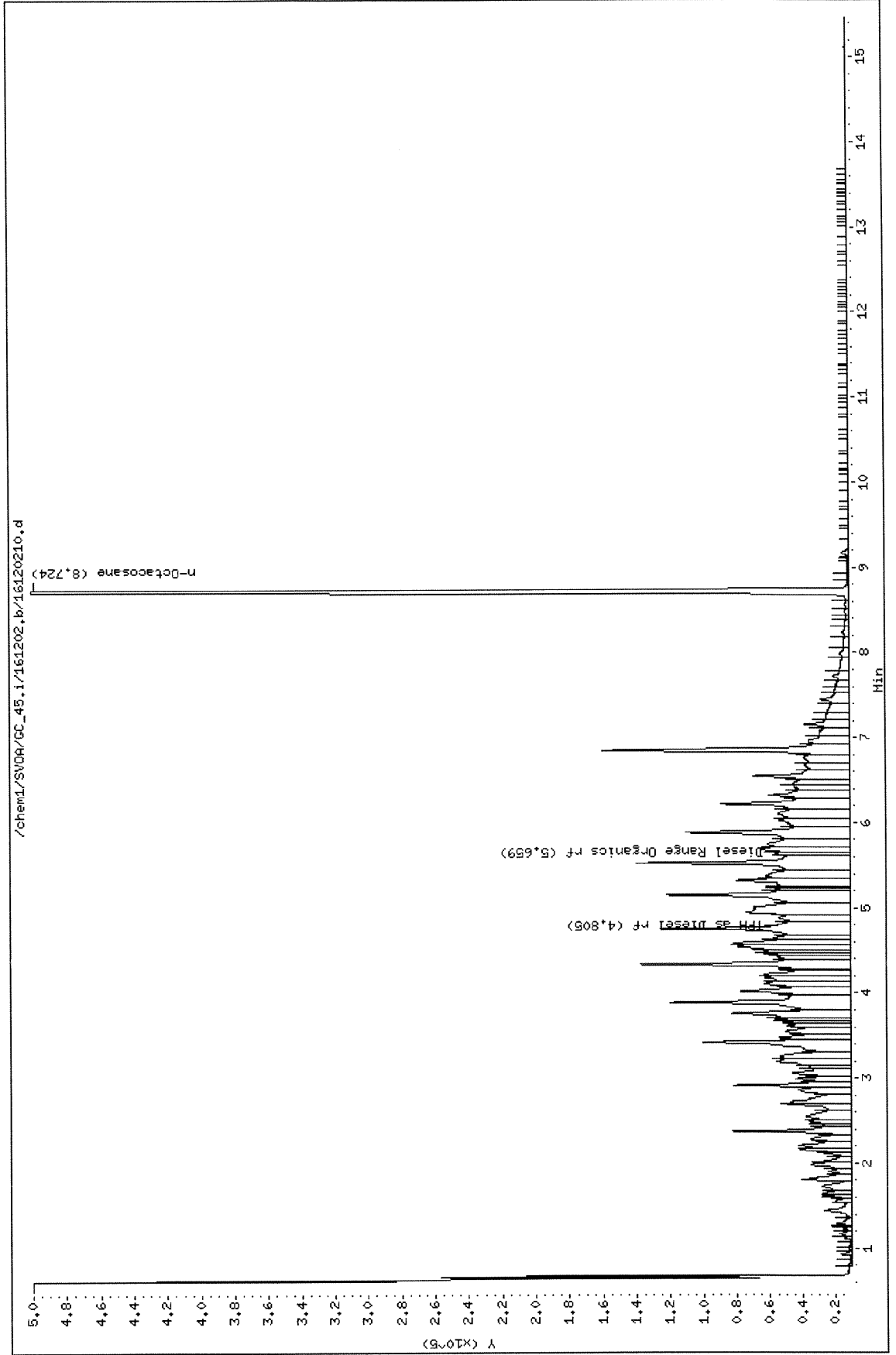
Sample Info: ICAL D800 C28 100 L102516E

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_45.i/161202.b/16120211.d
 Report Date: 05-Dec-2016 10:25

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/161202.b/16120211.d
 Lab Smp Id:
 Inj Date : 02-DEC-2016 21:09
 Operator : 682 Inst ID: GC_45.i
 Smp Info : ICAL D1600 C28 200 L102516F
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Meth Date : 05-Dec-2016 10:25 uj3k Quant Type: ESTD
 Cal Date : 03-DEC-2016 01:29 Cal File: 16120223.d
 Als bottle: 11 Calibration Sample, Level: 5
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: ICAL_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RT	RESPONSE	AMOUNTS	
						CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.666-8.944				134380116	1600.00	1679.295 (A)
S 27 Diesel Range Organics	2.374-8.944				123710939	1600.00	1652.927 (A)
\$ 93 n-Octacosane	8.737	8.737	0.000		17173600	200.000	199.089

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.



Data File: /chem1/SVDA/GC_45.i/161202.b/16120211.d

Date : 02-DEC-2016 21:09

Client ID:

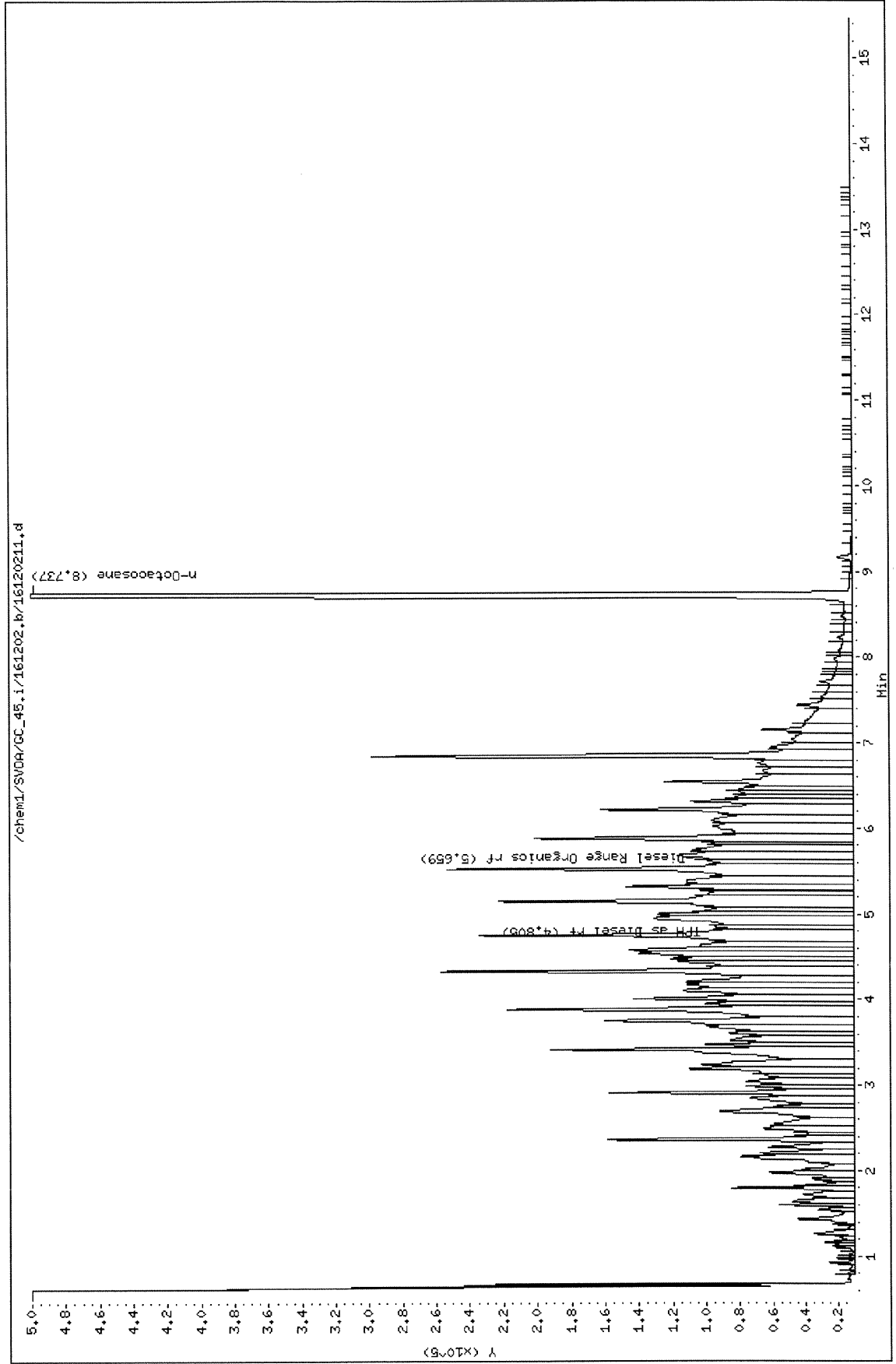
Sample Info: ICAL D1600 C28 200 L102516F

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_45.i/161202.b/16120212.d
 Report Date: 05-Dec-2016 10:25

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/161202.b/16120212.d
 Lab Smp Id:
 Inj Date : 02-DEC-2016 21:31
 Operator : 682 Inst ID: GC_45.i
 Smp Info : ICV D400 C28 50 L102516G
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/161202.b/8015d.m
 Meth Date : 05-Dec-2016 10:25 ujj3k Quant Type: ESTD
 Cal Date : 03-DEC-2016 01:29 Cal File: 16120223.d
 Als bottle: 12 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: ICAL_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

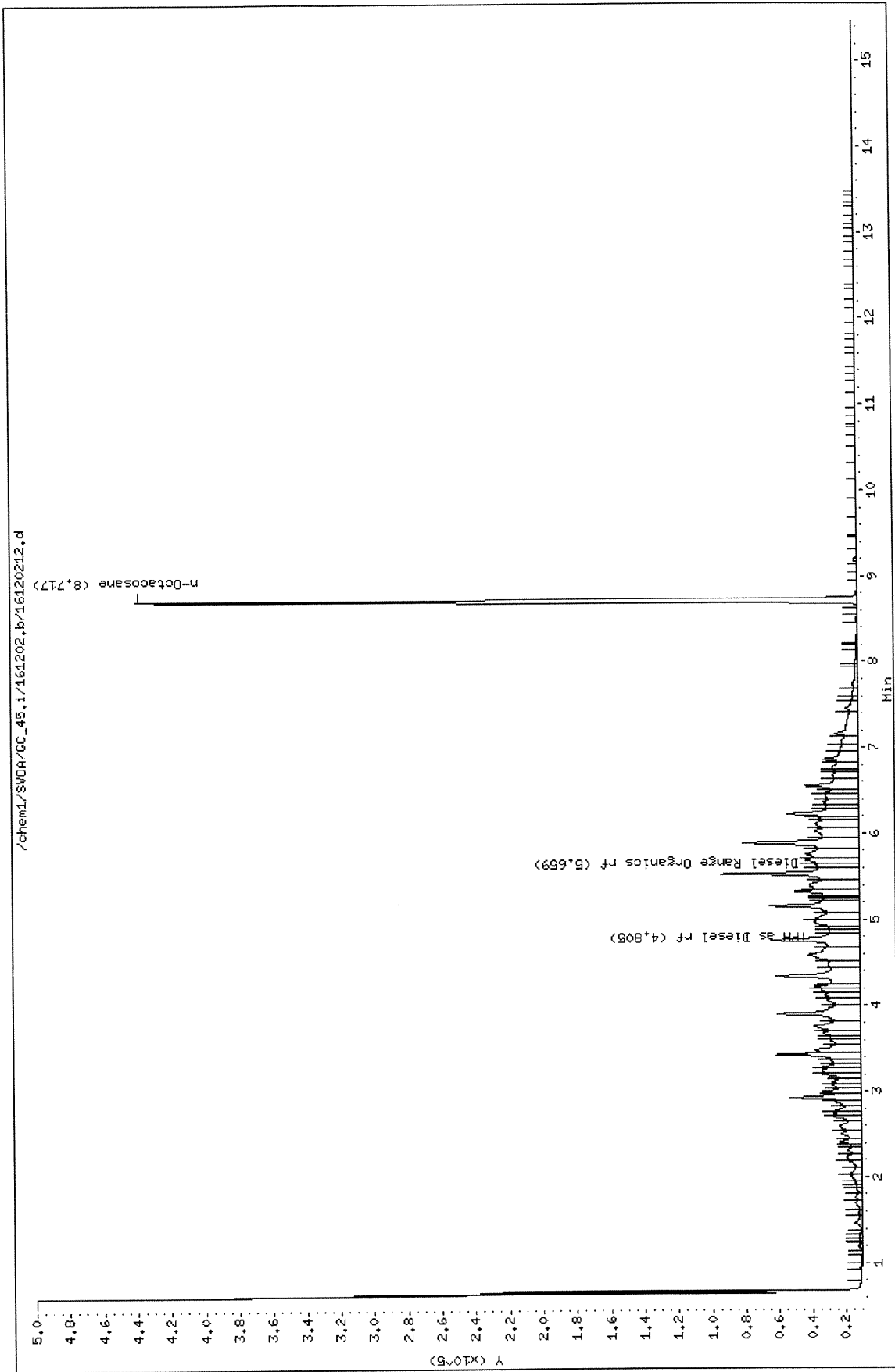
Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.666-8.944			32804272	400.000	409.942
S 27 Diesel Range Organics	2.374-8.944			31562730	400.000	421.716
\$ 93 n-Octacosane	8.717	8.717	0.000	4357915	50.0000	50.520



Data File: /chem1/SVDA/GC_45.i/161202.b/16120212.d
Date : 02-DEC-2016 21:31
Client ID:
Sample Info: ICV D400 C28 50 L102516G

Instrument: GC_45.i
Operator: 682
Column diameter: 2.00

Column phase:



=====
 External Standard Report
 =====

Data File Name : /chem1/SVOA/GC_45/161202/16120202.d
 Page Number :
 Operator : 682 Vial Number : Vial 2
 Instrument : GC 45 Injection Number : 2
 Sample Name : C6-C44 L110816A Sequence Line : 0
 Instrument Method: 8015d.m
 Acquired on : 02 DEC 16 18:40
 Report Created on: 05-DEC-16 10:26 Method Sublist : all
 Software Revision: Target 3.50

Sig. 1 in /chem1/SVOA/GC_45.i/161202.b/16120202.d

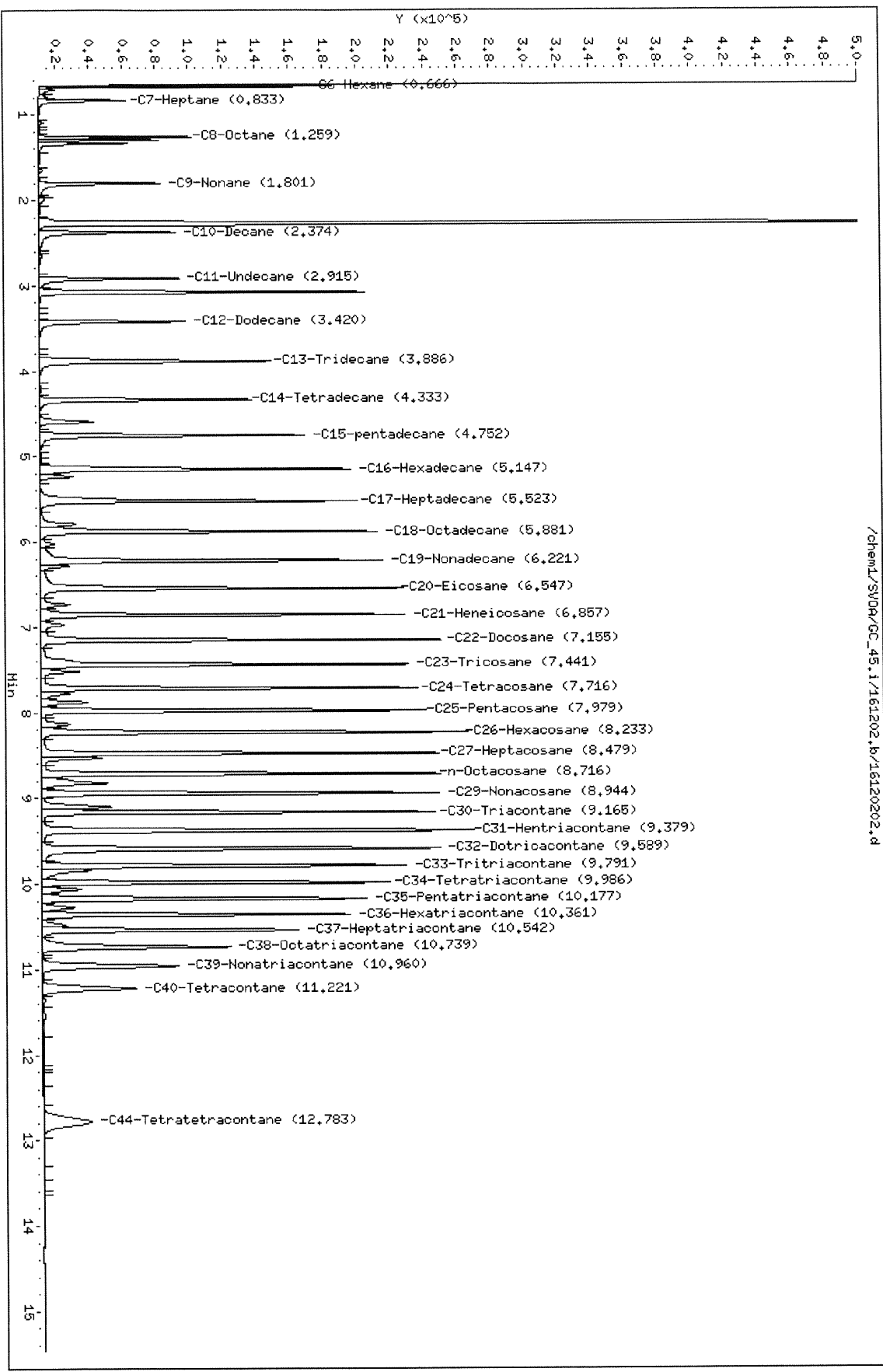
RT Range	Exp RT	DLT RT	Response	ppm	Compound
0.666	8.731	8.065	180018.00	0.00	C6-Hexane
0.833	8.731	7.898	363999.00	0.00	C7-Heptane
1.259	8.731	7.472	516823.00	0.00	C8-Octane
1.801	8.731	6.930	648211.00	0.00	C9-Nonane
2.374	8.731	6.357	819875.00	0.00	C10-Decane
2.915	8.731	5.816	821043.00	0.00	C11-Undecane
3.420	8.731	5.311	868861.00	0.00	C12-Dodecane
3.886	8.731	4.845	1699945.00	0.00	C13-Tridecane
4.333	8.731	4.398	1198646.00	0.00	C14-Tetradecane
4.752	8.731	3.979	1628623.00	0.00	C15-pentadecane
5.147	8.731	3.584	1935413.00	0.00	C16-Hexadecane
5.523	8.731	3.208	2181667.00	0.00	C17-Heptadecane
5.881	8.731	2.850	2092837.00	0.00	C18-Octadecane
6.221	8.731	2.510	2246859.00	0.00	C19-Nonadecane
6.547	8.731	2.184	2479179.00	0.00	C20-Eicosane
6.857	8.731	1.874	2195758.00	0.00	C21-Heneicosane
7.155	8.731	1.576	2470063.00	0.00	C22-Docosane
7.441	8.731	1.290	2524633.00	0.00	C23-Tricosane
7.716	8.731	1.015	2327237.00	0.00	C24-Tetracosane
7.979	8.731	0.752	2307901.00	0.00	C25-Pentacosane
8.233	8.731	0.498	2726070.00	0.00	C26-Hexacosane
8.479	8.731	0.252	2592723.00	0.00	C27-Heptacosane
8.716	8.731	0.015	2414138.00	1.63	n-Octacosane
8.944	8.731	-0.213	2391574.00	0.00	C29-Nonacosane
9.165	8.731	-0.434	2426307.00	0.00	C30-Triacontane
9.379	8.731	-0.648	2983404.00	0.00	C31-Hentriacontane
9.589	8.731	-0.858	2945350.00	0.00	C32-Dotriacontane
9.791	8.731	-1.060	2432008.00	0.00	C33-Tritriacontane
9.986	8.731	-1.255	2277735.00	0.00	C34-Tetratriacontane
10.177	8.731	-1.446	2179871.00	0.00	C35-Pentatriacontane
10.361	8.731	-1.630	2070018.00	0.00	C36-Hexatriacontane
10.542	8.731	-1.811	1782436.00	0.00	C37-Heptatriacontane
10.739	8.731	-2.008	1683811.00	0.00	C38-Octatriacontane
10.960	8.731	-2.229	1322621.00	0.00	C39-Nonatriacontane
11.221	8.731	-2.490	1024950.00	0.00	C40-Tetracontane
12.783	8.731	-4.052	1055033.00	0.00	C44-Tetratetracontane

End of File

Data File: /chem1/SVD9/CC_45.1/161202.b/16120202.d
 Date : 02-DEC-2016 18:40
 Client ID:
 Sample Info: C6-C44 L1108166
 Column phase:

Instrument: GC_45.1
 Operator: 682
 Column diameter: 2.00

/chem1/SVD9/CC_45.1/161202.b/16120202.d



[Return to Contents](#)

EPA 8015B (M) Diesel + Motor Oil

SAMPLE DATA

**RAW DATA SHEET
FOR METHOD: EPA 8015B (M)**

WORK ORDER: 17-03-1557
INSTRUMENT: GC 45
EXTRACTION: EPA 3550B
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 972
D/T ANALYZED: 2017-03-28 14:08
REVIEWED BY:
D/T REVIEWED:

DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032810.d\Report.txt17032810

25 **CLIENT SAMPLE NUMBER: B-DU1-ISM1-8**

LCS/MB BATCH: 170324B12 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 10.00 g / ACTUAL: 10.10 g
MS/MSD BATCH: 170324S12 **FINAL VOLUME / WEIGHT:** DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 0.99

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
TPH as Diesel	0.430	1.00	ND	5.0	
TPH as Motor Oil	0.700	1.00	ND	5.0	

```

=====
External Standard Report
=====

```

```

Data File Name : /chem1/SVOA/GC_45/170328/17032810.d
Page Number   :
Operator      : 682                               Vial Number    : Vial 10
Instrument    : GC 45                             Injection Number : 10
Sample Name   : 17-03-1557-25                    Sequence Line  : 0
                                                    Instrument Method: 8015d.m

Acquired on   : 28 MAR 17 14:08
Report Created on: 29-MAR-17 09:47               Compound Sublist : all
Software Revision: Target 3.50

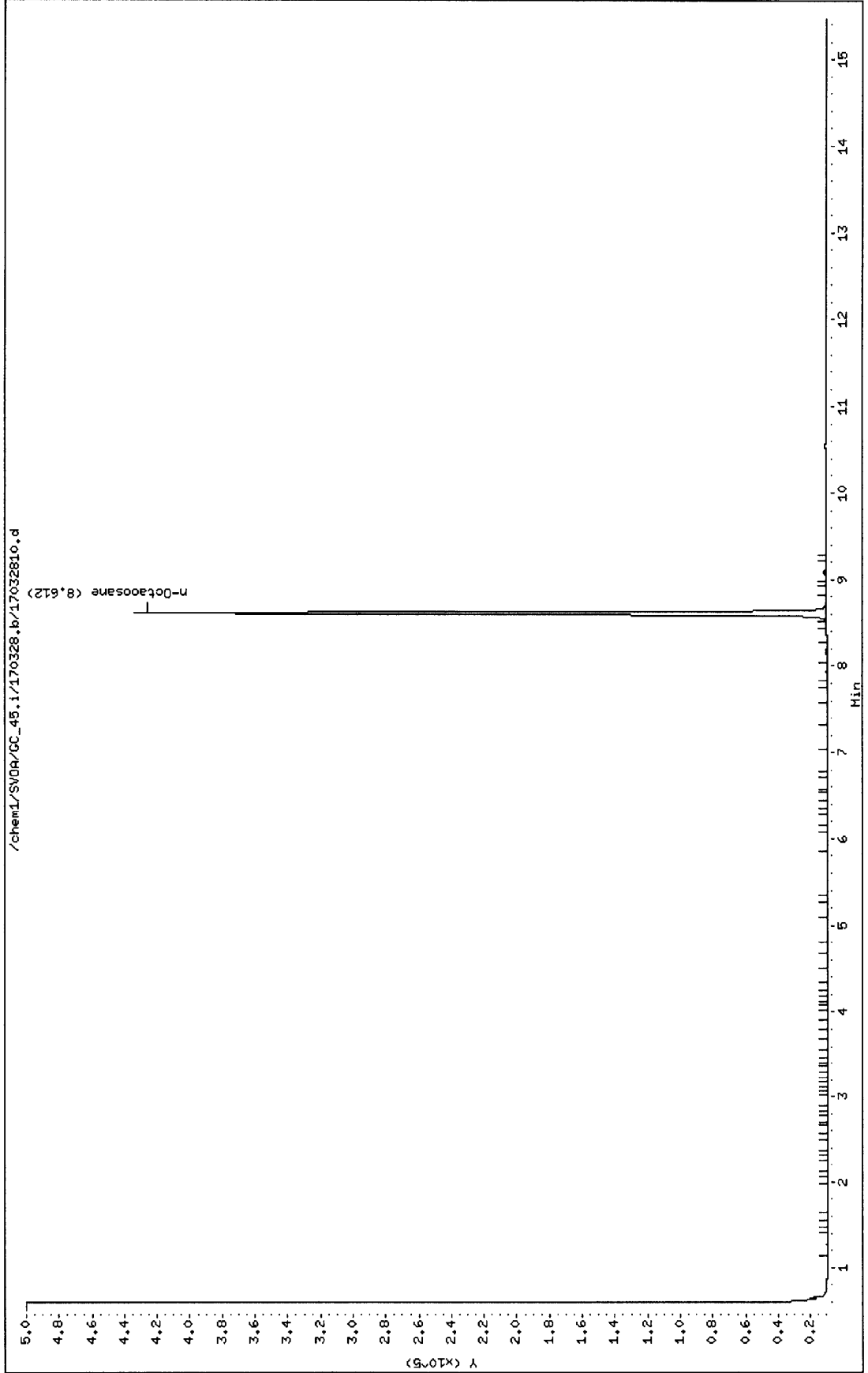
```

```

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032810.d
RT Range      Exp RT      DLT RT      Response      ppm      Compound
|-----|-----|-----|-----|-----|-----|
      8.612      8.628      0.017      4382608.00      50.81      n-Octacosane
2.300- 8.839      34198.42      0.43      TPH as Diesel
5.442-12.391      55892.46      0.70      TPH as Motor Oil
End of File

```

Data File: /chem1/SV0A/CC_45.i/170328.b/17032810.d
Date : 28-MAR-2017 14:08
Client ID:
Sample Info: 17-03-1557-25
Column phase:
Instrument: GC_45.i
Operator: 682
Column diameter: 2.00



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 17-03-1557
INSTRUMENT: GC 45
EXTRACTION: EPA 3550B
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 972
D/T ANALYZED: 2017-03-28 14:29
REVIEWED BY:
D/T REVIEWED:

DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032811.d\Report.txt17032811

26 **CLIENT SAMPLE NUMBER: B-DU1-ISM2-8**

LCS/MB BATCH: 170324B12 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 10.00 g / ACTUAL: 10.20 g
MS/MSD BATCH: 170324S12 **FINAL VOLUME / WEIGHT:** DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 0.98

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
TPH as Diesel	0.570	1.00	ND	4.9	
TPH as Motor Oil	0.940	1.00	ND	4.9	

External Standard Report

Data File Name : /chem1/SVOA/GC_45/170328/17032811.d
Page Number :
Operator : 682 Vial Number : Vial 11
Instrument : GC 45 Injection Number : 11
Sample Name : 17-03-1557-26 Sequence Line : 0
Instrument Method: 8015d.m
Acquired on : 28 MAR 17 14:29
Report Created on: 29-MAR-17 09:47 Compound Sublist : all
Software Revision: Target 3.50

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032811.d

RT Range	Exp RT	DLT RT	Response	ppm	Compound
8.611	8.628	0.017	4512502.00	52.31	n-Octacosane
2.300- 8.839			45521.73	0.57	TPH as Diesel
5.442-12.391			75515.93	0.94	TPH as Motor Oil

End of File

Page 1

Data File: /chem1/SV0A/GC_45.i/170328.b/17032811.d

Date : 28-MAR-2017 14:29

Client ID:

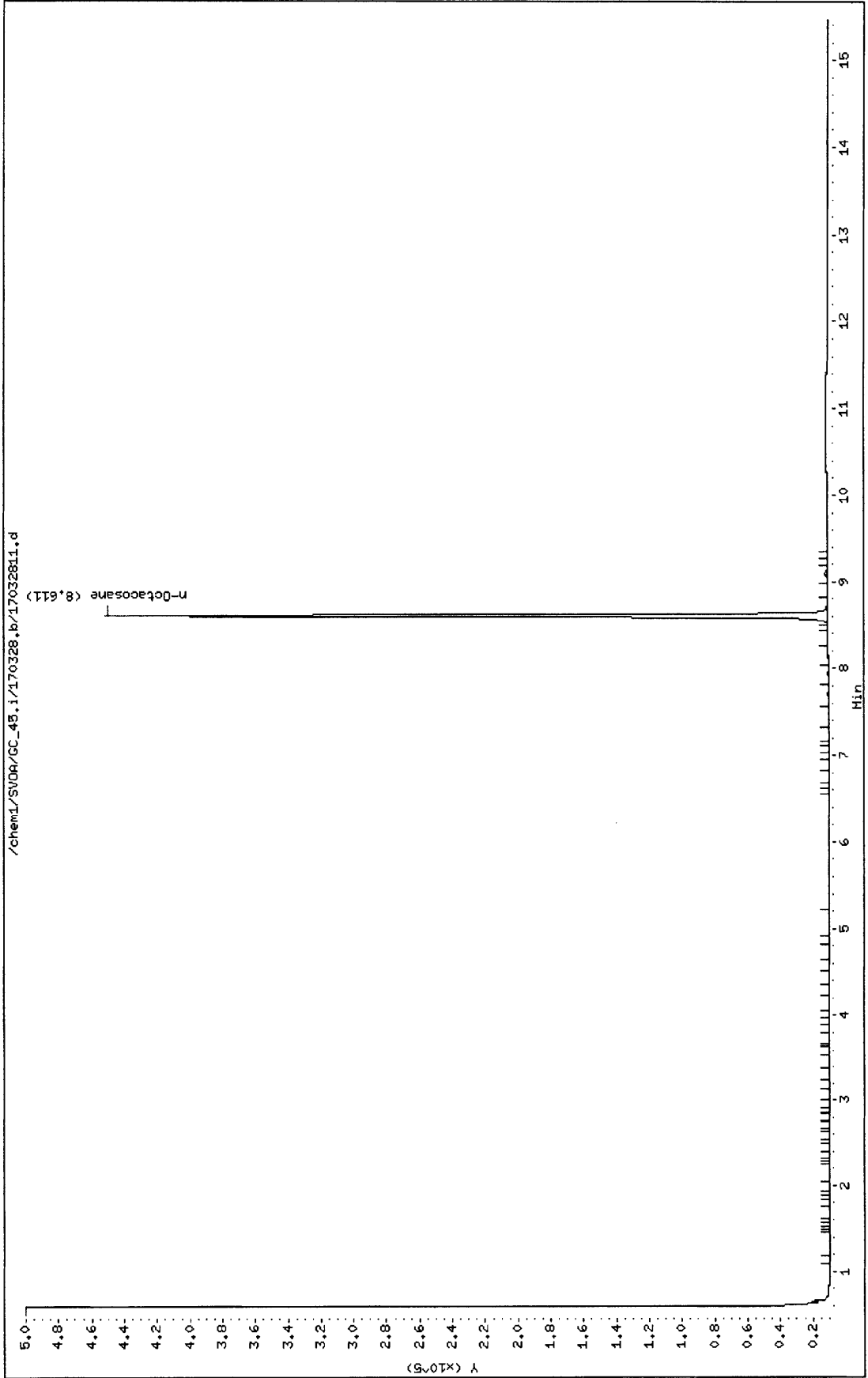
Sample Info: 17-03-1557-26

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 17-03-1557
INSTRUMENT: GC 45
EXTRACTION: EPA 3550B
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 972
D/T ANALYZED: 2017-03-28 14:51
REVIEWED BY:
D/T REVIEWED:

DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032812.d\Report.txt17032812

27 **CLIENT SAMPLE NUMBER: B-DU1-ISM3-8**

LCS/MB BATCH: 170324B12 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 10.00 g / ACTUAL: 10.10 g
MS/MSD BATCH: 170324S12 **FINAL VOLUME / WEIGHT:** DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 0.99

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
TPH as Diesel	0.450	1.00	ND	5.0	
TPH as Motor Oil	0.910	1.00	ND	5.0	

```

=====
External Standard Report
=====

```

```

Data File Name   : /chem1/SVOA/GC_45/170328/17032812.d
Page Number      :
Operator         : 682                               Vial Number      : Vial 12
Instrument       : GC 45                             Injection Number  : 12
Sample Name     : 17-03-1557-27                     Sequence Line    : 0
                                                    Instrument Method: 8015d.m

Acquired on     : 28 MAR 17 14:51
Report Created  : 29-MAR-17 09:48                   Compound Sublist : all
Software Revision: Target 3.50

```

```

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032812.d
RT Range      Exp RT    DLT RT    Response      ppm      Compound
|-----|-----|-----|-----|-----|-----|
      8.611      8.628      0.017      4345132.00    50.37    n-Octacosane
2.300- 8.839      35706.48      0.45    TPH as Diesel
5.442-12.391      73014.44      0.91    TPH as Motor Oil
End of File

```

Page 1

Data File: /chem1/SVOR/GC_45.i/170328.b/17032812.d

Date : 28-MAR-2017 14:51

Client ID:

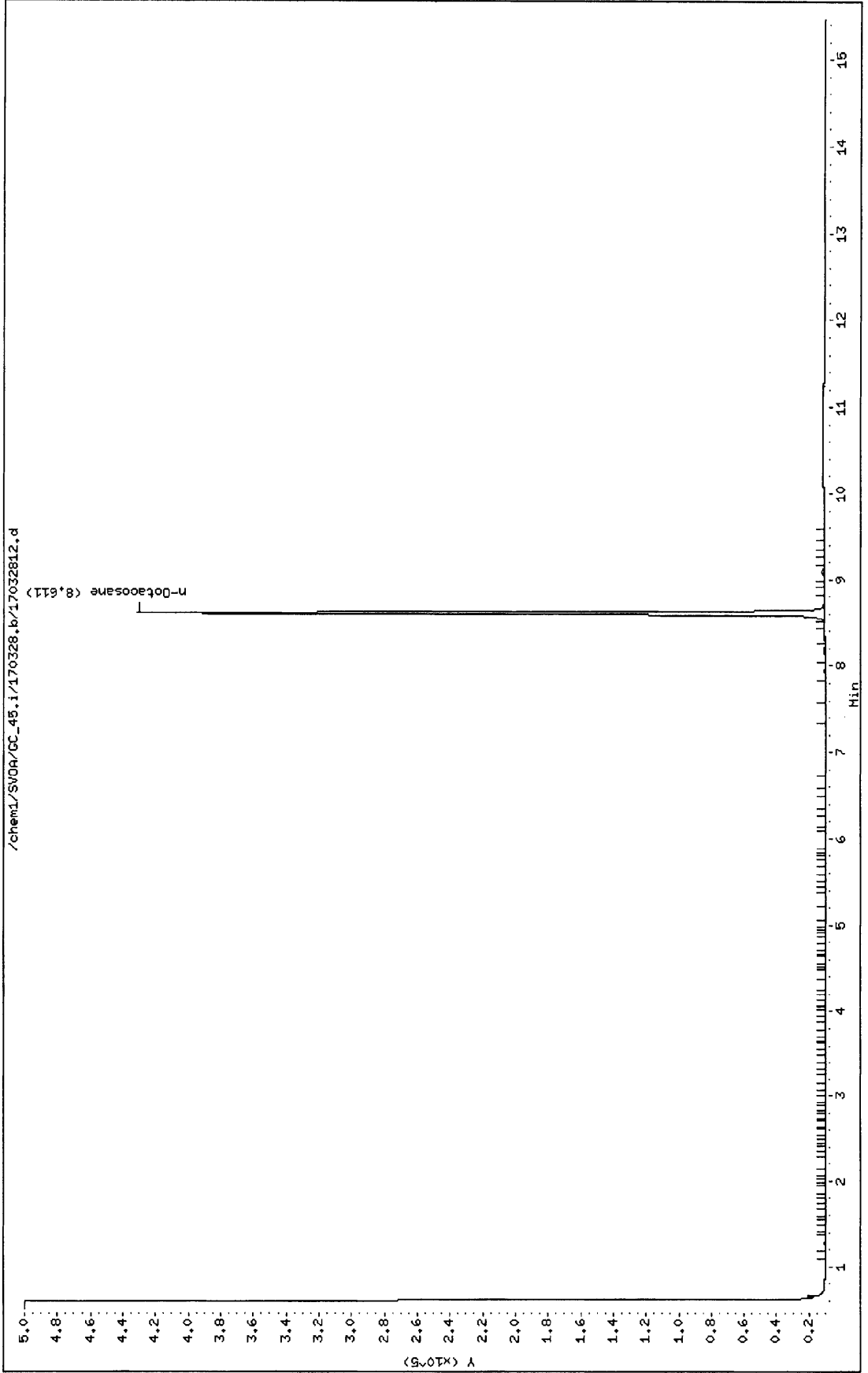
Sample Info: 17-03-1557-27

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



EPA 8015B (M)
Diesel + Motor Oil

QUALITY CONTROL

Method Blank
LCS/LCSD
MS/MSD

METHOD BLANK ASSOCIATION SUMMARY
FOR METHOD: EPA 8015B (M)

MB SAMPLE ID: 099-14-353-13
MB BATCH ID: 170324B12
INSTRUMENT: GC 45
EXTRACTION: EPA 3550B
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 972
D/T ANALYZED: 2017-03-28 11:16
REVIEWED BY:
D/T REVIEWED:
MATRIX: Soil

DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032806.d\Report.txt17032806

CLIENT WORK ORDER: 17-03-1557

<u>S#</u>	<u>RUN TYPE</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
25	B-DU1-ISM1-8		2017-03-28 14:08	T:\GC_45\GC_45_data\2017\170328\17032810.d\Report.txt17032810
26	B-DU1-ISM2-8		2017-03-28 14:29	T:\GC_45\GC_45_data\2017\170328\17032811.d\Report.txt17032811
27	B-DU1-ISM3-8		2017-03-28 14:51	T:\GC_45\GC_45_data\2017\170328\17032812.d\Report.txt17032812

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 099-14-353
INSTRUMENT: GC 45
EXTRACTION: EPA 3550B
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 972
D/T ANALYZED: 2017-03-28 11:16
REVIEWED BY:
D/T REVIEWED:

DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032806.d\Report.txt17032806

MB **CLIENT SAMPLE NUMBER: Method Blank**

LCS/MB BATCH: 170324B12 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 10.00 g / ACTUAL: 10.00 g
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
TPH as Diesel	0.360	1.00	ND	5.0	
TPH as Motor Oil	0.600	1.00	ND	5.0	



LCS QUALITY CONTROL SHEET FOR METHOD: EPA 8015B (M)

LCS SAMPLE ID: 099-14-353-13
LCS/MB BATCH ID: 170324B12
INSTRUMENT: GC 45

EXTRACTION: EPA 3550B
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 972
D/T ANALYZED: 2017-03-28 11:37
REVIEWED BY:
D/T REVIEWED:

DATA FILE: T:\GC_45\data\2017\170328\17032807.d\Report.txt\17032807

<u>COMPOUND NAME</u>	<u>CONC ADDED</u>	<u>CONC REC</u>	<u>%RECOVERY</u>	<u>%REC CONTROL LIMIT</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
TPH as Diesel	400.0	373.8	93	61-145	PASS	



Return to Contents

MATRIX SPIKE / MATRIX SPIKE DUPLICATE QUALITY CONTROL SHEET FOR METHOD: EPA 8015B (M)

SPIKED SAMPLE ID: 17-03-1557-27
MS/MSD BATCH: 170324S12

INSTRUMENTS:
SAMPLE: GC 45
MS: GC 45
MSD: GC 45

EXTRACTION: EPA 3550B
D/T EXTRACTED:

SAMPLE: 2017-03-24 00:00
MS: 2017-03-24 00:00
MSD: 2017-03-24 00:00

ANALYZED BY: 972

D/T ANALYZED:

SAMPLE: 2017-03-28 14:51
MS: 2017-03-28 12:42
MSD: 2017-03-28 13:03

REVIEWED BY:

D/T REVIEWED:

COMMENT:

COMPOUND NAME	SAMPLE	INITIAL	FINAL	MS CONC	% MS.REC	MSD CONC	% MSD.REC	% REC CL	RPD	RPD CL	STATUS	QUALIFIERS
TPH as Diesel	ND	400.0	400.0	379.0	95	393.3	98	61-145	4	0-25	PASS	

Data Files:

TYPE	DATA FILE	DATA FILE PATH
MS	17032808	T:\GC_45\GC_45_data\2017\170328\17032808.d\Report.txt
MSD	17032809	T:\GC_45\GC_45_data\2017\170328\17032809.d\Report.txt

SURROGATE RECOVERIES FOR METHOD: EPA 8015B (M)

WORK ORDER: 17-03-1557

BATCH ID:

REVIEWED BY:

D/T REVIEWED:

LCS/MB: 170324B12MS: 170324S12

EXTRACTION: EPA 3550B

25 **CLIENT SAMPLE NUMBER : B-DU1-ISM1-8**INSTRUMENT: GC 45ANALYZED BY: 972D/T EXTRACTED: 2017-03-24 00:00D/T ANALYZED 2017-03-28 14:08DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032810.d\Report.txt17032810COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	102	61-145	PASS	

26 **CLIENT SAMPLE NUMBER : B-DU1-ISM2-8**INSTRUMENT: GC 45ANALYZED BY: 972D/T EXTRACTED: 2017-03-24 00:00D/T ANALYZED 2017-03-28 14:29DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032811.d\Report.txt17032811COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	105	61-145	PASS	

27 **CLIENT SAMPLE NUMBER : B-DU1-ISM3-8**INSTRUMENT: GC 45ANALYZED BY: 972D/T EXTRACTED: 2017-03-24 00:00D/T ANALYZED 2017-03-28 14:51DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032812.d\Report.txt17032812COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	101	61-145	PASS	

MB **CLIENT SAMPLE NUMBER : Method Blank**INSTRUMENT: GC 45ANALYZED BY: 972D/T EXTRACTED: 2017-03-24 00:00D/T ANALYZED 2017-03-28 11:16DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032806.d\Report.txt17032806COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	103	61-145	PASS	

**SURROGATE RECOVERIES
FOR METHOD: EPA 8015B (M)**

WORK ORDER: 17-03-1557
BATCH ID:
LCS/MB: 170324B12
MS:
EXTRACTION: EPA 3550B

REVIEWED BY:
D/T REVIEWED:

LCS **CLIENT SAMPLE NUMBER :** Lab Control Sample

INSTRUMENT: GC 45 ANALYZED BY: 972
D/T EXTRACTED: 2017-03-24 00:00 D/T ANALYZED 2017-03-28 11:37
DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032807.d\Report.txt17032807

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	100	61-145	PASS	

MS **CLIENT SAMPLE NUMBER :** Matrix Spike

INSTRUMENT: GC 45 ANALYZED BY: 972
D/T EXTRACTED: 2017-03-24 00:00 D/T ANALYZED 2017-03-28 12:42
DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032808.d\Report.txt17032808

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	103	61-145	PASS	

MSD **CLIENT SAMPLE NUMBER :** Matrix Spike Duplicate

INSTRUMENT: GC 45 ANALYZED BY: 972
D/T EXTRACTED: 2017-03-24 00:00 D/T ANALYZED 2017-03-28 13:03
DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032809.d\Report.txt17032809

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
n-Octacosane	102	61-145	PASS	

Return to Contents

=====
External Standard Report
=====

Data File Name : /chem1/SVOA/GC_45/170328/17032806.d
Page Number :
Operator : 682 Vial Number : Vial 6
Instrument : GC 45 Injection Number : 6
Sample Name : MB 17032412 Sequence Line : 0
Instrument Method: 8015d.m
Acquired on : 28 MAR 17 11:16
Report Created on: 28-MAR-17 17:33 Compound Sublist : all
Software Revision: Target 3.50

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032806.d

RT Range	Exp RT	DLT RT	Response	ppm	Compound
8.611	8.628	0.017	4422114.00	51.26	n-Octacosane
2.300- 8.839			28906.63	0.36	TPH as Diesel
5.442-12.391			47888.98	0.60	TPH as Motor Oil

End of File

Data File: /chem1/SV0A/GC_45.i/170328.b/17032806.d

Date : 28-MAR-2017 11:16

Client ID:

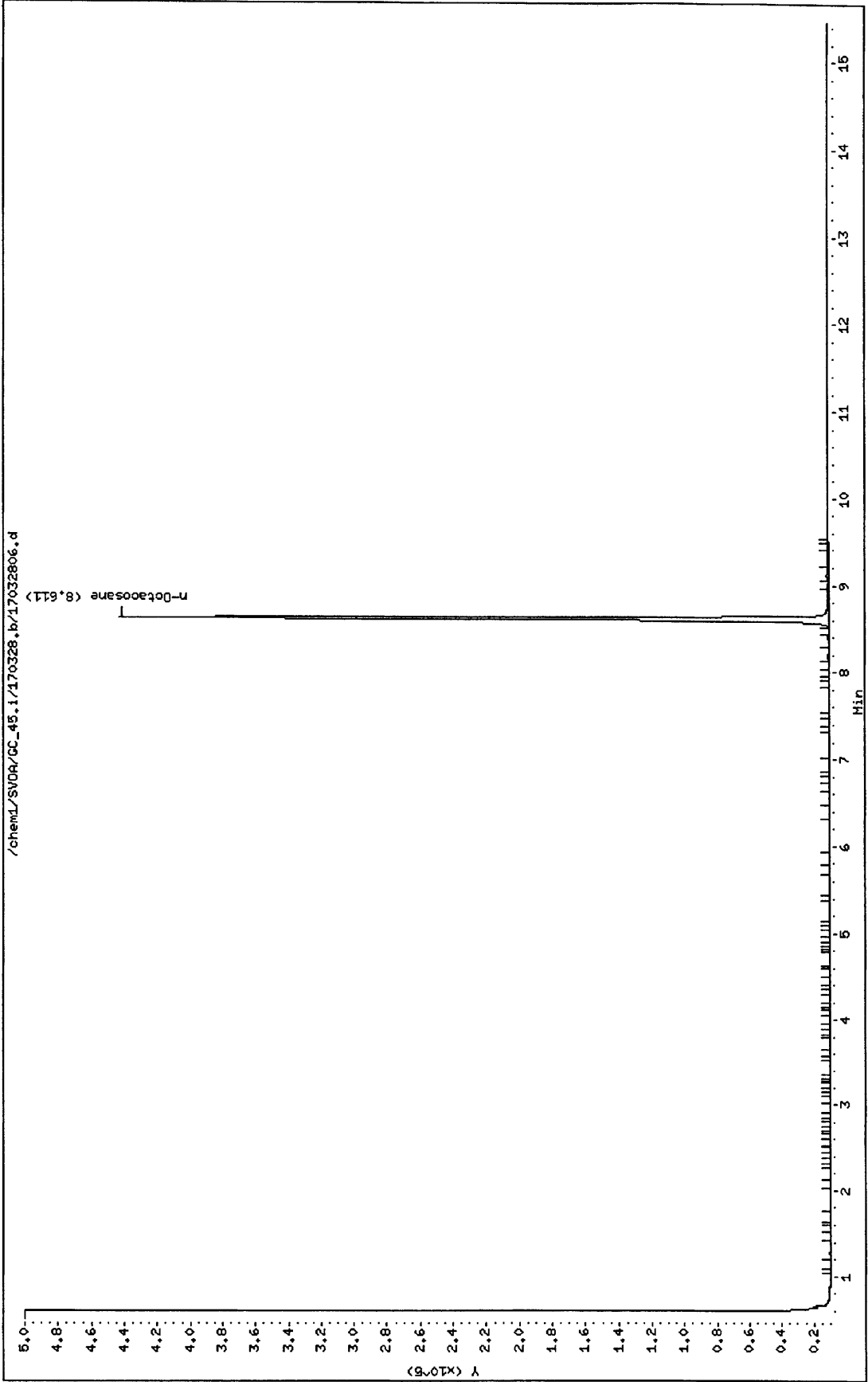
Sample Info: HB 17032412

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



```

=====
External Standard Report
=====

```

```

Data File Name   : /chem1/SVOA/GC_45/170328/17032807.d
Page Number      :
Operator         : 682                               Vial Number      : Vial 7
Instrument       : GC 45                             Injection Number : 7
Sample Name      : LCS 17032412                     Sequence Line    : 0
                                                    Instrument Method: 8015d.m

Acquired on      : 28 MAR 17 11:37
Report Created on: 28-MAR-17 17:33                 Compound Sublist : all
Software Revision: Target 3.50

```

```

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032807.d

```

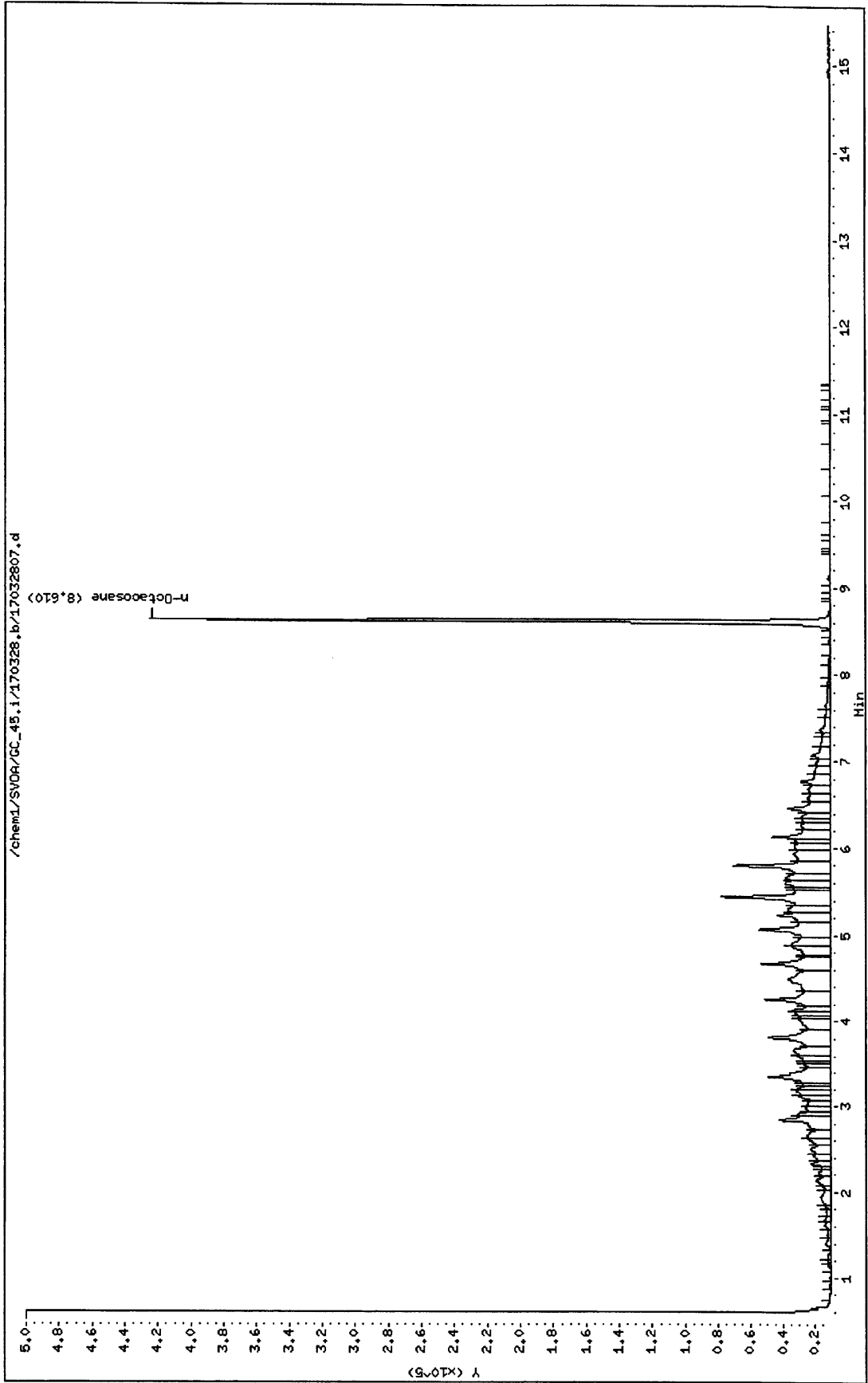
RT Range	Exp RT	DLT RT	Response	ppm	Compound
8.610	8.628	0.018	4297394.00	49.82	n-Octacosane
0.649- 8.839			29912924.71	373.81	TPH as Diesel
2.300- 8.839			28764552.27	384.33	Diesel Range Organics

End of File

Page 1

Data File: /chem1/SV09/GC_45.i/170328.b/17032807.d
Date : 28-MAR-2017 11:37
Client ID:
Sample Info: LCS 17032412

Instrument: GC_45.i
Operator: 682
Column diameter: 2.00



```

=====
External Standard Report
=====

```

```

Data File Name   : /chem1/SVOA/GC_45/170328/17032808.d
Page Number      :
Operator         : 682                               Vial Number      : Vial 8
Instrument       : GC 45                             Injection Number  : 8
Sample Name      : MS 17-03-1557-27                 Sequence Line    : 0
                                                    Instrument Method: 8015d.m

Acquired on      : 28 MAR 17 12:42
Report Created on: 29-MAR-17 09:47                   Compound Sublist : all
Software Revision: Target 3.50

```

```

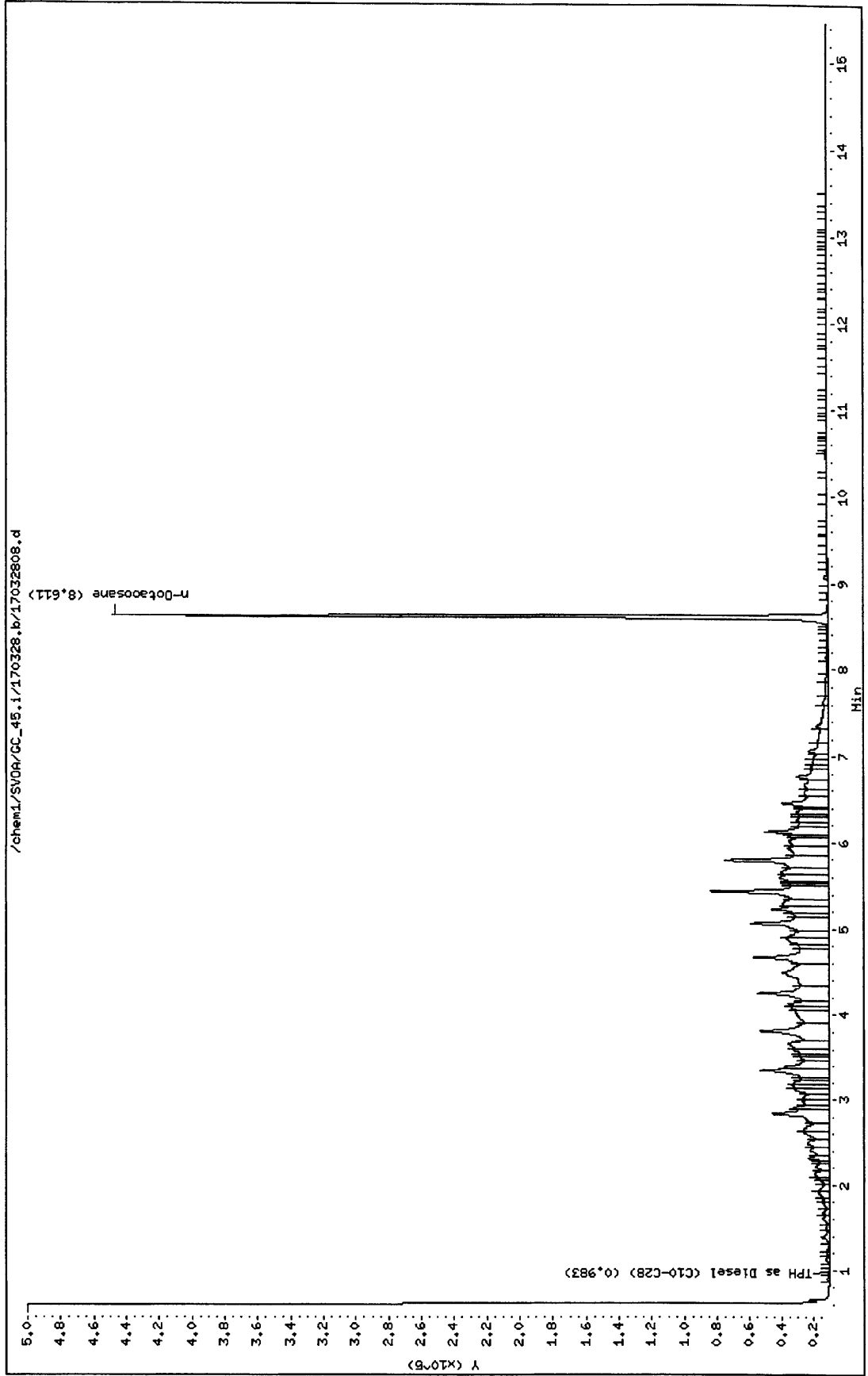
Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032808.d
RT Range      Exp RT   DLT RT   Response      ppm          Compound
|-----|-----|-----|-----|-----|-----|
      8.611      8.628      0.018      4450050.00    51.59    n-Octacosane
2.300- 8.839      30630369.62   382.78    TPH as Diesel
5.442-12.391     10370140.00   129.59    TPH as Motor Oil
End of File

```


Data File: /chem1/SV0A/GC_45.i/170328.b/17032808.d
Date : 28-MAR-2017 12:42
Client ID:
Sample Info: MS 17-03-1857-27

Instrument: GC_45.i
Operator: 682
Column diameter: 2.00

Column phase:



=====
External Standard Report
=====

Data File Name : /chem1/SVOA/GC_45/170328/17032809.d
Page Number :
Operator : 682 Vial Number : Vial 9
Instrument : GC 45 Injection Number : 9
Sample Name : MSD 17-03-1557-27 Sequence Line : 0
Instrument Method: 8015d.m
Acquired on : 28 MAR 17 13:03
Report Created on: 29-MAR-17 09:47 Compound Sublist : all
Software Revision: Target 3.50

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032809.d

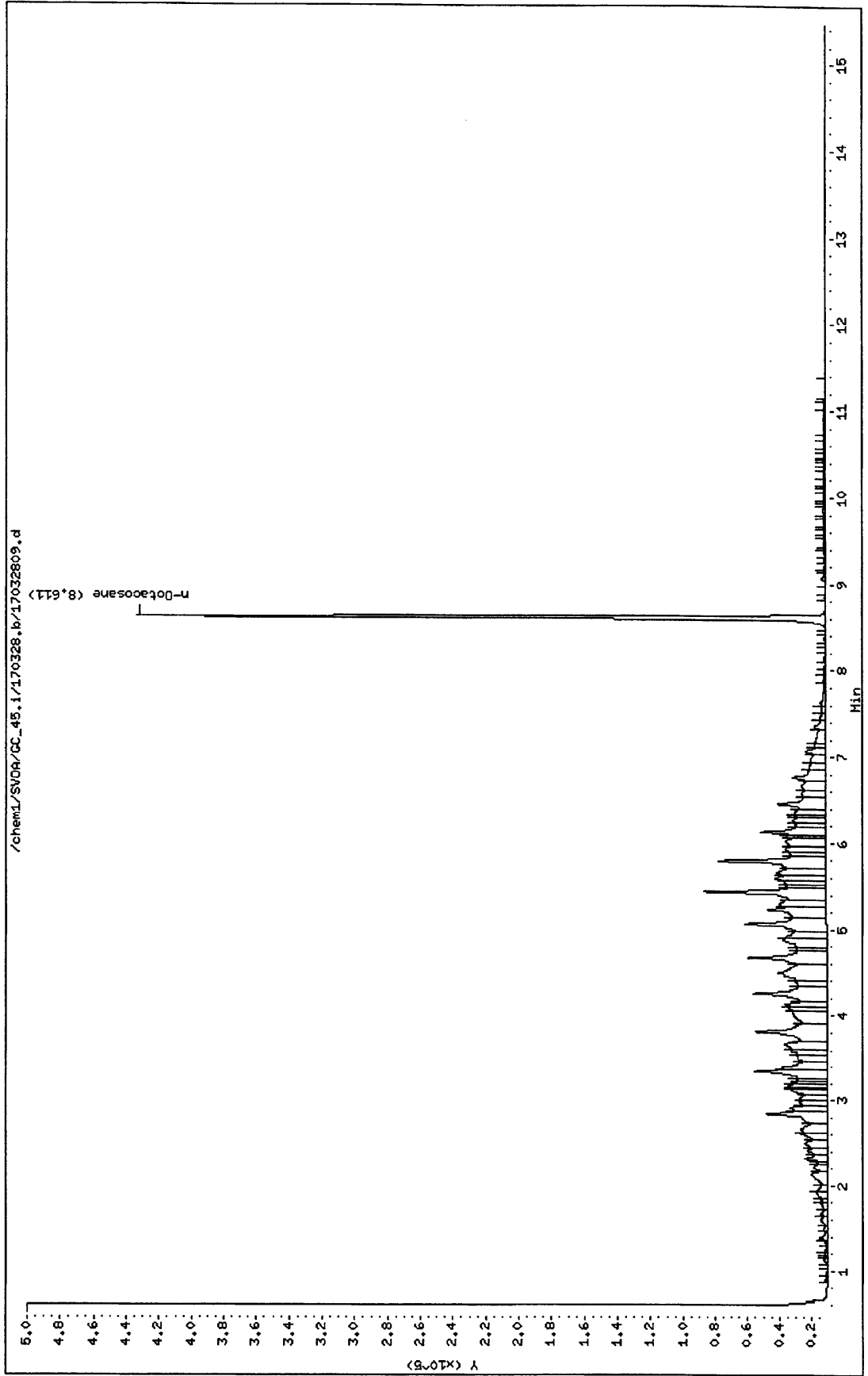
RT Range	Exp RT	DLT RT	Response	ppm	Compound
8.611	8.628	0.017	4386449.00	50.85	n-Octacosane
2.300- 8.839			31475683.35	393.34	TPH as Diesel
5.442-12.391			11132190.40	139.11	TPH as Motor Oil

End of File

Data File: /chem1/SVDA/GC_45.i/170328.b/17032809.d
Date : 28-MAR-2017 13:03
Client ID:
Sample Info: MSD 17-03-1557-27

Instrument: GC_45.i
Operator: 682
Column diameter: 2.00

Column phase:



EPA 8015B (M) Diesel + Motor Oil

CONTINUING CALIBRATION

CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8015B (M)

CCV WORK ORDER: 099-14-354-50-5901

BATCH ID:

1612021007

170328A031

GC 45

ANALYZED BY: 972

D/T ANALYZED:

INITIAL:

CCV:

REVIEWED BY:

D/T REVIEWED:

2016-12-02 19:43

2017-03-28 10:13

DATA FILE: T:\GC_45\data\20171170328\17032803.d\Report.txt\17032803

<u>COMPOUND NAME</u>	<u>TYPE</u>	<u>CALIB MODEL</u>	<u>MIN RF</u>	<u>AVG RF</u>	<u>CCV RE</u>	<u>AMOUNT</u>	<u>CCV CONC</u>	<u>CCV %D</u>	<u>CCV %D CL</u>	<u>STATUS</u>
TPH as Diesel	C	Avg Resp	0.00	80.022	79.568			1	0-20	PASS

MIN RF: Method Specified Minimum Response Factor

Data File: /chem1/SVOA/GC_45.i/170328.b/17032803.d
 Report Date: 28-Mar-2017 14:09

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/170328.b/17032803.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 10:13
 Operator : 682 Inst ID: GC_45.i
 Smp Info : CCV D400 C28 50 L102516D
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/170328.b/8015d.m
 Meth Date : 28-Mar-2017 14:09 umd6 Quant Type: ESTD
 Cal Date : 28-DEC-2016 19:01 Cal File: 16122828.d
 Als bottle: 3 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: CCV_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

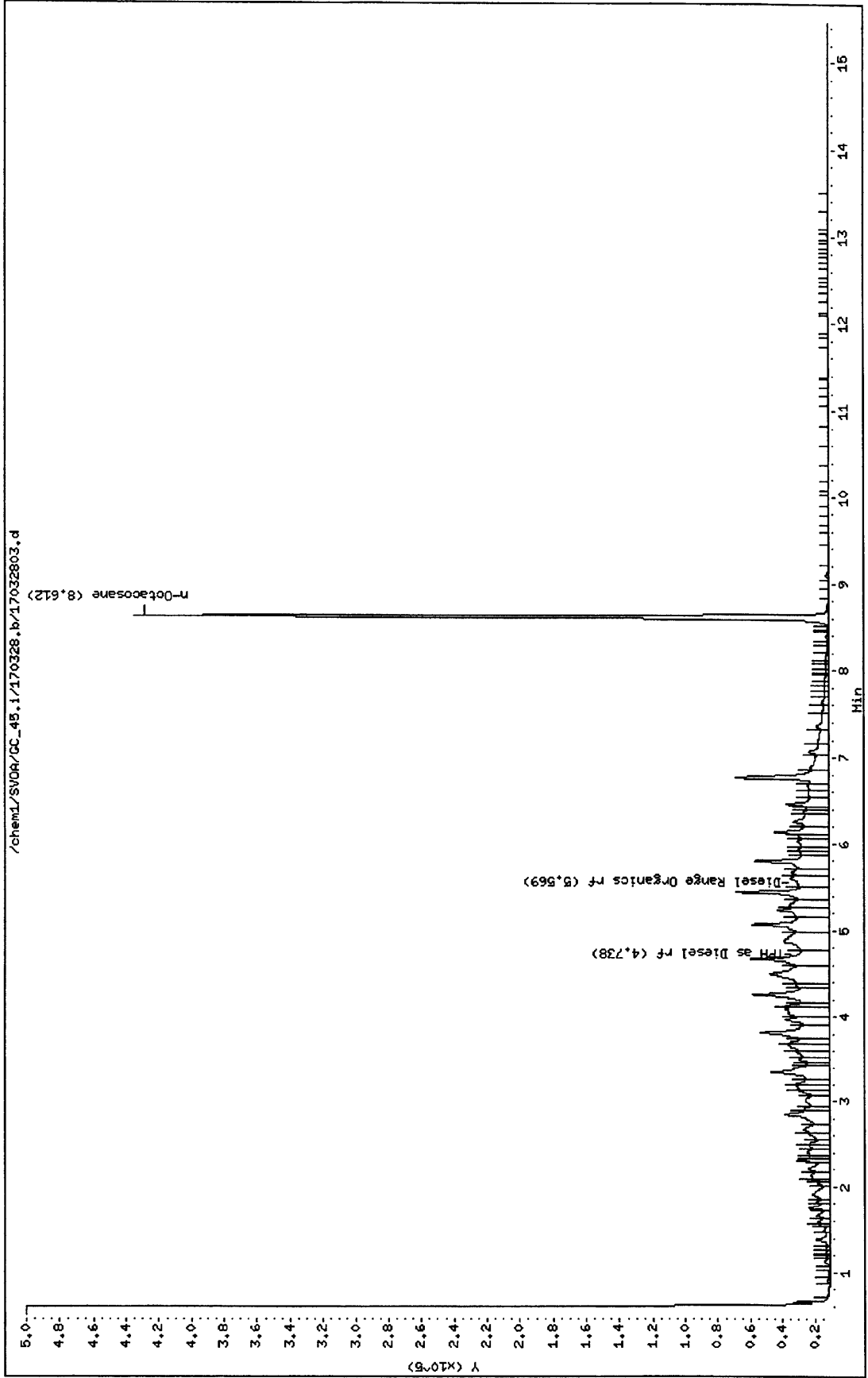
Compounds	RT	EXP RT	DLT RT	RT	RESPONSE	AMOUNTS	
						CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.649-8.839				31827054	400.000	397.730
S 27 Diesel Range Organics	2.300-8.839				29781149	400.000	397.912
\$ 93 n-Octacosane	8.612	8.612	0.000		4464923	50.0000	51.760

Page 1

Data File: /chem1/SVDA/CC_45.i/170328.b/17032803.d
Date : 28-MAR-2017 10:13
Client ID:
Sample Info: CCV D400 C28 50 L102516D

Instrument: GC_45.i
Operator: 682
Column diameter: 2.00

Column phase:



CCV ASSOCIATION SUMMARY FOR METHOD: EPA 8015B (M)

BATCH ID: 170328A055
INSTRUMENT: GC 45

ANALYZED BY: 972

WORK ORDER: 099-14-354
MATRIX: Water

REVIEWED BY: 1,027
D/T REVIEWED: 2017-03-29 11:48

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
51	Daily Calibration	2017-03-28 13:24	T:\GC_45\GC_45_data\2017\170328\17032829.d\Report.txt17032829

WORK ORDER: 17-03-1557
MATRIX: Soil

REVIEWED BY: 1,027
D/T REVIEWED: 2017-03-29 11:54

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
25	B-DU1-ISM1-8	2017-03-28 14:08	T:\GC_45\GC_45_data\2017\170328\17032810.d\Report.txt17032810
26	B-DU1-ISM2-8	2017-03-28 14:29	T:\GC_45\GC_45_data\2017\170328\17032811.d\Report.txt17032811
27	B-DU1-ISM3-8	2017-03-28 14:51	T:\GC_45\GC_45_data\2017\170328\17032812.d\Report.txt17032812

CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8015B (M)

CCV WORK ORDER: 099-14-354-51-5901

BATCH ID:

1612021007

170328A055

GC 45

ANALYZED BY: 972

D/T ANALYZED:

INITIAL:

CCV:

REVIEWED BY:

D/T REVIEWED:

2016-12-02 19:43

2017-03-28 13:24

DATA FILE: T:\GC_45\GC_45_data\2017\170328\17032829.d\Report.txt17032829

<u>COMPOUND NAME</u>	<u>TYPE</u>	<u>CALIB MODEL</u>	<u>MIN RF</u>	<u>AVG RF</u>	<u>CCV RF</u>	<u>AMOUNT</u>	<u>CCV CONC</u>	<u>CCV %D</u>	<u>CCV %D CL</u>	<u>STATUS</u>
TPH as Diesel	C	Avg Resp	0.00	80.022	81.349			-2	0-20	PASS

MIN RF: Method Specified Minimum Response Factor

Data File: /chem1/SVOA/GC_45.i/170328.b/17032829.d
 Report Date: 28-Mar-2017 17:22

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/170328.b/17032829.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 13:24
 Operator : 682 Inst ID: GC_45.i
 Smp Info : CCV D400 C28 50 L102516D
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/170328.b/8015d.m
 Meth Date : 28-Mar-2017 17:22 umd6 Quant Type: ESTD
 Cal Date : 28-DEC-2016 19:01 Cal File: 16122828.d
 Als bottle: 29 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: CCV_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.649-8.839			32539603	400.000	406.634
S 27 Diesel Range Organics	2.300-8.839			30028122	400.000	401.211
\$ 93 n-Octacosane	8.612	8.612	0.000	4662166	50.0000	54.047

Data File: /chem1/SVDR/GC_45.i/170328.b/17032829.d

Date : 28-MAR-2017 13:24

Client ID:

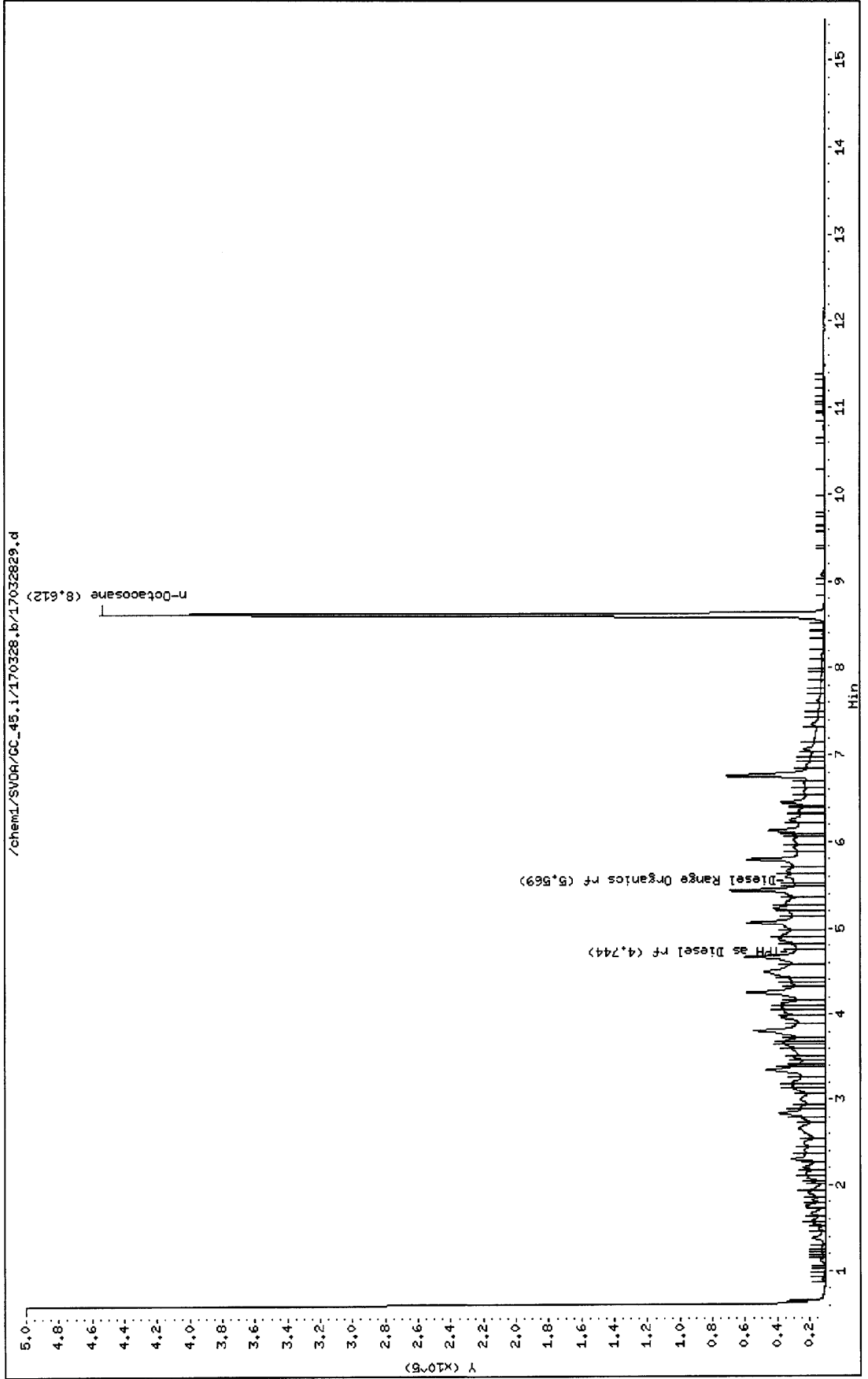
Sample Info: CCV D400 C28 90 L102516D

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_45.i/170328.b/17032819.d
 Report Date: 28-Mar-2017 18:06

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/170328.b/17032819.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 17:24
 Operator : 682 Inst ID: GC_45.i
 Smp Info : CCV D400 C28 50 L102516D
 Misc Info :
 Comment :
 Method : /chem1/SVOA/GC_45.i/170328.b/8015d.m
 Meth Date : 28-Mar-2017 18:06 umd6 Quant Type: ESTD
 Cal Date : 28-DEC-2016 19:01 Cal File: 16122828.d
 Als bottle: 19 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: CCV_D-DRO.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.649-8.839			33033904	400.000	412.811
S 27 Diesel Range Organics	2.300-8.839			30923192	400.000	413.171
\$ 93 n-Octacosane	8.614	8.614	0.000	4727662	50.0000	54.806



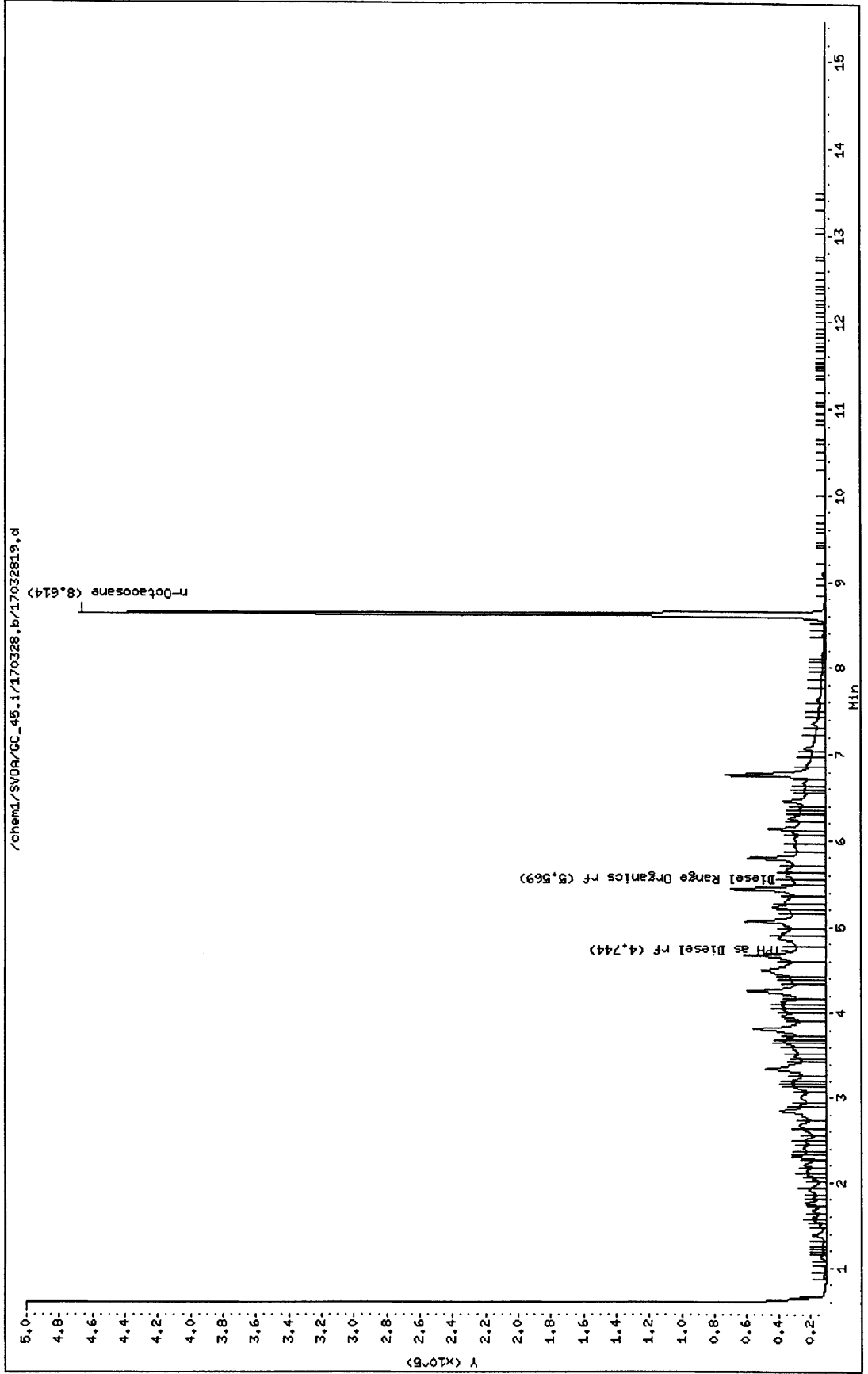
Data File: /chem1/SV0A/GC_45.i/170328.b/17032819.d
Date : 28-MAR-2017 17:24
Client ID:
Sample Info: CCV B400 C28 50 L102516D

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_45.i/170328.b/17032826.d
 Report Date: 03/29/2017 10:31

Eurofins CalScience
 Calibration Verification Report

Instrument ID: GC_45.i Injection Date and Time: 28-MAR-2017 19:54
 Sample Name: CCV D400 C28 50 L102516D Initial Calibration Date(s): 18-OCT-2016 29-MAR-2017
 Sublist used: CCV_D-DRO.sub Initial Calibration Time(s): 12:03 07:24
 Method used: /chem1/SVOA/GC_45.i/170328.b/8015d.m

Target Compounds	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift /Drift	Max%D	Curve Type
TPH as Diesel	80021.702	81209.305	0.00	-1 15	Averaged	
Diesel Range Organics	74843.538	75952.245	0.00	-1 15	Averaged	
Surrogate Standards	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift /Drift	Max%D	Curve Type
n-Octacosane	86260.794	93539.560	0.00	-8 20	Averaged	

page 1

Data File: /chem1/SVOA/GC_45.i/170328.b/17032826.d
Report Date: 29-Mar-2017 10:28

Page 1

Eurofins Calscience

EPA 8015B(M)

Data file : /chem1/SVOA/GC_45.i/170328.b/17032826.d
Lab Smp Id:
Inj Date : 28-MAR-2017 19:54
Operator : 682 Inst ID: GC_45.i
Smp Info : CCV D400 C28 50 L102516D
Misc Info :
Comment :
Method : /chem1/SVOA/GC_45.i/170328.b/8015d.m
Meth Date : 29-Mar-2017 10:27 umd6 Quant Type: ESTD
Cal Date : 29-MAR-2017 07:24 Cal File: 17032863.d
Als bottle: 26 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: HP Genie Compound Sublist: CCV_D-DRO.sub
Target Version: 3.50
Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RT	RESPONSE	AMOUNTS	
						CAL-AMT (ppm)	ON-COL (ppm)
S 15 TPH as Diesel	0.649-8.839				32483722	400.000	405.936
S 27 Diesel Range Organics	2.300-8.839				30380898	400.000	405.925
\$ 93 n-Octacosane	8.612	8.612	0.000		4676978	50.0000	54.219

Data File: /chem1/SV0A/GC_45.i/170328.b/17032826.d

Date : 28-MAR-2017 19:54

Client ID:

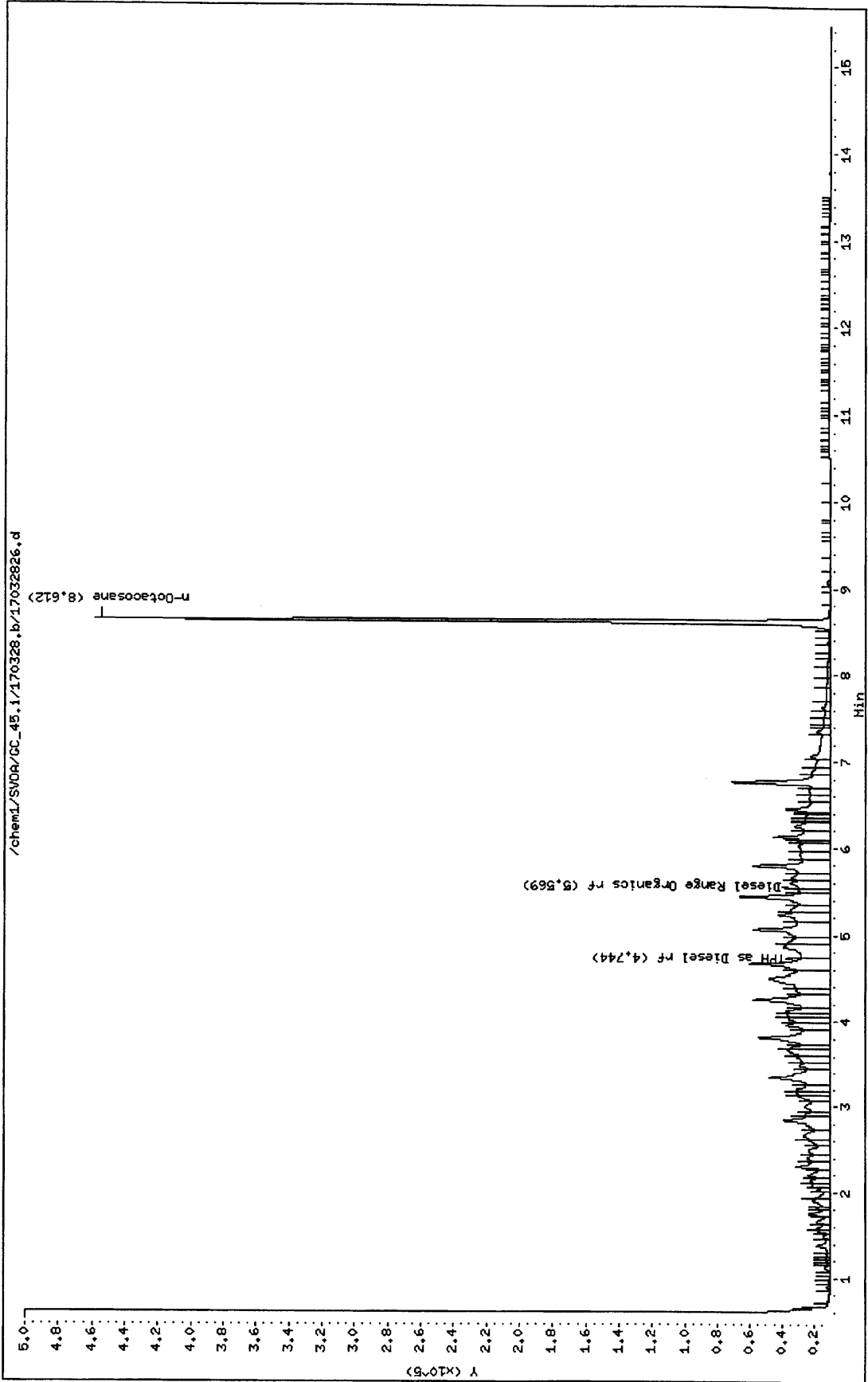
Sample Info: CCV D400 C28 50 L102516D

Instrument: GC_45.i

Operator: 682

Column diameter: 2.00

Column phase:



=====
 External Standard Report
 =====

Data File Name : /chem1/SVOA/GC_45/170328/17032802.d
 Page Number :
 Operator : 682 Vial Number : Vial 2
 Instrument : GC 45 Injection Number : 2
 Sample Name : C6-C44 L110816A Sequence Line : 0
 Instrument Method: 8015d.m
 Acquired on : 28 MAR 17 09:50
 Report Created on: 28-MAR-17 14:07 Compound Sublist : all
 Software Revision: Target 3.50

Sig. 1 in /chem1/SVOA/GC_45.i/170328.b/17032802.d

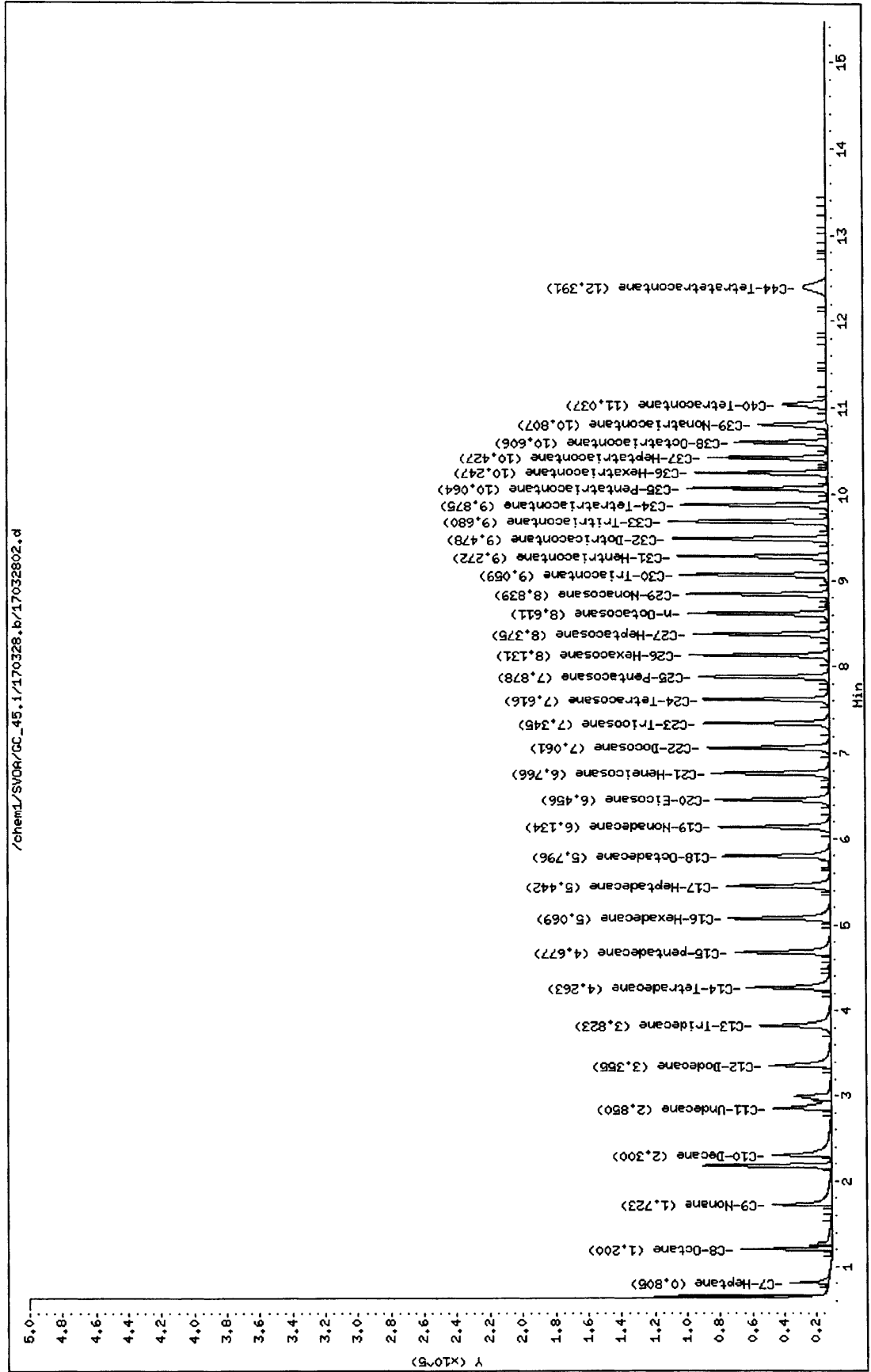
RT Range	Exp RT	DLT RT	Response	ppm	Compound
0.649	8.628	7.979	91768.00	0.00	C6-Hexane
0.805	8.628	7.823	194316.00	0.00	C7-Heptane
1.200	8.628	7.428	366057.00	0.00	C8-Octane
1.723	8.628	6.905	543918.00	0.00	C9-Nonane
2.300	8.628	6.328	649581.00	0.00	C10-Decane
2.850	8.628	5.778	531998.00	0.00	C11-Undecane
3.355	8.628	5.273	642332.00	0.00	C12-Dodecane
3.823	8.628	4.805	727422.00	0.00	C13-Tridecane
4.263	8.628	4.366	710780.00	0.00	C14-Tetradecane
4.677	8.628	3.952	758796.00	0.00	C15-pentadecane
5.069	8.628	3.559	781474.00	0.00	C16-Hexadecane
5.442	8.628	3.187	784796.00	0.00	C17-Heptadecane
5.796	8.628	2.832	788986.00	0.00	C18-Octadecane
6.134	8.628	2.495	800870.00	0.00	C19-Nonadecane
6.456	8.628	2.172	811307.00	0.00	C20-Eicosane
6.766	8.628	1.863	814573.00	0.00	C21-Heneicosane
7.061	8.628	1.568	827132.00	0.00	C22-Docosane
7.345	8.628	1.284	833698.00	0.00	C23-Tricosane
7.616	8.628	1.012	831196.00	0.00	C24-Tetracosane
7.878	8.628	0.750	840236.00	0.00	C25-Pentacosane
8.131	8.628	0.498	877188.00	0.00	C26-Hexacosane
8.375	8.628	0.254	871681.00	0.00	C27-Heptacosane
8.611	8.628	0.017	899905.00	10.43	n-Octacosane
8.839	8.628	-0.210	917053.00	0.00	C29-Nonacosane
9.059	8.628	-0.430	956005.00	0.00	C30-Triacontane
9.272	8.628	-0.644	974028.00	0.00	C31-Hentriacontane
9.478	8.628	-0.850	1008553.00	0.00	C32-Dotriacontane
9.680	8.628	-1.052	1023195.00	0.00	C33-Tritriacontane
9.875	8.628	-1.247	991845.00	0.00	C34-Tetratriacontane
10.064	8.628	-1.436	924018.00	0.00	C35-Pentatriacontane
10.247	8.628	-1.618	904871.00	0.00	C36-Hexatriacontane
10.427	8.628	-1.798	781874.00	0.00	C37-Heptatriacontane
10.606	8.628	-1.978	714520.00	0.00	C38-Octatriacontane
10.807	8.628	-2.178	582222.00	0.00	C39-Nonatriacontane
11.037	8.628	-2.408	472440.00	0.00	C40-Tetracontane
12.391	8.628	-3.762	479640.00	0.00	C44-Tetratetracontane

End of File

Data File: /chem1/SVDA/GC_45.i/170328.b/17032802.d
Date : 28-MAR-2017 09:50
Client ID:
Sample Info: C6-C44 L110816A

Instrument: GC_45.i
Operator: 682
Column diameter: 2.00

Column phase:



EPA 8015B (M) Diesel + Motor Oil

RUN LOGS

Line	Vial	File	Name	Method	Acquired
1	1	16120201	BLANK	8015D	02-Dec-16, 17:34:02
2	1	1612020102	BLANK	8015D	02-Dec-16, 17:55:52
3	1	1612020103	BLANK	8015D	02-Dec-16, 18:17:52
4	2	16120202	C6-C44 L110816A	8015D	02-Dec-16, 18:40:02
5	5	16120205	CCV D400 C28 50 L102516D	8015D	02-Dec-16, 19:00:47
6	6	16120206	CCV MO400 L090716D	8015D	02-Dec-16, 19:21:33
7	7	16120207	ICAL D5 C28 0.625 L102516B	8015D	02-Dec-16, 19:43:29
8	8	16120208	ICAL D200 C28 25 L102516C	8015D	02-Dec-16, 20:05:44
9	9	16120209	ICAL D400 C28 50 L102516D	8015D	02-Dec-16, 20:27:44
10	10	16120210	ICAL D800 C28 100 L102516E	8015D	02-Dec-16, 20:48:51
11	11	16120211	ICAL D1600 C28 200 L102516F	8015D	02-Dec-16, 21:09:46
12	12	16120212	ICV D400 C28 50 L102516G	8015D	02-Dec-16, 21:31:10
13	13	16120213	ICAL CO200 L111516C	8015D	02-Dec-16, 21:52:20
14	14	16120214	ICAL CO400 L111516D	8015D	02-Dec-16, 22:14:41
15	15	16120215	ICAL CO20 L111516B	8015D	02-Dec-16, 22:36:25
16	16	16120216	ICAL CO600 L111516E	8015D	02-Dec-16, 22:58:04
17	17	16120217	ICAL CO800 L111516F	8015D	02-Dec-16, 23:19:13
18	18	16120218	ICV CO400 L111516G	8015D	02-Dec-16, 23:40:58
19	19	16120219	ICAL GD5 L091416A	8015D	03-Dec-16, 00:02:10
20	20	16120220	ICAL GD200 L091416B	8015D	03-Dec-16, 00:23:54
21	21	16120221	ICAL GD400 L091416C	8015D	03-Dec-16, 00:46:08
22	22	16120222	ICAL GD800 L091416D	8015D	03-Dec-16, 01:08:26
23	23	16120223	ICAL GD1600 L091416E	8015D	03-Dec-16, 01:29:41
24	24	16120224	ICV GD400 L091416F	8015D	03-Dec-16, 01:50:38
25	2	16120225	C6-C44 L110816A	8015D	03-Dec-16, 02:11:22
26	26	16120226	CCV D400 C28 50 L102516D	8015D	03-Dec-16, 02:32:10
27	27	16120227	CCV GD400 L091416C	8015D	03-Dec-16, 02:53:05
28	28	16120228	MB 16120213	8015D	03-Dec-16, 03:13:44
29	29	16120229	GDLCS 16120213	8015D	03-Dec-16, 03:34:30
30	30	16120230	GDMS 16-12-0003-17	8015D	03-Dec-16, 03:55:45
31	31	16120231	GDMSD 16-12-0003-17	8015D	03-Dec-16, 04:17:45
32	32	16120232	16-12-0003-13	8015D	03-Dec-16, 04:39:44
33	33	16120233	16-12-0003-17	8015D	03-Dec-16, 05:02:19
34	34	16120234	16-12-0003-21	8015D	03-Dec-16, 05:24:56
35	35	16120235	CCV D400 C28 50 L102516D	8015D	03-Dec-16, 05:47:26
36	36	16120236	CCV GD400 L091416C	8015D	03-Dec-16, 06:09:53
37	37	16120237	MB 16120104	8015D	03-Dec-16, 06:31:28
38	38	16120238	LCS 16120104	8015D	03-Dec-16, 06:54:20
39	39	16120239	LCSD 16120104	8015D	03-Dec-16, 07:16:51
40	40	16120240	16-11-2665-1	8015D	03-Dec-16, 07:39:49
41	41	16120241	16-11-2665-2	8015D	03-Dec-16, 08:02:28
42	42	16120242	16-11-2665-3	8015D	03-Dec-16, 08:23:23
43	43	16120243	16-11-2665-4	8015D	03-Dec-16, 08:44:14
44	44	16120244	16-11-2665-5	8015D	03-Dec-16, 09:05:55
45	45	16120245	16-11-2665-6	8015D	03-Dec-16, 09:28:11
46	46	16120246	16-11-2665-7	8015D	03-Dec-16, 09:49:18
47	47	16120247	16-11-2665-8	8015D	03-Dec-16, 10:10:09
48	48	16120248	16-11-2665-9	8015D	03-Dec-16, 10:32:03
49	49	16120249	16-11-2606-1	8015D	03-Dec-16, 10:54:08
50	50	16120250	CCV D400 C28 50 L102516D	8015D	03-Dec-16, 11:16:34
51	51	16120251	16-11-2607-1	8015D	03-Dec-16, 11:38:13
52	52	16120252	16-11-2608-1	8015D	03-Dec-16, 11:59:20
53	2	16120253	C6-C44 L110816A	8015D	03-Dec-16, 12:21:29
54	54	16120254	CCV D400 C28 50 L102516D	8015D	03-Dec-16, 12:43:37
55	55	16120255	MB 16113008	8015D	03-Dec-16, 13:04:44
56	56	16120256	LCS 16113008	8015D	03-Dec-16, 13:25:48
57	57	16120257	LCSD 16113008	8015D	03-Dec-16, 13:48:06
58	58	16120258	16-11-2371-1	8015D	03-Dec-16, 14:10:26
59	59	16120259	16-11-2371-2	8015D	03-Dec-16, 14:33:03

10078

A085

A086

A087

A075

12/13/16
 42
 45
 8015
 94
 Logbook ID: 45

Return to Contents

Line	Vial	File	Name	Method	Acquired
1	1	17032800	BLANK	8015D	28-Mar-17, 09:06:52
2	1	17032801	BLANK	8015D	28-Mar-17, 09:28:38
3	2	17032802	C6-C44 L110816A	8015D	28-Mar-17, 09:50:34
4	3	17032803	CCV D400 C28 50 L102516D A031	8015D	28-Mar-17, 10:13:03
5	4	17032804	CCV MO400 L030317D	8015D	28-Mar-17, 10:34:41
6	5	17032805	17-03-1532-17 RB	8015D	28-Mar-17, 10:55:43
7	6	17032806	MB 17032412	8015D	28-Mar-17, 11:16:37
8	7	17032807	LCS 17032412	8015D	28-Mar-17, 11:37:35
9	27	17032827	17-03-1130-7 CONF	8015D	28-Mar-17, 11:59:52
10	28	17032828	17-03-1130-18 CONF	8015D	28-Mar-17, 12:21:29
11	8	17032808	MS 17-03-1557-27	8015D	28-Mar-17, 12:42:28
12	9	17032809	MSD 17-03-1557-27	8015D	28-Mar-17, 13:03:23
13	29	17032829	CCV D400 C28 50 L102516D A055	8015D	28-Mar-17, 13:24:59
14	42	17032842	17-03-1532-17 RB	8015D	28-Mar-17, 13:47:27
15	10	17032810	17-03-1557-25	8015D	28-Mar-17, 14:08:47
16	11	17032811	17-03-1557-26	8015D	28-Mar-17, 14:29:39
17	12	17032812	17-03-1557-27	8015D	28-Mar-17, 14:51:06
18	13	17032813	17-03-1556-25	8015D	28-Mar-17, 15:13:32
19	14	17032814	17-03-1556-26	8015D	28-Mar-17, 15:34:46
20	15	17032815	17-03-1556-27	8015D	28-Mar-17, 15:55:46
21	16	17032816	17-03-1418-25	8015D	28-Mar-17, 16:17:52
22	17	17032817	17-03-1418-26	8015D	28-Mar-17, 16:39:35
23	18	17032818	17-03-1418-27	8015D	28-Mar-17, 17:00:37
24	19	17032819	CCV D400 C28 50 L102516D A056	8015D	28-Mar-17, 17:24:04
25	20	17032820	17-03-1558-25	8015D	28-Mar-17, 17:46:42
26	21	17032821	17-03-1558-26	8015D	28-Mar-17, 18:07:57
27	22	17032822	17-03-1558-27	8015D	28-Mar-17, 18:28:51
28	23	17032823	17-03-1417-25	8015D	28-Mar-17, 18:51:02
29	24	17032824	17-03-1417-26	8015D	28-Mar-17, 19:12:37
30	25	17032825	17-03-1417-27	8015D	28-Mar-17, 19:33:39
31	26	17032826	CCV D400 C28 50 L102516D	8015D	28-Mar-17, 19:54:54
32	31	17032831	CCV D80 C28 10 L103116B A057	8015D	28-Mar-17, 20:17:21
33	32	17032832	CCV MO80 L030717D	8015D	28-Mar-17, 20:38:59
34	33	17032833	17-03-1588-3 RB	8015D	28-Mar-17, 20:59:52
35	34	17032834	CCV D80 C28 10 L103116B	8015D	28-Mar-17, 21:21:01
36	35	17032835	CCV MO80 L030717D	8015D	28-Mar-17, 21:43:26
37	36	17032836	ICAL MO25 L030317B	8015D	28-Mar-17, 22:04:48
38	37	17032837	ICAL MO200 L030317C	8015D	28-Mar-17, 22:25:51
39	38	17032838	ICAL MO400 L030317D	8015D	28-Mar-17, 22:47:24
40	39	17032839	ICAL MO600 L030317E	8015D	28-Mar-17, 23:09:31
41	40	17032840	ICAL MO800 L030317F	8015D	28-Mar-17, 23:31:49
42	41	17032841	ICV MO400 L030317G	8015D	28-Mar-17, 23:53:04
43	43	17032843	CCV D400 C28 50 L102516D A058	8015D	29-Mar-17, 00:14:00
44	44	17032844	CCV MO400 L030317D	8015D	29-Mar-17, 00:34:59
45	45	17032845	MB 17032803/04/05/06/14	8015D	29-Mar-17, 00:56:13
46	46	17032846	LCS 17032803/14	8015D	29-Mar-17, 01:18:26
47	47	17032847	LCSD 17032803/14	8015D	29-Mar-17, 01:40:28
48	48	17032848	MOLCS 17032804	8015D	29-Mar-17, 02:01:39
49	49	17032849	MOLCSD 17032804	8015D	29-Mar-17, 02:22:37
50	50	17032850	17-03-1973-1 5X	8015D	29-Mar-17, 02:43:46
51	51	17032851	17-03-1993-1	8015D	29-Mar-17, 03:05:20
52	52	17032852	17-03-1823-3	8015D	29-Mar-17, 03:27:25
53	53	17032853	CCV D400 C28 50 L102516D	8015D	29-Mar-17, 03:48:45
54	54	17032854	CCV MO400 L030317D	8015D	29-Mar-17, 04:09:39

Reviewed/Assign to Logbook Date: 3/29/17
Analysis: 8015 Chemist ID: 682
Lockbox Page: 04 Instrument ID: 45

I001

Return to Contents

EPA 8015B (M)
Diesel + Motor Oil

PREPARATION LOGS

Analysis Method (EPA Method): TPH 8015 NWTPH 8015
 Extraction Method (EPA Method): 3510 3511 3550
 Analyst ID#: Measuring Sample- 1112 Start Extraction- 605 Blow Down- Clean Up-
 Matrix: Solid Aqueous Oil Wipe/Filter Balance ID#: 15 Sand or Filter ID#: 68-19-20
 Extraction Start Date & Time: 3-24-17 18:00 Extractions End Date & Time: 3-24-17 20:00
 Drying Agent & ID#: Na₂SO₄ 66-23-20 Spike Added to: LCS LCSD MS MSD
 Surrogate Std ID# & Volume Added (mL): 2011717A 0.5
 Spike Std ID# & Volume Added (mL): D: 2102416B 0.2 MO:
 Extraction Solvent & ID#: MeCl₂ 507-58-1 Elution Solvent ID# & Volume (mL): 10
 Reverse Surrogate ID#: & Volume Added (mL):

Clean Up Start Date & Time: Clean Up End Date & Time:
 SGC Clean Up: 1. CEL 0.5g 2. 3630 - 2g 3. 3630 - 10g SPE Cartridge ID#:
 Clean Up Solvent ID#: Silica Gel ID#:

QC Batch #: 17032412	Test Analyte	Sample W (g) / V (mL)		SCG Clean Up	Comments
		Initial	Final		
MB	D+M	10.0	10	<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
LCS		10.0		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
MS 17-03-1557-27C		10.1		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
MSD		10.0		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
17-03-1417-25C	D+MO	9.86		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-26	10.2		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-27	9.86		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
17-03-1418-25C	D+MO	9.98		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-26	9.86		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-27	10.0		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
17-03-1419-25C	D+MO	10.0		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-26	9.80		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-27	9.87		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
17-03-1556-25C	D+MO	10.1		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-26	10.0		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-27	10.2		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
17-03-1557-25C	D+MO	10.1		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-26	10.2		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-27	10.1		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
17-03-1558-25C	D+MO	10.2		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-26	9.84		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
	-27	10.1		<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	
				<input type="checkbox"/> NA <input type="checkbox"/> 1 <input type="checkbox"/> 2 <input type="checkbox"/> 3	



EPA 6010B ICP Metals (Solid)

RAW DATA

EPA 6010B ICP Metals (Solid)

Initial Calibration

ICV/ICB
CCV/CCB
ICSA/B

Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte Name	Initial Calibration Verification				
	True	ICV-1		Control Limit	Comment
		Observed	%D		
Silver	0.500000	0.491810	2	+/-10	
Arsenic	5.000000	4.806104	4	+/-10	
Barium	1.000000	0.977623	2	+/-10	
Beryllium	0.500000	0.485431	3	+/-10	
Cadmium	1.500000	1.492932	0	+/-10	
Cobalt	1.000000	1.030879	-3	+/-10	
Chromium	0.400000	0.395313	1	+/-10	
Copper	1.000000	0.986354	1	+/-10	
Molybdenum	2.500000	2.376245	5	+/-10	
Nickel	0.400000	0.406362	-2	+/-10	
Lead	5.000000	5.071895	-1	+/-10	
Antimony	2.000000	1.903690	5	+/-10	
Selenium	2.000000	1.927283	4	+/-10	
Thallium	2.000000	1.940131	3	+/-10	
Vanadium	1.000000	0.976319	2	+/-10	
Zinc	1.500000	1.515325	-1	+/-10	

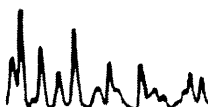
Report Time: 3/30/2017 6:02:51 PM

ICV-1 File: ICP-M072816C

Analysis Time: 3/29/2017 9:46:40 AM

Note: Note: %D= (True-Observed) / True x 100%

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

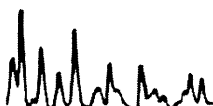
Analyte	Initial Calibration Blank		
	ICB-1	RL	Comment
Silver	-0.000620	0.005000	
Arsenic	0.003618	0.010000	
Barium	-0.000096	0.010000	
Beryllium	0.000030	0.010000	
Cadmium	0.000421	0.010000	
Cobalt	0.000260	0.010000	
Chromium	0.000106	0.010000	
Copper	-0.000241	0.010000	
Molybdenum	0.000250	0.010000	
Nickel	0.000954	0.010000	
Lead	0.000993	0.010000	
Antimony	-0.001530	0.015000	
Selenium	0.002268	0.015000	
Thallium	-0.001209	0.015000	
Vanadium	0.000021	0.010000	
Zinc	0.000192	0.010000	

Report Time: 3/30/2017 6:02:51 PM

ICB-1 File: ICB-R12091601

Analysis Time: 3/29/2017 9:47:39 AM

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

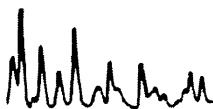
Concentration Unit: mg/L

Analyte	Interference Check						Comment
	ICS-A-1		ICS-AB-1			Control Limit	
	Observed	Control Limit	True	Observed	%D		
Silver	-0.000235	0.005000	0.300000	0.318524	-6	+/-20	
Arsenic	-0.004571	0.010000	1.000000	1.029879	-3	+/-20	
Barium	0.002481	0.010000	0.300000	0.310564	-4	+/-20	
Beryllium	-0.000002	0.010000	0.100000	0.105207	-5	+/-20	
Cadmium	0.002364	0.010000	0.300000	0.308962	-3	+/-20	
Cobalt	0.000896	0.010000	0.300000	0.313328	-4	+/-20	
Chromium	-0.001851	0.010000	0.300000	0.316153	-5	+/-20	
Copper	-0.001849	0.010000	0.300000	0.319465	-6	+/-20	
Molybdenum	0.001841	0.010000	0.300000	0.305124	-2	+/-20	
Nickel	0.002170	0.010000	0.300000	0.316239	-5	+/-20	
Lead	-0.005630	0.010000	1.000000	1.030376	-3	+/-20	
Antimony	-0.011601	0.015000	1.000000	0.984318	2	+/-20	
Selenium	-0.006483	0.015000	0.500000	0.515228	-3	+/-20	
Thallium	-0.005541	0.015000	1.000000	1.020640	-2	+/-20	
Vanadium	0.001300	0.010000	0.300000	0.313592	-5	+/-20	
Zinc	0.000752	0.010000	0.300000	0.318114	-6	+/-20	

Report Time: 3/30/2017 6:02:51 PM

ICS-A-1 File: ICS_A - M110116B Analysis Time: 3/29/2017 9:48:40 AM
ICS-AB-1 File: ICS_AB - M110116A Analysis Time: 3/29/2017 9:49:46 AM

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte	Continuing Calibration Verification						
	True	CCV-1		CCV-2		Control Limit	Comment
		Observed	%D	Observed	%D		
Silver	0.375000	0.364909	3	0.366493	2	+/-10	
Arsenic	3.750000	3.683143	2	3.748817	0	+/-10	
Barium	7.500000	7.466781	0	7.509775	0	+/-10	
Beryllium	0.562500	0.554684	1	0.552267	2	+/-10	
Cadmium	0.750000	0.765439	-2	0.772031	-3	+/-10	
Cobalt	1.875000	1.943705	-4	1.958211	-4	+/-10	
Chromium	0.600000	0.605527	-1	0.610239	-2	+/-10	
Copper	0.937500	0.914562	2	0.915045	2	+/-10	
Molybdenum	0.600000	0.590266	2	0.598323	0	+/-10	
Nickel	0.600000	0.605696	-1	0.612071	-2	+/-10	
Lead	3.750000	3.831537	-2	3.868642	-3	+/-10	
Antimony	4.500000	4.295910	5	4.356098	3	+/-10	
Selenium	1.500000	1.501677	0	1.536173	-2	+/-10	
Thallium	1.500000	1.503777	0	1.524691	-2	+/-10	
Vanadium	1.875000	1.884077	0	1.887622	-1	+/-10	
Zinc	2.500000	2.532855	-1	2.558497	-2	+/-10	

Report Time: 3/30/2017 6:02:51 PM

CCV-1 File: CCV= STD3x0.5

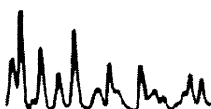
Analysis Time: 3/29/2017 4:56:49 PM

CCV-2 File: CCV= STD3x0.5

Analysis Time: 3/29/2017 5:14:09 PM

Note: Note: %D= (True-Observed) / True x 100%

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte	Continuing Calibration Blank			Qualifier
	CCB-1	CCB-2	RL	
Silver	0.000093	-0.000166	0.005000	
Arsenic	0.001411	0.003119	0.010000	
Barium	0.002932	0.003744	0.010000	
Beryllium	0.000202	0.000244	0.010000	
Cadmium	0.000400	0.000604	0.010000	
Cobalt	0.000682	0.001410	0.010000	
Chromium	0.000346	0.000524	0.010000	
Copper	-0.000270	0.000140	0.010000	
Molybdenum	0.000265	0.000834	0.010000	
Nickel	0.001092	0.001622	0.010000	
Lead	-0.000451	0.002461	0.010000	
Antimony	0.000419	0.002032	0.015000	
Selenium	0.004360	0.000890	0.015000	
Thallium	0.000833	0.003304	0.015000	
Vanadium	0.001293	0.001167	0.010000	
Zinc	-0.001586	-0.002259	0.010000	

Report Time: 3/30/2017 6:02:51 PM

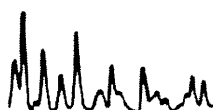
CCB-1 File: CCB-R12091601

Analysis Time: 3/29/2017 4:57:41 PM

CCB-2 File: CCB-R12091601

Analysis Time: 3/29/2017 5:15:02 PM

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte	Continuing Calibration Verification						Comment
	True	CCV-3		CCV-4		Control Limit	
		Observed	%D	Observed	%D		
Silver	0.375000	0.365890	2	0.364320	3	+/-10	
Arsenic	3.750000	3.639842	3	3.738350	0	+/-10	
Barium	7.500000	7.486822	0	7.490872	0	+/-10	
Beryllium	0.562500	0.552547	2	0.554388	1	+/-10	
Cadmium	0.750000	0.768964	-3	0.769795	-3	+/-10	
Cobalt	1.875000	1.953335	-4	1.946074	-4	+/-10	
Chromium	0.600000	0.605342	-1	0.606826	-1	+/-10	
Copper	0.937500	0.927333	1	0.913727	3	+/-10	
Molybdenum	0.600000	0.583597	3	0.601492	0	+/-10	
Nickel	0.600000	0.596226	1	0.612980	-2	+/-10	
Lead	3.750000	3.864061	-3	3.805985	-1	+/-10	
Antimony	4.500000	4.253002	5	4.379475	3	+/-10	
Selenium	1.500000	1.487794	1	1.520864	-1	+/-10	
Thallium	1.500000	1.479459	1	1.518761	-1	+/-10	
Vanadium	1.875000	1.881908	0	1.880491	0	+/-10	
Zinc	2.500000	2.584336	-3	2.548002	-2	+/-10	

Report Time: 3/30/2017 6:02:51 PM

CCV-3 File: CCV= STD3x0.5

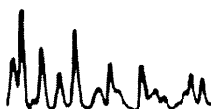
Analysis Time: 3/29/2017 5:25:20 PM

CCV-4 File: CCV= STD3x0.5

Analysis Time: 3/29/2017 6:49:54 PM

Note: Note: %D= (True-Observed) / True x 100%

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte	Continuing Calibration Blank			Qualifier
	CCB-3	CCB-4	RL	
Silver	0.000243	-0.001144	0.005000	
Arsenic	0.002115	0.006971	0.010000	
Barium	0.005343	0.006558	0.010000	
Beryllium	0.000338	0.000452	0.010000	
Cadmium	0.000642	0.000805	0.010000	
Cobalt	0.001688	0.002144	0.010000	
Chromium	0.000661	0.000733	0.010000	
Copper	0.007025	0.001044	0.010000	
Molybdenum	0.001400	0.001459	0.010000	
Nickel	0.001460	0.001553	0.010000	
Lead	0.002519	0.002238	0.010000	
Antimony	0.003017	0.009395	0.015000	
Selenium	-0.000616	0.001094	0.015000	
Thallium	0.000804	0.010618	0.015000	
Vanadium	0.001323	0.001626	0.010000	
Zinc	0.018746	-0.000224	0.010000	

Report Time: 3/30/2017 6:02:51 PM

CCB-3 File: CCB-R12091601

Analysis Time: 3/29/2017 5:26:13 PM

CCB-4 File: CCB-R12091601

Analysis Time: 3/29/2017 6:50:47 PM

01/22/2014 Revision



Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte	Continuing Calibration Verification						
	True	CCV-5				Control Limit	Comment
		Observed	%D	Observed	%D		
Silver	0.375000	0.365897	2			+/-10	
Arsenic	3.750000	3.766058	0			+/-10	
Barium	7.500000	7.590027	-1			+/-10	
Beryllium	0.562500	0.557685	1			+/-10	
Cadmium	0.750000	0.781077	-4			+/-10	
Cobalt	1.875000	1.977082	-5			+/-10	
Chromium	0.600000	0.612816	-2			+/-10	
Copper	0.937500	0.923631	1			+/-10	
Molybdenum	0.600000	0.598362	0			+/-10	
Nickel	0.600000	0.616309	-3			+/-10	
Lead	3.750000	3.912740	-4			+/-10	
Antimony	4.500000	4.368196	3			+/-10	
Selenium	1.500000	1.540234	-3			+/-10	
Thallium	1.500000	1.525337	-2			+/-10	
Vanadium	1.875000	1.890963	-1			+/-10	
Zinc	2.500000	2.601649	-4			+/-10	

Report Time: 3/30/2017 6:02:51 PM

CCV-5 File: CCV= STD3x0.5

Analysis Time: 3/29/2017 7:02:53 PM

Note: Note: %D= (True-Observed) / True x 100%

01/22/2014 Revision

Work Order No: 17-03-1557

Instrument ID: ICP 7300

Concentration Unit: mg/L

Analyte	Continuing Calibration Blank			Qualifier
	CCB-5		RL	
Silver	-0.000051		0.005000	
Arsenic	0.006579		0.010000	
Barium	0.006934		0.010000	
Beryllium	0.000482		0.010000	
Cadmium	0.000909		0.010000	
Cobalt	0.001929		0.010000	
Chromium	0.000660		0.010000	
Copper	0.000713		0.010000	
Molybdenum	0.001566		0.010000	
Nickel	0.001572		0.010000	
Lead	0.002269		0.010000	
Antimony	0.003698		0.015000	
Selenium	-0.001731		0.015000	
Thallium	0.007080		0.015000	
Vanadium	0.001911		0.010000	
Zinc	-0.000648		0.010000	

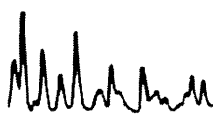
Report Time: 3/30/2017 6:02:51 PM

CCB-5 File: CCB-R12091601

Analysis Time: 3/29/2017 7:03:52 PM

01/22/2014 Revision

Return to Contents



=====
Analysis Begun

Start Time: 3/29/2017 9:42:08 AM
Logged In Analyst: Oscar Gomez 935
Spectrometer: Optima 7300 DV, S/N 77c8120401

Plasma On Time: 3/29/2017 9:19:17 AM
Technique: ICP Continuous
Autosampler: ESI

Sample Information File: C:\Documents and Settings\All Users\PerkinElmer\ICP\Data\Sample Information\
17032901.sif

Batch ID:
Results Data Set: 170329C 1
Results Library: W:\pe\7300\Results\results.mdb

=====
Sequence No.: 1
Sample ID: Cal blankR12091601_935
Analyst:
Initial Sample Wt:
Dilution:
Wash Time:
Autosampler Location: 1
Date Collected: 3/29/2017 9:42:24 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:

Mean Data: Cal blankR12091601_935

Analyte	Mean Corrected Intensity	Std.Dev.	RSD	Conc.	Units	Calib
Tb 384	79882.1	426.02	0.53%	100.0	%	
Tb 350	126132.3	1450.08	1.15%	100.0	%	
Ag 328.068*†	-1109.6	15.42	1.39%	[0.00]	mg/L	
Al 308.215*†	-2344.4	51.96	2.22%	[0.00]	mg/L	
As 188.979†	0.1	0.40	305.01%	[0.00]	mg/L	
As 193.696*†	-3.2	5.42	170.90%	[0.00]	mg/L	
B 249.677*†	-719.8	6.20	0.86%	[0.00]	mg/L	
Ba 233.527*†	-217.4	10.10	4.65%	[0.00]	mg/L	
Be 313.042*†	-398.3	118.04	29.63%	[0.00]	mg/L	
Ca 317.933*†	5.8	0.49	8.35%	[0.00]	mg/L	
Cd 226.502*†	-1.6	12.07	771.16%	[0.00]	mg/L	
Cd 228.802†	10.9	2.33	21.37%	[0.00]	mg/L	
Co 228.616*†	-67.4	9.47	14.05%	[0.00]	mg/L	
Cr 267.716*†	302.8	15.66	5.17%	[0.00]	mg/L	
Cu 324.752*†	1754.3	23.32	1.33%	[0.00]	mg/L	
Fe 273.955*†	-199.7	9.87	4.94%	[0.00]	mg/L	
K 766.490*†	1041.3	116.35	11.17%	[0.00]	mg/L	
Mg 279.077*†	-5115.6	30.88	0.60%	[0.00]	mg/L	
Mn 257.610*†	-49.5	19.18	38.75%	[0.00]	mg/L	
Mo 202.031*†	-33.8	5.89	17.42%	[0.00]	mg/L	
Na 589.592*†	755.1	78.40	10.38%	[0.00]	mg/L	
Ni 231.604*†	-71.1	12.46	17.53%	[0.00]	mg/L	
P 213.617*†	-105.6	4.60	4.36%	[0.00]	mg/L	
P 214.914†	-31.9	7.95	24.94%	[0.00]	mg/L	
Pb 220.353*†	-24.3	7.17	29.51%	[0.00]	mg/L	
Sb 206.836†	7.8	11.31	144.24%	[0.00]	mg/L	
Sb 217.582*†	-1.9	3.44	176.56%	[0.00]	mg/L	
Se 196.026*†	8.1	2.61	32.16%	[0.00]	mg/L	
Si 251.611*†	1405.3	3.05	0.22%	[0.00]	mg/L	
Sn 189.927*†	-64.3	2.08	3.24%	[0.00]	mg/L	
Sn 242.170†	-315.5	2.08	0.66%	[0.00]	mg/L	
Sr 407.771*†	60.3	7.88	13.08%	[0.00]	mg/L	
Ti 334.940†	27706.2	76.17	0.27%	[0.00]	mg/L	
Ti 336.121*†	-972.8	72.76	7.48%	[0.00]	mg/L	
Tl 190.801*†	-5.8	0.68	11.72%	[0.00]	mg/L	
V 292.402*†	132.7	20.06	15.12%	[0.00]	mg/L	
Zn 206.200*†	-34.9	0.39	1.11%	[0.00]	mg/L	
Zn 213.857*†	119.5	17.89	14.97%	[0.00]	mg/L	

User canceled analysis.

=====
Analysis BegunStart Time: 3/29/2017 9:43:25 AM
Logged In Analyst: Oscar Gomez 935
Spectrometer: Optima 7300 DV, S/N 77c8120401Plasma On Time: 3/29/2017 9:19:17 AM
Technique: ICP Continuous
Autosampler: ESISample Information File: C:\Documents and Settings\All Users\PerkinElmer\ICP\Data\Sample Information\
17032901.sifBatch ID:
Results Data Set: 170329C 1
Results Library: W:\pe\7300\Results\results.mdb=====
Sequence No.: 2
Sample ID: STD3-M111116A_935_ICP7300
Analyst:
Initial Sample Wt:
Dilution:
Wash Time:
Autosampler Location: 2
Date Collected: 3/29/2017 9:43:26 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:-----
Mean Data: STD3-M111116A_935_ICP7300

Analyte	Mean Corrected	Std.Dev.	RSD	Conc.	Units	Calib
Tb 384	70310.9	175.42	0.25%	88.02	%	
Tb 350	114006.6	129.66	0.11%	90.39	%	
Ag 328.068*†	132760.8	506.60	0.38%	[0.75]	mg/L	
Al 308.215*†	417970.9	1264.07	0.30%	[27.0]	mg/L	
As 188.979†	13072.6	17.28	0.13%	[7.50]	mg/L	
As 193.696*†	9184.5	5.51	0.06%	[7.50]	mg/L	
B 249.677*†	331734.3	3157.74	0.95%	[7.50]	mg/L	
Ba 233.527*†	2067399.5	35890.20	1.74%	[15.0]	mg/L	
Be 313.042*†	4061144.7	55774.45	1.37%	[1.125]	mg/L	
Ca 317.933*†	132985.6	2989.41	2.25%	[60.0]	mg/L	
Cd 226.502*†	95218.2	56.21	0.06%	[1.50]	mg/L	
Cd 228.802†	49766.0	136.18	0.27%	[1.50]	mg/L	
Co 228.616*†	89541.7	116.69	0.13%	[3.75]	mg/L	
Cr 267.716*†	121151.8	259.85	0.21%	[1.20]	mg/L	
Cu 324.752*†	451519.6	842.06	0.19%	[1.875]	mg/L	
Fe 273.955*†	154634.3	29.78	0.02%	[7.50]	mg/L	
K 766.490*†	181803.1	3712.57	2.04%	[54.0]	mg/L	
Mg 279.077*†	231349.7	410.69	0.18%	[15.0]	mg/L	
Mn 257.610*†	883422.8	14802.15	1.68%	[1.50]	mg/L	
Mo 202.031*†	10066.3	11.91	0.12%	[1.20]	mg/L	
Na 589.592*†	373282.3	7409.12	1.98%	[72.0]	mg/L	
Ni 231.604*†	25478.2	161.58	0.63%	[1.20]	mg/L	
P 213.617*†	19555.6	1.92	0.01%	[12.0]	mg/L	
P 214.914†	12138.0	23.23	0.19%	[12.0]	mg/L	
Pb 220.353*†	48071.5	334.82	0.70%	[7.50]	mg/L	
Sb 206.836†	15257.3	69.05	0.45%	[9.0]	mg/L	
Sb 217.582*†	15357.9	28.80	0.19%	[9.0]	mg/L	
Se 196.026*†	5293.3	26.37	0.50%	[3.0]	mg/L	
Si 251.611*†	421514.9	3936.53	0.93%	[12.0]	mg/L	
Sn 189.927*†	30377.8	53.27	0.18%	[6.0]	mg/L	
Sn 242.170†	9669.3	63.12	0.65%	[6.0]	mg/L	
Sr 407.771*†	222734.4	5110.22	2.29%	[0.60]	mg/L	
Ti 334.940†	975572.3	13341.42	1.37%	[1.20]	mg/L	
Ti 336.121*†	673724.1	1066.79	0.16%	[1.20]	mg/L	
Tl 190.801*†	4756.7	7.82	0.16%	[3.0]	mg/L	
V 292.402*†	468425.9	1116.44	0.24%	[3.75]	mg/L	
Zn 206.200*†	161669.7	701.58	0.43%	[5.0]	mg/L	
Zn 213.857*†	298383.1	319.81	0.11%	[5.0]	mg/L	

User canceled analysis.

=====
Analysis Begun

Start Time: 3/29/2017 9:46:38 AM Plasma On Time: 3/29/2017 9:19:17 AM
 Logged In Analyst: Oscar Gomez 935 Technique: ICP Continuous
 Spectrometer: Optima 7300 DV, S/N 77c8120401 Autosampler: ESI

Sample Information File: C:\Documents and Settings\All Users\PerkinElmer\ICP\Data\Sample Information\17032901.sif

Batch ID:
 Results Data Set: 170329C 1
 Results Library: W:\pe\7300\Results\results.mdb

=====
 Sequence No.: 3 Autosampler Location: 10
 Sample ID: ICV-M072816C Date Collected: 3/29/2017 9:46:40 AM
 Analyst: 935 icp 7300 Data Type: Original
 Initial Sample Wt: Initial Sample Vol:
 Dilution: Sample Prep Vol:
 Wash Time: 20 Auto Dilution Factor: 1

 Mean Data: ICV-M072816C

Analyte	Mean Corrected		Calib.		Sample		RSD
	Intensity	Conc.	Units	Std.Dev.	Conc.	Std.Dev.	
Tb 384	74252.1	92.95	%	1.178			1.27%
Tb 350	120370.3	95.43	%	0.983			1.03%
Ag 328.068*†	87057.5	0.4918	mg/L	0.00376	0.4918 mg/L	0.00376	0.76%
	QC value within limits for Ag 328.068* Recovery = 98.36%						
Al 308.215*†	62137.3	4.014	mg/L	0.0023	4.014 mg/L	0.0023	0.06%
	QC value within limits for Al 308.215* Recovery = 100.35%						
As 188.979†	8467.2	4.858	mg/L	0.0687	4.858 mg/L	0.0687	1.41%
	QC value within limits for As 188.979 Recovery = 97.16%						
As 193.696*†	5885.6	4.806	mg/L	0.0697	4.806 mg/L	0.0697	1.45%
	QC value within limits for As 193.696* Recovery = 96.12%						
B 249.677*†	108210.5	2.446	mg/L	0.0327	2.446 mg/L	0.0327	1.34%
	QC value within limits for B 249.677* Recovery = 97.86%						
Ba 233.527*†	134742.5	0.9776	mg/L	0.00166	0.9776 mg/L	0.00166	0.17%
	QC value within limits for Ba 233.527* Recovery = 97.76%						
Be 313.042*†	1752361.6	0.4854	mg/L	0.00056	0.4854 mg/L	0.00056	0.12%
	QC value within limits for Be 313.042* Recovery = 97.09%						
Ca 317.933*†	40919.5	18.46	mg/L	0.137	18.46 mg/L	0.137	0.74%
	QC value within limits for Ca 317.933* Recovery = 92.31%						
Cd 226.502*†	94769.5	1.493	mg/L	0.0050	1.493 mg/L	0.0050	0.33%
	QC value within limits for Cd 226.502* Recovery = 99.53%						
Cd 228.802†	49016.9	1.477	mg/L	0.0045	1.477 mg/L	0.0045	0.30%
Co 228.616*†	24615.1	1.031	mg/L	0.0170	1.031 mg/L	0.0170	1.65%
	QC value within limits for Co 228.616* Recovery = 103.09%						
Cr 267.716*†	39910.8	0.3953	mg/L	0.00152	0.3953 mg/L	0.00152	0.38%
	QC value within limits for Cr 267.716* Recovery = 98.83%						
Cu 324.752*†	237524.2	0.9864	mg/L	0.00299	0.9864 mg/L	0.00299	0.30%
	QC value within limits for Cu 324.752* Recovery = 98.64%						
Fe 273.955*†	2084449.0	101.1	mg/L	0.54	101.1 mg/L	0.54	0.54%
	QC value within limits for Fe 273.955* Recovery = 101.10%						
K 766.490*†	25136.2	7.466	mg/L	0.2347	7.466 mg/L	0.2347	3.14%
	QC value within limits for K 766.490* Recovery = 93.33%						
Mg 279.077*†	152956.5	9.917	mg/L	0.0222	9.917 mg/L	0.0222	0.22%
	QC value within limits for Mg 279.077* Recovery = 99.17%						
Mn 257.610*†	577658.6	0.9808	mg/L	0.00144	0.9808 mg/L	0.00144	0.15%
	QC value within limits for Mn 257.610* Recovery = 98.08%						
Mo 202.031*†	19933.3	2.376	mg/L	0.0406	2.376 mg/L	0.0406	1.71%
	QC value within limits for Mo 202.031* Recovery = 95.05%						
Na 589.592*†	270521.4	52.18	mg/L	0.492	52.18 mg/L	0.492	0.94%
	QC value within limits for Na 589.592* Recovery = 96.63%						
Ni 231.604*†	8627.8	0.4064	mg/L	0.00598	0.4064 mg/L	0.00598	1.47%
	QC value within limits for Ni 231.604* Recovery = 101.59%						
P 213.617*†	7685.4	4.716	mg/L	0.0525	4.716 mg/L	0.0525	1.11%
	QC value within limits for P 213.617* Recovery = 94.32%						
P 214.914†	4932.4	4.876	mg/L	0.0578	4.876 mg/L	0.0578	1.19%
Pb 220.353*†	32508.5	5.072	mg/L	0.0101	5.072 mg/L	0.0101	0.20%

Return to Contents

	QC value within limits for Pb 220.353*	Recovery = 101.44%				
Sb	206.836†	3192.6	1.883 mg/L	0.0399	1.883 mg/L	0.0399 2.12%
	QC value within limits for Sb 206.836	Recovery = 94.16%				
Sb	217.582*†	3248.5	1.904 mg/L	0.0259	1.904 mg/L	0.0259 1.36%
	QC value within limits for Sb 217.582*	Recovery = 95.18%				
Se	196.026*†	3400.5	1.927 mg/L	0.0389	1.927 mg/L	0.0389 2.02%
	QC value within limits for Se 196.026*	Recovery = 96.36%				
Si	251.611*†	327242.2	9.316 mg/L	0.0398	9.316 mg/L	0.0398 0.43%
	QC value within limits for Si 251.611*	Recovery = 93.16%				
Sn	189.927*†	12344.0	2.438 mg/L	0.0383	2.438 mg/L	0.0383 1.57%
	QC value within limits for Sn 189.927*	Recovery = 97.52%				
Sn	242.170†	4361.7	2.707 mg/L	0.0428	2.707 mg/L	0.0428 1.58%
Sr	407.771*†	71013.0	0.1913 mg/L	0.00178	0.1913 mg/L	0.00178 0.93%
	QC value within limits for Sr 407.771*	Recovery = 95.65%				
Ti	334.940†	3828522.5	4.709 mg/L	0.0185	4.709 mg/L	0.0185 0.39%
Ti	336.121*†	2698233.6	4.806 mg/L	0.0203	4.806 mg/L	0.0203 0.42%
	QC value within limits for Ti 336.121*	Recovery = 96.12%				
Tl	190.801*†	3076.2	1.940 mg/L	0.0300	1.940 mg/L	0.0300 1.55%
	QC value within limits for Tl 190.801*	Recovery = 97.01%				
V	292.402*†	122162.7	0.9763 mg/L	0.00045	0.9763 mg/L	0.00045 0.05%
	QC value within limits for V 292.402*	Recovery = 97.63%				
Zn	206.200*†	48996.4	1.515 mg/L	0.0094	1.515 mg/L	0.0094 0.62%
	QC value within limits for Zn 206.200*	Recovery = 101.02%				
Zn	213.857*†	90376.6	1.506 mg/L	0.0014	1.506 mg/L	0.0014 0.09%
	QC value within limits for Zn 213.857*	Recovery = 100.40%				

All analyte(s) passed QC.

Sequence No.: 4

Sample ID: ICB-R12091601

Analyst: 935 icp 7300

Initial Sample Wt:

Dilution:

Wash Time: 20

Autosampler Location: 1

Date Collected: 3/29/2017 9:47:39 AM

Data Type: Original

Initial Sample Vol:

Sample Prep Vol:

Auto Dilution Factor: 1

Mean Data: ICB-R12091601

Analyte	Mean Corrected		Calib. Conc. Units	Std.Dev.	Sample		Std.Dev.	RSD
	Intensity	Conc.			Conc.	Units		
Tb 384	80062.2	100.2	%	1.32				1.31%
Tb 350	126192.1	100.0	%	0.73				0.73%
Ag 328.068*†	-109.7	-0.0006	mg/L	0.00085	-0.0006	mg/L	0.00085	136.57%
Al 308.215*†	-27.1	-0.0018	mg/L	0.00237	-0.0018	mg/L	0.00237	135.61%
As 188.979†	2.5	0.0015	mg/L	0.00654	0.0015	mg/L	0.00654	448.28%
As 193.696*†	4.4	0.0036	mg/L	0.00058	0.0036	mg/L	0.00058	15.92%
QC value within limits for As 193.696* Recovery = Not calculated								
B 249.677*†	1487.2	0.0336	mg/L	0.00158	0.0336	mg/L	0.00158	4.70%
QC value within limits for B 249.677* Recovery = Not calculated								
Ba 233.527*†	-13.2	-0.0001	mg/L	0.00001	-0.0001	mg/L	0.00001	9.43%
Be 313.042*†	107.8	0.0000	mg/L	0.00001	0.0000	mg/L	0.00001	40.92%
Ca 317.933*†	-10.7	-0.0048	mg/L	0.00563	-0.0048	mg/L	0.00563	116.82%
Cd 226.502*†	26.7	0.0004	mg/L	0.00005	0.0004	mg/L	0.00005	12.58%
Cd 228.802†	9.4	0.0003	mg/L	0.00030	0.0003	mg/L	0.00030	106.23%
Co 228.616*†	6.2	0.0003	mg/L	0.00044	0.0003	mg/L	0.00044	170.02%
Cr 267.716*†	10.7	0.0001	mg/L	0.00030	0.0001	mg/L	0.00030	281.33%
Cu 324.752*†	-58.0	-0.0002	mg/L	0.00029	-0.0002	mg/L	0.00029	119.06%
Fe 273.955*†	-3.3	-0.0002	mg/L	0.00019	-0.0002	mg/L	0.00019	117.82%
K 766.490*†	-219.3	-0.0651	mg/L	0.09670	-0.0651	mg/L	0.09670	148.43%
Mg 279.077*†	33.0	0.0021	mg/L	0.00730	0.0021	mg/L	0.00730	340.90%
Mn 257.610*†	-47.6	-0.0001	mg/L	0.00002	-0.0001	mg/L	0.00002	27.17%
Mo 202.031*†	2.1	0.0002	mg/L	0.00053	0.0002	mg/L	0.00053	214.34%
Na 589.592*†	-65.9	-0.0127	mg/L	0.02287	-0.0127	mg/L	0.02287	179.88%
Ni 231.604*†	20.3	0.0010	mg/L	0.00083	0.0010	mg/L	0.00083	86.80%
P 213.617*†	1.5	0.0009	mg/L	0.00987	0.0009	mg/L	0.00987	>999.9%
P 214.914†	6.8	0.0067	mg/L	0.00567	0.0067	mg/L	0.00567	84.20%
Pb 220.353*†	6.4	0.0010	mg/L	0.00019	0.0010	mg/L	0.00019	19.07%
QC value within limits for Pb 220.353* Recovery = Not calculated								
Sb 206.836†	9.6	0.0056	mg/L	0.00309	0.0056	mg/L	0.00309	54.85%
Sb 217.582*†	-2.6	-0.0015	mg/L	0.00049	-0.0015	mg/L	0.00049	32.11%
QC value within limits for Sb 217.582* Recovery = Not calculated								
Se 196.026*†	4.0	0.0023	mg/L	0.00539	0.0023	mg/L	0.00539	237.50%
QC value within limits for Se 196.026* Recovery = Not calculated								
Si 251.611*†	-12.4	-0.0004	mg/L	0.00026	-0.0004	mg/L	0.00026	74.63%
QC value within limits for Si 251.611* Recovery = Not calculated								
Sn 189.927*†	3.0	0.0006	mg/L	0.00118	0.0006	mg/L	0.00118	197.30%
Sn 242.170†	27.5	0.0171	mg/L	0.01715	0.0171	mg/L	0.01715	100.55%
Sr 407.771*†	3.1	0.0000	mg/L	0.00003	0.0000	mg/L	0.00003	388.36%
Ti 334.940†	580.3	0.0007	mg/L	0.00020	0.0007	mg/L	0.00020	28.61%
Ti 336.121*†	497.4	0.0009	mg/L	0.00024	0.0009	mg/L	0.00024	26.60%
Tl 190.801*†	-1.9	-0.0012	mg/L	0.00109	-0.0012	mg/L	0.00109	90.05%
QC value within limits for Tl 190.801* Recovery = Not calculated								
V 292.402*†	2.6	0.0000	mg/L	0.00057	0.0000	mg/L	0.00057	>999.9%
Zn 206.200*†	6.2	0.0002	mg/L	0.00042	0.0002	mg/L	0.00042	218.81%
Zn 213.857*†	4.5	0.0001	mg/L	0.00027	0.0001	mg/L	0.00027	359.95%

All analyte(s) passed QC.

Sequence No.: 5

Sample ID: ICS_A - M110116B

Analyst: 935 icp 7300

Initial Sample Wt:

Dilution:

Wash Time: 15

Autosampler Location: 8

Date Collected: 3/29/2017 9:48:40 AM

Data Type: Original

Initial Sample Vol:

Sample Prep Vol:

Auto Dilution Factor: 1

Mean Data: ICS_A - M110116B

Analyte	Mean Corrected Intensity	Conc. Units	Calib. Units	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	73445.6	91.94	%	1.573			1.71%
Tb 350	118669.9	94.08	%	1.683			1.79%
Ag 328.068*†	-41.5	-0.0002	mg/L	0.00014	-0.0002 mg/L	0.00014	60.69%
Al 308.215*†	364506.1	23.55	mg/L	0.374	23.55 mg/L	0.374	1.59%
As 188.979†	10.7	0.0061	mg/L	0.00615	0.0061 mg/L	0.00615	100.24%
As 193.696*†	-5.6	-0.0046	mg/L	0.00786	-0.0046 mg/L	0.00786	171.99%
B 249.677*†	865.8	0.0196	mg/L	0.00061	0.0196 mg/L	0.00061	3.12%
Ba 233.527*†	341.9	0.0025	mg/L	0.00008	0.0025 mg/L	0.00008	3.11%
Be 313.042*†	-6.9	-0.0000	mg/L	0.00000	-0.0000 mg/L	0.00000	55.15%
Ca 317.933*†	245949.8	111.0	mg/L	3.11	111.0 mg/L	3.11	2.80%
Cd 226.502*†	150.1	0.0024	mg/L	0.00022	0.0024 mg/L	0.00022	9.12%
Cd 228.802†	13.1	0.0004	mg/L	0.00018	0.0004 mg/L	0.00018	45.94%
Co 228.616*†	21.4	0.0009	mg/L	0.00024	0.0009 mg/L	0.00024	26.67%
Cr 267.716*†	-186.9	-0.0019	mg/L	0.00008	-0.0019 mg/L	0.00008	4.56%
Cu 324.752*†	-445.2	-0.0018	mg/L	0.00058	-0.0018 mg/L	0.00058	31.17%
Fe 273.955*†	1877600.3	91.07	mg/L	1.379	91.07 mg/L	1.379	1.51%
K 766.490*†	284.4	0.0845	mg/L	0.05387	0.0845 mg/L	0.05387	63.77%
Mg 279.077*†	881903.1	57.18	mg/L	0.889	57.18 mg/L	0.889	1.55%
Mn 257.610*†	39.5	0.0001	mg/L	0.00018	0.0001 mg/L	0.00018	271.60%
Mo 202.031*†	15.4	0.0018	mg/L	0.00148	0.0018 mg/L	0.00148	80.19%
Na 589.592*†	103624.1	19.99	mg/L	0.504	19.99 mg/L	0.504	2.52%
Ni 231.604*†	46.1	0.0022	mg/L	0.00044	0.0022 mg/L	0.00044	20.44%
P 213.617*†	-175.0	-0.1074	mg/L	0.00670	-0.1074 mg/L	0.00670	6.24%
P 214.914†	44.5	0.0440	mg/L	0.00351	0.0440 mg/L	0.00351	7.98%
Pb 220.353*†	-36.1	-0.0056	mg/L	0.00068	-0.0056 mg/L	0.00068	12.00%
Sb 206.836†	13.6	0.0081	mg/L	0.00665	0.0081 mg/L	0.00665	82.58%
Sb 217.582*†	-19.8	-0.0116	mg/L	0.00039	-0.0116 mg/L	0.00039	3.33%
Se 196.026*†	-11.4	-0.0065	mg/L	0.00011	-0.0065 mg/L	0.00011	1.75%
Si 251.611*†	-112.0	-0.0032	mg/L	0.00078	-0.0032 mg/L	0.00078	24.45%
Sn 189.927*†	-30.7	-0.0061	mg/L	0.00213	-0.0061 mg/L	0.00213	35.13%
Sn 242.170†	383.3	0.2379	mg/L	0.01939	0.2379 mg/L	0.01939	8.15%
Sr 407.771*†	901.8	0.0024	mg/L	0.00022	0.0024 mg/L	0.00022	8.97%
Ti 334.940†	-1941.7	-0.0024	mg/L	0.00000	-0.0024 mg/L	0.00000	0.18%
Ti 336.121*†	-727.7	-0.0013	mg/L	0.00001	-0.0013 mg/L	0.00001	0.47%
Tl 190.801*†	-8.8	-0.0055	mg/L	0.00484	-0.0055 mg/L	0.00484	87.29%
V 292.402*†	349.0	0.0013	mg/L	0.00021	0.0013 mg/L	0.00021	16.02%
Zn 206.200*†	24.3	0.0008	mg/L	0.00025	0.0008 mg/L	0.00025	32.85%
Zn 213.857*†	436.2	-0.0003	mg/L	0.00017	-0.0003 mg/L	0.00017	66.49%

User canceled analysis.

=====
Analysis Begun

Start Time: 3/29/2017 9:49:44 AM Plasma On Time: 3/29/2017 9:19:17 AM
 Logged In Analyst: Oscar Gomez 935 Technique: ICP Continuous
 Spectrometer: Optima 7300 DV, S/N 77c8120401 Autosampler: ESI

Sample Information File: C:\Documents and Settings\All Users\PerkinElmer\ICP\Data\Sample Information\17032901.sif

Batch ID:
 Results Data Set: 170329C 1
 Results Library: W:\pe\7300\Results\results.mdb

=====
 Sequence No.: 6 Autosampler Location: 9
 Sample ID: ICS_AB - M110116A Date Collected: 3/29/2017 9:49:46 AM
 Analyst: 935 icp 7300 Data Type: Original
 Initial Sample Wt: Initial Sample Vol:
 Dilution: Sample Prep Vol:
 Wash Time:

 Mean Data: ICS_AB - M110116A

Analyte	Mean Corrected Intensity	Conc. Units	Calib.	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	69213.3	86.64	%	0.628			0.73%
Tb 350	111741.1	88.59	%	0.558			0.63%
Ag 328.068*†	56383.3	0.3185	mg/L	0.00094	0.3185 mg/L	0.00094	0.29%
Al 308.215*†	386172.1	24.95	mg/L	0.043	24.95 mg/L	0.043	0.17%
As 188.979†	1815.2	1.041	mg/L	0.0159	1.041 mg/L	0.0159	1.52%
As 193.696*†	1261.2	1.030	mg/L	0.0158	1.030 mg/L	0.0158	1.54%
B 249.677*†	22608.1	0.5111	mg/L	0.00130	0.5111 mg/L	0.00130	0.25%
Ba 233.527*†	42804.0	0.3106	mg/L	0.00161	0.3106 mg/L	0.00161	0.52%
Be 313.042*†	379786.1	0.1052	mg/L	0.00024	0.1052 mg/L	0.00024	0.22%
Ca 317.933*†	259231.0	117.0	mg/L	1.61	117.0 mg/L	1.61	1.38%
Cd 226.502*†	19612.6	0.3090	mg/L	0.00254	0.3090 mg/L	0.00254	0.82%
Cd 228.802†	10291.7	0.3102	mg/L	0.00244	0.3102 mg/L	0.00244	0.79%
Co 228.616*†	7481.6	0.3133	mg/L	0.00140	0.3133 mg/L	0.00140	0.45%
Cr 267.716*†	31918.8	0.3162	mg/L	0.00263	0.3162 mg/L	0.00263	0.83%
Cu 324.752*†	76930.6	0.3195	mg/L	0.00068	0.3195 mg/L	0.00068	0.21%
Fe 273.955*†	2004522.9	97.22	mg/L	0.297	97.22 mg/L	0.297	0.31%
K 766.490*†	71628.9	21.28	mg/L	0.299	21.28 mg/L	0.299	1.41%
Mg 279.077*†	933465.3	60.52	mg/L	0.221	60.52 mg/L	0.221	0.37%
Mn 257.610*†	121136.8	0.2057	mg/L	0.00018	0.2057 mg/L	0.00018	0.09%
Mo 202.031*†	2559.6	0.3051	mg/L	0.00361	0.3051 mg/L	0.00361	1.18%
Na 589.592*†	108252.2	20.88	mg/L	0.238	20.88 mg/L	0.238	1.14%
Ni 231.604*†	6714.3	0.3162	mg/L	0.00284	0.3162 mg/L	0.00284	0.90%
P 213.617*†	-179.6	-0.1102	mg/L	0.01262	-0.1102 mg/L	0.01262	11.45%
P 214.914†	53.0	0.0524	mg/L	0.00286	0.0524 mg/L	0.00286	5.47%
Pb 220.353*†	6604.2	1.030	mg/L	0.0063	1.030 mg/L	0.0063	0.61%
Sb 206.836†	1683.3	0.9929	mg/L	0.00751	0.9929 mg/L	0.00751	0.76%
Sb 217.582*†	1679.7	0.9843	mg/L	0.01766	0.9843 mg/L	0.01766	1.79%
Se 196.026*†	909.1	0.5152	mg/L	0.00282	0.5152 mg/L	0.00282	0.55%
Si 251.611*†	7843.2	0.2233	mg/L	0.00499	0.2233 mg/L	0.00499	2.23%
Sn 189.927*†	-46.3	-0.0091	mg/L	0.00084	-0.0091 mg/L	0.00084	9.23%
Sn 242.170†	427.6	0.2653	mg/L	0.00089	0.2653 mg/L	0.00089	0.34%
Sr 407.771*†	1019.1	0.0027	mg/L	0.00013	0.0027 mg/L	0.00013	4.64%
Ti 334.940†	820110.1	1.009	mg/L	0.0018	1.009 mg/L	0.0018	0.18%
Ti 336.121*†	576771.8	1.027	mg/L	0.0013	1.027 mg/L	0.0013	0.12%
Tl 190.801*†	1618.3	1.021	mg/L	0.0032	1.021 mg/L	0.0032	0.32%
V 292.402*†	39371.0	0.3136	mg/L	0.00014	0.3136 mg/L	0.00014	0.04%
Zn 206.200*†	10285.9	0.3181	mg/L	0.00327	0.3181 mg/L	0.00327	1.03%
Zn 213.857*†	19555.3	0.3196	mg/L	0.00270	0.3196 mg/L	0.00270	0.85%

Sequence No.: 4

Autosampler Location: 3

Sample ID: CCV= STD3x0.5

Date Collected: 3/29/2017 4:56:49 PM

Analyst: 935 icp 7300

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time: 15

Auto Dilution Factor: 1

Mean Data: CCV= STD3x0.5

Analyte	Mean Corrected Intensity	Conc.	Calib. Units	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	79501.9	99.52	%	1.057			1.06%
Tb 350	129845.7	102.9	%	1.53			1.49%
Ag 328.068*†	64594.2	0.3649	mg/L	0.00125	0.3649 mg/L	0.00125	0.34%
	QC value within limits for Ag 328.068* Recovery = 97.31%						
Al 308.215*†	203459.0	13.14	mg/L	0.016	13.14 mg/L	0.016	0.12%
	QC value within limits for Al 308.215* Recovery = 97.36%						
As 188.979†	6495.3	3.727	mg/L	0.0684	3.727 mg/L	0.0684	1.84%
	QC value within limits for As 188.979 Recovery = 99.37%						
As 193.696*†	4510.4	3.683	mg/L	0.0643	3.683 mg/L	0.0643	1.74%
	QC value within limits for As 193.696* Recovery = 98.22%						
B 249.677*†	167196.1	3.780	mg/L	0.0642	3.780 mg/L	0.0642	1.70%
	QC value within limits for B 249.677* Recovery = 100.80%						
Ba 233.527*†	1029121.3	7.467	mg/L	0.0022	7.467 mg/L	0.0022	0.03%
	QC value within limits for Ba 233.527* Recovery = 99.56%						
Be 313.042*†	2002359.1	0.5547	mg/L	0.00157	0.5547 mg/L	0.00157	0.28%
	QC value within limits for Be 313.042* Recovery = 98.61%						
Ca 317.933*†	63340.7	28.58	mg/L	0.373	28.58 mg/L	0.373	1.31%
	QC value within limits for Ca 317.933* Recovery = 95.26%						
Cd 226.502*†	48589.2	0.7654	mg/L	0.00203	0.7654 mg/L	0.00203	0.27%
	QC value within limits for Cd 226.502* Recovery = 102.06%						
Cd 228.802†	25174.3	0.7588	mg/L	0.00116	0.7588 mg/L	0.00116	0.15%
Co 228.616*†	46411.4	1.944	mg/L	0.0056	1.944 mg/L	0.0056	0.29%
	QC value within limits for Co 228.616* Recovery = 103.66%						
Cr 267.716*†	61133.9	0.6055	mg/L	0.00200	0.6055 mg/L	0.00200	0.33%
	QC value within limits for Cr 267.716* Recovery = 100.92%						
Cu 324.752*†	220236.1	0.9146	mg/L	0.00000	0.9146 mg/L	0.00000	0.00%
	QC value within limits for Cu 324.752* Recovery = 97.55%						
Fe 273.955*†	78942.0	3.829	mg/L	0.0125	3.829 mg/L	0.0125	0.33%
	QC value within limits for Fe 273.955* Recovery = 102.10%						
K 766.490*†	85604.5	25.43	mg/L	0.446	25.43 mg/L	0.446	1.75%
	QC value within limits for K 766.490* Recovery = 94.17%						
Mg 279.077*†	119014.6	7.717	mg/L	0.0180	7.717 mg/L	0.0180	0.23%
	QC value within limits for Mg 279.077* Recovery = 102.89%						
Mn 257.610*†	436371.8	0.7409	mg/L	0.00034	0.7409 mg/L	0.00034	0.05%
	QC value within limits for Mn 257.610* Recovery = 98.79%						
Mo 202.031*†	4951.5	0.5903	mg/L	0.01250	0.5903 mg/L	0.01250	2.12%
	QC value within limits for Mo 202.031* Recovery = 98.38%						
Na 589.592*†	179849.3	34.69	mg/L	0.475	34.69 mg/L	0.475	1.37%
	QC value within limits for Na 589.592* Recovery = 96.36%						
Ni 231.604*†	12860.0	0.6057	mg/L	0.01357	0.6057 mg/L	0.01357	2.24%
	QC value within limits for Ni 231.604* Recovery = 100.95%						
P 213.617*†	9433.6	5.789	mg/L	0.1357	5.789 mg/L	0.1357	2.35%
	QC value within limits for P 213.617* Recovery = 96.48%						
P 214.914†	6032.9	5.964	mg/L	0.1051	5.964 mg/L	0.1051	1.76%
Pb 220.353*†	24558.4	3.832	mg/L	0.0072	3.832 mg/L	0.0072	0.19%
	QC value within limits for Pb 220.353* Recovery = 102.17%						
Sb 206.836†	7354.7	4.338	mg/L	0.0727	4.338 mg/L	0.0727	1.68%
	QC value within limits for Sb 206.836 Recovery = 96.41%						
Sb 217.582*†	7330.7	4.296	mg/L	0.0861	4.296 mg/L	0.0861	2.00%
	QC value within limits for Sb 217.582* Recovery = 95.46%						
Se 196.026*†	2649.6	1.502	mg/L	0.0216	1.502 mg/L	0.0216	1.44%
	QC value within limits for Se 196.026* Recovery = 100.11%						
Si 251.611*†	204763.4	5.829	mg/L	0.0354	5.829 mg/L	0.0354	0.61%
	QC value within limits for Si 251.611* Recovery = 97.16%						
Sn 189.927*†	15221.0	3.006	mg/L	0.0546	3.006 mg/L	0.0546	1.82%
	QC value within limits for Sn 189.927* Recovery = 100.21%						
Sn 242.170†	4737.4	2.940	mg/L	0.0233	2.940 mg/L	0.0233	0.79%
Sr 407.771*†	108135.4	0.2913	mg/L	0.00424	0.2913 mg/L	0.00424	1.46%
	QC value within limits for Sr 407.771* Recovery = 97.10%						

Ti 334.940†	474413.1	0.5836 mg/L	0.00142	0.5836 mg/L	0.00142	0.24%
Ti 336.121*†	334591.0	0.5960 mg/L	0.00096	0.5960 mg/L	0.00096	0.16%
QC value within limits for Ti 336.121* Recovery = 99.33%						
Tl 190.801*†	2384.3	1.504 mg/L	0.0274	1.504 mg/L	0.0274	1.83%
QC value within limits for Tl 190.801* Recovery = 100.25%						
V 292.402*†	235354.6	1.884 mg/L	0.0041	1.884 mg/L	0.0041	0.22%
QC value within limits for V 292.402* Recovery = 100.48%						
Zn 206.200*†	81897.2	2.533 mg/L	0.0117	2.533 mg/L	0.0117	0.46%
QC value within limits for Zn 206.200* Recovery = 101.31%						
Zn 213.857*†	152244.9	2.551 mg/L	0.0007	2.551 mg/L	0.0007	0.03%
QC value within limits for Zn 213.857* Recovery = 102.03%						

All analyte(s) passed QC.

Sequence No.: 5

Autosampler Location: 1

Sample ID: CCB-R12091601

Date Collected: 3/29/2017 4:57:41 PM

Analyst: 935 icp 7300

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time: 15

Auto Dilution Factor: 1

Mean Data: CCB-R12091601

Analyte	Mean Corrected		Calib. Units	Std.Dev.	Sample		Std.Dev.	RSD
	Intensity	Conc.			Conc.	Units		
Tb 384	86105.4	107.8	%	1.44				1.33%
Tb 350	138375.3	109.7	%	1.32				1.20%
Ag 328.068*†	16.5	0.0001	mg/L	0.00070	0.0001	mg/L	0.00070	758.37%
QC value within limits for Ag 328.068* Recovery = Not calculated								
Al 308.215*†	-66.0	-0.0043	mg/L	0.00210	-0.0043	mg/L	0.00210	49.12%
As 188.979†	5.7	0.0032	mg/L	0.00024	0.0032	mg/L	0.00024	7.51%
As 193.696*†	1.7	0.0014	mg/L	0.00184	0.0014	mg/L	0.00184	130.69%
QC value within limits for As 193.696* Recovery = Not calculated								
B 249.677*†	1691.8	0.0382	mg/L	0.00259	0.0382	mg/L	0.00259	6.77%
Ba 233.527*†	404.1	0.0029	mg/L	0.00014	0.0029	mg/L	0.00014	4.77%
Be 313.042*†	728.5	0.0002	mg/L	0.00001	0.0002	mg/L	0.00001	5.03%
Ca 317.933*†	52.4	0.0236	mg/L	0.00026	0.0236	mg/L	0.00026	1.09%
Cd 226.502*†	25.4	0.0004	mg/L	0.00003	0.0004	mg/L	0.00003	7.92%
Cd 228.802†	24.8	0.0007	mg/L	0.00006	0.0007	mg/L	0.00006	8.63%
Co 228.616*†	16.3	0.0007	mg/L	0.00062	0.0007	mg/L	0.00062	90.23%
Cr 267.716*†	34.9	0.0003	mg/L	0.00021	0.0003	mg/L	0.00021	60.61%
Cu 324.752*†	-65.1	-0.0003	mg/L	0.00008	-0.0003	mg/L	0.00008	28.49%
Fe 273.955*†	214.7	0.0104	mg/L	0.00072	0.0104	mg/L	0.00072	6.88%
K 766.490*†	-252.9	-0.0751	mg/L	0.07945	-0.0751	mg/L	0.07945	105.75%
Mg 279.077*†	344.7	0.0224	mg/L	0.00143	0.0224	mg/L	0.00143	6.38%
Mn 257.610*†	315.7	0.0005	mg/L	0.00000	0.0005	mg/L	0.00000	0.88%
Mo 202.031*†	2.2	0.0003	mg/L	0.00084	0.0003	mg/L	0.00084	317.17%
Na 589.592*†	27.7	0.0053	mg/L	0.01470	0.0053	mg/L	0.01470	275.36%
Ni 231.604*†	23.2	0.0011	mg/L	0.00053	0.0011	mg/L	0.00053	48.23%
P 213.617*†	9.6	0.0059	mg/L	0.00079	0.0059	mg/L	0.00079	13.49%
P 214.914†	5.9	0.0059	mg/L	0.00091	0.0059	mg/L	0.00091	15.49%
Pb 220.353*†	-2.9	-0.0005	mg/L	0.00036	-0.0005	mg/L	0.00036	79.94%
Sb 206.836†	25.5	0.0151	mg/L	0.00201	0.0151	mg/L	0.00201	13.34%
Sb 217.582*†	0.7	0.0004	mg/L	0.00118	0.0004	mg/L	0.00118	281.37%
QC value within limits for Sb 217.582* Recovery = Not calculated								
Se 196.026*†	7.7	0.0044	mg/L	0.00097	0.0044	mg/L	0.00097	22.25%
QC value within limits for Se 196.026* Recovery = Not calculated								
Si 251.611*†	-68.7	-0.0020	mg/L	0.00389	-0.0020	mg/L	0.00389	198.84%
Sn 189.927*†	14.5	0.0029	mg/L	0.00152	0.0029	mg/L	0.00152	53.32%
Sn 242.170†	12.8	0.0079	mg/L	0.01786	0.0079	mg/L	0.01786	224.66%
Sr 407.771*†	82.8	0.0002	mg/L	0.00003	0.0002	mg/L	0.00003	12.39%
Ti 334.940†	707.6	0.0009	mg/L	0.00030	0.0009	mg/L	0.00030	34.83%
Ti 336.121*†	628.9	0.0011	mg/L	0.00030	0.0011	mg/L	0.00030	26.44%
Tl 190.801*†	1.3	0.0008	mg/L	0.00034	0.0008	mg/L	0.00034	41.34%
V 292.402*†	161.5	0.0013	mg/L	0.00146	0.0013	mg/L	0.00146	112.55%
Zn 206.200*†	-51.3	-0.0016	mg/L	0.00060	-0.0016	mg/L	0.00060	37.62%
QC value within limits for Zn 206.200* Recovery = Not calculated								
Zn 213.857*†	-43.8	-0.0007	mg/L	0.00019	-0.0007	mg/L	0.00019	25.42%
QC value within limits for Zn 213.857* Recovery = Not calculated								

All analyte(s) passed QC.

Sequence No.: 11

Autosampler Location: 3

Sample ID: CCV= STD3x0.5

Date Collected: 3/29/2017 5:14:09 PM

Analyst: 935 icp 7300

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time: 15

Auto Dilution Factor: 1

Mean Data: CCV= STD3x0.5

Analyte	Mean Corrected		Calib.	Std.Dev.	Sample		Std.Dev.	RSD
	Intensity	Conc.			Units	Conc.		
Tb 384	79519.1	99.55	%	0.022				0.02%
Tb 350	131160.1	104.0	%	0.11				0.10%
Ag 328.068*†	64874.6	0.3665	mg/L	0.00021	0.3665	mg/L	0.00021	0.06%
	QC value within limits for Ag 328.068* Recovery = 97.73%							
Al 308.215*†	204627.2	13.22	mg/L	0.021	13.22	mg/L	0.021	0.16%
	QC value within limits for Al 308.215* Recovery = 97.91%							
As 188.979†	6594.7	3.784	mg/L	0.0224	3.784	mg/L	0.0224	0.59%
	QC value within limits for As 188.979 Recovery = 100.89%							
As 193.696*†	4590.8	3.749	mg/L	0.0185	3.749	mg/L	0.0185	0.49%
	QC value within limits for As 193.696* Recovery = 99.97%							
B 249.677*†	167474.5	3.786	mg/L	0.0544	3.786	mg/L	0.0544	1.44%
	QC value within limits for B 249.677* Recovery = 100.97%							
Ba 233.527*†	1035047.1	7.510	mg/L	0.0045	7.510	mg/L	0.0045	0.06%
	QC value within limits for Ba 233.527* Recovery = 100.13%							
Be 313.042*†	1993630.9	0.5523	mg/L	0.00033	0.5523	mg/L	0.00033	0.06%
	QC value within limits for Be 313.042* Recovery = 98.18%							
Ca 317.933*†	65347.7	29.48	mg/L	0.091	29.48	mg/L	0.091	0.31%
	QC value within limits for Ca 317.933* Recovery = 98.28%							
Cd 226.502*†	49007.6	0.7720	mg/L	0.00153	0.7720	mg/L	0.00153	0.20%
	QC value within limits for Cd 226.502* Recovery = 102.94%							
Cd 228.802†	25246.5	0.7610	mg/L	0.00301	0.7610	mg/L	0.00301	0.40%
Co 228.616*†	46757.8	1.958	mg/L	0.0073	1.958	mg/L	0.0073	0.37%
	QC value within limits for Co 228.616* Recovery = 104.44%							
Cr 267.716*†	61609.6	0.6102	mg/L	0.00041	0.6102	mg/L	0.00041	0.07%
	QC value within limits for Cr 267.716* Recovery = 101.71%							
Cu 324.752*†	220352.4	0.9150	mg/L	0.00010	0.9150	mg/L	0.00010	0.01%
	QC value within limits for Cu 324.752* Recovery = 97.60%							
Fe 273.955*†	81168.6	3.937	mg/L	0.0177	3.937	mg/L	0.0177	0.45%
	QC value within limits for Fe 273.955* Recovery = 104.98%							
K 766.490*†	86754.1	25.77	mg/L	0.056	25.77	mg/L	0.056	0.22%
	QC value within limits for K 766.490* Recovery = 95.44%							
Mg 279.077*†	119999.3	7.780	mg/L	0.0015	7.780	mg/L	0.0015	0.02%
	QC value within limits for Mg 279.077* Recovery = 103.74%							
Mn 257.610*†	439771.5	0.7467	mg/L	0.00181	0.7467	mg/L	0.00181	0.24%
	QC value within limits for Mn 257.610* Recovery = 99.56%							
Mo 202.031*†	5019.1	0.5983	mg/L	0.00255	0.5983	mg/L	0.00255	0.43%
	QC value within limits for Mo 202.031* Recovery = 99.72%							
Na 589.592*†	182050.8	35.11	mg/L	0.007	35.11	mg/L	0.007	0.02%
	QC value within limits for Na 589.592* Recovery = 97.54%							
Ni 231.604*†	12995.4	0.6121	mg/L	0.00179	0.6121	mg/L	0.00179	0.29%
	QC value within limits for Ni 231.604* Recovery = 102.01%							
P 213.617*†	9532.3	5.849	mg/L	0.0674	5.849	mg/L	0.0674	1.15%
	QC value within limits for P 213.617* Recovery = 97.49%							
P 214.914†	6151.3	6.081	mg/L	0.0305	6.081	mg/L	0.0305	0.50%
Pb 220.353*†	24796.2	3.869	mg/L	0.0250	3.869	mg/L	0.0250	0.65%
	QC value within limits for Pb 220.353* Recovery = 103.16%							
Sb 206.836†	7437.0	4.387	mg/L	0.0116	4.387	mg/L	0.0116	0.26%
	QC value within limits for Sb 206.836 Recovery = 97.49%							
Sb 217.582*†	7433.4	4.356	mg/L	0.0213	4.356	mg/L	0.0213	0.49%
	QC value within limits for Sb 217.582* Recovery = 96.80%							
Se 196.026*†	2710.5	1.536	mg/L	0.0078	1.536	mg/L	0.0078	0.51%
	QC value within limits for Se 196.026* Recovery = 102.41%							
Si 251.611*†	207337.9	5.903	mg/L	0.0448	5.903	mg/L	0.0448	0.76%
	QC value within limits for Si 251.611* Recovery = 98.38%							
Sn 189.927*†	15430.2	3.048	mg/L	0.0216	3.048	mg/L	0.0216	0.71%
	QC value within limits for Sn 189.927* Recovery = 101.59%							
Sn 242.170†	4821.1	2.992	mg/L	0.0395	2.992	mg/L	0.0395	1.32%
Sr 407.771*†	109828.3	0.2959	mg/L	0.00005	0.2959	mg/L	0.00005	0.02%
	QC value within limits for Sr 407.771* Recovery = 98.62%							

Ti 334.940†	482106.2	0.5930 mg/L	0.00020	0.5930 mg/L	0.00020	0.03%
Ti 336.121*†	340660.6	0.6068 mg/L	0.00031	0.6068 mg/L	0.00031	0.05%
QC value within limits for Ti 336.121* Recovery = 101.13%						
Tl 190.801*†	2417.5	1.525 mg/L	0.0115	1.525 mg/L	0.0115	0.76%
QC value within limits for Tl 190.801* Recovery = 101.65%						
V 292.402*†	235797.7	1.888 mg/L	0.0022	1.888 mg/L	0.0022	0.12%
QC value within limits for V 292.402* Recovery = 100.67%						
Zn 206.200*†	82726.3	2.558 mg/L	0.0217	2.558 mg/L	0.0217	0.85%
QC value within limits for Zn 206.200* Recovery = 102.34%						

All analyte(s) passed QC.

Sequence No.: 12
 Sample ID: CCB-R12091601
 Analyst: 935 icp 7300
 Initial Sample Wt:
 Dilution:
 Wash Time: 15

Autosampler Location: 1
 Date Collected: 3/29/2017 5:15:02 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol:
 Auto Dilution Factor: 1

Mean Data: CCB-R12091601

Analyte	Mean Corrected Intensity	Conc. Units	Calib.	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	85492.7	107.0 %		0.27			0.25%
Tb 350	138055.2	109.5 %		0.60			0.55%
Ag 328.068*†	-29.4	-0.0002 mg/L		0.00099	-0.0002 mg/L	0.00099	599.31%
Al 308.215*†	12.3	0.0008 mg/L		0.00200	0.0008 mg/L	0.00200	250.17%
As 188.979†	10.2	0.0059 mg/L		0.00304	0.0059 mg/L	0.00304	51.95%
As 193.696*†	3.8	0.0031 mg/L		0.00286	0.0031 mg/L	0.00286	91.85%
QC value within limits for As 193.696* Recovery = Not calculated							
B 249.677*†	1583.3	0.0358 mg/L		0.00164	0.0358 mg/L	0.00164	4.59%
Ba 233.527*†	516.1	0.0037 mg/L		0.00005	0.0037 mg/L	0.00005	1.24%
Be 313.042*†	882.7	0.0002 mg/L		0.00003	0.0002 mg/L	0.00003	11.53%
Ca 317.933*†	42.6	0.0192 mg/L		0.00319	0.0192 mg/L	0.00319	16.60%
Cd 226.502*†	38.4	0.0006 mg/L		0.00002	0.0006 mg/L	0.00002	3.97%
Cd 228.802†	24.0	0.0007 mg/L		0.00016	0.0007 mg/L	0.00016	22.27%
Co 228.616*†	33.7	0.0014 mg/L		0.00033	0.0014 mg/L	0.00033	23.45%
Cr 267.716*†	52.8	0.0005 mg/L		0.00003	0.0005 mg/L	0.00003	5.70%
Cu 324.752*†	33.6	0.0001 mg/L		0.00061	0.0001 mg/L	0.00061	437.36%
Fe 273.955*†	230.7	0.0112 mg/L		0.00027	0.0112 mg/L	0.00027	2.41%
K 766.490*†	-90.9	-0.0270 mg/L		0.05264	-0.0270 mg/L	0.05264	195.05%
Mg 279.077*†	99.6	0.0065 mg/L		0.00078	0.0065 mg/L	0.00078	12.08%
Mn 257.610*†	383.7	0.0007 mg/L		0.00002	0.0007 mg/L	0.00002	2.82%
Mo 202.031*†	7.0	0.0008 mg/L		0.00047	0.0008 mg/L	0.00047	56.36%
Na 589.592*†	-64.9	-0.0125 mg/L		0.03335	-0.0125 mg/L	0.03335	266.57%
Ni 231.604*†	34.4	0.0016 mg/L		0.00053	0.0016 mg/L	0.00053	32.75%
P 213.617*†	24.3	0.0149 mg/L		0.00462	0.0149 mg/L	0.00462	31.00%
P 214.914†	7.7	0.0077 mg/L		0.00525	0.0077 mg/L	0.00525	68.51%
Pb 220.353*†	15.8	0.0025 mg/L		0.00069	0.0025 mg/L	0.00069	27.93%
Sb 206.836†	25.2	0.0149 mg/L		0.00252	0.0149 mg/L	0.00252	16.96%
Sb 217.582*†	3.5	0.0020 mg/L		0.00000	0.0020 mg/L	0.00000	0.02%
QC value within limits for Sb 217.582* Recovery = Not calculated							
Se 196.026*†	1.6	0.0009 mg/L		0.01087	0.0009 mg/L	0.01087	>999.9%
QC value within limits for Se 196.026* Recovery = Not calculated							
Si 251.611*†	215.4	0.0061 mg/L		0.00111	0.0061 mg/L	0.00111	18.04%
Sn 189.927*†	17.6	0.0035 mg/L		0.00377	0.0035 mg/L	0.00377	108.28%
Sn 242.170†	12.9	0.0080 mg/L		0.01297	0.0080 mg/L	0.01297	162.37%
Sr 407.771*†	78.5	0.0002 mg/L		0.00001	0.0002 mg/L	0.00001	7.02%
Ti 334.940†	1179.7	0.0015 mg/L		0.00002	0.0015 mg/L	0.00002	1.19%
Ti 336.121*†	574.8	0.0010 mg/L		0.00024	0.0010 mg/L	0.00024	23.63%
Tl 190.801*†	5.2	0.0033 mg/L		0.00462	0.0033 mg/L	0.00462	139.79%
QC value within limits for Tl 190.801* Recovery = Not calculated							
V 292.402*†	145.8	0.0012 mg/L		0.00026	0.0012 mg/L	0.00026	22.48%
Zn 206.200*†	-73.1	-0.0023 mg/L		0.00034	-0.0023 mg/L	0.00034	14.97%
QC value within limits for Zn 206.200* Recovery = Not calculated							

All analyte(s) passed QC.

Return to Contents

Sequence No.: 23

Autosampler Location: 3

Sample ID: CCV= STD3x0.5

Date Collected: 3/29/2017 5:25:20 PM

Analyst: 935 icp 7300

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time: 15

Auto Dilution Factor: 1

Mean Data: CCV= STD3x0.5

Analyte	Mean Corrected		Calib. Units	Std.Dev.	Sample		Std.Dev.	RSD
	Intensity	Conc.			Conc.	Units		
Tb 384	81435.6	101.9	%	0.40				0.40%
Tb 350	135223.7	107.2	%	0.48				0.44%
Ag 328.068*†	64767.7	0.3659	mg/L	0.00231	0.3659	mg/L	0.00231	0.63%
QC value within limits for Ag 328.068* Recovery = 97.57%								
Al 308.215*†	203883.0	13.17	mg/L	0.055	13.17	mg/L	0.055	0.41%
QC value within limits for Al 308.215* Recovery = 97.56%								
As 188.979†	6388.3	3.665	mg/L	0.0016	3.665	mg/L	0.0016	0.04%
QC value within limits for As 188.979 Recovery = 97.74%								
As 193.696*†	4457.4	3.640	mg/L	0.0044	3.640	mg/L	0.0044	0.12%
QC value within limits for As 193.696* Recovery = 97.06%								
B 249.677*†	168270.0	3.804	mg/L	0.0629	3.804	mg/L	0.0629	1.65%
QC value within limits for B 249.677* Recovery = 101.45%								
Ba 233.527*†	1031883.5	7.487	mg/L	0.0363	7.487	mg/L	0.0363	0.48%
QC value within limits for Ba 233.527* Recovery = 99.82%								
Be 313.042*†	1994642.8	0.5525	mg/L	0.00361	0.5525	mg/L	0.00361	0.65%
QC value within limits for Be 313.042* Recovery = 98.23%								
Ca 317.933*†	64567.9	29.13	mg/L	0.189	29.13	mg/L	0.189	0.65%
QC value within limits for Ca 317.933* Recovery = 97.11%								
Cd 226.502*†	48812.9	0.7690	mg/L	0.00345	0.7690	mg/L	0.00345	0.45%
QC value within limits for Cd 226.502* Recovery = 102.53%								
Cd 228.802†	25357.6	0.7643	mg/L	0.00768	0.7643	mg/L	0.00768	1.00%
Co 228.616*†	46641.3	1.953	mg/L	0.0152	1.953	mg/L	0.0152	0.78%
QC value within limits for Co 228.616* Recovery = 104.18%								
Cr 267.716*†	61115.2	0.6053	mg/L	0.00335	0.6053	mg/L	0.00335	0.55%
QC value within limits for Cr 267.716* Recovery = 100.89%								
Cu 324.752*†	223311.5	0.9273	mg/L	0.00292	0.9273	mg/L	0.00292	0.31%
QC value within limits for Cu 324.752* Recovery = 98.92%								
Fe 273.955*†	80306.7	3.895	mg/L	0.0292	3.895	mg/L	0.0292	0.75%
QC value within limits for Fe 273.955* Recovery = 103.87%								
K 766.490*†	85015.0	25.25	mg/L	0.135	25.25	mg/L	0.135	0.53%
QC value within limits for K 766.490* Recovery = 93.52%								
Mg 279.077*†	119994.5	7.780	mg/L	0.0674	7.780	mg/L	0.0674	0.87%
QC value within limits for Mg 279.077* Recovery = 103.73%								
Mn 257.610*†	437996.9	0.7437	mg/L	0.00391	0.7437	mg/L	0.00391	0.53%
QC value within limits for Mn 257.610* Recovery = 99.16%								
Mo 202.031*†	4895.6	0.5836	mg/L	0.00110	0.5836	mg/L	0.00110	0.19%
QC value within limits for Mo 202.031* Recovery = 97.27%								
Na 589.592*†	179185.8	34.56	mg/L	0.022	34.56	mg/L	0.022	0.06%
QC value within limits for Na 589.592* Recovery = 96.01%								
Ni 231.604*†	12659.0	0.5962	mg/L	0.00327	0.5962	mg/L	0.00327	0.55%
QC value within limits for Ni 231.604* Recovery = 99.37%								
P 213.617*†	9320.1	5.719	mg/L	0.0266	5.719	mg/L	0.0266	0.47%
QC value within limits for P 213.617* Recovery = 95.32%								
P 214.914†	5980.7	5.913	mg/L	0.0117	5.913	mg/L	0.0117	0.20%
Pb 220.353*†	24766.8	3.864	mg/L	0.0321	3.864	mg/L	0.0321	0.83%
QC value within limits for Pb 220.353* Recovery = 103.04%								
Sb 206.836†	7258.7	4.282	mg/L	0.0115	4.282	mg/L	0.0115	0.27%
QC value within limits for Sb 206.836 Recovery = 95.15%								
Sb 217.582*†	7257.5	4.253	mg/L	0.0259	4.253	mg/L	0.0259	0.61%
QC value within limits for Sb 217.582* Recovery = 94.51%								
Se 196.026*†	2625.1	1.488	mg/L	0.0028	1.488	mg/L	0.0028	0.19%
QC value within limits for Se 196.026* Recovery = 99.19%								
Si 251.611*†	207433.5	5.905	mg/L	0.0705	5.905	mg/L	0.0705	1.19%
QC value within limits for Si 251.611* Recovery = 98.42%								
Sn 189.927*†	15032.8	2.969	mg/L	0.0001	2.969	mg/L	0.0001	0.00%
QC value within limits for Sn 189.927* Recovery = 98.97%								
Sn 242.170†	4688.7	2.909	mg/L	0.0194	2.909	mg/L	0.0194	0.67%
Sr 407.771*†	108181.2	0.2914	mg/L	0.00101	0.2914	mg/L	0.00101	0.35%
QC value within limits for Sr 407.771* Recovery = 97.14%								

Return to Contents

Ti 334.940†	476182.1	0.5857 mg/L	0.00273	0.5857 mg/L	0.00273	0.47%
Ti 336.121*†	336509.2	0.5994 mg/L	0.00320	0.5994 mg/L	0.00320	0.53%
QC value within limits for Ti 336.121* Recovery = 99.90%						
Tl 190.801*†	2345.8	1.479 mg/L	0.0005	1.479 mg/L	0.0005	0.03%
QC value within limits for Tl 190.801* Recovery = 98.63%						
V 292.402*†	235083.9	1.882 mg/L	0.0110	1.882 mg/L	0.0110	0.59%
QC value within limits for V 292.402* Recovery = 100.37%						
Zn 206.200*†	83561.8	2.584 mg/L	0.0138	2.584 mg/L	0.0138	0.53%
QC value within limits for Zn 206.200* Recovery = 103.37%						

All analyte(s) passed QC.

Sequence No.: 24
Sample ID: CCB-R12091601
Analyst: 935 icp 7300
Initial Sample Wt:
Dilution:
Wash Time: 15

Autosampler Location: 1
Date Collected: 3/29/2017 5:26:13 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1

Mean Data: CCB-R12091601

Table with 9 columns: Analyte, Mean Corrected Intensity, Conc., Calib. Units, Std.Dev., Sample Conc. Units, Std.Dev., RSD. Lists various elements like Tb, Ag, Al, As, B, Ba, Be, Ca, Cd, Co, Cr, Cu, Fe, K, Mg, Mn, Mo, Na, Ni, P, Pb, Sb, Se, Si, Sn, Sr, Ti, Tl, V, Zn with their respective values and recovery percentages.

Sequence No.: 25
Sample ID: CCB-R12091601
Analyst: 935 icp 7300
Initial Sample Wt:
Dilution:
Wash Time: 15

Handwritten note: 1012 3/20/17

Autosampler Location: 1
Date Collected: 3/29/2017 5:26:59 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1

Mean Data: CCB-R12091601

Table with 9 columns: Analyte, Mean Corrected Intensity, Conc., Calib. Units, Std.Dev., Sample Conc. Units, Std.Dev., RSD. Lists elements Tb, Ag with their respective values and recovery percentages.

Return to Contents (vertical arrow)

Sequence No.: 11

Autosampler Location: 3

Sample ID: CCV= STD3x0.5

Date Collected: 3/29/2017 6:49:54 PM

Analyst: 935 icp 7300

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time: 20

Auto Dilution Factor: 1

Mean Data: CCV= STD3x0.5

Analyte	Mean Corrected Intensity	Conc.	Calib. Units	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	79358.9	99.35	%	0.593			0.60%
Tb 350	130127.7	103.2	%	0.73			0.71%
Ag 328.068*†	64489.9	0.3643	mg/L	0.00130	0.3643 mg/L	0.00130	0.36%
QC value within limits for Ag 328.068* Recovery = 97.15%							
Al 308.215*†	203857.8	13.17	mg/L	0.094	13.17 mg/L	0.094	0.71%
QC value within limits for Al 308.215* Recovery = 97.55%							
As 188.979†	6610.2	3.792	mg/L	0.0048	3.792 mg/L	0.0048	0.13%
QC value within limits for As 188.979 Recovery = 101.13%							
As 193.696*†	4578.0	3.738	mg/L	0.0173	3.738 mg/L	0.0173	0.46%
QC value within limits for As 193.696* Recovery = 99.69%							
B 249.677*†	166568.8	3.766	mg/L	0.0262	3.766 mg/L	0.0262	0.70%
QC value within limits for B 249.677* Recovery = 100.42%							
Ba 233.527*†	1032441.6	7.491	mg/L	0.0382	7.491 mg/L	0.0382	0.51%
QC value within limits for Ba 233.527* Recovery = 99.88%							
Be 313.042*†	2001290.4	0.5544	mg/L	0.00154	0.5544 mg/L	0.00154	0.28%
QC value within limits for Be 313.042* Recovery = 98.56%							
Ca 317.933*†	64772.7	29.22	mg/L	0.788	29.22 mg/L	0.788	2.70%
QC value within limits for Ca 317.933* Recovery = 97.41%							
Cd 226.502*†	48865.7	0.7698	mg/L	0.00321	0.7698 mg/L	0.00321	0.42%
QC value within limits for Cd 226.502* Recovery = 102.64%							
Cd 228.802†	25050.9	0.7551	mg/L	0.00753	0.7551 mg/L	0.00753	1.00%
Co 228.616*†	46467.9	1.946	mg/L	0.0203	1.946 mg/L	0.0203	1.04%
QC value within limits for Co 228.616* Recovery = 103.79%							
Cr 267.716*†	61265.1	0.6068	mg/L	0.00167	0.6068 mg/L	0.00167	0.27%
QC value within limits for Cr 267.716* Recovery = 101.14%							
Cu 324.752*†	220035.0	0.9137	mg/L	0.00469	0.9137 mg/L	0.00469	0.51%
QC value within limits for Cu 324.752* Recovery = 97.46%							
Fe 273.955*†	79800.1	3.870	mg/L	0.0221	3.870 mg/L	0.0221	0.57%
QC value within limits for Fe 273.955* Recovery = 103.21%							
K 766.490*†	85973.8	25.54	mg/L	0.588	25.54 mg/L	0.588	2.30%
QC value within limits for K 766.490* Recovery = 94.58%							
Mg 279.077*†	119653.1	7.758	mg/L	0.0007	7.758 mg/L	0.0007	0.01%
QC value within limits for Mg 279.077* Recovery = 103.44%							
Mn 257.610*†	437872.1	0.7435	mg/L	0.00372	0.7435 mg/L	0.00372	0.50%
QC value within limits for Mn 257.610* Recovery = 99.13%							
Mo 202.031*†	5045.7	0.6015	mg/L	0.00475	0.6015 mg/L	0.00475	0.79%
QC value within limits for Mo 202.031* Recovery = 100.25%							
Na 589.592*†	180334.2	34.78	mg/L	0.730	34.78 mg/L	0.730	2.10%
QC value within limits for Na 589.592* Recovery = 96.62%							
Ni 231.604*†	13014.7	0.6130	mg/L	0.00535	0.6130 mg/L	0.00535	0.87%
QC value within limits for Ni 231.604* Recovery = 102.16%							
P 213.617*†	9599.6	5.891	mg/L	0.0166	5.891 mg/L	0.0166	0.28%
QC value within limits for P 213.617* Recovery = 98.18%							
P 214.914†	6163.2	6.093	mg/L	0.0358	6.093 mg/L	0.0358	0.59%
Pb 220.353*†	24394.6	3.806	mg/L	0.0014	3.806 mg/L	0.0014	0.04%
QC value within limits for Pb 220.353* Recovery = 101.49%							
Sb 206.836†	7490.0	4.418	mg/L	0.0137	4.418 mg/L	0.0137	0.31%
QC value within limits for Sb 206.836 Recovery = 98.18%							
Sb 217.582*†	7473.3	4.379	mg/L	0.0313	4.379 mg/L	0.0313	0.71%
QC value within limits for Sb 217.582* Recovery = 97.32%							
Se 196.026*†	2683.4	1.521	mg/L	0.0042	1.521 mg/L	0.0042	0.28%
QC value within limits for Se 196.026* Recovery = 101.39%							
Si 251.611*†	207523.4	5.908	mg/L	0.0097	5.908 mg/L	0.0097	0.16%
QC value within limits for Si 251.611* Recovery = 98.47%							
Sn 189.927*†	15486.6	3.059	mg/L	0.0106	3.059 mg/L	0.0106	0.35%
QC value within limits for Sn 189.927* Recovery = 101.96%							
Sn 242.170†	4811.9	2.986	mg/L	0.0001	2.986 mg/L	0.0001	0.00%
Sr 407.771*†	108771.6	0.2930	mg/L	0.00698	0.2930 mg/L	0.00698	2.38%
QC value within limits for Sr 407.771* Recovery = 97.67%							

Return to Contents

Ti 334.940†	475601.3	0.5850 mg/L	0.00425	0.5850 mg/L	0.00425	0.73%
Ti 336.121*†	335088.3	0.5968 mg/L	0.00268	0.5968 mg/L	0.00268	0.45%
QC value within limits for Ti 336.121* Recovery = 99.47%						
Tl 190.801*†	2408.1	1.519 mg/L	0.0018	1.519 mg/L	0.0018	0.12%
QC value within limits for Tl 190.801* Recovery = 101.25%						
V 292.402*†	234906.7	1.880 mg/L	0.0083	1.880 mg/L	0.0083	0.44%
QC value within limits for V 292.402* Recovery = 100.29%						
Zn 206.200*†	82386.9	2.548 mg/L	0.0021	2.548 mg/L	0.0021	0.08%
QC value within limits for Zn 206.200* Recovery = 101.92%						
Zn 213.857*†	152037.2	2.547 mg/L	0.0010	2.547 mg/L	0.0010	0.04%
QC value within limits for Zn 213.857* Recovery = 101.89%						

All analyte(s) passed QC.

Sequence No.: 12

Autosampler Location: 1

Sample ID: CCB-R12091601

Date Collected: 3/29/2017 6:50:47 PM

Analyst: 935 icp 7300

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time: 15

Auto Dilution Factor: 1

Mean Data: CCB-R12091601

Analyte	Mean Corrected Intensity	Conc.	Calib. Units	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	87734.0	109.8	%	0.53			0.48%
Tb 350	142073.9	112.6	%	0.01			0.01%
Ag 328.068*†	-202.4	-0.0011	mg/L	0.00060	-0.0011 mg/L	0.00060	52.66%
QC value within limits for Ag 328.068* Recovery = Not calculated							
Al 308.215*†	55.8	0.0036	mg/L	0.00301	0.0036 mg/L	0.00301	83.41%
As 188.979†	9.0	0.0052	mg/L	0.00180	0.0052 mg/L	0.00180	34.80%
As 193.696*†	8.5	0.0070	mg/L	0.00103	0.0070 mg/L	0.00103	14.84%
QC value within limits for As 193.696* Recovery = Not calculated							
B 249.677*†	1804.5	0.0408	mg/L	0.00235	0.0408 mg/L	0.00235	5.76%
Ba 233.527*†	903.8	0.0066	mg/L	0.00021	0.0066 mg/L	0.00021	3.15%
Be 313.042*†	1629.8	0.0005	mg/L	0.00003	0.0005 mg/L	0.00003	5.56%
Ca 317.933*†	100.4	0.0453	mg/L	0.00273	0.0453 mg/L	0.00273	6.02%
Cd 226.502*†	51.1	0.0008	mg/L	0.00005	0.0008 mg/L	0.00005	6.22%
Cd 228.802†	31.5	0.0010	mg/L	0.00002	0.0010 mg/L	0.00002	2.34%
Co 228.616*†	51.2	0.0021	mg/L	0.00006	0.0021 mg/L	0.00006	2.86%
Cr 267.716*†	74.0	0.0007	mg/L	0.00014	0.0007 mg/L	0.00014	18.73%
Cu 324.752*†	251.3	0.0010	mg/L	0.00016	0.0010 mg/L	0.00016	15.22%
Fe 273.955*†	304.0	0.0147	mg/L	0.00001	0.0147 mg/L	0.00001	0.04%
K 766.490*†	-265.0	-0.0787	mg/L	0.08257	-0.0787 mg/L	0.08257	104.91%
Mg 279.077*†	314.3	0.0204	mg/L	0.00559	0.0204 mg/L	0.00559	27.42%
Mn 257.610*†	481.5	0.0008	mg/L	0.00007	0.0008 mg/L	0.00007	8.30%
Mo 202.031*†	12.2	0.0015	mg/L	0.00051	0.0015 mg/L	0.00051	34.63%
Na 589.592*†	82.2	0.0159	mg/L	0.00070	0.0159 mg/L	0.00070	4.39%
Ni 231.604*†	33.0	0.0016	mg/L	0.00001	0.0016 mg/L	0.00001	0.68%
P 213.617*†	16.8	0.0103	mg/L	0.00318	0.0103 mg/L	0.00318	30.86%
P 214.914†	5.0	0.0049	mg/L	0.00242	0.0049 mg/L	0.00242	49.24%
Pb 220.353*†	14.3	0.0022	mg/L	0.00144	0.0022 mg/L	0.00144	64.25%
Sb 206.836†	23.0	0.0136	mg/L	0.00106	0.0136 mg/L	0.00106	7.78%
Sb 217.582*†	16.0	0.0094	mg/L	0.00502	0.0094 mg/L	0.00502	53.48%
QC value within limits for Sb 217.582* Recovery = Not calculated							
Se 196.026*†	1.9	0.0011	mg/L	0.00363	0.0011 mg/L	0.00363	332.02%
QC value within limits for Se 196.026* Recovery = Not calculated							
Si 251.611*†	115.3	0.0033	mg/L	0.00079	0.0033 mg/L	0.00079	23.98%
Sn 189.927*†	25.3	0.0050	mg/L	0.00074	0.0050 mg/L	0.00074	14.85%
Sn 242.170†	23.5	0.0146	mg/L	0.00024	0.0146 mg/L	0.00024	1.64%
Sr 407.771*†	151.1	0.0004	mg/L	0.00002	0.0004 mg/L	0.00002	4.64%
Ti 334.940†	1313.5	0.0016	mg/L	0.00009	0.0016 mg/L	0.00009	5.83%
Ti 336.121*†	789.6	0.0014	mg/L	0.00000	0.0014 mg/L	0.00000	0.19%
Tl 190.801*†	16.8	0.0106	mg/L	0.00149	0.0106 mg/L	0.00149	14.01%
V 292.402*†	203.2	0.0016	mg/L	0.00014	0.0016 mg/L	0.00014	8.59%
Zn 206.200*†	-7.2	-0.0002	mg/L	0.00011	-0.0002 mg/L	0.00011	50.11%
QC value within limits for Zn 206.200* Recovery = Not calculated							
Zn 213.857*†	28.0	0.0005	mg/L	0.00026	0.0005 mg/L	0.00026	54.66%
QC value within limits for Zn 213.857* Recovery = Not calculated							

All analyte(s) passed QC.

=====
Analysis Begun

Start Time: 3/29/2017 7:02:35 PM
Logged In Analyst: Oscar Gomez 935
Spectrometer: Optima 7300 DV, S/N 77c8120401

Plasma On Time: 3/29/2017 4:49:39 PM
Technique: ICP Continuous
Autosampler: ESI

Sample Information File: C:\Documents and Settings\All Users\PerkinElmer\ICP\Data\Sample Information\
17032902.sif

Batch ID:
Results Data Set: 170329C 1
Results Library: W:\pe\7300\Results\results.mdb

=====
Sequence No.: 1
Sample ID: CCV= STD3x0.5
Analyst: 935 icp 7300
Initial Sample Wt:
Dilution:
Wash Time:
Autosampler Location: 3
Date Collected: 3/29/2017 7:02:53 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:

Mean Data: CCV= STD3x0.5

Analyte	Mean Corrected		Calib.	Std.Dev.	Sample		RSD
	Intensity	Conc.			Units	Std.Dev.	
Tb 384	79586.6	99.63	%	1.006			1.01%
Tb 350	132281.0	104.9	%	0.79			0.75%
Ag 328.068*†	64769.0	0.3659	mg/L	0.00156	0.3659	mg/L	0.00156 0.43%
	QC value within limits for Ag 328.068* Recovery = 97.57%						
Al 308.215*†	206702.8	13.35	mg/L	0.035	13.35	mg/L	0.035 0.26%
	QC value within limits for Al 308.215* Recovery = 98.91%						
As 188.979†	6628.4	3.803	mg/L	0.0266	3.803	mg/L	0.0266 0.70%
	QC value within limits for As 188.979 Recovery = 101.41%						
As 193.696*†	4611.9	3.766	mg/L	0.0197	3.766	mg/L	0.0197 0.52%
	QC value within limits for As 193.696* Recovery = 100.43%						
B 249.677*†	169490.4	3.832	mg/L	0.0216	3.832	mg/L	0.0216 0.56%
	QC value within limits for B 249.677* Recovery = 102.18%						
Ba 233.527*†	1046107.9	7.590	mg/L	0.0449	7.590	mg/L	0.0449 0.59%
	QC value within limits for Ba 233.527* Recovery = 101.20%						
Be 313.042*†	2013190.4	0.5577	mg/L	0.00408	0.5577	mg/L	0.00408 0.73%
	QC value within limits for Be 313.042* Recovery = 99.14%						
Ca 317.933*†	63373.3	28.59	mg/L	0.142	28.59	mg/L	0.142 0.50%
	QC value within limits for Ca 317.933* Recovery = 95.31%						
Cd 226.502*†	49581.9	0.7811	mg/L	0.00335	0.7811	mg/L	0.00335 0.43%
	QC value within limits for Cd 226.502* Recovery = 104.14%						
Cd 228.802†	25482.2	0.7681	mg/L	0.00480	0.7681	mg/L	0.00480 0.62%
Co 228.616*†	47208.3	1.977	mg/L	0.0215	1.977	mg/L	0.0215 1.09%
	QC value within limits for Co 228.616* Recovery = 105.44%						
Cr 267.716*†	61869.8	0.6128	mg/L	0.00375	0.6128	mg/L	0.00375 0.61%
	QC value within limits for Cr 267.716* Recovery = 102.14%						
Cu 324.752*†	222419.9	0.9236	mg/L	0.00146	0.9236	mg/L	0.00146 0.16%
	QC value within limits for Cu 324.752* Recovery = 98.52%						
Fe 273.955*†	81797.6	3.967	mg/L	0.0015	3.967	mg/L	0.0015 0.04%
	QC value within limits for Fe 273.955* Recovery = 105.79%						
K 766.490*†	83769.4	24.88	mg/L	0.138	24.88	mg/L	0.138 0.55%
	QC value within limits for K 766.490* Recovery = 92.15%						
Mg 279.077*†	121598.4	7.884	mg/L	0.0339	7.884	mg/L	0.0339 0.43%
	QC value within limits for Mg 279.077* Recovery = 105.12%						
Mn 257.610*†	444131.6	0.7541	mg/L	0.00477	0.7541	mg/L	0.00477 0.63%
	QC value within limits for Mn 257.610* Recovery = 100.55%						
Mo 202.031*†	5019.4	0.5984	mg/L	0.00311	0.5984	mg/L	0.00311 0.52%
	QC value within limits for Mo 202.031* Recovery = 99.73%						
Na 589.592*†	175700.8	33.89	mg/L	0.219	33.89	mg/L	0.219 0.64%
	QC value within limits for Na 589.592* Recovery = 94.14%						
Ni 231.604*†	13085.4	0.6163	mg/L	0.00407	0.6163	mg/L	0.00407 0.66%
	QC value within limits for Ni 231.604* Recovery = 102.72%						
P 213.617*†	9658.1	5.927	mg/L	0.0659	5.927	mg/L	0.0659 1.11%
	QC value within limits for P 213.617* Recovery = 98.78%						
P 214.914†	6201.9	6.131	mg/L	0.0637	6.131	mg/L	0.0637 1.04%
Pb 220.353*†	25078.8	3.913	mg/L	0.0044	3.913	mg/L	0.0044 0.11%
	QC value within limits for Pb 220.353* Recovery = 104.34%						

Sb 206.836†	7472.2	4.408 mg/L	0.0302	4.408 mg/L	0.0302	0.68%
QC value within limits for Sb 206.836 Recovery = 97.95%						
Sb 217.582*†	7454.0	4.368 mg/L	0.0376	4.368 mg/L	0.0376	0.86%
QC value within limits for Sb 217.582* Recovery = 97.07%						
Se 196.026*†	2717.6	1.540 mg/L	0.0261	1.540 mg/L	0.0261	1.69%
QC value within limits for Se 196.026* Recovery = 102.68%						
Si 251.611*†	210785.1	6.001 mg/L	0.0030	6.001 mg/L	0.0030	0.05%
QC value within limits for Si 251.611* Recovery = 100.01%						
Sn 189.927*†	15533.5	3.068 mg/L	0.0333	3.068 mg/L	0.0333	1.08%
QC value within limits for Sn 189.927* Recovery = 102.27%						
Sn 242.170†	4812.8	2.986 mg/L	0.0384	2.986 mg/L	0.0384	1.29%
Sr 407.771*†	105832.0	0.2851 mg/L	0.00107	0.2851 mg/L	0.00107	0.37%
QC value within limits for Sr 407.771* Recovery = 95.03%						
Ti 334.940†	498627.4	0.6133 mg/L	0.00059	0.6133 mg/L	0.00059	0.10%
Ti 336.121*†	351352.1	0.6258 mg/L	0.00036	0.6258 mg/L	0.00036	0.06%
QC value within limits for Ti 336.121* Recovery = 104.30%						
Tl 190.801*†	2418.5	1.525 mg/L	0.0154	1.525 mg/L	0.0154	1.01%
QC value within limits for Tl 190.801* Recovery = 101.69%						
V 292.402*†	236215.1	1.891 mg/L	0.0111	1.891 mg/L	0.0111	0.59%
QC value within limits for V 292.402* Recovery = 100.85%						
Zn 206.200*†	84121.6	2.602 mg/L	0.0160	2.602 mg/L	0.0160	0.61%
QC value within limits for Zn 206.200* Recovery = 104.07%						
Zn 213.857*†	154374.6	2.587 mg/L	0.0121	2.587 mg/L	0.0121	0.47%
QC value within limits for Zn 213.857* Recovery = 103.46%						

All analyte(s) passed QC.

Sequence No.: 2
 Sample ID: CCB-R12091601
 Analyst: 935 icp 7300
 Initial Sample Wt:
 Dilution:
 Wash Time: 15

Autosampler Location: 1
 Date Collected: 3/29/2017 7:03:52 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol:
 Auto Dilution Factor: 1

Mean Data: CCB-R12091601

Analyte	Mean Corrected Intensity	Conc.	Calib. Units	Std.Dev.	Sample Conc. Units	Std.Dev.	RSD
Tb 384	86528.6	108.3	%	1.11			1.03%
Tb 350	140513.5	111.4	%	0.62			0.56%
Ag 328.068*†	-9.1	-0.0001	mg/L	0.00175	-0.0001 mg/L	0.00175	>999.9%
QC value within limits for Ag 328.068* Recovery = Not calculated							
Al 308.215*†	102.9	0.0066	mg/L	0.00182	0.0066 mg/L	0.00182	27.33%
As 188.979†	12.3	0.0071	mg/L	0.00204	0.0071 mg/L	0.00204	28.92%
As 193.696*†	8.1	0.0066	mg/L	0.00949	0.0066 mg/L	0.00949	144.29%
QC value within limits for As 193.696* Recovery = Not calculated							
B 249.677*†	1567.3	0.0354	mg/L	0.00230	0.0354 mg/L	0.00230	6.50%
Ba 233.527*†	955.7	0.0069	mg/L	0.00030	0.0069 mg/L	0.00030	4.32%
Be 313.042*†	1739.5	0.0005	mg/L	0.00003	0.0005 mg/L	0.00003	6.31%
Ca 317.933*†	94.7	0.0427	mg/L	0.00574	0.0427 mg/L	0.00574	13.43%
Cd 226.502*†	57.7	0.0009	mg/L	0.00021	0.0009 mg/L	0.00021	22.84%
Cd 228.802†	27.7	0.0008	mg/L	0.00013	0.0008 mg/L	0.00013	15.14%
Co 228.616*†	46.1	0.0019	mg/L	0.00021	0.0019 mg/L	0.00021	10.71%
Cr 267.716*†	66.7	0.0007	mg/L	0.00010	0.0007 mg/L	0.00010	15.18%
Cu 324.752*†	171.7	0.0007	mg/L	0.00002	0.0007 mg/L	0.00002	2.54%
Fe 273.955*†	251.3	0.0122	mg/L	0.00137	0.0122 mg/L	0.00137	11.27%
K 766.490*†	28.5	0.0085	mg/L	0.02904	0.0085 mg/L	0.02904	342.66%
Mg 279.077*†	266.7	0.0173	mg/L	0.00666	0.0173 mg/L	0.00666	38.55%
Mn 257.610*†	466.6	0.0008	mg/L	0.00002	0.0008 mg/L	0.00002	2.32%
Mo 202.031*†	13.1	0.0016	mg/L	0.00061	0.0016 mg/L	0.00061	38.97%
Na 589.592*†	70.3	0.0136	mg/L	0.00083	0.0136 mg/L	0.00083	6.09%
Ni 231.604*†	33.4	0.0016	mg/L	0.00028	0.0016 mg/L	0.00028	18.00%
P 213.617*†	18.5	0.0114	mg/L	0.02398	0.0114 mg/L	0.02398	210.96%
P 214.914†	12.2	0.0120	mg/L	0.00803	0.0120 mg/L	0.00803	66.82%
Pb 220.353*†	14.5	0.0023	mg/L	0.00283	0.0023 mg/L	0.00283	124.58%
Sb 206.836†	30.8	0.0182	mg/L	0.00217	0.0182 mg/L	0.00217	11.97%
Sb 217.582*†	6.3	0.0037	mg/L	0.00046	0.0037 mg/L	0.00046	12.40%
QC value within limits for Sb 217.582* Recovery = Not calculated							
Se 196.026*†	-3.1	-0.0017	mg/L	0.00163	-0.0017 mg/L	0.00163	94.41%
QC value within limits for Se 196.026* Recovery = Not calculated							
Si 251.611*†	183.7	0.0052	mg/L	0.00061	0.0052 mg/L	0.00061	11.75%
Sn 189.927*†	32.2	0.0064	mg/L	0.00113	0.0064 mg/L	0.00113	17.82%
Sn 242.170†	15.5	0.0096	mg/L	0.00526	0.0096 mg/L	0.00526	54.63%
Sr 407.771*†	143.0	0.0004	mg/L	0.00001	0.0004 mg/L	0.00001	2.23%
Ti 334.940†	2050.8	0.0025	mg/L	0.00051	0.0025 mg/L	0.00051	20.24%
Ti 336.121*†	1392.1	0.0025	mg/L	0.00041	0.0025 mg/L	0.00041	16.46%
Tl 190.801*†	11.2	0.0071	mg/L	0.00106	0.0071 mg/L	0.00106	14.95%
V 292.402*†	238.7	0.0019	mg/L	0.00003	0.0019 mg/L	0.00003	1.54%
Zn 206.200*†	-20.9	-0.0006	mg/L	0.00015	-0.0006 mg/L	0.00015	23.12%
QC value within limits for Zn 206.200* Recovery = Not calculated							
Zn 213.857*†	7.3	0.0001	mg/L	0.00003	0.0001 mg/L	0.00003	21.42%
QC value within limits for Zn 213.857* Recovery = Not calculated							

All analyte(s) passed QC.

EPA 6010B ICP Metals (Solid)

Sample Data

RAW DATA SHEET FOR METHOD: EPA 6010B

WORK ORDER: 17-03-1557
INSTRUMENT: ICP 7300
EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 18:56
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\ICP-DATA\170329C 1\17-03-1557-28.icp

28 **CLIENT SAMPLE NUMBER: B-DU1-ISM1-8**

LCS/MB BATCH: 170329L05 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 2.00 g / ACTUAL: 2.00 g
MS/MSD BATCH: 170329S05 **FINAL VOLUME / WEIGHT:** DEFAULT: 100.00 ml / ACTUAL: 100.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Antimony	0.00788	0.985	ND	0.739	
Arsenic	0.0278	0.985	1.37	0.739	
Barium	1.24	0.985	61.1	0.493	
Beryllium	0.00296	0.985	ND	0.246	
Cadmium	0.00500	0.985	ND	0.493	
Chromium	0.0984	0.985	4.85	0.246	
Cobalt	0.0999	0.985	4.92	0.246	
Copper	0.141	0.985	6.93	0.493	
Lead	0.0892	0.985	4.39	0.493	
Molybdenum	0.00171	0.985	ND	0.246	
Nickel	0.0875	0.985	4.31	0.246	
Selenium	-0.00306	0.985	ND	0.739	
Silver	-0.00316	0.985	ND	0.246	
Thallium	-0.00717	0.985	ND	0.739	
Vanadium	0.335	0.985	16.5	0.246	
Zinc	0.493	0.985	24.3	0.985	



Sequence No.: 3
 Sample ID: 17-03-1557-28
 Analyst: 1030 icp 7300
 Initial Sample Wt: 2.03 g
 Dilution:
 Wash Time: 15

Autosampler Location: 134
 Date Collected: 3/29/2017 6:56:44 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Mean Data: 17-03-1557-28

Analyte	Mean Corrected		Calib.	Std.Dev.	Sample		Std.Dev.	RSD
	Intensity	Conc. Units			Conc. Units	Std.Dev.		
Tb 384	67749.9	84.81 %	%	0.497				0.59%
Tb 350	126349.2	100.2 %	%	0.88				0.88%
Ag 328.068*†	-560.0	-0.0032	mg/L	0.00060	-0.1559	mg/kg	0.02970	19.05%
Al 308.215*†	1713264.2	110.7	mg/L	0.14	5452	mg/kg	6.74	0.12%
As 188.979†	51.1	0.0293	mg/L	0.00283	1.445	mg/kg	0.1396	9.66%
As 193.696*†	34.0	0.0278	mg/L	0.00528	1.368	mg/kg	0.2602	19.03%
B 249.677*†	747.5	0.0169	mg/L	0.00601	0.8325	mg/kg	0.29591	35.55%
Ba 233.527*†	170926.8	1.240	mg/L	0.0041	61.09	mg/kg	0.204	0.33%
Be 313.042*†	10688.1	0.0030	mg/L	0.00001	0.1459	mg/kg	0.00041	0.28%
Ca 317.933*†	133803.9	60.37	mg/L	1.311	2974	mg/kg	64.59	2.17%
Cd 226.502*†	317.2	0.0050	mg/L	0.00000	0.2462	mg/kg	0.00017	0.07%
Cd 228.802†	33.0	0.0010	mg/L	0.00009	0.0490	mg/kg	0.00422	8.60%
Co 228.616*†	2384.6	0.0999	mg/L	0.00082	4.920	mg/kg	0.0402	0.82%
Cr 267.716*†	9931.3	0.0984	mg/L	0.00123	4.846	mg/kg	0.0607	1.25%
Cu 324.752*†	33860.0	0.1406	mg/L	0.00006	6.927	mg/kg	0.0031	0.05%
Fe 273.955*†	3544151.7	171.9	mg/L	4.87	8468	mg/kg	240.01	2.83%
K 766.490*†	107408.9	31.90	mg/L	0.603	1572	mg/kg	29.73	1.89%
Mg 279.077*†	873986.6	56.67	mg/L	0.339	2791	mg/kg	16.70	0.60%
Mn 257.610*†	1625277.9	2.760	mg/L	0.0087	135.9	mg/kg	0.43	0.32%
Mo 202.031*†	14.4	0.0017	mg/L	0.00082	0.0844	mg/kg	0.04048	47.98%
Na 589.592*†	19653.9	3.791	mg/L	0.0849	186.7	mg/kg	4.18	2.24%
Ni 231.604*†	1857.3	0.0875	mg/L	0.00109	4.309	mg/kg	0.0539	1.25%
P 213.617*†	16660.7	10.22	mg/L	0.093	503.6	mg/kg	4.56	0.91%
P 214.914†	11053.6	10.93	mg/L	0.096	538.3	mg/kg	4.73	0.88%
Pb 220.353*†	571.6	0.0892	mg/L	0.00202	4.393	mg/kg	0.0997	2.27%
Sb 206.836†	13.3	0.0078	mg/L	0.00038	0.3865	mg/kg	0.01896	4.91%
Sb 217.582*†	13.4	0.0079	mg/L	0.00251	0.3882	mg/kg	0.12353	31.82%
Se 196.026*†	-5.4	-0.0031	mg/L	0.00191	-0.1509	mg/kg	0.09426	62.48%
Si 251.611*†	311712.8	8.874	mg/L	0.1160	437.1	mg/kg	5.71	1.31%
Sn 189.927*†	-363.6	-0.0718	mg/L	0.00283	-3.538	mg/kg	0.1392	3.93%
Sn 242.170†	693.0	0.4300	mg/L	0.00465	21.18	mg/kg	0.229	1.08%
Sr 407.771*†	275485.0	0.7421	mg/L	0.01644	36.56	mg/kg	0.810	2.22%
Ti 334.940†	8744139.9	10.76	mg/L	0.259	529.8	mg/kg	12.78	2.41%
Ti 336.121*†	6232628.4	11.10	mg/L	0.272	546.9	mg/kg	13.38	2.45%
Tl 190.801*†	-11.4	-0.0072	mg/L	0.00764	-0.3531	mg/kg	0.37611	106.51%
V 292.402*†	42231.6	0.3353	mg/L	0.00458	16.52	mg/kg	0.226	1.37%
Zn 206.200*†	15942.2	0.4930	mg/L	0.00386	24.29	mg/kg	0.190	0.78%

RAW DATA SHEET FOR METHOD: EPA 6010B

WORK ORDER: 17-03-1557
INSTRUMENT: ICP 7300
EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 18:57
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\ICP-DATA\170329C 1\17-03-1557-29.icp

29 **CLIENT SAMPLE NUMBER: B-DU1-ISM2-8**

LCS/MB BATCH: 170329L05 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 2.00 g / ACTUAL: 2.00 g
MS/MSD BATCH: 170329S05 **FINAL VOLUME / WEIGHT:** DEFAULT: 100.00 ml / ACTUAL: 100.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Antimony	-0.00144	0.980	ND	0.735	
Arsenic	0.00514	0.980	ND	0.735	
Barium	1.36	0.980	66.9	0.490	
Beryllium	0.00274	0.980	ND	0.245	
Cadmium	0.00469	0.980	ND	0.490	
Chromium	0.0966	0.980	4.74	0.245	
Cobalt	0.0969	0.980	4.75	0.245	
Copper	0.137	0.980	6.71	0.490	
Lead	0.0705	0.980	3.46	0.490	
Molybdenum	0.00150	0.980	ND	0.245	
Nickel	0.0777	0.980	3.81	0.245	
Selenium	0.00322	0.980	ND	0.735	
Silver	-0.00251	0.980	ND	0.245	
Thallium	-0.00798	0.980	ND	0.735	
Vanadium	0.322	0.980	15.8	0.245	
Zinc	0.496	0.980	24.3	0.980	

Return to Contents

Sequence No.: 4
 Sample ID: 17-03-1557-29
 Analyst: 1030 icp 7300
 Initial Sample Wt: 2.04 g
 Dilution:
 Wash Time: 15

Autosampler Location: 135
 Date Collected: 3/29/2017 6:57:37 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Mean Data: 17-03-1557-29

Analyte	Mean Corrected		Calib.		Sample			
	Intensity	Conc.	Units	Std.Dev.	Conc.	Units	Std.Dev.	RSD
Tb 384	67797.0	84.87	%	0.497				0.59%
Tb 350	127607.7	101.2	%	1.32				1.30%
Ag 328.068*†	-444.4	-0.0025	mg/L	0.00113	-0.1231	mg/kg	0.05523	44.88%
Al 308.215*†	1686580.8	108.9	mg/L	0.03	5341	mg/kg	1.44	0.03%
As 188.979†	56.5	0.0324	mg/L	0.00236	1.589	mg/kg	0.1155	7.27%
As 193.696*†	6.3	0.0051	mg/L	0.00281	0.2519	mg/kg	0.13764	54.65%
B 249.677*†	397.3	0.0090	mg/L	0.00235	0.4403	mg/kg	0.11499	26.11%
Ba 233.527*†	187982.4	1.364	mg/L	0.0046	66.86	mg/kg	0.225	0.34%
Be 313.042*†	9908.1	0.0027	mg/L	0.00005	0.1345	mg/kg	0.00269	2.00%
Ca 317.933*†	130585.5	58.92	mg/L	1.096	2888	mg/kg	53.71	1.86%
Cd 226.502*†	297.7	0.0047	mg/L	0.00000	0.2299	mg/kg	0.00002	0.01%
Cd 228.802†	6.2	0.0002	mg/L	0.00017	0.0092	mg/kg	0.00855	93.14%
Co 228.616*†	2313.4	0.0969	mg/L	0.00082	4.749	mg/kg	0.0401	0.84%
Cr 267.716*†	9755.5	0.0966	mg/L	0.00104	4.737	mg/kg	0.0512	1.08%
Cu 324.752*†	32958.7	0.1369	mg/L	0.00049	6.709	mg/kg	0.0240	0.36%
Fe 273.955*†	3477701.5	168.7	mg/L	2.67	8268	mg/kg	130.64	1.58%
K 766.490*†	108959.5	32.36	mg/L	0.692	1586	mg/kg	33.92	2.14%
Mg 279.077*†	883463.9	57.28	mg/L	0.003	2808	mg/kg	0.16	0.01%
Mn 257.610*†	1605043.5	2.725	mg/L	0.0010	133.6	mg/kg	0.05	0.04%
Mo 202.031*†	12.6	0.0015	mg/L	0.00090	0.0733	mg/kg	0.04420	60.26%
Na 589.592*†	19951.6	3.848	mg/L	0.0934	188.6	mg/kg	4.58	2.43%
Ni 231.604*†	1649.8	0.0777	mg/L	0.00112	3.809	mg/kg	0.0550	1.44%
P 213.617*†	17071.7	10.48	mg/L	0.148	513.5	mg/kg	7.27	1.42%
P 214.914†	11295.0	11.17	mg/L	0.108	547.4	mg/kg	5.29	0.97%
Pb 220.353*†	451.9	0.0705	mg/L	0.00178	3.456	mg/kg	0.0871	2.52%
Sb 206.836†	13.5	0.0080	mg/L	0.00749	0.3913	mg/kg	0.36725	93.86%
Sb 217.582*†	-2.5	-0.0014	mg/L	0.00309	-0.0704	mg/kg	0.15138	214.89%
Se 196.026*†	5.7	0.0032	mg/L	0.00643	0.1578	mg/kg	0.31533	199.84%
Si 251.611*†	295745.4	8.420	mg/L	0.1392	412.7	mg/kg	6.82	1.65%
Sn 189.927*†	-333.6	-0.0659	mg/L	0.00189	-3.230	mg/kg	0.0928	2.87%
Sn 242.170†	675.3	0.4191	mg/L	0.02941	20.54	mg/kg	1.442	7.02%
Sr 407.771*†	239896.2	0.6462	mg/L	0.01298	31.68	mg/kg	0.636	2.01%
Ti 334.940†	8644708.4	10.63	mg/L	0.136	521.2	mg/kg	6.66	1.28%
Ti 336.121*†	6155086.7	10.96	mg/L	0.145	537.4	mg/kg	7.09	1.32%
Tl 190.801*†	-12.6	-0.0080	mg/L	0.01013	-0.3909	mg/kg	0.49636	126.97%
V 292.402*†	40553.4	0.3219	mg/L	0.00188	15.78	mg/kg	0.092	0.58%
Zn 206.200*†	16043.5	0.4962	mg/L	0.00573	24.32	mg/kg	0.281	1.16%

RAW DATA SHEET FOR METHOD: EPA 6010B

WORK ORDER: 17-03-1557
INSTRUMENT: ICP 7300
EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 18:58
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\ICP-DATA\170329C 1\17-03-1557-30.icp

30 **CLIENT SAMPLE NUMBER: B-DU1-ISM3-8**

LCS/MB BATCH: 170329L05 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 2.00 g / ACTUAL: 2.00 g
MS/MSD BATCH: 170329S05 **FINAL VOLUME / WEIGHT:** DEFAULT: 100.00 ml / ACTUAL: 100.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Antimony	0.00207	0.980	ND	0.735	
Arsenic	0.0194	0.980	0.950	0.735	
Barium	1.16	0.980	56.9	0.490	
Beryllium	0.00258	0.980	ND	0.245	
Cadmium	0.00405	0.980	ND	0.490	
Chromium	0.0841	0.980	4.12	0.245	
Cobalt	0.0862	0.980	4.22	0.245	
Copper	0.125	0.980	6.12	0.490	
Lead	0.0763	0.980	3.74	0.490	
Molybdenum	0.000970	0.980	ND	0.245	
Nickel	0.0709	0.980	3.48	0.245	
Selenium	0.00118	0.980	ND	0.735	
Silver	-0.00248	0.980	ND	0.245	
Thallium	-0.0116	0.980	ND	0.735	
Vanadium	0.273	0.980	13.4	0.245	
Zinc	0.473	0.980	23.2	0.980	

Return to Contents

Sequence No.: 5
 Sample ID: 17-03-1557-30
 Analyst: 1030 icp 7300
 Initial Sample Wt: 2.04 g
 Dilution:
 Wash Time: 15

Autosampler Location: 136
 Date Collected: 3/29/2017 6:58:29 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Mean Data: 17-03-1557-30

Analyte	Mean Corrected		Calib.		Sample		Std.Dev.	RSD
	Intensity	Conc.	Units	Std.Dev.	Conc.	Units		
Tb 384	70482.6	88.23	%	0.824				0.93%
Tb 350	127648.3	101.2	%	0.65				0.64%
Ag 328.068*†	-439.6	-0.0025	mg/L	0.00056	-0.1217	mg/kg	0.02747	22.56%
Al 308.215*†	1552436.6	100.3	mg/L	0.37	4916	mg/kg	18.15	0.37%
As 188.979†	40.2	0.0231	mg/L	0.00107	1.131	mg/kg	0.0527	4.66%
As 193.696*†	23.7	0.0194	mg/L	0.00920	0.9502	mg/kg	0.45094	47.46%
B 249.677*†	394.2	0.0089	mg/L	0.00117	0.4368	mg/kg	0.05752	13.17%
Ba 233.527*†	159924.0	1.160	mg/L	0.0044	56.88	mg/kg	0.216	0.38%
Be 313.042*†	9311.1	0.0026	mg/L	0.00001	0.1264	mg/kg	0.00067	0.53%
Ca 317.933*†	116209.7	52.43	mg/L	0.333	2570	mg/kg	16.32	0.63%
Cd 226.502*†	257.3	0.0041	mg/L	0.00002	0.1987	mg/kg	0.00120	0.60%
Cd 228.802†	11.1	0.0003	mg/L	0.00041	0.0164	mg/kg	0.01988	120.89%
Co 228.616*†	2057.4	0.0862	mg/L	0.00106	4.224	mg/kg	0.0520	1.23%
Cr 267.716*†	8490.0	0.0841	mg/L	0.00053	4.122	mg/kg	0.0261	0.63%
Cu 324.752*†	30054.0	0.1248	mg/L	0.00006	6.118	mg/kg	0.0030	0.05%
Fe 273.955*†	3012378.6	146.1	mg/L	4.24	7162	mg/kg	207.69	2.90%
K 766.490*†	101256.7	30.08	mg/L	0.214	1474	mg/kg	10.48	0.71%
Mg 279.077*†	841817.7	54.58	mg/L	0.057	2676	mg/kg	2.80	0.10%
Mn 257.610*†	1521876.8	2.584	mg/L	0.0046	126.7	mg/kg	0.22	0.18%
Mo 202.031*†	8.1	0.0010	mg/L	0.00079	0.0476	mg/kg	0.03870	81.39%
Na 589.592*†	17775.5	3.429	mg/L	0.0223	168.1	mg/kg	1.09	0.65%
Ni 231.604*†	1505.5	0.0709	mg/L	0.00126	3.476	mg/kg	0.0619	1.78%
P 213.617*†	15893.1	9.753	mg/L	0.0324	478.1	mg/kg	1.59	0.33%
P 214.914†	10561.0	10.44	mg/L	0.062	511.8	mg/kg	3.02	0.59%
Pb 220.353*†	489.2	0.0763	mg/L	0.00034	3.741	mg/kg	0.0166	0.44%
Sb 206.836†	14.0	0.0082	mg/L	0.00743	0.4036	mg/kg	0.36398	90.19%
Sb 217.582*†	3.5	0.0021	mg/L	0.00313	0.1016	mg/kg	0.15361	151.22%
Se 196.026*†	2.1	0.0012	mg/L	0.00786	0.0577	mg/kg	0.38521	667.23%
Si 251.611*†	303503.4	8.640	mg/L	0.0583	423.5	mg/kg	2.86	0.67%
Sn 189.927*†	-300.8	-0.0594	mg/L	0.00260	-2.913	mg/kg	0.1274	4.37%
Sn 242.170†	602.5	0.3739	mg/L	0.01277	18.33	mg/kg	0.626	3.41%
Sr 407.771*†	234567.1	0.6319	mg/L	0.00556	30.97	mg/kg	0.273	0.88%
Ti 334.940†	7789122.6	9.581	mg/L	0.2312	469.7	mg/kg	11.33	2.41%
Ti 336.121*†	5554103.8	9.893	mg/L	0.2456	484.9	mg/kg	12.04	2.48%
Tl 190.801*†	-18.4	-0.0116	mg/L	0.00614	-0.5675	mg/kg	0.30099	53.04%
V 292.402*†	34434.2	0.2733	mg/L	0.00185	13.40	mg/kg	0.091	0.68%
Zn 206.200*†	15281.9	0.4726	mg/L	0.00082	23.17	mg/kg	0.040	0.17%

EPA 6010B ICP Metals (Solid)

Quality Control

Method Blank
LCS/LCSD
MS/MSD

**METHOD BLANK ASSOCIATION SUMMARY
FOR METHOD: EPA 6010B**

MB SAMPLE ID: 097-01-002-24537
MB BATCH ID: 170329L05
INSTRUMENT: ICP 7300
EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 17:07
REVIEWED BY:
D/T REVIEWED:
MATRIX: Soil

DATA FILE: W:\ICP-DATA\170329C 1\170329-b-05__376.icp

CLIENT WORK ORDER: 17-03-1557

<u>S#</u>	<u>RUN TYPE</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
28	B-DU1-ISM1-8		2017-03-29 18:56	W:\ICP-DATA\170329C 1\17-03-1557-28.icp
29	B-DU1-ISM2-8		2017-03-29 18:57	W:\ICP-DATA\170329C 1\17-03-1557-29.icp
30	B-DU1-ISM3-8		2017-03-29 18:58	W:\ICP-DATA\170329C 1\17-03-1557-30.icp

RAW DATA SHEET FOR METHOD: EPA 6010B

WORK ORDER: 097-01-002
INSTRUMENT: ICP 7300
EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 17:07
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\ICP-DATA\170329C 1\170329-b-05__376.icp

MB **CLIENT SAMPLE NUMBER:** Method Blank

LCS/MB BATCH: 170329L05 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 2.00 g / ACTUAL: 2.00 g
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 100.00 ml / ACTUAL: 100.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Antimony	0.00945	0.976	ND	0.732	
Arsenic	0.00243	0.976	ND	0.732	
Barium	0.000101	0.976	ND	0.488	
Beryllium	0.000126	0.976	ND	0.244	
Cadmium	0.000351	0.976	ND	0.488	
Chromium	0.000619	0.976	ND	0.244	
Cobalt	0.000590	0.976	ND	0.244	
Copper	0.000285	0.976	ND	0.488	
Lead	-0.000160	0.976	ND	0.488	
Molybdenum	0.000532	0.976	ND	0.244	
Nickel	0.000974	0.976	ND	0.244	
Selenium	0.00485	0.976	ND	0.732	
Silver	-0.000838	0.976	ND	0.244	
Thallium	-0.000864	0.976	ND	0.732	
Vanadium	0.000318	0.976	ND	0.244	
Zinc	-0.00205	0.976	ND	0.976	

Return to Contents

LCS QUALITY CONTROL SHEET FOR METHOD: EPA 6010B

LCS SAMPLE ID: 097-01-002-24537
LCS/MB BATCH ID: 170329L05
INSTRUMENT: ICP 7300

EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 17:08
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\ICP-DATA\170329C 1170329-I05_377.icp

<u>COMPOUND</u>	<u>CONC</u>	<u>CONC REC</u>	<u>%REC</u>	<u>%REC CL</u>	<u>ME CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Antimony	25.00	20.99	84	80-120	73-127	PASS	
Arsenic	25.00	22.47	90	80-120	73-127	PASS	
Barium	25.00	25.38	102	80-120	73-127	PASS	
Beryllium	25.00	22.33	89	80-120	73-127	PASS	
Cadmium	25.00	25.54	102	80-120	73-127	PASS	
Chromium	25.00	24.93	100	80-120	73-127	PASS	
Cobalt	25.00	26.08	104	80-120	73-127	PASS	
Copper	25.00	24.17	97	80-120	73-127	PASS	
Lead	25.00	25.61	102	80-120	73-127	PASS	
Molybdenum	25.00	23.03	92	80-120	73-127	PASS	
Nickel	25.00	24.87	99	80-120	73-127	PASS	
Phosphorus	25.00	22.97	92	80-120	73-127	PASS	
Selenium	25.00	23.31	93	80-120	73-127	PASS	
Silver	12.50	11.95	96	80-120	73-127	PASS	
Thallium	25.00	24.20	97	80-120	73-127	PASS	
Vanadium	25.00	23.74	95	80-120	73-127	PASS	
Zinc	25.00	25.45	102	80-120	73-127	PASS	
Aluminum	25.00	23.68	95	80-120	73-127	PASS	
Calcium	25.00	23.08	92	80-120	73-127	PASS	
Iron	25.00	24.89	100	80-120	73-127	PASS	
Magnesium	25.00	25.96	104	80-120	73-127	PASS	
Manganese	25.00	24.29	97	80-120	73-127	PASS	
Potassium	250.0	218.0	87	80-120	73-127	PASS	
Sodium	250.0	225.7	90	80-120	73-127	PASS	
Strontium	25.00	23.68	95	80-120	73-127	PASS	
Tin	25.00	23.49	94	80-120	73-127	PASS	
Titanium	25.00	23.67	95	80-120	73-127	PASS	
Boron	25.00	23.78	95	80-120	73-127	PASS	
Silicon	25.00	23.33	93	80-120	73-127	PASS	

LCS QUALITY CONTROL SHEET FOR METHOD: EPA 6010B

LCS SAMPLE ID: **097-01-002-24537**
LCS/MB BATCH ID: 170329L05
INSTRUMENT: ICP 7300

EXTRACTION: EPA 3050B
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED: 2017-03-29 17:08
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\ICP-DATA\170329C 1\170329-I-05__377.icp

<u>COMPOUND</u>	<u>CONC</u>	<u>CONC REC</u>	<u>%REC</u>	<u>%REC CL</u>	<u>ME CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
-----------------	-------------	-----------------	-------------	----------------	--------------	---------------	-------------------

Total number of LCS compounds: 29
 Total number of ME compounds: 0
 Total number of ME compounds allowed: 1
 LCS ME CL validation result: Pass



MATRIX SPIKE / MATRIX SPIKE DUPLICATE QUALITY CONTROL SHEET FOR METHOD: EPA 6010B

SPIKED SAMPLE ID: 17-03-1813-1
MS/MSD BATCH: 170329S05

INSTRUMENTS:

SAMPLE: ICP 7300
MS: ICP 7300
MSD: ICP 7300

EXTRACTION: EPA 3050B
D/T EXTRACTED:

SAMPLE: 2017-03-29 00:00
MS: 2017-03-29 00:00
MSD: 2017-03-29 00:00

ANALYZED BY: 935

D/T ANALYZED:

SAMPLE: 2017-03-29 17:12
MS: 2017-03-29 17:13
MSD: 2017-03-29 17:16

REVIEWED BY:

D/T REVIEWED:

COMMENT:

COMPOUND NAME	SAMPLE	INITIAL	FINAL	MS CONC	%MS.REC	MSD CONC	%MSD.REC	%REC CL	RPD	RPD CL	STATUS	QUALIFIERS
Antimony	ND	0.5000	25.00	11.32	45	11.05	44	50-115	2	0-20	FAIL	3G
Arsenic	2.141	0.5000	25.00	28.57	106	26.28	97	75-125	8	0-20	PASS	
Barium	65.88	0.5000	25.00	95.97	120	78.14	49	75-125	20	0-20	FAIL	3G
Beryllium	0.3817	0.5000	25.00	26.81	106	25.14	99	75-125	6	0-20	PASS	
Cadmium	ND	0.5000	25.00	28.11	112	26.45	106	75-125	6	0-20	PASS	
Chromium	12.69	0.5000	25.00	39.67	108	35.64	92	75-125	11	0-20	PASS	
Cobalt	7.098	0.5000	25.00	35.02	112	32.08	100	75-125	9	0-20	PASS	
Copper	10.42	0.5000	25.00	37.00	106	33.11	91	75-125	11	0-20	PASS	
Lead	3.987	0.5000	25.00	32.73	115	30.39	106	75-125	7	0-20	PASS	
Molybdenum	ND	0.5000	25.00	24.76	99	23.57	94	75-125	5	0-20	PASS	
Nickel	8.515	0.5000	25.00	34.90	106	31.89	93	75-125	9	0-20	PASS	
Phosphorus	343.9	0.5000	25.00	340.2	4x	303.6	4x	75-125	4x	0-20	PASS	Q
Selenium	ND	0.5000	25.00	27.32	109	25.65	103	75-125	6	0-20	PASS	
Silver	ND	0.2500	12.50	13.18	105	12.25	98	75-125	7	0-20	PASS	
Thallium	ND	0.5000	25.00	25.36	101	24.63	99	75-125	3	0-20	PASS	
Vanadium	24.51	0.5000	25.00	50.22	103	43.83	77	75-125	14	0-20	PASS	
Zinc	29.81	0.5000	25.00	57.58	111	49.16	77	75-125	16	0-20	PASS	
Aluminum	7204	0.5000	25.00	7385	4x	5915	4x	75-125	4x	0-20	PASS	Q
Calcium	2062	0.5000	25.00	1931	4x	1604	4x	75-125	4x	0-20	PASS	Q
Iron	11500	0.5000	25.00	11170	4x	8906	4x	75-125	4x	0-20	PASS	Q
Magnesium	3039	0.5000	25.00	2843	4x	2242	4x	75-125	4x	0-20	PASS	Q
Manganese	272.5	0.5000	25.00	287.3	4x	228.9	4x	75-125	4x	0-20	PASS	Q
Potassium	2239	5.000	250.0	2385	4x	1876	4x	75-125	4x	0-20	PASS	Q
Sodium	115.9	5.000	250.0	380.6	106	346.5	92	75-125	9	0-20	PASS	
Strontium	24.23	0.5000	25.00	49.64	102	43.61	78	75-125	13	0-20	PASS	
Tin	ND	0.5000	25.00	22.89	92	21.88	88	75-125	5	0-20	PASS	
Titanium	465.0	0.5000	25.00	531.3	4x	435.8	4x	75-125	4x	0-20	PASS	Q
Boron	ND	0.5000	25.00	25.76	103	25.29	101	75-125	2	0-20	PASS	

MATRIX SPIKE / MATRIX SPIKE DUPLICATE QUALITY CONTROL SHEET FOR METHOD: EPA 6010B

SPIKED SAMPLE ID: 17-03-1813-1
MS/MSD BATCH: 170329S05
INSTRUMENTS:
SAMPLE: ICP 7300
MS: ICP 7300
MSD: ICP 7300

EXTRACTION: EPA 3050B
D/T EXTRACTED:
SAMPLE: 2017-03-29 00:00
MS: 2017-03-29 00:00
MSD: 2017-03-29 00:00

ANALYZED BY: 935
D/T ANALYZED:
SAMPLE: 2017-03-29 17:12
MS: 2017-03-29 17:13
MSD: 2017-03-29 17:16
REVIEWED BY:
D/T REVIEWED:

COMMENT:

COMPOUND NAME	SAMPLE	INITIAL	FINAL	MS CONC	% MS.REC	MSD CONC	% MSD.REC	% REC CL	RPD	RPD CL	STATUS	QUALIFIERS
Silicon	290.2	0.5000	25.00	621.6	4x	634.2	4x	75-125	4x	0-20	PASS	Q

Data Files:

TYPE	DATA FILE	DATA FILE PATH
MS	17-03-1813-1 ms.icp	W:\ICP-DATA\170329C 1\
MSD	17-03-1813-1 msd.icp	W:\ICP-DATA\170329C 1\

Sequence No.: 3
Sample ID: 170329-b-05
Analyst: 1030 icp 7300
Initial Sample Wt: 2.05 g
Dilution:
Wash Time: 15

Autosampler Location: 103
Date Collected: 3/29/2017 5:07:11 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Mean Data: 170329-b-05

Analyte	Mean Corrected		Calib. Conc. Units	Std.Dev.	Sample		RSD
	Intensity				Conc. Units	Std.Dev.	
Tb 384	85764.3		107.4 %	3.74			3.48%
Tb 350	138557.8		109.9 %	4.15			3.78%
Ag 328.068*†	-148.3	-0.0008	mg/L	0.00141	-0.0409	mg/kg	0.06886 168.54%
Al 308.215*†	-205.2	-0.0133	mg/L	0.00337	-0.6467	mg/kg	0.16421 25.39%
As 188.979†	2.7	0.0016	mg/L	0.00035	0.0758	mg/kg	0.01727 22.76%
As 193.696*†	3.0	0.0024	mg/L	0.00407	0.1185	mg/kg	0.19856 167.63%
B 249.677*†	413.6	0.0094	mg/L	0.00011	0.4562	mg/kg	0.00518 1.14%
Ba 233.527*†	13.9	0.0001	mg/L	0.00008	0.0049	mg/kg	0.00382 77.51%
Be 313.042*†	456.3	0.0001	mg/L	0.00001	0.0062	mg/kg	0.00033 5.37%
Ca 317.933*†	6.0	0.0027	mg/L	0.00077	0.1310	mg/kg	0.03769 28.78%
Cd 226.502*†	22.3	0.0004	mg/L	0.00001	0.0171	mg/kg	0.00049 2.89%
Cd 228.802†	7.0	0.0002	mg/L	0.00030	0.0103	mg/kg	0.01444 139.79%
Co 228.616*†	14.1	0.0006	mg/L	0.00029	0.0288	mg/kg	0.01400 48.67%
Cr 267.716*†	62.5	0.0006	mg/L	0.00002	0.0302	mg/kg	0.00106 3.50%
Cu 324.752*†	68.6	0.0003	mg/L	0.00006	0.0139	mg/kg	0.00279 20.05%
Fe 273.955*†	71.3	0.0035	mg/L	0.00014	0.1686	mg/kg	0.00659 3.91%
K 766.490*†	-194.2	-0.0577	mg/L	0.07207	-2.814	mg/kg	3.5154 124.92%
Mg 279.077*†	148.1	0.0096	mg/L	0.01779	0.4685	mg/kg	0.86801 185.28%
Mn 257.610*†	824.0	0.0014	mg/L	0.00001	0.0682	mg/kg	0.00044 0.65%
Mo 202.031*†	4.5	0.0005	mg/L	0.00037	0.0259	mg/kg	0.01805 69.58%
Na 589.592*†	-115.1	-0.0222	mg/L	0.02283	-1.083	mg/kg	1.1138 102.84%
Ni 231.604*†	20.7	0.0010	mg/L	0.00121	0.0475	mg/kg	0.05887 123.89%
P 213.617*†	1.1	0.0007	mg/L	0.02091	0.0324	mg/kg	1.01980 >999.9%
P 214.914†	3.7	0.0037	mg/L	0.00245	0.1795	mg/kg	0.11967 66.67%
Pb 220.353*†	-1.0	-0.0002	mg/L	0.00129	-0.0078	mg/kg	0.06269 805.04%
Sb 206.836†	27.5	0.0162	mg/L	0.00325	0.7919	mg/kg	0.15848 20.01%
Sb 217.582*†	16.1	0.0095	mg/L	0.00009	0.4611	mg/kg	0.00434 0.94%
Se 196.026*†	8.6	0.0048	mg/L	0.00445	0.2365	mg/kg	0.21695 91.74%
Si 251.611*†	-375.5	-0.0107	mg/L	0.00060	-0.5215	mg/kg	0.02938 5.63%
Sn 189.927*†	-24.5	-0.0048	mg/L	0.00103	-0.2359	mg/kg	0.05011 21.24%
Sn 242.170†	-11.7	-0.0073	mg/L	0.00834	-0.3549	mg/kg	0.40663 114.57%
Sr 407.771*†	69.9	0.0002	mg/L	0.00004	0.0092	mg/kg	0.00206 22.45%
Ti 334.940†	358.2	0.0004	mg/L	0.00041	0.0215	mg/kg	0.02020 94.01%
Ti 336.121*†	474.1	0.0008	mg/L	0.00019	0.0412	mg/kg	0.00915 22.22%
Tl 190.801*†	-1.4	-0.0009	mg/L	0.00545	-0.0421	mg/kg	0.26582 630.77%
V 292.402*†	39.7	0.0003	mg/L	0.00060	0.0155	mg/kg	0.02932 189.10%
Zn 206.200*†	-66.3	-0.0021	mg/L	0.00063	-0.1001	mg/kg	0.03070 30.67%

Sequence No.: 4
 Sample ID: 170329-1-05
 Analyst: 1030 icp 7300
 Initial Sample Wt: 2.08 g
 Dilution:
 Wash Time: 15

Autosampler Location: 104
 Date Collected: 3/29/2017 5:08:12 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Mean Data: 170329-1-05

Analyte	Mean Corrected		Calib.		Sample		RSD
	Intensity	Conc.	Units	Std.Dev.	Conc.	Units	
Tb 384	86898.2	108.8	%	0.65			0.60%
Tb 350	142895.2	113.3	%	0.71			0.63%
Ag 328.068*†	44003.8	0.2486	mg/L	0.00008	11.95	mg/kg	0.004 0.03%
Al 308.215*†	7624.2	0.4925	mg/L	0.00806	23.68	mg/kg	0.387 1.64%
As 188.979†	841.8	0.4829	mg/L	0.00677	23.22	mg/kg	0.325 1.40%
As 193.696*†	572.4	0.4674	mg/L	0.00113	22.47	mg/kg	0.054 0.24%
B 249.677*†	21880.2	0.4947	mg/L	0.00221	23.78	mg/kg	0.106 0.45%
Ba 233.527*†	72768.0	0.5280	mg/L	0.00145	25.38	mg/kg	0.070 0.27%
Be 313.042*†	1676315.9	0.4644	mg/L	0.00999	22.33	mg/kg	0.481 2.15%
Ca 317.933*†	1064.0	0.4801	mg/L	0.01466	23.08	mg/kg	0.705 3.05%
Cd 226.502*†	33723.8	0.5313	mg/L	0.00069	25.54	mg/kg	0.033 0.13%
Cd 228.802†	17095.4	0.5153	mg/L	0.00366	24.77	mg/kg	0.176 0.71%
Co 228.616*†	12954.0	0.5425	mg/L	0.00087	26.08	mg/kg	0.042 0.16%
Cr 267.716*†	52354.3	0.5186	mg/L	0.00085	24.93	mg/kg	0.041 0.16%
Cu 324.752*†	121051.2	0.5027	mg/L	0.00236	24.17	mg/kg	0.113 0.47%
Fe 273.955*†	10674.8	0.5177	mg/L	0.00049	24.89	mg/kg	0.023 0.09%
K 766.490*†	15262.8	4.533	mg/L	0.0101	218.0	mg/kg	0.48 0.22%
Mg 279.077*†	8329.4	0.5401	mg/L	0.00147	25.96	mg/kg	0.070 0.27%
Mn 257.610*†	297594.5	0.5053	mg/L	0.00039	24.29	mg/kg	0.019 0.08%
Mo 202.031*†	4017.9	0.4790	mg/L	0.00285	23.03	mg/kg	0.137 0.60%
Na 589.592*†	24338.1	4.694	mg/L	0.0024	225.7	mg/kg	0.12 0.05%
Ni 231.604*†	10985.2	0.5174	mg/L	0.00782	24.87	mg/kg	0.376 1.51%
P 213.617*†	778.7	0.4778	mg/L	0.00461	22.97	mg/kg	0.222 0.97%
P 214.914†	471.8	0.4664	mg/L	0.00125	22.42	mg/kg	0.060 0.27%
Pb 220.353*†	3413.7	0.5326	mg/L	0.00057	25.61	mg/kg	0.027 0.11%
Sb 206.836†	758.6	0.4475	mg/L	0.00739	21.51	mg/kg	0.355 1.65%
Sb 217.582*†	745.1	0.4366	mg/L	0.00937	20.99	mg/kg	0.450 2.15%
Se 196.026*†	855.3	0.4848	mg/L	0.00019	23.31	mg/kg	0.009 0.04%
Si 251.611*†	17045.8	0.4853	mg/L	0.00031	23.33	mg/kg	0.015 0.06%
Sn 189.927*†	2473.6	0.4886	mg/L	0.00173	23.49	mg/kg	0.083 0.35%
Sn 242.170†	765.3	0.4749	mg/L	0.00080	22.83	mg/kg	0.038 0.17%
Sr 407.771*†	182846.7	0.4926	mg/L	0.00347	23.68	mg/kg	0.167 0.70%
Ti 334.940†	378838.1	0.4660	mg/L	0.00912	22.40	mg/kg	0.438 1.96%
Ti 336.121*†	276396.0	0.4923	mg/L	0.00008	23.67	mg/kg	0.004 0.02%
Tl 190.801*†	798.0	0.5033	mg/L	0.00213	24.20	mg/kg	0.102 0.42%
V 292.402*†	61688.5	0.4938	mg/L	0.00012	23.74	mg/kg	0.006 0.02%
Zn 206.200*†	17114.6	0.5293	mg/L	0.00309	25.45	mg/kg	0.149 0.58%

Sequence No.: 10
 Sample ID: 17-03-1813-1 ms
 Analyst: 1030 icp 7300
 Initial Sample Wt: 1.96 g
 Dilution:
 Wash Time: 15

Autosampler Location: 106
 Date Collected: 3/29/2017 5:13:18 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Mean Data: 17-03-1813-1 ms

Analyte	Mean Corrected		Calib.	Std.Dev.	Sample		RSD
	Intensity	Conc. Units			Conc. Units	Std.Dev.	
Tb 384	69206.7	86.64 %		1.576			1.82%
Tb 350	127784.2	101.3 %		1.06			1.05%
Ag 328.068*†	45731.6	0.2583 mg/L		0.00065	13.18 mg/kg	0.033	0.25%
Al 308.215*†	2240725.0	144.7 mg/L		0.36	7385 mg/kg	18.58	0.25%
As 188.979†	1016.2	0.5830 mg/L		0.00268	29.74 mg/kg	0.136	0.46%
As 193.696*†	685.8	0.5600 mg/L		0.00098	28.57 mg/kg	0.050	0.17%
B 249.677*†	22330.0	0.5048 mg/L		0.00464	25.76 mg/kg	0.237	0.92%
Ba 233.527*†	259257.6	1.881 mg/L		0.0004	95.97 mg/kg	0.021	0.02%
Be 313.042*†	1896769.4	0.5254 mg/L		0.00120	26.81 mg/kg	0.061	0.23%
Ca 317.933*†	83871.9	37.84 mg/L		0.123	1931 mg/kg	6.27	0.32%
Cd 226.502*†	34976.5	0.5510 mg/L		0.00304	28.11 mg/kg	0.155	0.55%
Cd 228.802†	18083.4	0.5451 mg/L		0.00280	27.81 mg/kg	0.143	0.51%
Co 228.616*†	16389.6	0.6864 mg/L		0.00241	35.02 mg/kg	0.123	0.35%
Cr 267.716*†	78491.4	0.7775 mg/L		0.00076	39.67 mg/kg	0.039	0.10%
Cu 324.752*†	174659.4	0.7253 mg/L		0.00041	37.01 mg/kg	0.021	0.06%
Fe 273.955*†	4515453.9	219.0 mg/L		3.66	11170 mg/kg	186.72	1.67%
K 766.490*†	157382.8	46.75 mg/L		0.223	2385 mg/kg	11.39	0.48%
Mg 279.077*†	859499.9	55.73 mg/L		0.120	2843 mg/kg	6.13	0.22%
Mn 257.610*†	3315944.7	5.630 mg/L		0.0040	287.3 mg/kg	0.20	0.07%
Mo 202.031*†	4071.6	0.4854 mg/L		0.00147	24.76 mg/kg	0.075	0.30%
Na 589.592*†	38672.9	7.459 mg/L		0.0602	380.6 mg/kg	3.07	0.81%
Ni 231.604*†	14522.6	0.6840 mg/L		0.00326	34.90 mg/kg	0.166	0.48%
P 213.617*†	10866.0	6.668 mg/L		0.0253	340.2 mg/kg	1.29	0.38%
P 214.914†	7356.9	7.273 mg/L		0.0531	371.1 mg/kg	2.71	0.73%
Pb 220.353*†	4111.8	0.6415 mg/L		0.00241	32.73 mg/kg	0.123	0.38%
Sb 206.836†	405.4	0.2391 mg/L		0.01082	12.20 mg/kg	0.552	4.53%
Sb 217.582*†	378.6	0.2219 mg/L		0.00360	11.32 mg/kg	0.184	1.62%
Se 196.026*†	945.0	0.5356 mg/L		0.00395	27.32 mg/kg	0.201	0.74%
Si 251.611*†	427954.9	12.18 mg/L		0.041	621.6 mg/kg	2.07	0.33%
Sn 189.927*†	2271.6	0.4487 mg/L		0.00864	22.89 mg/kg	0.441	1.93%
Sn 242.170†	1717.2	1.066 mg/L		0.0102	54.37 mg/kg	0.521	0.96%
Sr 407.771*†	361183.2	0.9730 mg/L		0.00639	49.64 mg/kg	0.326	0.66%
Ti 334.940†	8198521.5	10.08 mg/L		0.143	514.5 mg/kg	7.30	1.42%
Ti 336.121*†	5846264.4	10.41 mg/L		0.150	531.3 mg/kg	7.64	1.44%
Tl 190.801*†	788.1	0.4971 mg/L		0.00898	25.36 mg/kg	0.458	1.81%
V 292.402*†	123410.1	0.9844 mg/L		0.00432	50.22 mg/kg	0.220	0.44%
Zn 206.200*†	36489.6	1.129 mg/L		0.0029	57.58 mg/kg	0.147	0.25%

Sequence No.: 13
 Sample ID: 17-03-1813-1 msd
 Analyst: 1030 icp 7300
 Initial Sample Wt: 2.04 g
 Dilution:
 Wash Time: 15

Autosampler Location: 107
 Date Collected: 3/29/2017 5:16:02 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Mean Data: 17-03-1813-1 msd

Analyte	Mean Corrected		Calib.		Sample		RSD
	Intensity	Conc.	Units	Std.Dev.	Conc.	Units	
Tb 384	70700.4	88.51	%	1.621			1.83%
Tb 350	126817.2	100.5	%	1.34			1.33%
Ag 328.068*†	44228.6	0.2499	mg/L	0.00135	12.25	mg/kg	0.066 0.54%
Al 308.215*†	1868023.3	120.7	mg/L	1.55	5915	mg/kg	76.19 1.29%
As 188.979†	976.9	0.5605	mg/L	0.00458	27.47	mg/kg	0.224 0.82%
As 193.696*†	656.4	0.5360	mg/L	0.01203	26.28	mg/kg	0.590 2.24%
B 249.677*†	22816.7	0.5159	mg/L	0.00877	25.29	mg/kg	0.430 1.70%
Ba 233.527*†	219705.9	1.594	mg/L	0.0203	78.14	mg/kg	0.993 1.27%
Be 313.042*†	1851346.2	0.5129	mg/L	0.00616	25.14	mg/kg	0.302 1.20%
Ca 317.933*†	72521.1	32.72	mg/L	1.111	1604	mg/kg	54.48 3.40%
Cd 226.502*†	34249.1	0.5395	mg/L	0.00546	26.45	mg/kg	0.267 1.01%
Cd 228.802†	17779.5	0.5359	mg/L	0.00695	26.27	mg/kg	0.341 1.30%
Co 228.616*†	15627.4	0.6545	mg/L	0.00701	32.08	mg/kg	0.344 1.07%
Cr 267.716*†	73405.0	0.7271	mg/L	0.00667	35.64	mg/kg	0.327 0.92%
Cu 324.752*†	162668.7	0.6755	mg/L	0.00769	33.11	mg/kg	0.377 1.14%
Fe 273.955*†	3745721.1	181.7	mg/L	4.03	8906	mg/kg	197.42 2.22%
K 766.490*†	128876.3	38.28	mg/L	1.011	1876	mg/kg	49.58 2.64%
Mg 279.077*†	705514.8	45.74	mg/L	0.519	2242	mg/kg	25.44 1.13%
Mn 257.610*†	2750197.6	4.670	mg/L	0.0509	228.9	mg/kg	2.49 1.09%
Mo 202.031*†	4033.2	0.4808	mg/L	0.00286	23.57	mg/kg	0.140 0.60%
Na 589.592*†	36649.3	7.069	mg/L	0.1906	346.5	mg/kg	9.34 2.70%
Ni 231.604*†	13812.4	0.6506	mg/L	0.00174	31.89	mg/kg	0.085 0.27%
P 213.617*†	10091.8	6.193	mg/L	0.0187	303.6	mg/kg	0.92 0.30%
P 214.914†	6802.8	6.725	mg/L	0.0352	329.7	mg/kg	1.73 0.52%
Pb 220.353*†	3974.1	0.6200	mg/L	0.00301	30.39	mg/kg	0.147 0.49%
Sb 206.836†	412.5	0.2433	mg/L	0.00052	11.93	mg/kg	0.026 0.21%
Sb 217.582*†	384.8	0.2255	mg/L	0.00141	11.05	mg/kg	0.069 0.63%
Se 196.026*†	923.3	0.5233	mg/L	0.00718	25.65	mg/kg	0.352 1.37%
Si 251.611*†	454462.0	12.94	mg/L	0.206	634.2	mg/kg	10.10 1.59%
Sn 189.927*†	2260.3	0.4464	mg/L	0.00044	21.88	mg/kg	0.021 0.10%
Sn 242.170†	1574.4	0.9769	mg/L	0.00278	47.89	mg/kg	0.136 0.28%
Sr 407.771*†	330261.8	0.8897	mg/L	0.02465	43.61	mg/kg	1.208 2.77%
Ti 334.940†	7001799.4	8.613	mg/L	0.1470	422.2	mg/kg	7.21 1.71%
Ti 336.121*†	4991082.4	8.890	mg/L	0.1529	435.8	mg/kg	7.50 1.72%
Tl 190.801*†	796.6	0.5024	mg/L	0.00918	24.63	mg/kg	0.450 1.83%
V 292.402*†	112053.1	0.8941	mg/L	0.01133	43.83	mg/kg	0.555 1.27%
Zn 206.200*†	32425.5	1.003	mg/L	0.0075	49.16	mg/kg	0.367 0.75%

Sequence No.: 9
Sample ID: 17-03-1813-1
Analyst: 1030 icp 7300
Initial Sample Wt: 2.02 g
Dilution:
Wash Time: 15

Autosampler Location: 105
Date Collected: 3/29/2017 5:12:24 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Mean Data: 17-03-1813-1

Analyte	Mean Corrected		Calib.		Sample		Std.Dev.	RSD
	Intensity	Conc.	Units	Std.Dev.	Conc.	Units		
Tb 384	69133.0	86.54	%	0.100				0.12%
Tb 350	129419.1	102.6	%	0.42				0.41%
Ag 328.068*†	-445.9	-0.0025	mg/L	0.00020	-0.1247	mg/kg	0.00994	7.97%
Al 308.215*†	2252863.3	145.5	mg/L	0.51	7204	mg/kg	25.49	0.35%
As 188.979†	99.1	0.0569	mg/L	0.00576	2.815	mg/kg	0.2851	10.13%
As 193.696*†	53.0	0.0433	mg/L	0.00276	2.141	mg/kg	0.1365	6.37%
B 249.677*†	183.7	0.0042	mg/L	0.00218	0.2056	mg/kg	0.10787	52.46%
Ba 233.527*†	183408.8	1.331	mg/L	0.0061	65.88	mg/kg	0.300	0.46%
Be 313.042*†	27830.3	0.0077	mg/L	0.00005	0.3817	mg/kg	0.00256	0.67%
Ca 317.933*†	92329.7	41.66	mg/L	0.070	2062	mg/kg	3.47	0.17%
Cd 226.502*†	354.2	0.0056	mg/L	0.00014	0.2763	mg/kg	0.00685	2.48%
Cd 228.802†	-1.0	-0.0000	mg/L	0.00003	-0.0015	mg/kg	0.00162	106.92%
Co 228.616*†	3423.7	0.1434	mg/L	0.00046	7.098	mg/kg	0.0229	0.32%
Cr 267.716*†	25871.6	0.2563	mg/L	0.00136	12.69	mg/kg	0.067	0.53%
Cu 324.752*†	50674.9	0.2104	mg/L	0.00063	10.42	mg/kg	0.031	0.30%
Fe 273.955*†	4788115.1	232.2	mg/L	1.44	11500	mg/kg	71.24	0.62%
K 766.490*†	152241.2	45.22	mg/L	0.133	2239	mg/kg	6.56	0.29%
Mg 279.077*†	946792.7	61.39	mg/L	0.417	3039	mg/kg	20.64	0.68%
Mn 257.610*†	3242286.8	5.505	mg/L	0.0345	272.5	mg/kg	1.71	0.63%
Mo 202.031*†	-10.0	-0.0012	mg/L	0.00096	-0.0588	mg/kg	0.04764	81.04%
Na 589.592*†	12134.0	2.340	mg/L	0.0057	115.9	mg/kg	0.28	0.24%
Ni 231.604*†	3652.1	0.1720	mg/L	0.00107	8.515	mg/kg	0.0532	0.62%
P 213.617*†	11320.1	6.946	mg/L	0.0376	343.9	mg/kg	1.86	0.54%
P 214.914†	7791.9	7.703	mg/L	0.0526	381.4	mg/kg	2.61	0.68%
Pb 220.353*†	516.2	0.0805	mg/L	0.00051	3.987	mg/kg	0.0253	0.63%
Sb 206.836†	26.9	0.0159	mg/L	0.00662	0.7865	mg/kg	0.32768	41.66%
Sb 217.582*†	0.1	0.0000	mg/L	0.01088	0.0023	mg/kg	0.53872	>999.9%
Se 196.026*†	15.4	0.0087	mg/L	0.00202	0.4313	mg/kg	0.09988	23.16%
Si 251.611*†	205893.7	5.862	mg/L	0.3338	290.2	mg/kg	16.53	5.70%
Sn 189.927*†	-275.1	-0.0543	mg/L	0.00223	-2.690	mg/kg	0.1106	4.11%
Sn 242.170†	973.9	0.6043	mg/L	0.02232	29.92	mg/kg	1.105	3.69%
Sr 407.771*†	181720.7	0.4895	mg/L	0.00055	24.23	mg/kg	0.027	0.11%
Ti 334.940†	7396041.7	9.097	mg/L	0.0541	450.4	mg/kg	2.68	0.59%
Ti 336.121*†	5273784.8	9.393	mg/L	0.0558	465.0	mg/kg	2.76	0.59%
Tl 190.801*†	-34.7	-0.0219	mg/L	0.00799	-1.082	mg/kg	0.3958	36.58%
V 292.402*†	62333.4	0.4952	mg/L	0.00465	24.52	mg/kg	0.230	0.94%
Zn 206.200*†	19469.7	0.6021	mg/L	0.00403	29.81	mg/kg	0.200	0.67%

EPA 6010B ICP Metals (Solid)

Run Logs

170329C 1

M006-032-05 0.050 ml

10 ml

INT STD M120716A

PDS/PDSD

R.B. R12091602

M006-032-06 0.050 ml

Carrier/wash sol R12091604/R12091603

No.	File Name	Date	Time	Analyst Name	A/S	Location
1	Cal blankR12091601_935	3/29/2017	9:40:37 AM	935 icp 7300	1	
2	Cal blankR12091601_935	3/29/2017	9:42:24 AM	935 icp 7300	1	
3	STD3-M11116A_935_ICP7300	3/29/2017	9:43:26 AM	935 icp 7300	2	
4	ICV-M072816C	3/29/2017	9:44:25 AM	935 icp 7300	10	<i>out. Renew</i>
5	ICB-R12091601	3/29/2017	9:45:44 AM	935 icp 7300	1	
6	ICV-M072816C	3/29/2017	9:46:40 AM	935 icp 7300	10	
7	ICB-R12091601	3/29/2017	9:47:39 AM	935 icp 7300	1	
8	ICS_A - M110116B	3/29/2017	9:48:40 AM	935 icp 7300	8	
9	ICS_AB - M110116A	3/29/2017	9:49:46 AM	935 icp 7300	9	
10	LLCV-M010617A	3/29/2017	9:50:36 AM	935 icp 7300	101	<i>\$ out</i>
11	LLCV--M010617A	3/29/2017	9:51:52 AM	935 icp 7300	102	
12	CCV= STD3x0.5	3/29/2017	9:52:52 AM	935 icp 7300	3	
13	CCB-R12091601	3/29/2017	9:53:44 AM	935 icp 7300	1	
14	17-03-1526-4 pds	3/29/2017	9:57:10 AM	935 icp 7300	359	
15	LLCV-M010617A	3/29/2017	9:58:13 AM	935 icp 7300	101	<i>B, Sb, Se out</i>
16	ICS_A - M110116B	3/29/2017	9:59:15 AM	935 icp 7300	8	
17	ICS_AB - M110116A	3/29/2017	10:00:05 AM	935 icp 7300	9	
18	CCV= STD3x0.5	3/29/2017	10:00:55 AM	935 icp 7300	3	
19	CCB-R12091601	3/29/2017	10:01:47 AM	935 icp 7300	1	
20	170328-b-07	3/29/2017	10:07:01 AM	935 icp 7300	103	
21	170328-l-07	3/29/2017	10:08:08 AM	935 icp 7300	104	
22	170328-b-08	3/29/2017	10:09:09 AM	935 icp 7300	105	
23	170328-l-08	3/29/2017	10:10:08 AM	935 icp 7300	106	
24	170328-b-02	3/29/2017	10:10:59 AM	935 icp 7300	107	

Reviewed/Assign to Logbook Date: 3/30/17
 Analysis: 200-11610 Chemist ID: 1012
 Logbook Page: 69 Instrument ID: 1017

No.	File Name	Date	Time	Analyst Name	A/S	Location
345	CCV= STD3x0.5	3/29/2017	4:14:19 PM	935 icp 7300		3
346	CCB-R12091601	3/29/2017	4:15:11 PM	935 icp 7300		1
347	170329-b-02	3/29/2017	4:31:05 PM	935 icp 7300		125
348	170329-l-02	3/29/2017	4:32:12 PM	935 icp 7300		126
349	17-03-1625-1 pds	3/29/2017	4:33:03 PM	935 icp 7300		137
350	17-03-1625-1 pdsd	3/29/2017	4:33:47 PM	935 icp 7300		138
351	17-03-2113-1 ms	3/29/2017	4:34:30 PM	935 icp 7300		139
352	17-03-2113-1 msd	3/29/2017	4:35:22 PM	935 icp 7300		140
353	17-03-2113-1	3/29/2017	4:36:15 PM	935 icp 7300		141
354	17-03-1667-1	3/29/2017	4:37:08 PM	935 icp 7300		127
355	17-03-1667-1 ms	3/29/2017	4:38:01 PM	935 icp 7300		128
356	17-03-1667-1 msd	3/29/2017	4:38:43 PM	935 icp 7300		129
357	CCV= STD3x0.5	3/29/2017	4:39:24 PM	935 icp 7300		3
358	CCB-R12091601	3/29/2017	4:40:17 PM	935 icp 7300		1
359	17-03-1667-3	3/29/2017	4:41:18 PM	935 icp 7300		130
360	17-03-1667-5	3/29/2017	4:42:12 PM	935 icp 7300		131
361	17-03-1667-7	3/29/2017	4:43:04 PM	935 icp 7300		132
362	17-03-1667-9	3/29/2017	4:43:56 PM	935 icp 7300		133
363	blank	3/29/2017	4:44:47 PM	935 icp 7300		5
364	blank	3/29/2017	4:45:49 PM	935 icp 7300		6
365	LLCV-M010617A	3/29/2017	4:46:48 PM	935 icp 7300		101
366	LLCV-M010617A	3/29/2017	4:50:06 PM	935 icp 7300		101
367	CCV= STD3x0.5	3/29/2017	4:52:11 PM	935 icp 7300		3
368	CCB-R12091601	3/29/2017	4:53:10 PM	935 icp 7300		1
369	LLCV-M010617A	3/29/2017	4:54:02 PM	935 icp 7300		101
370	ICS_A - M110116B	3/29/2017	4:55:09 PM	935 icp 7300		8
371	ICS_AB - M110116A	3/29/2017	4:55:59 PM	935 icp 7300		9
372	CCV= STD3x0.5	3/29/2017	4:56:49 PM	935 icp 7300		3
373	CCB-R12091601	3/29/2017	4:57:41 PM	935 icp 7300		1
374	1703239-b-03	3/29/2017	5:05:12 PM	935 icp 7300		108
375	1703239-l-03	3/29/2017	5:06:20 PM	935 icp 7300		109
376	170329-b-05	3/29/2017	5:07:11 PM	935 icp 7300		103

*No solution
Pb, Se, Zn out*

Reviewed/Assign to Logbook Date: 3/30/17
 Analysis: 2071690 Chemist ID: 1012
 Logbook Page: 80 Instrument ID: 1017

No.	File Name	Date	Time	Analyst Name	A/S	Location
377	170329-I-05	3/29/2017	5:08:12 PM	935 icp 7300		104
378	17-03-2139-1	3/29/2017	5:09:03 PM	935 icp 7300		144
379	17-03-2139-2	3/29/2017	5:09:47 PM	935 icp 7300		145
380	17-03-2139-3	3/29/2017	5:10:40 PM	935 icp 7300		146
381	17-03-2139-4	3/29/2017	5:11:32 PM	935 icp 7300		147
382	17-03-1813-1	3/29/2017	5:12:24 PM	935 icp 7300		105
383	17-03-1813-1 ms	3/29/2017	5:13:18 PM	935 icp 7300		106
384	CCV= STD3x0.5	3/29/2017	5:14:09 PM	935 icp 7300		3
385	CCB-R12091601	3/29/2017	5:15:02 PM	935 icp 7300		1
386	17-03-1813-1 msd	3/29/2017	5:16:02 PM	935 icp 7300		107
387	17-03-1672-1	3/29/2017	5:16:55 PM	935 icp 7300		110
388	17-03-1672-1 ms	3/29/2017	5:17:54 PM	935 icp 7300		111
389	17-03-1672-1 msd	3/29/2017	5:18:53 PM	935 icp 7300		112
390	17-03-1679-1	3/29/2017	5:19:52 PM	935 icp 7300		113
391	17-03-1679-2	3/29/2017	5:20:48 PM	935 icp 7300		114
392	17-03-1679-3	3/29/2017	5:21:42 PM	935 icp 7300		115
393	17-03-1679-4	3/29/2017	5:22:39 PM	935 icp 7300		116
394	17-03-1680-1	3/29/2017	5:23:38 PM	935 icp 7300		117
395	17-03-1680-2	3/29/2017	5:24:29 PM	935 icp 7300		118
396	CCV= STD3x0.5	3/29/2017	5:25:20 PM	935 icp 7300		3
397	CCB-R12091601	3/29/2017	5:26:13 PM	935 icp 7300		1
398	CCB-R12091601	3/29/2017	5:26:59 PM	935 icp 7300		1
399	CCB-R12091601	3/29/2017	5:27:47 PM	935 icp 7300		1
400	17-03-1680-3	3/29/2017	5:28:50 PM	935 icp 7300		119
401	17-03-1680-4	3/29/2017	5:29:43 PM	935 icp 7300		120
402	17-03-1680-5	3/29/2017	5:30:34 PM	935 icp 7300		121
403	17-03-1680-6	3/29/2017	5:31:25 PM	935 icp 7300		122
404	17-03-1680-7	3/29/2017	5:32:16 PM	935 icp 7300		123
405	17-03-1681-1	3/29/2017	5:33:07 PM	935 icp 7300		124
406	CCV= STD3x0.5	3/29/2017	5:33:58 PM	935 icp 7300		3
407	CCB-R12091601	3/29/2017	5:34:50 PM	935 icp 7300		1
408	CCV= STD3x0.5	3/29/2017	6:39:14 PM	935 icp 7300		3

207 RL
}

Reviewed/Assign to Logbook Date: 3/30/17
 Analysis: 2007/6010 Chemist ID: 1012
 Logbook Page: 81 Instrument ID: 1017

No.	File Name	Date	Time	Analyst Name	A/S	Location
409	CCB-R12091601	3/29/2017	6:40:13 PM	935 icp 7300		1
410	17-03-1835-1 pds	3/29/2017	6:41:05 PM	935 icp 7300		142
411	17-03-1835-1 pdsd	3/29/2017	6:42:08 PM	935 icp 7300		143
412	17-03-1780x5-1	3/29/2017	6:43:04 PM	935 icp 7300		148
413	17-03-1780-1 pds	3/29/2017	6:43:53 PM	935 icp 7300		149
414	17-03-1965x5-1	3/29/2017	6:44:49 PM	935 icp 7300		150
415	17-03-1965-1 pds	3/29/2017	6:45:36 PM	935 icp 7300		151
416	17-03-1973x5-1	3/29/2017	6:46:31 PM	935 icp 7300		152
417	17-03-1973-1 pds	3/29/2017	6:47:19 PM	935 icp 7300		153
418	17-03-1959x5-f-2	3/29/2017	6:48:10 PM	935 icp 7300		154
419	17-03-1959-f-2 pds	3/29/2017	6:48:58 PM	935 icp 7300		155
420	CCV= STD3x0.5	3/29/2017	6:49:54 PM	935 icp 7300		3
421	CCB-R12091601	3/29/2017	6:50:47 PM	935 icp 7300		1
422	17-03-1959x5-9	3/29/2017	6:51:47 PM	935 icp 7300		156
423	17-03-1959-9 pds	3/29/2017	6:52:42 PM	935 icp 7300		157
424	170328-la-4	3/29/2017	6:53:39 PM	935 icp 7300		160
425	17-03-1667x5-1	3/29/2017	6:54:54 PM	935 icp 7300		158
426	17-03-1667-1 pds	3/29/2017	6:55:52 PM	935 icp 7300		159
427	17-03-1557-28	3/29/2017	6:56:44 PM	935 icp 7300		134
428	17-03-1557-29	3/29/2017	6:57:37 PM	935 icp 7300		135
429	17-03-1557-30	3/29/2017	6:58:29 PM	935 icp 7300		136
430	CCV= STD3x0.5	3/29/2017	7:02:53 PM	935 icp 7300		3
431	CCB-R12091601	3/29/2017	7:03:52 PM	935 icp 7300		1
432	LLCV-M010617A	3/29/2017	7:04:43 PM	935 icp 7300		101
433	LLCV-M010617A	3/29/2017	7:06:31 PM	935 icp 7300		101
434	ICS_A - M110116B	3/29/2017	7:07:40 PM	935 icp 7300		8
435	ICS_AB - M110116A	3/29/2017	7:08:30 PM	935 icp 7300		9
436	blank	3/29/2017	7:09:20 PM	935 icp 7300		5
437	blank	3/29/2017	7:10:21 PM	935 icp 7300		6
438	CCV= STD3x0.5	3/29/2017	7:11:20 PM	935 icp 7300		3
439	CCB-R12091601	3/29/2017	7:12:12 PM	935 icp 7300		1
440	170329-b-06	3/29/2017	7:38:03 PM	935 icp 7300		301

Ag, As, B, Sb, Se, Zn out

Reviewed/Assign to Logbook Date: 3/30/17
 Analysis: 20071600 Chemist ID: R12
 Logbook Page: 82 Instrument ID: 1017

EPA 6010B ICP Metals (Solid)

Preparation Logs

Metals Sample Preparation Logbook (Solid / Other)

METHOD	MATRIX	EQUIPMENT ID #	REAGENT ID #	STANDARD ID #												
<input checked="" type="checkbox"/> EPA 3050B <input type="checkbox"/> EPA 200.7 <input type="checkbox"/> EPA 200.8	<input checked="" type="checkbox"/> Solid <input type="checkbox"/> Other (Specify)	Thermometer 1711427 (CF 0.0 °C) Block Digester J Pipetter / Dispenser P-0730-276-028	HNO ₃ R03271702 20 mL HCl R03271701 15 mL H ₂ O ₂ M00604073 3 mL	Spike 1 M02217A Spike 2 M020617C Spike 3												
BATCH NUMBER		SUPPLY LOT #	BALANCE ID #	QUALITY SYSTEM MATRIX ID #	SAMPLE HANDLING											
MS/MSD 170329-505 (Specify)		Tube / Container 170104 Filter	65	Teflon Chip M006-035-115 (Specify)	1 = Composite 2 = Subsample 3 = Homogenize 4 = None											
DIGESTION																
DATE	START			END			SAMPLE HANDLING	ECL ID #	ANALYTE(S)	SAMPLE		SPIKE STANDARD				
	TIME	TEMP W/O CF (°C)	PREP TECH ID #	TIME	TEMP W/O CF (°C)	PREP TECH ID #				INITIAL (g)	FINAL (mL)	1 (µL)	2 (µL)	3 (µL)		
03/29/17	11:40	95	70	13:10	95	70	4	MS 17-03-1813-1 ^A MSD ↓ LCS 170329-105 LCSD / MB 805 17-03-1813-1 ^A 17-03-157-28 ^A 28 ^B 28 ^C 28 ^D 28 ^E 29 ^A 29 ^B 29 ^C 29 ^D 29 ^E 30 ^A 30 ^B 30 ^C 30 ^D 30 ^E	T ₁₂	1.96 2.04 2.08 2.05 2.02 2.02 2.03 1.98 2.01 2.09 2.07 1.95 2.06 2.06 2.05 2.04 1.98 2.02 2.08 2.07	1.00	500	500			

COMMENTS:

EPA 7471A Mercury (Solid)

RAW DATA

EPA 7471A Mercury (Solid)

Initial Calibration

ICV/ICB

CCV/CCB

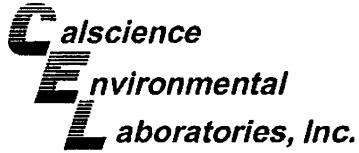
Sample Data

Quality Control

Method Blank

LCS/LCSD

MS/MSD



**EPA Method 7471A
Initial Calibration Verification**



Work Order No.: 17-03-1557

Instrument ID: HG 8 (H)

Concentration Unit: µg/L

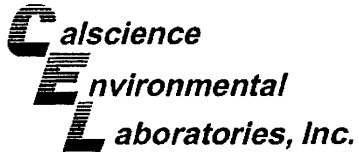
Test Method: EPA 7471A

Analyte	Initial Calibration Verification			
	True	ICV-1		Control Limit
		Observed	%D	
Mercury	5.000000	4.964654	1	0 - 10

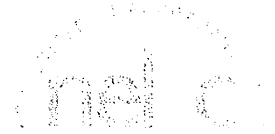
03/29/2017 17:09

ICV-1 File: ICV M030617B 03/29/2017 11:45:01 AM

Return to Contents



EPA Method 7471A
Continuing Calibration Verification



Work Order No.: 17-03-1557

Instrument ID: HG 8 (H)

Concentration Unit: µg/L

Test Method: EPA 7471A

Analyte	Continuing Calibration Verification									
	True	CCV-1		CCV-2		CCV-3		CCV-4		Control Limit
		Observed	%D	Observed	%D	Observed	%D	Observed	%D	
Mercury	2.000000	1.998235	0	1.742626	13	1.713826	14	1.786031	11	0 - 20

03/29/2017 17:10

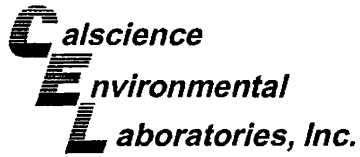
CCV-1 File: CCV 0.2x10ppb 03/29/2017 02:19:44 PM

CCV-2 File: CCV 0.2x10ppb 03/29/2017 03:15:49 PM

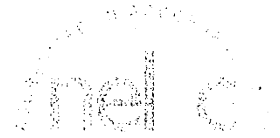
CCV-3 File: CCV 0.2x10ppb 03/29/2017 03:50:04 PM

CCV-4 File: CCV 0.2x10ppb 03/29/2017 04:17:32 PM

Return to Contents



EPA Method 7471A
Initial and Continuing Calibration Blanks

Work Order No.: 17-03-1557Instrument ID: HG 8 (H)Concentration Unit: µg/LTest Method: EPA 7471A

Analyte	Initial and Continuing Calibration Blanks					
	ICB-1	CCB-1	CCB-2	CCB-3	CCB-4	RL (No PF)
Mercury	-0.027629	-0.018645	-0.031176	-0.017623	-0.030838	0.500000

03/29/2017 17:11

ICB-1 File: ICB 03/29/2017 11:47:15 AM

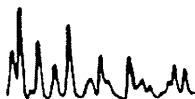
CCB-1 File: CCB 03/29/2017 02:48:04 PM

CCB-2 File: CCB 03/29/2017 03:18:06 PM

CCB-3 File: CCB 03/29/2017 03:52:21 PM

CCB-4 File: CCB 03/29/2017 04:19:49 PM

Note: Preparation factor (PF) = 167 L/kg



**RAW DATA SHEET
FOR METHOD: EPA 7471A**

WORK ORDER: 17-03-1557
INSTRUMENT: Mercury 08
EXTRACTION : EPA 7471A Total
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED: 2017-03-29 15:54
REVIEWED BY: 309
D/T REVIEWED: 2017-03-29 17:04

DATA FILE: W:\MERCURY_DATA\FINAL\170329H1\17-03-1557-25.icp

25 **CLIENT SAMPLE NUMBER:** B-DU1-ISM1-8

<u>LCS/MB BATCH:</u> 170329L01	<u>SAMPLE VOLUME / WEIGHT:</u> DEFAULT: 0.60 g / ACTUAL: 0.63 g
<u>MS/MSD BATCH:</u> 170329S01	<u>FINAL VOLUME / WEIGHT:</u> DEFAULT: 100.00 ml
<u>UNITS:</u> mg/kg	<u>ADJUSTMENT RATIO TO PF:</u> 0.95

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Mercury	0.000120	1.00	ND	0.0794	

**RAW DATA SHEET
FOR METHOD: EPA 7471A**

WORK ORDER: 17-03-1557
INSTRUMENT: Mercury 08
EXTRACTION: EPA 7471A Total
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED: 2017-03-29 15:56
REVIEWED BY: 309
D/T REVIEWED: 2017-03-29 17:04

DATA FILE: W:\MERCURY_DATA\FINAL\170329H1\17-03-1557-26.icp

26 **CLIENT SAMPLE NUMBER:** B-DU1-ISM2-8

<u>LCS/MB BATCH:</u> 170329L01	<u>SAMPLE VOLUME / WEIGHT:</u> DEFAULT: 0.60 g / ACTUAL: 0.62 g
<u>MS/MSD BATCH:</u> 170329S01	<u>FINAL VOLUME / WEIGHT:</u> DEFAULT: 100.00 ml
<u>UNITS:</u> mg/kg	<u>ADJUSTMENT RATIO TO PF:</u> 0.97

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Mercury	0.000359	1.00	ND	0.0806	

**RAW DATA SHEET
FOR METHOD: EPA 7471A**

WORK ORDER: 17-03-1557
INSTRUMENT: Mercury 08
EXTRACTION : EPA 7471A Total
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED: 2017-03-29 15:59
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\MERCURY_DATA\FINAL\170329H1\170329H1_1557\17-03-1557-27.icp

27 **CLIENT SAMPLE NUMBER:** B-DU1-ISM3-8

LCS/MB BATCH: 170329L01 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 0.60 g / ACTUAL: 0.60 g
MS/MSD BATCH: 170329S01 **FINAL VOLUME / WEIGHT:** DEFAULT: 100.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Mercury	0.0000575	1.00	ND	0.0833	

METHOD BLANK ASSOCIATION SUMMARY FOR METHOD: EPA 7471A

MB SAMPLE ID: 099-16-272-2910
MB BATCH ID: 170329L01
INSTRUMENT: Mercury 08
EXTRACTION: EPA 7471A Total
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED: 2017-03-29 15:08
REVIEWED BY: 309
D/T REVIEWED: 2017-03-29 17:00
MATRIX: Soil

DATA FILE: W:\MERCURY_DATA\FINAL\170329H1\170329-B-01.icp

CLIENT WORK ORDER: 17-03-1557

<u>S#</u>	<u>RUN TYPE</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
25	B-DU1-ISM1-8		2017-03-29 15:54	W:\MERCURY_DATA\FINAL\170329H1\17-03-1557-25.icp
26	B-DU1-ISM2-8		2017-03-29 15:56	W:\MERCURY_DATA\FINAL\170329H1\17-03-1557-26.icp
27	B-DU1-ISM3-8		2017-03-29 15:59	W:\MERCURY_DATA\FINAL\170329H1\170329H1_1557\17-03-1557-27.icp

RAW DATA SHEET FOR METHOD: EPA 7471A

WORK ORDER: 099-16-272
INSTRUMENT: Mercury 08
EXTRACTION : EPA 7471A Total
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED: 2017-03-29 15:08
REVIEWED BY: 309
D/T REVIEWED: 2017-03-29 17:00

DATA FILE: W:\MERCURY_DATA\FINAL\170329H1\170329-B-01.icp

MB **CLIENT SAMPLE NUMBER:** Method Blank

LCS/MB BATCH: 170329L01 SAMPLE VOLUME / WEIGHT: DEFAULT: 0.60 g / ACTUAL: 0.60 g
MS/MSD BATCH: FINAL VOLUME / WEIGHT: DEFAULT: 100.00 ml
UNITS: mg/kg ADJUSTMENT RATIO TO PF: 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Mercury	-0.00000955	1.00	ND	0.0833	



LCS QUALITY CONTROL SHEET FOR METHOD: EPA 7471A

LCS SAMPLE ID: 099-16-272- 2910
LCS/MB BATCH ID: 170329L01
INSTRUMENT: Mercury 08

EXTRACTION: EPA 7471A Total
D/T EXTRACTED: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED: 2017-03-29 15:11
REVIEWED BY: 309
D/T REVIEWED: 2017-03-29 17:00

DATA FILE: W:\MERCURY_DATA\FINAL\170329H1\170329-L-01.icp

<u>COMPOUND NAME</u>	<u>CONC ADDED</u>	<u>CONC REC</u>	<u>%RECOVERY</u>	<u>%REC CONTROL LIMIT</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Mercury	0.8350	0.7369	88	85-121	PASS	

**MATRIX SPIKE / MATRIX SPIKE DUPLICATE QUALITY CONTROL SHEET
FOR METHOD: EPA 7471A**

SPIKED SAMPLE ID: 17-03-1569-1
MS/MSD BATCH: 170329S01
INSTRUMENTS:
SAMPLE: Mercury 08
MS: Mercury 08
MSD: Mercury 08

EXTRACTION: EPA 7471A Total
D/T EXTRACTED:
SAMPLE: 2017-03-29 00:00
MS: 2017-03-29 00:00
MSD: 2017-03-29 00:00

ANALYZED BY: 868
D/T ANALYZED:
SAMPLE: 2017-03-29 15:13
MS: 2017-03-29 15:20
MSD: 2017-03-29 15:22
REVIEWED BY: 309
D/T REVIEWED: 2017-03-29 16:59

COMMENT:

COMPOUND NAME	SAMPLE	INITIAL	FINAL	MS CONC	% MS.REC	MSD CONC	% MSD.REC	% REC CL	RPD	RPD CL	STATUS	QUALIFIERS
Mercury	0.06019	0.005000	0.8350	0.7189	79	0.8023	89	76-136	11	0-16	PASS	

Data Files:

TYPE	DATA FILE	DATA FILE PATH
MS	17-03-1569-1 MS.icp	W:\MERCURY_DATA\FINAL\170329H1\
MSD	17-03-1569-1 MSD.icp	W:\MERCURY_DATA\FINAL\170329H1\



=====
Analysis Begun

Logged In Analyst: US26_SVC_INSTRUMENT
Spectrometer: FIMS-400, S/N B050-9560

Technique: AA FIMS-MHS
Autosampler: S10

Sample Information File: C:\Users\Public\PerkinElmer Syngistix\AA\Data\Sample Information\
170329H1.sifx

Batch ID:
Results Data Set: 170329H1
Results Library: U:\MERCURY_8\Data\Results\results.mdb

=====
Sequence No.: 1
Sample ID: Calib blank_868
Analyst: 868
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 1
Date Collected: 3/29/2017 11:23:11 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1

Replicate Data: Calib blank_868
Analyte: Hg 253.7
Repl SampleConc StndConc BlnkCorr Peak Peak Time Peak
mg/L ug/L Signal Area Height Stored
1 [0.00] 0.0000 -0.0002 0.0000 11:24:15 AM Yes
2 [0.00] 0.0000 0.0000 0.0000 11:25:00 AM Yes
Mean: [0.00] 0.0000
SD: 0.0000 0.0000
%RSD: 0.00% 37.15
Auto-zero performed.

=====
Sequence No.: 2
Sample ID: 0.025ppb 0.005x5ppb
Analyst: 868
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 2
Date Collected: 3/29/2017 11:25:25 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1

Replicate Data: 0.025ppb 0.005x5ppb
Analyte: Hg 253.7
Repl SampleConc StndConc BlnkCorr Peak Peak Time Peak
mg/L ug/L Signal Area Height Stored
1 [0.025] 0.0002 0.0000 0.0000 11:26:30 AM Yes
2 [0.025] 0.0002 -0.0008 0.0002 11:27:14 AM Yes
Mean: [0.025] 0.0002
SD: 0.00000 0.0000
%RSD: 0.00% 11.59
Standard number 1 applied. [0.025]
Correlation Coef.: 1.000000 Slope: 0.00769 Intercept: 0.00000

=====
Sequence No.: 3
Sample ID: 0.10ppb M030617AX0.0001
Analyst: 868
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 3
Date Collected: 3/29/2017 11:27:40 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1

Replicate Data: 0.10ppb M030617AX0.0001
Analyte: Hg 253.7
Repl SampleConc StndConc BlnkCorr Peak Peak Time Peak
mg/L ug/L Signal Area Height Stored
1 [0.100] 0.0010 0.0027 0.0011 11:28:44 AM Yes
2 [0.100] 0.0010 0.0022 0.0010 11:29:30 AM Yes
Mean: [0.100] 0.0010
SD: 0.00000 0.0001
%RSD: 0.00% 5.16
Standard number 2 applied. [0.100]
Correlation Coef.: 0.998204 Slope: 0.01015 Intercept: -0.00003



```

=====
Sequence No.: 4                               Autosampler Location: 4
Sample ID: 1.00ppb M030617AX0.001           Date Collected: 3/29/2017 11:29:56 AM
Analyst: 868                                  Data Type: Original
Initial Sample Wt:                             Initial Sample Vol:
Dilution:                                     Sample Prep Vol:
Wash Time (before sample): 0                 Auto Dilution Factor: 1
=====

```

```

-----
Replicate Data: 1.00ppb M030617AX0.001       Analyte: Hg 253.7
Repl  SampleConc  StndConc  BlnkCorr  Peak    Peak    Time    Peak
#      mg/L        ug/L      Signal   Area   Height  Time    Stored
1      [1.000]      0.0103   0.0301   0.0103 11:31:01 AM Yes
2      [1.000]      0.0095   0.0253   0.0095 11:31:45 AM Yes
Mean:  [1.000]      0.0099
SD:     0.00000     0.0005
%RSD:   0.00%        5.55
Standard number 3 applied. [1.000]
Correlation Coef.: 0.999982  Slope: 0.00988  Intercept: -0.00002
=====

```

```

=====
Sequence No.: 5                               Autosampler Location: 5
Sample ID: 2.00ppb M030617AX0.002           Date Collected: 3/29/2017 11:32:12 AM
Analyst: 868                                  Data Type: Original
Initial Sample Wt:                             Initial Sample Vol:
Dilution:                                     Sample Prep Vol:
Wash Time (before sample): 0                 Auto Dilution Factor: 1
=====

```

```

-----
Replicate Data: 2.00ppb M030617AX0.002       Analyte: Hg 253.7
Repl  SampleConc  StndConc  BlnkCorr  Peak    Peak    Time    Peak
#      mg/L        ug/L      Signal   Area   Height  Time    Stored
1      [2.000]      0.0198   0.0618   0.0199 11:33:17 AM Yes
2      [2.000]      0.0181   0.0484   0.0182 11:34:02 AM Yes
Mean:  [2.000]      0.0190
SD:     0.00000     0.0012
%RSD:   0.00%        6.37
Standard number 4 applied. [2.000]
Correlation Coef.: 0.999807  Slope: 0.00954  Intercept: 0.00005
=====

```

```

=====
Sequence No.: 6                               Autosampler Location: 6
Sample ID: 5.00ppb M030617AX0.005           Date Collected: 3/29/2017 11:34:30 AM
Analyst: 868                                  Data Type: Original
Initial Sample Wt:                             Initial Sample Vol:
Dilution:                                     Sample Prep Vol:
Wash Time (before sample): 0                 Auto Dilution Factor: 1
=====

```

```

-----
Replicate Data: 5.00ppb M030617AX0.005       Analyte: Hg 253.7
Repl  SampleConc  StndConc  BlnkCorr  Peak    Peak    Time    Peak
#      mg/L        ug/L      Signal   Area   Height  Time    Stored
1      [5.000]      0.0494   0.1503   0.0494 11:35:33 AM Yes
2      [5.000]      0.0447   0.1195   0.0448 11:36:18 AM Yes
Mean:  [5.000]      0.0471
SD:     0.00000     0.0033
%RSD:   0.00%        7.01
Standard number 5 applied. [5.000]
Correlation Coef.: 0.999949  Slope: 0.00941  Intercept: 0.00011
=====

```

```

=====
Sequence No.: 7                               Autosampler Location: 7
Sample ID: 10.0ppb M030617AX0.01            Date Collected: 3/29/2017 11:36:44 AM
Analyst: 868                                  Data Type: Original
Initial Sample Wt:                             Initial Sample Vol:
Dilution:                                     Sample Prep Vol:
Wash Time (before sample): 0                 Auto Dilution Factor: 1
=====

```

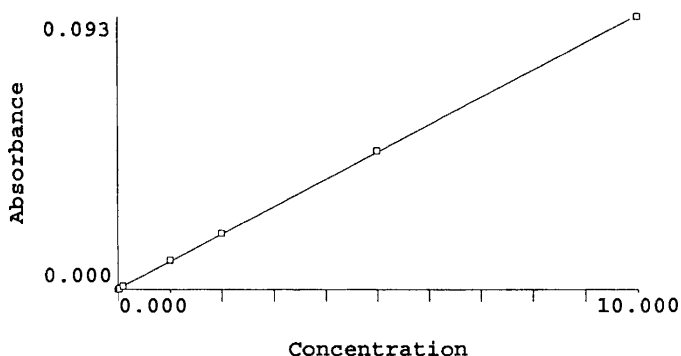
```

-----
Replicate Data: 10.0ppb M030617AX0.01        Analyte: Hg 253.7
Repl  SampleConc  StndConc  BlnkCorr  Peak    Peak    Time    Peak
#      mg/L        ug/L      Signal   Area   Height  Time    Stored
=====

```



1 [10.00] 0.0972 0.3015 0.0972 11:37:48 AM Yes
 2 [10.00] 0.0882 0.2366 0.0882 11:38:33 AM Yes
 Mean: [10.00] 0.0927
 SD: 0.0000 0.0064
 %RSD: 0.00% 6.87
 Standard number 6 applied. [10.00]
 Correlation Coef.: 0.999956 Slope: 0.00927 Intercept: 0.00025



Calibration data for Hg 253.7

Equation: Linear, Calculated Intercept

ID	Mean Signal (Abs)	Entered Conc. ug/L	Calculated Conc. ug/L	Standard Deviation	%RSD
Calib blank_868	0.0000	0	-0.026433	0.00	37.15
0.025ppb 0.005x5ppb	0.0002	0.025	-0.005694	0.00	11.59
0.10ppb M030617AX0.0001	0.0010	0.100	0.081189	0.00	5.16
1.00ppb M030617AX0.001	0.0099	1.000	1.037536	0.00	5.55
2.00ppb M030617AX0.002	0.0190	2.000	2.021603	0.00	6.37
5.00ppb M030617AX0.005	0.0471	5.000	5.049216	0.00	7.01
10.0ppb M030617AX0.01	0.0927	10.00	9.967583	0.01	6.87
Correlation Coef.: 0.999956		Slope: 0.00927		Intercept: 0.00025	

=====
Analysis Begun

Logged In Analyst: US26_SVC_INSTRUMENT
Spectrometer: FIMS-400, S/N B050-9560

Technique: AA FIMS-MHS
Autosampler: S10

Sample Information File: C:\Users\Public\PerkinElmer Syngistix\AA\Data\Sample Information\
170329H1.sifx

Batch ID:
Results Data Set: 170329H1
Results Library: U:\MERCURY_8\Data\Results\results.mdb

=====
Sequence No.: 1
Sample ID: ICV M030617B
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 8
Date Collected: 3/29/2017 11:43:13 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: ICV M030617B				Analyte: Hg 253.7			
Repl #	SampleConc mg/L	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00510	5.10	0.0475	0.1395	0.0475	11:44:16 AM	Yes
2	0.00483	4.83	0.0451	0.1197	0.0451	11:45:01 AM	Yes
Mean:	0.00496	4.96	0.0463				
SD:	0.000185	0.185	0.0017				
%RSD:	3.72%	3.72%	3.70				

QC value within limits for Hg 253.7 Recovery = 99.29%
All analyte(s) passed QC.

=====
Sequence No.: 2
Sample ID: ICB
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 1
Date Collected: 3/29/2017 11:45:27 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: ICB				Analyte: Hg 253.7			
Repl #	SampleConc mg/L	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	-0.000027	-0.0270	-0.0000	-0.0005	0.0000	11:46:30 AM	Yes
2	-0.000028	-0.0283	-0.0000	-0.0003	0.0000	11:47:15 AM	Yes
Mean:	-0.000028	-0.0276	-0.0000				
SD:	0.0000009	0.00091	0.0000				
%RSD:	3.28%	3.28%	75.78				

QC value within limits for Hg 253.7 Recovery = Not calculated
All analyte(s) passed QC.

=====
Sequence No.: 3
Sample ID: CRQL 0.25
Analyst: 868 HG-8
Initial Sample Wt: 0.6 g
Dilution:
Wash Time (before sample): 0
Autosampler Location: 9
Date Collected: 3/29/2017 11:47:41 AM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: CRQL 0.25				Analyte: Hg 253.7			
Repl #	SampleConc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0429	0.258	0.0026	0.0071	0.0027	11:48:45 AM	Yes
2	0.0392	0.235	0.0024	0.0053	0.0025	11:49:31 AM	Yes
Mean:	0.0411	0.246	0.0025				
SD:	0.00265	0.0159	0.0001				
%RSD:	6.45%	6.45%	5.82				



=====
Analysis Begun

Logged In Analyst: US26_SVC_INSTRUMENT
Spectrometer: FIMS-400, S/N B050-9560

Technique: AA FIMS-MHS
Autosampler: S10

Sample Information File: C:\Users\Public\PerkinElmer Syngistix\AA\Data\Sample Information\
170329H1.sifx

Batch ID:
Results Data Set: 170329H1
Results Library: U:\MERCURY_8\Data\Results\results.mdb

=====
Sequence No.: 1
Sample ID: CCV 0.2x10ppb
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 5
Date Collected: 3/29/2017 2:15:37 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: CCV 0.2x10ppb Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00140	1.40	0.0132	-0.0159	0.0132	2:16:42 PM	Yes
2	0.00150	1.50	0.0142	0.0062	0.0142	2:17:27 PM	Yes
Mean:	0.00145	1.45	0.0137				
SD:	0.000073	0.073	0.0007				
%RSD:	5.01%	5.01%	4.92				

QC value less than the lower limit for Hg 253.7 Recovery = 72.50%
QC Failed. Retry.

=====
Sequence No.: 2
Sample ID: CCV 0.2x10ppb
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 5
Date Collected: 3/29/2017 2:17:54 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: CCV 0.2x10ppb Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00207	2.07	0.0194	0.0634	0.0194	2:18:59 PM	Yes
2	0.00193	1.93	0.0182	0.0556	0.0182	2:19:44 PM	Yes
Mean:	0.00200	2.00	0.0188				
SD:	0.000095	0.095	0.0009				
%RSD:	4.73%	4.73%	4.67				

QC value within limits for Hg 253.7 Recovery = 99.91%
All analyte(s) passed QC.

=====
Sequence No.: 3
Sample ID: CCB
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0
Autosampler Location: 1
Date Collected: 3/29/2017 2:20:11 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: CCB Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.000256	0.256	0.0026	-0.0040	0.0027	2:21:15 PM	Yes
2	0.00250	2.50	0.0235	0.2001	0.0235	2:22:01 PM	Yes
Mean:	0.00138	1.38	0.0130				
SD:	0.001590	1.590	0.0147				
%RSD:	115.23%	115.23%	113.07				

QC value greater than the upper limit for Hg 253.7 Recovery = Not calculated
QC Failed. Stop the analysis.

Replicate Data: 17-03-1667-9

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00574	0.0350	0.0006	0.0021	0.0006	3:03:39 PM	Yes
2	0.00543	0.0331	0.0006	0.0020	0.0006	3:04:25 PM	Yes
Mean:	0.00559	0.0341	0.0006				
SD:	0.000216	0.00132	0.0000				
%RSD:	3.87%	3.87%	2.18				

=====

Sequence No.: 8	Autosampler Location: 9
Sample ID: CRQL 0.25	Date Collected: 3/29/2017 3:04:52 PM
Analyst: 868 HG-8	Data Type: Original
Initial Sample Wt: 0.6 g	Initial Sample Vol:
Dilution:	Sample Prep Vol: 100 mL
Wash Time (before sample): 0	Auto Dilution Factor: 1

Replicate Data: CRQL 0.25

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0363	0.218	0.0023	0.0092	0.0023	3:05:57 PM	Yes
2	0.00115	0.00691	0.0003	0.0027	0.0003	3:06:42 PM	Yes
Mean:	0.0187	0.112	0.0013				
SD:	0.02487	0.1492	0.0014				
%RSD:	132.73%	132.73%	107.46				

=====

Sequence No.: 9	Autosampler Location: 40
Sample ID: 170329-B-01	Date Collected: 3/29/2017 3:07:09 PM
Analyst: 868 HG-8	Data Type: Original
Initial Sample Wt: 0.6 g	Initial Sample Vol:
Dilution:	Sample Prep Vol: 100 mL
Wash Time (before sample): 0	Auto Dilution Factor: 1

Replicate Data: 170329-B-01

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	-0.00239	-0.0144	0.0001	0.0019	0.0001	3:08:13 PM	Yes
2	-0.000788	-0.00473	0.0002	0.0027	0.0002	3:08:58 PM	Yes
Mean:	-0.00159	-0.00955	0.0002				
SD:	0.001136	0.006815	0.0001				
%RSD:	71.39%	71.39%	40.36				

=====

Sequence No.: 10	Autosampler Location: 41
Sample ID: 170329-L-01	Date Collected: 3/29/2017 3:09:25 PM
Analyst: 868 HG-8	Data Type: Original
Initial Sample Wt: 0.6 g	Initial Sample Vol:
Dilution:	Sample Prep Vol: 100 mL
Wash Time (before sample): 0	Auto Dilution Factor: 1

Replicate Data: 170329-L-01

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.755	4.53	0.0422	0.1503	0.0423	3:10:29 PM	Yes
2	0.719	4.31	0.0402	0.1233	0.0403	3:11:15 PM	Yes
Mean:	0.737	4.42	0.0412				
SD:	0.0251	0.151	0.0014				
%RSD:	3.41%	3.41%	3.39				

=====

Sequence No.: 11	Autosampler Location: 42
Sample ID: 17-03-1569-1	Date Collected: 3/29/2017 3:11:41 PM
Analyst: 868 HG-8	Data Type: Original
Initial Sample Wt: 0.61 g	Initial Sample Vol:
Dilution:	Sample Prep Vol: 100 mL
Wash Time (before sample): 0	Auto Dilution Factor: 1

Replicate Data: 17-03-1569-1

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0616	0.376	0.0037	0.0135	0.0038	3:12:46 PM	Yes
2	0.0588	0.359	0.0036	0.0113	0.0036	3:13:31 PM	Yes
Mean:	0.0602	0.367	0.0036				
SD:	0.00194	0.0118	0.0001				
%RSD:	3.22%	3.22%	3.01				

Sequence No.: 12

Autosampler Location: 5

Sample ID: CCV 0.2x10ppb

Date Collected: 3/29/2017 3:13:57 PM

Analyst: 868 HG-8

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time (before sample): 0

Auto Dilution Factor: 1.0000

Replicate Data: CCV 0.2x10ppb

Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00179	1.79	0.0169	0.0598	0.0169	3:15:03 PM	Yes
2	0.00169	1.69	0.0159	0.0483	0.0160	3:15:49 PM	Yes
Mean:	0.00174	1.74	0.0164				
SD:	0.000073	0.073	0.0007				
%RSD:	4.18%	4.18%	4.11				

QC value within limits for Hg 253.7 Recovery = 87.13%

All analyte(s) passed QC.

Sequence No.: 13

Autosampler Location: 1

Sample ID: CCB

Date Collected: 3/29/2017 3:16:16 PM

Analyst: 868 HG-8

Data Type: Original

Initial Sample Wt:

Initial Sample Vol:

Dilution:

Sample Prep Vol:

Wash Time (before sample): 0

Auto Dilution Factor: 1.0000

Replicate Data: CCB

Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	-0.000031	-0.0311	-0.0000	-0.0021	-0.0000	3:17:21 PM	Yes
2	-0.000031	-0.0312	-0.0000	-0.0010	-0.0000	3:18:06 PM	Yes
Mean:	-0.000031	-0.0312	-0.0000				
SD:	0.0000001	0.00007	0.0000				
%RSD:	0.24%	0.24%	1.56				

QC value within limits for Hg 253.7 Recovery = Not calculated

All analyte(s) passed QC.

Sequence No.: 14

Autosampler Location: 43

Sample ID: 17-03-1569-1 MS

Date Collected: 3/29/2017 3:18:32 PM

Analyst: 868 HG-8

Data Type: Original

Initial Sample Wt: 0.61 g

Initial Sample Vol:

Dilution:

Sample Prep Vol: 100 mL

Wash Time (before sample): 0

Auto Dilution Factor: 1

Replicate Data: 17-03-1569-1 MS

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.731	4.46	0.0416	0.1532	0.0416	3:19:36 PM	Yes
2	0.707	4.31	0.0402	0.1216	0.0402	3:20:22 PM	Yes
Mean:	0.719	4.39	0.0409				
SD:	0.0173	0.106	0.0010				
%RSD:	2.41%	2.41%	2.39				

Sequence No.: 15

Autosampler Location: 44

Sample ID: 17-03-1569-1 MSD

Date Collected: 3/29/2017 3:20:48 PM

Analyst: 868 HG-8

Data Type: Original

Return to Contents

Initial Sample Wt: 0.59 g
Dilution:
Wash Time (before sample): 0

Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1569-1 MSD

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.821	4.84	0.0451	0.1638	0.0452	3:21:53 PM	Yes
2	0.784	4.63	0.0431	0.1293	0.0432	3:22:39 PM	Yes
Mean:	0.802	4.73	0.0441				
SD:	0.0260	0.154	0.0014				
%RSD:	3.25%	3.25%	3.23				

=====

Sequence No.: 16
Sample ID: 17-03-1569-2
Analyst: 868 HG-8
Initial Sample Wt: 0.61 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 45
Date Collected: 3/29/2017 3:23:05 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1569-2

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0817	0.499	0.0049	0.0184	0.0049	3:24:10 PM	Yes
2	0.0782	0.477	0.0047	0.0149	0.0047	3:24:56 PM	Yes
Mean:	0.0800	0.488	0.0048				
SD:	0.00250	0.0153	0.0001				
%RSD:	3.13%	3.13%	2.97				

=====

Sequence No.: 17
Sample ID: 17-03-1569-3
Analyst: 868 HG-8
Initial Sample Wt: 0.59 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 46
Date Collected: 3/29/2017 3:25:23 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1569-3

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.119	0.699	0.0067	0.0235	0.0068	3:26:28 PM	Yes
2	0.114	0.671	0.0065	0.0198	0.0065	3:27:14 PM	Yes
Mean:	0.116	0.685	0.0066				
SD:	0.0034	0.0199	0.0002				
%RSD:	2.90%	2.90%	2.80				

=====

Sequence No.: 18
Sample ID: 17-03-1569-4
Analyst: 868 HG-8
Initial Sample Wt: 0.6 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 47
Date Collected: 3/29/2017 3:27:40 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1569-4

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0924	0.554	0.0054	0.0187	0.0054	3:28:46 PM	Yes
2	0.0900	0.540	0.0052	0.0160	0.0053	3:29:31 PM	Yes
Mean:	0.0912	0.547	0.0053				
SD:	0.00171	0.0103	0.0001				
%RSD:	1.88%	1.88%	1.79				

=====

Sequence No.: 19
Sample ID: 17-03-1569-5
Analyst: 868 HG-8

Autosampler Location: 48
Date Collected: 3/29/2017 3:29:58 PM
Data Type: Original

Initial Sample Wt: 0.61 g
Dilution:
Wash Time (before sample): 0

Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1569-5

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlnkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.107	0.656	0.0063	0.0219	0.0064	3:31:04 PM	Yes
2	0.101	0.614	0.0059	0.0188	0.0060	3:31:50 PM	Yes
Mean:	0.104	0.635	0.0061				
SD:	0.0048	0.0294	0.0003				
%RSD:	4.62%	4.62%	4.44				

Sequence No.: 20
Sample ID: 17-03-1569-6
Analyst: 868 HG-8
Initial Sample Wt: 0.62 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 49
Date Collected: 3/29/2017 3:32:17 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1569-6

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlnkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0380	0.236	0.0024	0.0089	0.0025	3:33:23 PM	Yes
2	0.0356	0.221	0.0023	0.0067	0.0023	3:34:08 PM	Yes
Mean:	0.0368	0.228	0.0024				
SD:	0.00170	0.0105	0.0001				
%RSD:	4.60%	4.60%	4.12				

Sequence No.: 21
Sample ID: 17-03-1556-25
Analyst: 868 HG-8
Initial Sample Wt: 0.61 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 50
Date Collected: 3/29/2017 3:34:35 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1556-25

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlnkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0103	0.0628	0.0008	0.0029	0.0009	3:35:40 PM	Yes
2	0.00997	0.0608	0.0008	0.0020	0.0008	3:36:26 PM	Yes
Mean:	0.0101	0.0618	0.0008				
SD:	0.00023	0.00143	0.0000				
%RSD:	2.31%	2.31%	1.62				

Sequence No.: 22
Sample ID: 17-03-1556-26
Analyst: 868 HG-8
Initial Sample Wt: 0.58 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 51
Date Collected: 3/29/2017 3:36:52 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1556-26

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StdConc ug/L	BlnkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00923	0.0535	0.0007	0.0023	0.0008	3:37:57 PM	Yes
2	0.00943	0.0547	0.0008	0.0019	0.0008	3:38:43 PM	Yes
Mean:	0.00933	0.0541	0.0007				
SD:	0.000144	0.00084	0.0000				
%RSD:	1.54%	1.54%	1.04				

Sequence No.: 23
Sample ID: 17-03-1556-27
Analyst: 868 HG-8

Autosampler Location: 52
Date Collected: 3/29/2017 3:39:09 PM
Data Type: Original

Initial Sample Wt: 0.6 g
Dilution:
Wash Time (before sample): 0

Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1556-27

Analyte: Hg 253.7

Table with 8 columns: Repl #, SampleConc mg/kg, StndConc ug/L, BlnkCorr Signal, Peak Area, Peak Height, Time, Peak Stored. Contains 2 replicate rows and summary statistics (Mean, SD, %RSD).

Sequence No.: 24
Sample ID: CCV 0.2x10ppb
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0

Autosampler Location: 5
Date Collected: 3/29/2017 3:41:26 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: CCV 0.2x10ppb

Analyte: Hg 253.7

Table with 8 columns: Repl #, SampleConc mg/L, StndConc ug/L, BlnkCorr Signal, Peak Area, Peak Height, Time, Peak Stored. Contains 2 replicate rows and summary statistics (Mean, SD, %RSD).

QC value less than the lower limit for Hg 253.7 Recovery = 75.41%
QC Failed. Stop the analysis.

Analysis Begun

Logged In Analyst: US26_SVC_INSTRUMENT
Spectrometer: FIMS-400, S/N B050-9560

Technique: AA FIMS-MHS
Autosampler: S10

Sample Information File: C:\Users\Public\PerkinElmer Syngistix\AA\Data\Sample Information\170329H1.sifx

Batch ID:
Results Data Set: 170329H1
Results Library: U:\MERCURY_8\Data\Results\results.mdb

Sequence No.: 24

Autosampler Location: 5

Sample ID: CCV 0.2x10ppb
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0

Date Collected: 3/29/2017 3:48:12 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: CCV 0.2x10ppb

Analyte: Hg 253.7

Table with 8 columns: Repl #, SampleConc mg/L, StndConc ug/L, BlnkCorr Signal, Peak Area, Peak Height, Time, Peak Stored. Contains 2 replicate rows and summary statistics (Mean, SD, %RSD).

QC value within limits for Hg 253.7 Recovery = 85.69%
All analyte(s) passed QC.

Sequence No.: 25
Sample ID: CCB
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:

Autosampler Location: 1
Date Collected: 3/29/2017 3:50:31 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:

Wash Time (before sample): 0

Auto Dilution Factor: 1.0000

Replicate Data: CCB

Analyte: Hg 253.7

Repl #	Sample Conc mg/L	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	-0.000021	-0.0206	0.0001	0.0010	0.0001	3:51:35 PM	Yes
2	-0.000015	-0.0147	0.0001	0.0014	0.0001	3:52:21 PM	Yes
Mean:	-0.000018	-0.0176	0.0001				
SD:	0.000042	0.00420	0.0000				
%RSD:	23.82%	23.82%	47.64				

QC value within limits for Hg 253.7 Recovery = Not calculated
All analyte(s) passed QC.

Sequence No.: 26

Autosampler Location: 53

Sample ID: 17-03-1557-25

Date Collected: 3/29/2017 3:52:46 PM

Analyst: 868 HG-8

Data Type: Original

Initial Sample Wt: 0.63 g

Initial Sample Vol:

Dilution:

Sample Prep Vol: 100 mL

Wash Time (before sample): 0

Auto Dilution Factor: 1

Replicate Data: 17-03-1557-25

Analyte: Hg 253.7

Repl #	Sample Conc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0195	0.123	0.0014	0.0054	0.0014	3:53:51 PM	Yes
2	0.0185	0.116	0.0013	0.0048	0.0014	3:54:37 PM	Yes
Mean:	0.0190	0.120	0.0014				
SD:	0.00075	0.0047	0.0000				
%RSD:	3.93%	3.93%	3.22				

Sequence No.: 27

Autosampler Location: 54

Sample ID: 17-03-1557-26

Date Collected: 3/29/2017 3:55:03 PM

Analyst: 868 HG-8

Data Type: Original

Initial Sample Wt: 0.62 g

Initial Sample Vol:

Dilution:

Sample Prep Vol: 100 mL

Wash Time (before sample): 0

Auto Dilution Factor: 1

Replicate Data: 17-03-1557-26

Analyte: Hg 253.7

Repl #	Sample Conc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0593	0.368	0.0037	0.0132	0.0037	3:56:08 PM	Yes
2	0.0564	0.350	0.0035	0.0112	0.0035	3:56:54 PM	Yes
Mean:	0.0578	0.359	0.0036				
SD:	0.00206	0.0128	0.0001				
%RSD:	3.57%	3.57%	3.32				

Sequence No.: 28

Autosampler Location: 55

Sample ID: 17-03-1557-27

Date Collected: 3/29/2017 3:57:20 PM

Analyst: 868 HG-8 *0.6*

Data Type: Original

Initial Sample Wt: *0.5* g

Initial Sample Vol:

Dilution:

Sample Prep Vol: 100 mL

Wash Time (before sample): 0

Auto Dilution Factor: 1

PF = 166.6666670425

Replicate Data: 17-03-1557-27

Analyte: Hg 253.7

Repl #	Sample Conc mg/kg	StdConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0113	0.0565	0.0008	0.0010	0.0008	3:58:25 PM	Yes
2	0.0117	0.0586	0.0008	0.0014	0.0008	3:59:11 PM	Yes
Mean:	0.0115	0.0575	0.0008				
SD:	0.00029	0.00146	0.0000				
%RSD:	2.53%	2.53%	1.73				

Sequence No.: 29

Autosampler Location: 56

Sample ID: 17-03-1679-1

Date Collected: 3/29/2017 3:59:38 PM

Analyst: 868 HG-8

Data Type: Original

Initial Sample Wt: 0.6 g
Dilution:
Wash Time (before sample): 0

Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 170329-B-02

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	-0.00428	-0.0257	0.0000	0.0000	0.0000	4:09:53 PM	Yes
2	-0.00430	-0.0258	0.0000	-0.0000	0.0000	4:10:39 PM	Yes
Mean:	-0.00429	-0.0257	0.0000				
SD:	0.000019	0.00011	0.0000				
%RSD:	0.44%	0.44%	16.23				

Sequence No.: 34
Sample ID: 170329-L-02
Analyst: 868 HG-8
Initial Sample Wt: 0.6 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 61
Date Collected: 3/29/2017 4:11:06 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 170329-L-02

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.743	4.46	0.0416	0.1463	0.0416	4:12:11 PM	Yes
2	0.703	4.22	0.0393	0.1210	0.0394	4:12:56 PM	Yes
Mean:	0.723	4.34	0.0404				
SD:	0.0284	0.170	0.0016				
%RSD:	3.93%	3.93%	3.91				

Sequence No.: 35
Sample ID: 17-03-1672-1
Analyst: 868 HG-8
Initial Sample Wt: 0.6 g
Dilution:
Wash Time (before sample): 0

Autosampler Location: 62
Date Collected: 3/29/2017 4:13:23 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol: 100 mL
Auto Dilution Factor: 1

Replicate Data: 17-03-1672-1

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0777	0.466	0.0046	0.0165	0.0046	4:14:29 PM	Yes
2	0.0722	0.433	0.0043	0.0131	0.0043	4:15:14 PM	Yes
Mean:	0.0750	0.450	0.0044				
SD:	0.00391	0.0235	0.0002				
%RSD:	5.22%	5.22%	4.93				

Sequence No.: 36
Sample ID: CCV 0.2x10ppb
Analyst: 868 HG-8
Initial Sample Wt:
Dilution:
Wash Time (before sample): 0

Autosampler Location: 5
Date Collected: 3/29/2017 4:15:41 PM
Data Type: Original
Initial Sample Vol:
Sample Prep Vol:
Auto Dilution Factor: 1.0000

Replicate Data: CCV 0.2x10ppb

Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	BlkCorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.00183	1.83	0.0172	0.0618	0.0172	4:16:47 PM	Yes
2	0.00174	1.74	0.0164	0.0503	0.0164	4:17:32 PM	Yes
Mean:	0.00179	1.79	0.0168				
SD:	0.000061	0.061	0.0006				
%RSD:	3.42%	3.42%	3.37				

QC value within limits for Hg 253.7 Recovery = 89.30%
All analyte(s) passed QC.

Sequence No.: 37

Autosampler Location: 1



Method: EPA 7470A+7471A-Hg-8

Page 11

Date: 3/29/2017 4:27:08 PM

Sample ID: CCB
 Analyst: 868 HG-8
 Initial Sample Wt:
 Dilution:
 Wash Time (before sample): 0

Date Collected: 3/29/2017 4:17:59 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol:
 Auto Dilution Factor: 1.0000

Replicate Data: CCB

Analyte: Hg 253.7

Repl #	SampleConc mg/L	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	-0.000032	-0.0324	-0.0001	-0.0016	-0.0000	4:19:04 PM	Yes
2	-0.000029	-0.0293	-0.0000	-0.0019	0.0000	4:19:49 PM	Yes
Mean:	-0.000031	-0.0308	-0.0000				
SD:	0.000021	0.00215	0.0000				
%RSD:	6.96%	6.96%	48.69				

QC value within limits for Hg 253.7 Recovery = Not calculated
 All analyte(s) passed QC.

=====

Sequence No.: 38
 Sample ID: 17-03-1672-1 MS
 Analyst: 868 HG-8
 Initial Sample Wt: 0.61 g
 Dilution:
 Wash Time (before sample): 0

Autosampler Location: 63
 Date Collected: 3/29/2017 4:20:15 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Replicate Data: 17-03-1672-1 MS

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.705	4.30	0.0401	0.1413	0.0401	4:21:20 PM	Yes
2	0.678	4.13	0.0386	0.1198	0.0386	4:22:06 PM	Yes
Mean:	0.691	4.22	0.0393				
SD:	0.0190	0.116	0.0011				
%RSD:	2.75%	2.75%	2.73				

=====

Sequence No.: 39
 Sample ID: 17-03-1672-1 MSD
 Analyst: 868 HG-8
 Initial Sample Wt: 0.62 g
 Dilution:
 Wash Time (before sample): 0

Autosampler Location: 64
 Date Collected: 3/29/2017 4:22:33 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Replicate Data: 17-03-1672-1 MSD

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.712	4.41	0.0412	0.1456	0.0412	4:23:38 PM	Yes
2	0.681	4.22	0.0394	0.1232	0.0394	4:24:23 PM	Yes
Mean:	0.696	4.32	0.0403				
SD:	0.0220	0.137	0.0013				
%RSD:	3.16%	3.16%	3.14				

=====

Sequence No.: 40
 Sample ID: 17-03-1681-1
 Analyst: 868 HG-8
 Initial Sample Wt: 0.63 g
 Dilution:
 Wash Time (before sample): 0

Autosampler Location: 65
 Date Collected: 3/29/2017 4:24:50 PM
 Data Type: Original
 Initial Sample Vol:
 Sample Prep Vol: 100 mL
 Auto Dilution Factor: 1

Replicate Data: 17-03-1681-1

Analyte: Hg 253.7

Repl #	SampleConc mg/kg	StndConc ug/L	Blncorr Signal	Peak Area	Peak Height	Time	Peak Stored
1	0.0262	0.165	0.0018	0.0065	0.0018	4:25:55 PM	Yes
2	0.0228	0.143	0.0016	0.0043	0.0016	4:26:41 PM	Yes
Mean:	0.0245	0.154	0.0017				
SD:	0.00241	0.0152	0.0001				
%RSD:	9.85%	9.85%	8.41				

EPA 7471A Mercury (Solid)

Run Logs

170329H1

Carrier solution R07141602

Reducing Agent R07141603

Sample ID	Analyst Name	Sample Name	Analyte Name	Date	Time	Conc (Calib)	Units (Calib)	Conc (Samp)	Units (Samp)	Corr Coef
Calib blank_868	868 HG-8		Hg 253.7	3/29/2017	11:25:00 AM	ug/L			mg/L	
0.025ppb 0.005x5ppb	868 HG-8		Hg 253.7	3/29/2017	11:27:14 AM	ug/L			mg/L	
0.10ppb M030617AX0.0001	868 HG-8		Hg 253.7	3/29/2017	11:29:30 AM	ug/L			mg/L	
1.00ppb M030617AX0.001	868 HG-8		Hg 253.7	3/29/2017	11:31:45 AM	ug/L			mg/L	
2.00ppb M030617AX0.002	868 HG-8		Hg 253.7	3/29/2017	11:34:02 AM	ug/L			mg/L	
5.00ppb M030617AX0.005	868 HG-8		Hg 253.7	3/29/2017	11:36:18 AM	ug/L			mg/L	
10.0ppb M030617AX0.01	868 HG-8		Hg 253.7	3/29/2017	11:38:33 AM	ug/L			mg/L	
ICV M030617B	868 HG-8		Hg 253.7	3/29/2017	11:45:01 AM	4.964654 ug/L		0.004965 mg/L		0.999956
ICB	868 HG-8		Hg 253.7	3/29/2017	11:47:15 AM	-0.02763 ug/L		-2.76E-05 mg/L		0.999956
CRQL 0.25	868 HG-8	0.6	Hg 253.7	3/29/2017	11:49:31 AM	0.246378 ug/L		0.041063 mg/kg		0.999956
17-03-2051-1	868 HG-8	0.61	Hg 253.7	3/29/2017	11:51:46 AM	0.229835 ug/L		0.037678 mg/kg		0.999956
17-03-2051-3	868 HG-8	0.6	Hg 253.7	3/29/2017	11:54:01 AM	0.290939 ug/L		0.04849 mg/kg		0.999956
CCV 0.2x10ppb	868 HG-8		Hg 253.7	3/29/2017	11:56:18 AM	1.99546 ug/L		0.001995 mg/L		0.999956
CCB	868 HG-8		Hg 253.7	3/29/2017	11:58:33 AM	-0.02897 ug/L		-2.90E-05 mg/L		0.999956
CCV 0.2x10ppb } *	868 HG-8		Hg 253.7	3/29/2017	2:17:27 PM	1.450055 ug/L		0.00145 mg/L		0.999956
CCV 0.2x10ppb }	868 HG-8		Hg 253.7	3/29/2017	2:19:44 PM	1.998235 ug/L		0.001998 mg/L		0.999956
CCB } **	868 HG-8		Hg 253.7	3/29/2017	2:22:01 PM	1.379949 ug/L		0.00138 mg/L		0.999956
CCB }	868 HG-8		Hg 253.7	3/29/2017	2:48:04 PM	-0.01864 ug/L		-1.86E-05 mg/L		0.999956
17-03-2069-1	868 HG-8	0.63	Hg 253.7	3/29/2017	2:50:20 PM	0.037432 ug/L		0.005942 mg/kg		0.999956
17-03-1667-1	868 HG-8	0.61	Hg 253.7	3/29/2017	2:52:38 PM	0.055516 ug/L		0.009101 mg/kg		0.999956
17-03-1667-3	868 HG-8	0.62	Hg 253.7	3/29/2017	2:54:55 PM	0.038839 ug/L		0.006264 mg/kg		0.999956
17-03-1667-5	868 HG-8	0.62	Hg 253.7	3/29/2017	2:59:51 PM	0.054284 ug/L		0.008755 mg/kg		0.999956
17-03-1667-7	868 HG-8	0.59	Hg 253.7	3/29/2017	3:02:09 PM	0.040424 ug/L		0.006852 mg/kg		0.999956
17-03-1667-9	868 HG-8	0.61	Hg 253.7	3/29/2017	3:04:25 PM	0.034079 ug/L		0.005587 mg/kg		0.999956
CRQL 0.25	868 HG-8	0.6	Hg 253.7	3/29/2017	3:06:42 PM	0.112428 ug/L		0.018738 mg/kg		0.999956
170329-B-01	868 HG-8	0.6	Hg 253.7	3/29/2017	3:08:58 PM	-0.00955 ug/L		-0.00159 mg/kg		0.999956

* time gap, CCV did not pass, re-run - passed
 ** CCB failed, re-run - passed

Reviewed/Assign to Logbook Date: 03-29-17
 Analysis Hg Chemist ID: 309
 Logbook Page: 85 Instrument ID: Hg-8

Sample ID	Analyst Name	Sample Wt	Analyte Name	Date	Time	Conc		Units		Corr Coef
						(Calib)	(Samp)	(Calib)	(Samp)	
170329-L01	868 HG-8	0.6	Hg 253.7	3/29/2017	3:11:15 PM	4.42118	0.736863	ug/L	mg/kg	0.999956
17-03-1569-1	868 HG-8	0.61	Hg 253.7	3/29/2017	3:13:31 PM	0.367148	0.060188	ug/L	mg/kg	0.999956
CCV 0.2x10ppb	868 HG-8		Hg 253.7	3/29/2017	3:15:49 PM	1.742626	0.001743	ug/L	mg/L	0.999956
CCB	868 HG-8		Hg 253.7	3/29/2017	3:18:06 PM	-0.03118	-3.12E-05	ug/L	mg/L	0.999956
17-03-1569-1 MS	868 HG-8	0.61	Hg 253.7	3/29/2017	3:20:22 PM	4.385517	0.718937	ug/L	mg/kg	0.999956
17-03-1569-1 MSD	868 HG-8	0.59	Hg 253.7	3/29/2017	3:22:39 PM	4.733645	0.802313	ug/L	mg/kg	0.999956
17-03-1569-2	868 HG-8	0.61	Hg 253.7	3/29/2017	3:24:56 PM	0.487886	0.079981	ug/L	mg/kg	0.999956
17-03-1569-3	868 HG-8	0.59	Hg 253.7	3/29/2017	3:27:14 PM	0.685259	0.116146	ug/L	mg/kg	0.999956
17-03-1569-4	868 HG-8	0.6	Hg 253.7	3/29/2017	3:29:31 PM	0.547062	0.091177	ug/L	mg/kg	0.999956
17-03-1569-5	868 HG-8	0.61	Hg 253.7	3/29/2017	3:31:50 PM	0.634961	0.104092	ug/L	mg/kg	0.999956
17-03-1569-6	868 HG-8	0.62	Hg 253.7	3/29/2017	3:34:08 PM	0.228437	0.036845	ug/L	mg/kg	0.999956
17-03-1556-25	868 HG-8	0.61	Hg 253.7	3/29/2017	3:36:26 PM	0.061837	0.010137	ug/L	mg/kg	0.999956
17-03-1556-26	868 HG-8	0.58	Hg 253.7	3/29/2017	3:38:43 PM	0.054118	0.009331	ug/L	mg/kg	0.999956
17-03-1556-27	868 HG-8	0.6	Hg 253.7	3/29/2017	3:40:59 PM	0.088094	0.014682	ug/L	mg/kg	0.999956
CCV 0.2x10ppb	868 HG-8		Hg 253.7	3/29/2017	3:43:17 PM	1.508247	0.001508	ug/L	mg/L	0.999956
CCV 0.2x10ppb	868 HG-8		Hg 253.7	3/29/2017	3:50:04 PM	1.713826	0.001714	ug/L	mg/L	0.999956
CCB	868 HG-8		Hg 253.7	3/29/2017	3:52:21 PM	-0.01762	-1.76E-05	ug/L	mg/L	0.999956
17-03-1557-25	868 HG-8	0.63	Hg 253.7	3/29/2017	3:54:37 PM	0.119648	0.018992	ug/L	mg/kg	0.999956
17-03-1557-26	868 HG-8	0.62	Hg 253.7	3/29/2017	3:56:54 PM	0.358583	0.057836	ug/L	mg/kg	0.999956
17-03-1557-27	868 HG-8	0.6	Hg 253.7	3/29/2017	3:59:11 PM	0.057546	0.009591	ug/L	mg/kg	0.999956
17-03-1679-1	868 HG-8	0.6	Hg 253.7	3/29/2017	4:01:28 PM	-0.01356	-0.00226	ug/L	mg/kg	0.999956
17-03-1679-2	868 HG-8	0.63	Hg 253.7	3/29/2017	4:03:46 PM	-0.0119	-0.00189	ug/L	mg/kg	0.999956
17-03-1679-3	868 HG-8	0.63	Hg 253.7	3/29/2017	4:06:02 PM	-3.80E-05	-6.03E-06	ug/L	mg/kg	0.999956
17-03-1679-4	868 HG-8	0.63	Hg 253.7	3/29/2017	4:08:20 PM	0.095148	0.015103	ug/L	mg/kg	0.999956
170329-B-02	868 HG-8	0.6	Hg 253.7	3/29/2017	4:10:39 PM	-0.02574	-0.00429	ug/L	mg/kg	0.999956
170329-L-02	868 HG-8	0.6	Hg 253.7	3/29/2017	4:12:56 PM	4.335919	0.722653	ug/L	mg/kg	0.999956
17-03-1672-1	868 HG-8	0.6	Hg 253.7	3/29/2017	4:15:14 PM	0.449749	0.074958	ug/L	mg/kg	0.999956
CCV 0.2x10ppb	868 HG-8		Hg 253.7	3/29/2017	4:17:32 PM	1.786031	0.001786	ug/L	mg/L	0.999956
CCB	868 HG-8		Hg 253.7	3/29/2017	4:19:49 PM	-0.03084	-3.08E-05	ug/L	mg/L	0.999956

* CCV failed, re-run - passed

Reviewed/Assign to Logbook Date:	03-29-17
Analysis	Hg
Chemist ID:	309
Logbook Page:	86
Instrument ID:	Hg-8

EPA 7471A Mercury (Solid)

Preparation Logs

Mercury Sample Preparation Logbook (Solid / Other)

METHOD		MATRIX		EQUIPMENT ID #			REAGENT ID #			STANDARD ID #				
<input type="checkbox"/> EPA 7471A	<input checked="" type="checkbox"/> Solid			Thermometer	GT-4 (CF-2.0°C)		Aqua Regia	R04291601 10 mL		Spike M030617 A				
<input type="checkbox"/> EPA 7471B	<input type="checkbox"/> Other (Specify)			Block Digester	3		5% KMnO ₄	R10061601		IC A				
				Pipetter / Dispenser	P071/*		NaCl-H ₂ NO-HCl	R10261601 6 mL		ICV B				
BATCH NUMBER			SUPPLY LOT #			BALANCE ID #			QUALITY SYSTEM MATRIX ID #			SAMPLE HANDLING		
MS/MSD 170329.5(1)			Tube / Container 146977			36			Teflon Chip M066-35-15			1 = Composite 2 = Subsample		
(Specify)			Filter						(Specify)			3 = Homogenize 4 = None		
DIGESTION													SPIKE OR IC/ICV	
DATE	START			END			SAMPLE HANDLING	ECI ID #	SAMPLE		5% KMNO ₄ V (mL)	SPIKE OR IC/ICV V (μL)		
	TIME	TEMP W/O CF (°C)	PREP TECH ID #	TIME	TEMP W/O CF (°C)	PREP TECH ID #			INITIAL (g)	FINAL (mL)				
3/29/17	10:00	95	805	10:45	95	805	3	MS 17-03-1569-1AA	0.61	100	15	500		
							3	MSD	0.59					
							4	LCS 170329-LV1	0.60					
							4	LCS/MB 170329-B01	0.60					
							3	17-03-1569-1AA	0.61					
								2	0.61					
								3	0.59					
								4	0.60					
								5	0.61					
								6	0.62					
							4	17-03-1556-25B	0.61					
								26	0.58					
								27	0.60					
								17-03-1557-25B	0.63					
								26	0.62					
								27	0.60					
								17-03-1617-1A	0.60					
								17-03-1619-1A	0.60					
								2A	0.63					
								868 25/29/12 3A	0.63					
								17-03-1619-1A 4A	0.63					
								IC						
								ICV						
								CB						

COMMENTS: *D003 / MD043 / D058

EPA METHOD 8082 PCB

RAW DATA

EPA METHOD 8082 PCB

Initial Calibration

INITIAL CALIBRATION QUALITY CONTROL SUMMARY FOR METHOD: EPA 8082

ICAL WORK ORDER: 099-12-532-9351-5154
 ICAL BATCH ID: 1702221003
 INSTRUMENT: GC 66

ANALYZED BY: 669
 ICAL D/T ANALYZED: 2017-02-22 17:27
 REVIEWED BY: 27
 D/T REVIEWED: 2017-02-24 16:33

COMPOUND	COMP. TYPE	CALIB. MODEL	1	2	3	4	5	6	7	8	9	Avg. RF	Min. RF	%RSD CL	R. or R ² CL	R. or R ² CL	STATUS
Aroclor-1016	C		975,287	941,045	882,730	864,861	802,600					893.304	0.00	8	0-20		PASS
Aroclor-1260	C		1,012,414	995,847	939,916	914,296	876,433					947.781	0.00	6	0-20		PASS

Data Files:

Level #	D/T Analyzed	Data File
1	2017-02-22 17:27	/chem1/SVOA/GC_66/170222A/b1702220117022201
2	2017-02-22 17:45	/chem1/SVOA/GC_66/170222A/b1702220217022202
3	2017-02-22 18:03	/chem1/SVOA/GC_66/170222A/b1702220317022203
4	2017-02-22 18:21	/chem1/SVOA/GC_66/170222A/b1702220417022204
5	2017-02-22 18:38	/chem1/SVOA/GC_66/170222A/b1702220517022205

LR - E: Linear Regression (Equal Weight)
 Avg RF: Average Response Factor
 LR - IC: Linear Regression (Inverse Concentration Weight)
 QR - E: Quadratic Regression (Equal Weight)
 LR - ISC: Linear Regression (Inverse Square Concentration Weight)



INITIAL CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8082

ICV WORK ORDER: 099-12-532-9351-5154

INITIAL BATCH: 170222I003

INSTRUMENT: GC 66

ANALYZED BY: 669

D/T ANALYZED:

INITIAL:

ICV:

REVIEWED BY: 27

D/T REVIEWED:

2017-02-22 17:27

2017-02-22 18:56

2017-02-24 16:33

DATA FILE: /chem1/SVOA/GC_66/170222A/b1702220617022206

<u>COMPOUND NAME</u>	<u>COMP TYPE</u>	<u>CALIB MODEL</u>	<u>MIN RF</u>	<u>AVG RF</u>	<u>ICV RF</u>	<u>AMOUNT</u>	<u>ICV CONC</u>	<u>ICV %D</u>	<u>ICV %D CL</u>	<u>STATUS</u>
Atroclor-1016	C	Avg Resp	0.00	893304.438	1007962.398			-13	0-15	PASS
Atroclor-1260	C	Avg Resp	0.00	947781.003	1078849.128			-14	0-15	PASS

MIN RF: Method Specified Minimum Response Factor

Report Date : 03-Mar-2017 10:49

Page 1

Eurofins Calscience
INITIAL CALIBRATION DATA

Start Cal Date : 22-FEB-2017 17:27
 End Cal Date : 22-FEB-2017 20:07
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Cal Date : 03-Mar-2017 10:49 uhhn
 Curve Type : Average

Calibration File Names:

Level 1: /chem1/SVOA/GC_66.i/170222A.b/b17022201.d
 Level 2: /chem1/SVOA/GC_66.i/170222A.b/b17022202.d
 Level 3: /chem1/SVOA/GC_66.i/170222A.b/b17022210.d
 Level 4: /chem1/SVOA/GC_66.i/170222A.b/b17022204.d
 Level 5: /chem1/SVOA/GC_66.i/170222A.b/b17022205.d

Compound	100.000	250.000	500.000	750.000	2000.000	RRF	% RSD
====	====	====	====	====	====	====	====
M 2 Aroclor-1016	975287	941045	882730	864861	802600	893304	8
3 Aroclor 1016 (1)	101963	96557	88889	85902	76867	90036	11
4 Aroclor 1016 (2)	177940	169399	157230	152953	139805	159465	9
5 Aroclor 1016 (3)	368139	367391	353292	350925	334872	354924	4
6 Aroclor 1016 (4)	185303	173132	158540	153765	139752	162098	11
7 Aroclor 1016 (5)	141942	134567	124778	121316	111304	126781	9
M 8 Aroclor-1260	1012414	995847	939916	914296	876433	947781	6
9 Aroclor 1260 (1)	295733	284622	266899	258701	245467	270285	7
10 Aroclor 1260 (2)	233964	225940	211260	204056	194563	213956	7
11 Aroclor 1260 (3)	244538	243699	232324	225188	219576	233065	5
12 Aroclor 1260 (4)	84774	85815	81119	85251	78652	83122	4
13 Aroclor 1260 (5)	153405	155772	148314	141098	138176	147353	5
M 14 Aroclor-1221	++++	++++	293209	++++	++++	293209	0
15 Aroclor 1221 (1)	++++	++++	43149	++++	++++	43149	0
16 Aroclor 1221 (2)	++++	++++	55578	++++	++++	55578	0
17 Aroclor 1221 (3)	++++	++++	34870	++++	++++	34870	0
18 Aroclor 1221 (4)	++++	++++	138535	++++	++++	138535	0
19 Aroclor 1221 (5)	++++	++++	21077	++++	++++	21077	0
M 20 Aroclor-1232	++++	++++	461201	++++	++++	461201	0
21 Aroclor 1232 (1)	++++	++++	93661	++++	++++	93661	0
22 Aroclor 1232 (2)	++++	++++	75982	++++	++++	75982	0
23 Aroclor 1232 (3)	++++	++++	160461	++++	++++	160461	0

Report Date : 03-Mar-2017 10:49

Page 2

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 22-FEB-2017 17:27
 End Cal Date : 22-FEB-2017 20:07
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Cal Date : 03-Mar-2017 10:49 uhhn
 Curve Type : Average

Compound	100.000 Level 1	250.000 Level 2	500.000 Level 3	750.000 Level 4	2000.000 Level 5	RRF	% RSD
24 Aroclor 1232 (4)	++++	++++	74085	++++	++++	74085	0
25 Aroclor 1232 (5)	++++	++++	57013	++++	++++	57013	0
M 26 Aroclor-1242	++++	++++	721392	++++	++++	721392	0
27 Aroclor 1242 (1)	++++	++++	71280	++++	++++	71280	0
28 Aroclor 1242 (2)	++++	++++	129507	++++	++++	129507	0
29 Aroclor 1242 (3)	++++	++++	287006	++++	++++	287006	0
30 Aroclor 1242 (4)	++++	++++	130903	++++	++++	130903	0
31 Aroclor 1242 (5)	++++	++++	102696	++++	++++	102696	0
M 32 Aroclor-1248	++++	++++	698713	++++	++++	698713	0
33 Aroclor 1248 (1)	++++	++++	185195	++++	++++	185195	0
34 Aroclor 1248 (2)	++++	++++	103145	++++	++++	103145	0
35 Aroclor 1248 (3)	++++	++++	76812	++++	++++	76812	0
36 Aroclor 1248 (4)	++++	++++	152023	++++	++++	152023	0
37 Aroclor 1248 (5)	++++	++++	181538	++++	++++	181538	0
M 38 Aroclor-1254	++++	++++	1318858	++++	++++	1318858	0
39 Aroclor 1254 (1)	++++	++++	265379	++++	++++	265379	0
40 Aroclor 1254 (2)	++++	++++	186148	++++	++++	186148	0
41 Aroclor 1254 (3)	++++	++++	350152	++++	++++	350152	0
42 Aroclor 1254 (4)	++++	++++	239737	++++	++++	239737	0
43 Aroclor 1254 (5)	++++	++++	277442	++++	++++	277442	0
M 44 Aroclor-1262	++++	++++	1059164	++++	++++	1059164	0
45 Aroclor 1262 (1)	++++	++++	199559	++++	++++	199559	0
46 Aroclor 1262 (2)	++++	++++	285769	++++	++++	285769	0
47 Aroclor 1262 (3)	++++	++++	264628	++++	++++	264628	0
48 Aroclor 1262 (4)	++++	++++	89771	++++	++++	89771	0
49 Aroclor 1262 (5)	++++	++++	219437	++++	++++	219437	0
M 50 Aroclor-1268	++++	++++	4264652	++++	++++	4264652	0
51 Aroclor 1268 (1)	++++	++++	707347	++++	++++	707347	0
52 Aroclor 1268 (2)	++++	++++	709040	++++	++++	709040	0

Report Date : 03-Mar-2017 10:49

Page 3

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 22-FEB-2017 17:27
 End Cal Date : 22-FEB-2017 20:07
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Cal Date : 03-Mar-2017 10:49 uhn
 Curve Type : Average

Compound	100.000 Level 1	250.000 Level 2	500.000 Level 3	750.000 Level 4	2000.000 Level 5	RRF	% RSD
53 Aroclor 1268 (3)	+++++	+++++	580421	+++++	+++++	580421	0
54 Aroclor 1268 (4)	+++++	+++++	255277	+++++	+++++	255277	0
55 Aroclor 1268 (5)	+++++	+++++	2012567	+++++	+++++	2012567	0
\$ 1 2,4,5,6-Tetrachloro-m-xylene	5417720	5823446	5836705	5879203	5752252	5741865	3
\$ 56 Decachlorobiphenyl	5382862	5603807	5539339	5351948	5420179	5459627	2

Report Date : 03-Mar-2017 10:49

Page 4

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 22-FEB-2017 17:27
End Cal Date : 22-FEB-2017 20:07
Quant Method : ESTD
Origin : Disabled
Target Version : 3.50
Integrator : HP Genie
Method file : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
Cal Date : 03-Mar-2017 10:49 uhhn
Curve Type : Average

```
|Average %RSD Results. |  
|=====|  
|Calculated Average %RSD = 6.53024 |  
|Maximun Average %RSD = 20.00000 |  
|* Passed Average %RSD Test. |  
|_____|
```

Eurofins CalScience
 Calibration Verification Report

Instrument ID: GC_66.i Injection Date and Time: 22-FEB-2017 18:56
 Sample Name: PCB ICV P021517H 500PPB Initial Calibration Date(s): 17-OCT-2016 22-FEB-2017
 Sublist used: p1016_1260.sub Initial Calibration Time(s): 20:04 20:07
 Method used: /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m

Target Compounds	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift	Max%D / Drift	Curve Type		
Aroclor 1260 (1)	270284.532	318239.382	0.01	-18	15	Averaged	<-Failed	
Aroclor 1260 (2)	213956.488	237968.342	0.01	-11	15	Averaged		
Aroclor 1260 (3)	233064.808	267435.304	0.01	-15	15	Averaged		
Aroclor 1260 (4)	83122.167	83559.856	0.01	-1	15	Averaged		
Aroclor-1260	947781.003	1078849.128	0.01	-14	15	Averaged		
Aroclor-1016	893304.438	1007962.398	0.01	-13	15	Averaged		
Aroclor 1016 (1)	90035.625	99898.782	0.01	-11	15	Averaged		
Aroclor 1016 (2)	159465.395	177206.146	0.01	-11	15	Averaged		
Aroclor 1016 (3)	354923.864	406965.954	0.01	-15	15	Averaged		
Aroclor 1016 (4)	162098.269	181154.002	0.01	-12	15	Averaged		
Aroclor 1016 (5)	126781.283	142737.514	0.01	-13	15	Averaged		
Aroclor 1260 (5)	147353.008	171646.244	0.01	-16	15	Averaged		<-Failed
Surrogate Standards	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift	Max%D / Drift	Curve Type		
2,4,5,6-Tetrachloro-m-xylene	5741865.170	5554897.950	0.01	3	15	Averaged		
Decachlorobiphenyl	5459627.040	5505849.720	0.01	-1	15	Averaged		

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022201.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 17:27
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB ICAL1 P021517F 100PPB
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 22-FEB-2017 17:27 Cal File: b17022201.d
 Als bottle: 1 Calibration Sample, Level: 1
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.694	4.694	0.000	108354401	20.0000	18.9(a)
M 2 Aroclor-1016				97528686	100.000	109
3 Aroclor 1016 (1)	5.539	5.533	0.006	10196308	100.000	113
4 Aroclor 1016 (2)	6.168	6.165	0.003	17794023	100.000	112
5 Aroclor 1016 (3)	6.852	6.844	0.008	36813879	100.000	104
6 Aroclor 1016 (4)	7.050	7.043	0.007	18530301	100.000	114
7 Aroclor 1016 (5)	7.204	7.194	0.010	14194175	100.000	112
M 8 Aroclor-1260				101241371	100.000	107
9 Aroclor 1260 (1)	9.753	9.749	0.004	29573320	100.000	109
10 Aroclor 1260 (2)	10.352	10.348	0.004	23396360	100.000	109
11 Aroclor 1260 (3)	10.745	10.742	0.003	24453781	100.000	105
12 Aroclor 1260 (4)	12.119	12.115	0.004	8477403	100.000	102
13 Aroclor 1260 (5)	12.340	12.337	0.003	15340507	100.000	104
\$ 56 Decachlorobiphenyl	13.107	13.103	0.004	107657233	20.0000	19.7(a)

QC Flag Legend

a - Target compound detected but, quantitated amount
 Below Limit Of Quantitation(BLOQ).

Data File: /chem1/SV004/GC_66.i/170222A.b/b17022201.d

Date : 22-FEB-2017 17:27

Client ID:

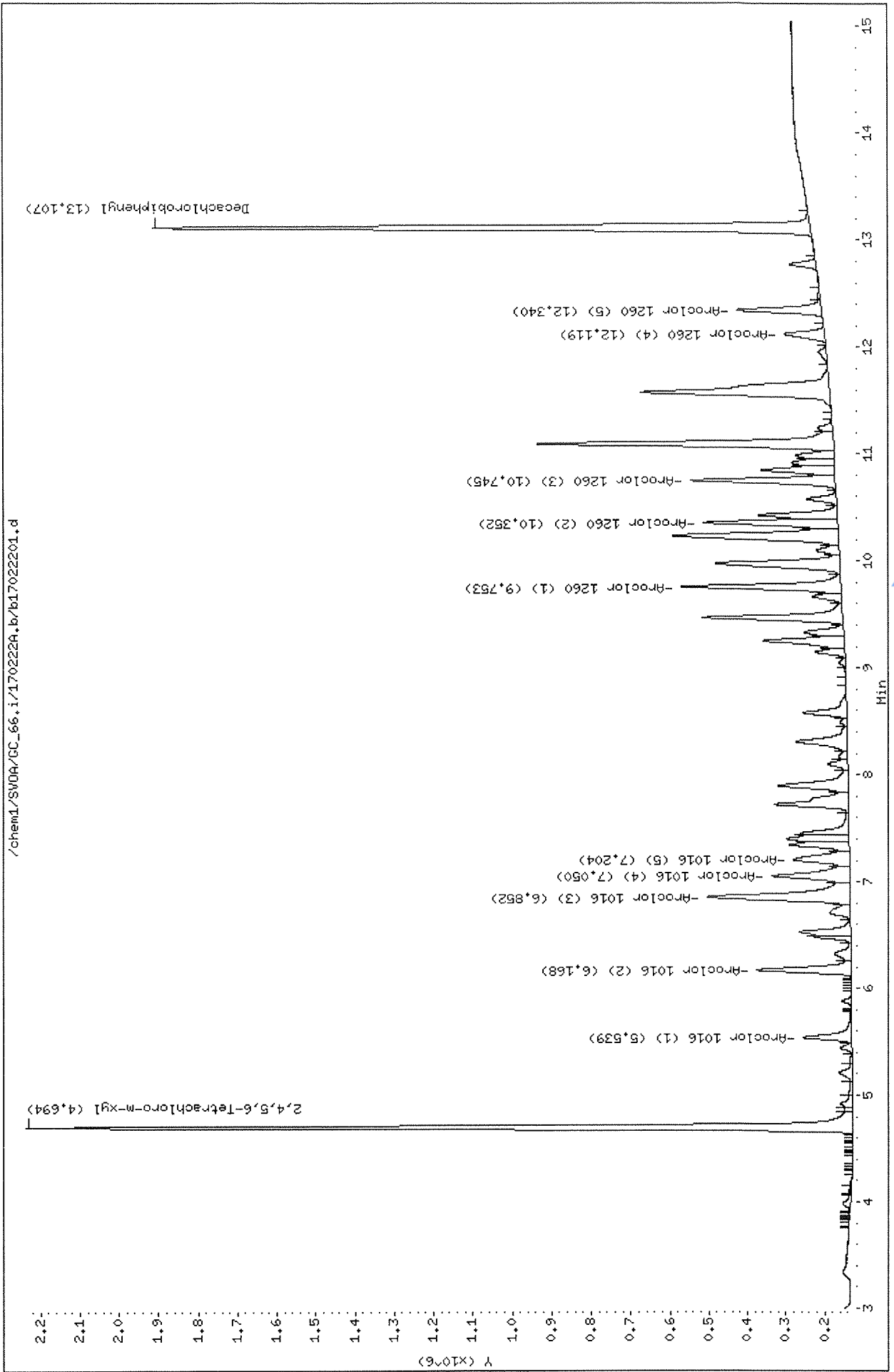
Sample Info: PCB ICAL1 F021517F 100PPB

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022202.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 17:45
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB ICAL2 P021517E 250PPB
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 22-FEB-2017 17:45 Cal File: b17022202.d
 Als bottle: 2 Calibration Sample, Level: 2
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.693	4.694	-0.001	291172299	50.0000	50.7
M 2 Aroclor-1016				235261233	250.000	263
3 Aroclor 1016 (1)	5.536	5.539	-0.003	24139170	250.000	268
4 Aroclor 1016 (2)	6.166	6.168	-0.002	42349726	250.000	266
5 Aroclor 1016 (3)	6.848	6.852	-0.004	91847697	250.000	259
6 Aroclor 1016 (4)	7.046	7.050	-0.004	43282979	250.000	267
7 Aroclor 1016 (5)	7.198	7.204	-0.006	33641661	250.000	265
M 8 Aroclor-1260				248961695	250.000	263
9 Aroclor 1260 (1)	9.751	9.753	-0.002	71155515	250.000	263
10 Aroclor 1260 (2)	10.349	10.352	-0.003	56484954	250.000	264
11 Aroclor 1260 (3)	10.743	10.745	-0.002	60924633	250.000	261
12 Aroclor 1260 (4)	12.116	12.119	-0.003	21453649	250.000	258
13 Aroclor 1260 (5)	12.339	12.340	-0.001	38942944	250.000	264
\$ 56 Decachlorobiphenyl	13.105	13.107	-0.002	280190360	50.0000	51.3

Data File: /chem1/SV04/GC_66.i/170222A.b/17022202.d

Date : 22-FEB-2017 17:45

Client ID:

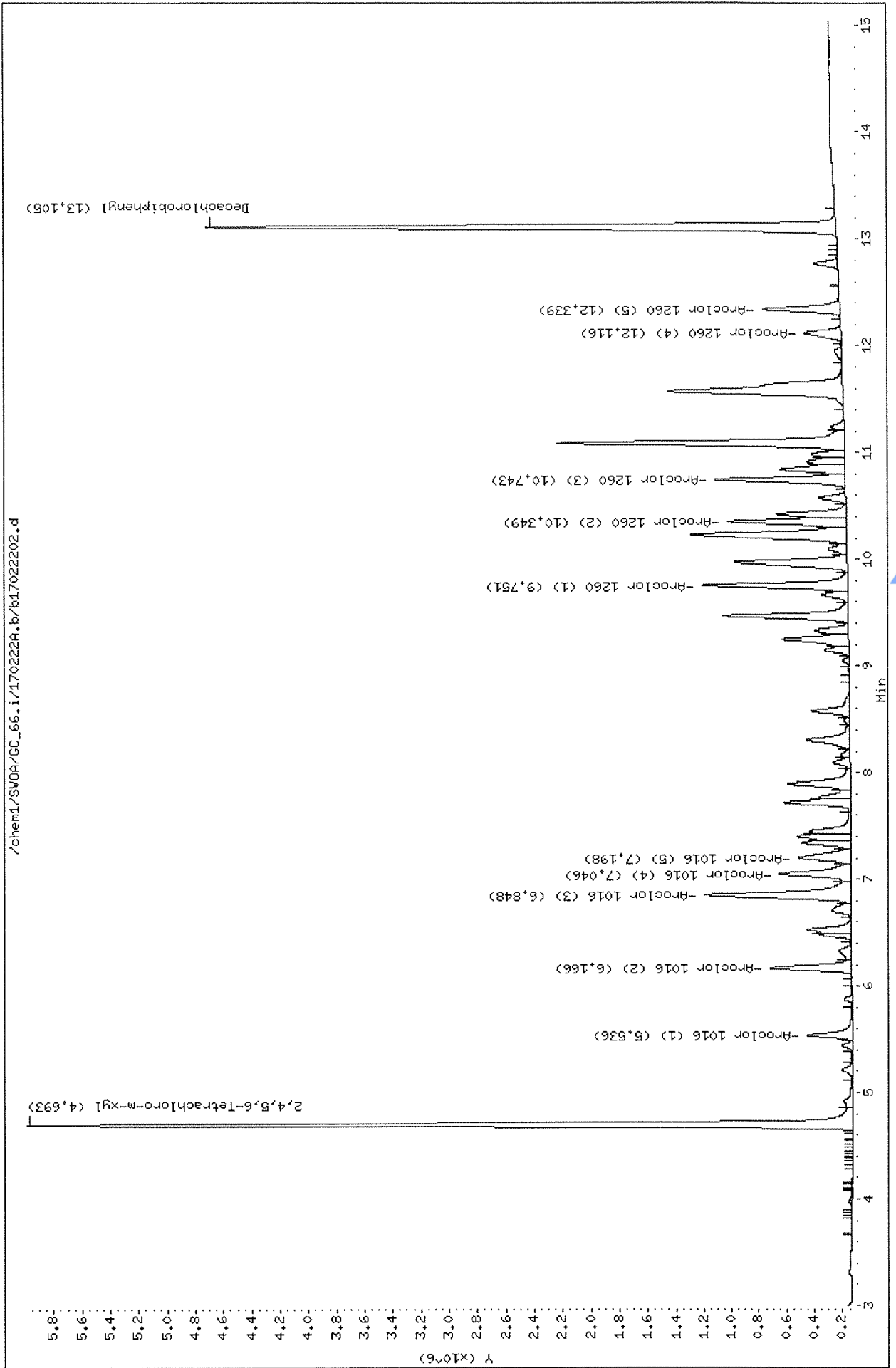
Sample Info: PCB ICAL2 P021517E 250PPB

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022203.d
Lab Smp Id:
Inj Date : 22-FEB-2017 18:03
Operator : 944 Inst ID: GC_66.i
Smp Info : PCB ICAL3 P021517D 500PPB
Misc Info :
Comment : Rtx-CLPesticide II
Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
Cal Date : 27-JAN-2017 13:18 Cal File: b17012710.d
Als bottle: 3 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: HP Genie Compound Sublist: p1016_1260.sub
Target Version: 3.50
Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
=====	==	=====	=====	=====	=====	=====
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.692	4.693	-0.001	583670480	100.000	102
M 2 Aroclor-1016				441364860	500.000	494
3 Aroclor 1016 (1)	5.533	5.536	-0.003	44444648	500.000	494
4 Aroclor 1016 (2)	6.164	6.166	-0.002	78615211	500.000	493
5 Aroclor 1016 (3)	6.844	6.848	-0.004	176646127	500.000	498
6 Aroclor 1016 (4)	7.043	7.046	-0.003	79269778	500.000	489
7 Aroclor 1016 (5)	7.195	7.198	-0.003	62389096	500.000	492
M 8 Aroclor-1260				469957779	500.000	496
9 Aroclor 1260 (1)	9.749	9.751	-0.002	133449409	500.000	494
10 Aroclor 1260 (2)	10.348	10.349	-0.001	105630017	500.000	494
11 Aroclor 1260 (3)	10.742	10.743	-0.001	116161904	500.000	498
12 Aroclor 1260 (4)	12.115	12.116	-0.001	40559632	500.000	488
13 Aroclor 1260 (5)	12.337	12.339	-0.002	74156817	500.000	503
\$ 56 Decachlorobiphenyl	13.105	13.105	0.000	553933911	100.000	101

Data File: /chem1/SW04/GC_66.i/170222A.b/b17022203.d

Date : 22-FEB-2017 18:03

Client ID:

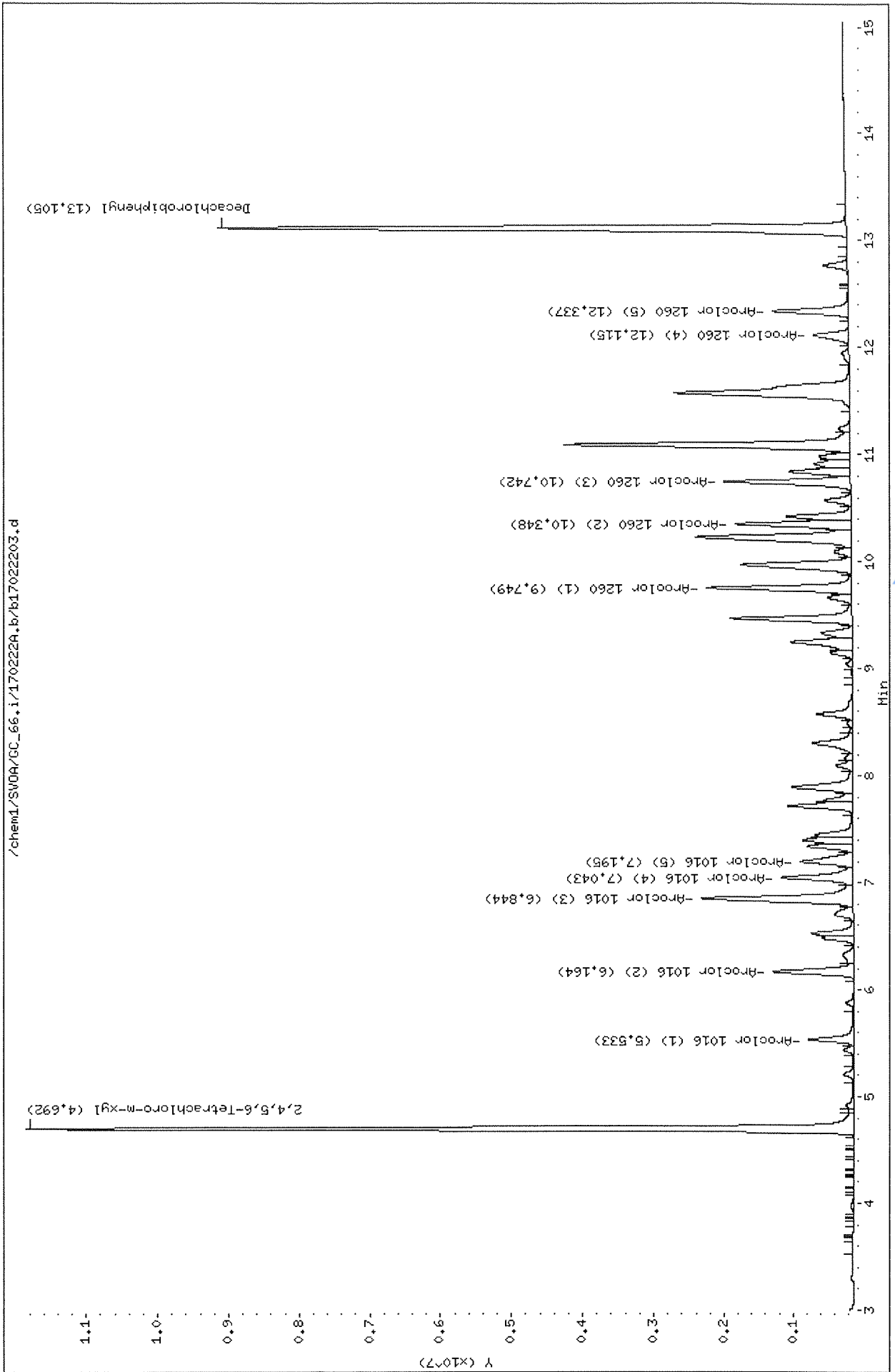
Sample Info: PCB ICAL3 P021517D 500PPB

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022204.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 18:21
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB ICAL4 P021517C 750PPB
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 22-FEB-2017 18:21 Cal File: b17022204.d
 Als bottle: 4 Calibration Sample, Level: 4
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.691	4.692	-0.001	881880457	150.000	154
M 2 Aroclor-1016				648645686	750.000	726
3 Aroclor 1016 (1)	5.530	5.533	-0.003	64426684	750.000	716
4 Aroclor 1016 (2)	6.163	6.164	-0.001	114714592	750.000	719
5 Aroclor 1016 (3)	6.842	6.844	-0.002	263194053	750.000	742
6 Aroclor 1016 (4)	7.041	7.043	-0.002	115323393	750.000	711
7 Aroclor 1016 (5)	7.192	7.195	-0.003	90986964	750.000	718
M 8 Aroclor-1260				685721676	750.000	724
9 Aroclor 1260 (1)	9.748	9.749	-0.001	194025997	750.000	718
10 Aroclor 1260 (2)	10.348	10.348	0.000	153042314	750.000	715
11 Aroclor 1260 (3)	10.742	10.742	0.000	168891068	750.000	725
12 Aroclor 1260 (4)	12.113	12.115	-0.002	63938532	750.000	769
13 Aroclor 1260 (5)	12.336	12.337	-0.001	105823765	750.000	718
\$ 56 Decachlorobiphenyl	13.104	13.105	-0.001	802792169	150.000	147



Data File: /chem1/SV00A/GC_66.i/170222A.b/b17022204.d

Date : 22-FEB-2017 18:21

Client ID:

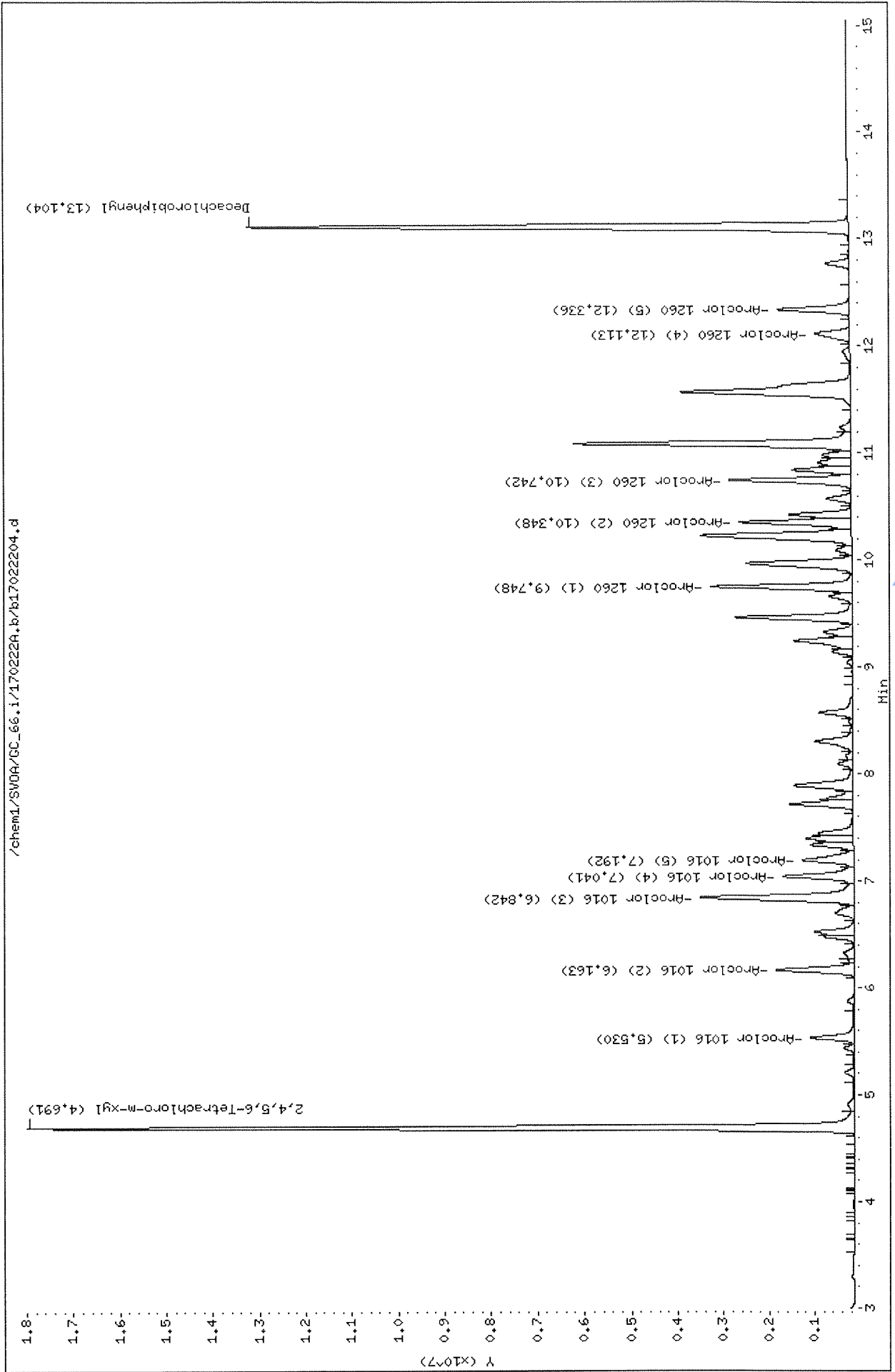
Sample Info: PCB ICAL4 P021517C 750PPB

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022205.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 18:38
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB ICAL5 P021517B 2000PPB
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 22-FEB-2017 18:38 Cal File: b17022205.d
 Als bottle: 5 Calibration Sample, Level: 5
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.693	4.691	0.002	2300900796	400.000	401(A)
M 2 Aroclor-1016				1605199523	2000.00	1800
3 Aroclor 1016 (1)	5.530	5.530	0.000	153733649	2000.00	1710
4 Aroclor 1016 (2)	6.163	6.163	0.000	279609263	2000.00	1750
5 Aroclor 1016 (3)	6.839	6.842	-0.003	669744171	2000.00	1890
6 Aroclor 1016 (4)	7.041	7.041	0.000	279504681	2000.00	1720
7 Aroclor 1016 (5)	7.190	7.192	-0.002	222607759	2000.00	1760
M 8 Aroclor-1260				1752866795	2000.00	1850
9 Aroclor 1260 (1)	9.747	9.748	-0.001	490934501	2000.00	1820
10 Aroclor 1260 (2)	10.347	10.348	-0.001	389125144	2000.00	1820
11 Aroclor 1260 (3)	10.742	10.742	0.000	439151601	2000.00	1880
12 Aroclor 1260 (4)	12.114	12.113	0.001	157303133	2000.00	1890
13 Aroclor 1260 (5)	12.336	12.336	0.000	276352416	2000.00	1880
\$ 56 Decachlorobiphenyl	13.104	13.104	0.000	2168071770	400.000	397

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

Data File: /chem1/SV0A/GC_66.i/170222A.b/b17022205.d

Date : 22-FEB-2017 18:38

Client ID:

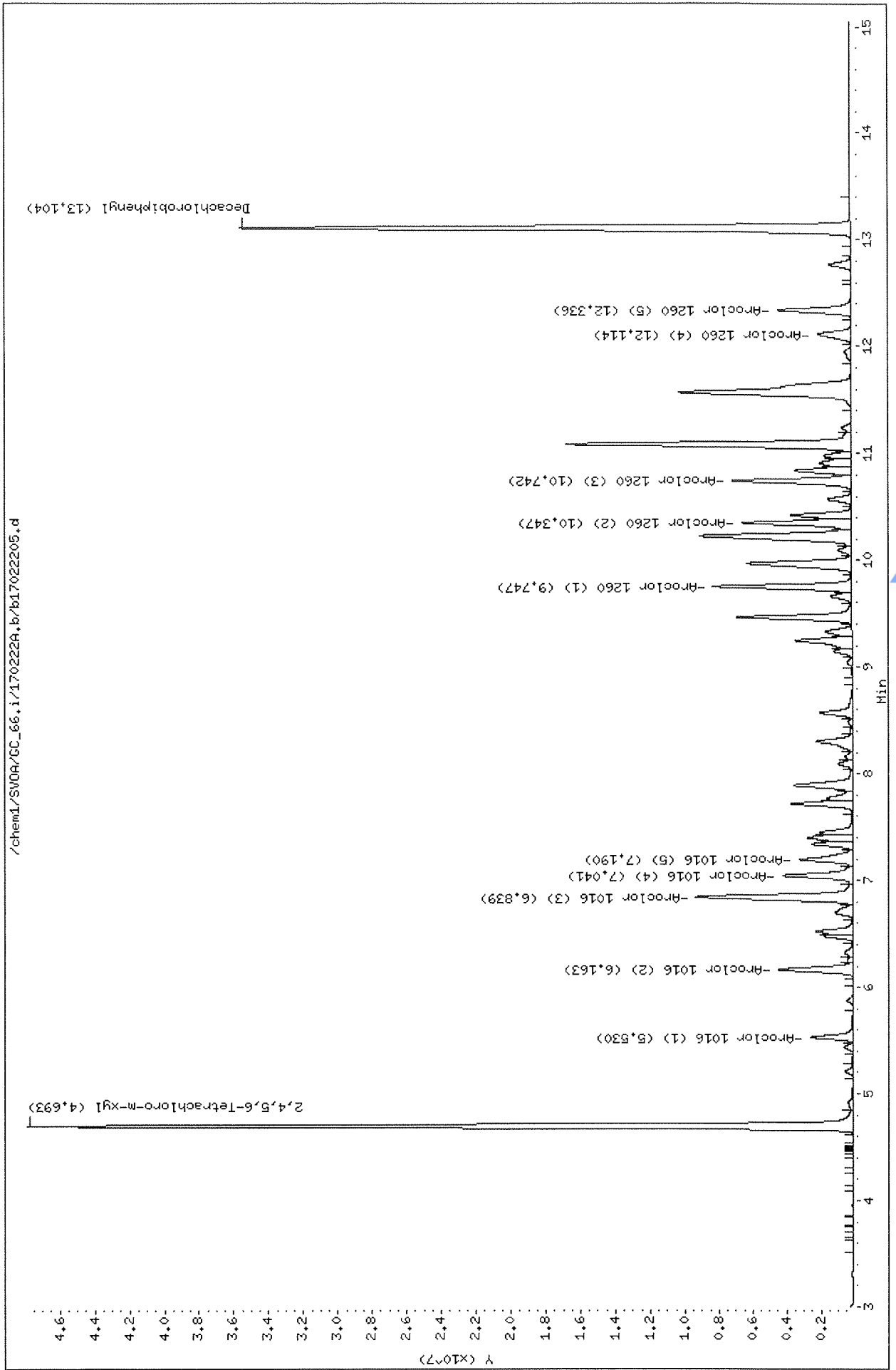
Sample Info: PCB ICALS P021517B 2000FPB

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022206.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 18:56
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB ICV P021517H 500PPB
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 22-FEB-2017 18:38 Cal File: b17022205.d
 Als bottle: 6 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.694	4.693	0.001	555489795	100.000	96.7
M 2 Aroclor-1016				503981199	500.000	564
3 Aroclor 1016 (1)	5.533	5.530	0.003	49949391	500.000	555
4 Aroclor 1016 (2)	6.165	6.163	0.002	88603073	500.000	556
5 Aroclor 1016 (3)	6.844	6.839	0.005	203482977	500.000	573
6 Aroclor 1016 (4)	7.043	7.041	0.002	90577001	500.000	559
7 Aroclor 1016 (5)	7.194	7.190	0.004	71368757	500.000	563
M 8 Aroclor-1260				539424564	500.000	569
9 Aroclor 1260 (1)	9.749	9.747	0.002	159119691	500.000	589
10 Aroclor 1260 (2)	10.348	10.347	0.001	118984171	500.000	556
11 Aroclor 1260 (3)	10.742	10.742	0.000	133717652	500.000	574
12 Aroclor 1260 (4)	12.115	12.114	0.001	41779928	500.000	503
13 Aroclor 1260 (5)	12.337	12.336	0.001	85823122	500.000	582
\$ 56 Decachlorobiphenyl	13.103	13.104	-0.001	550584972	100.000	101



Data File: /chem1/SV00A/GC_66.i/170222A,b/b17022206.d

Date : 22-FEB-2017 18:56

Client ID:

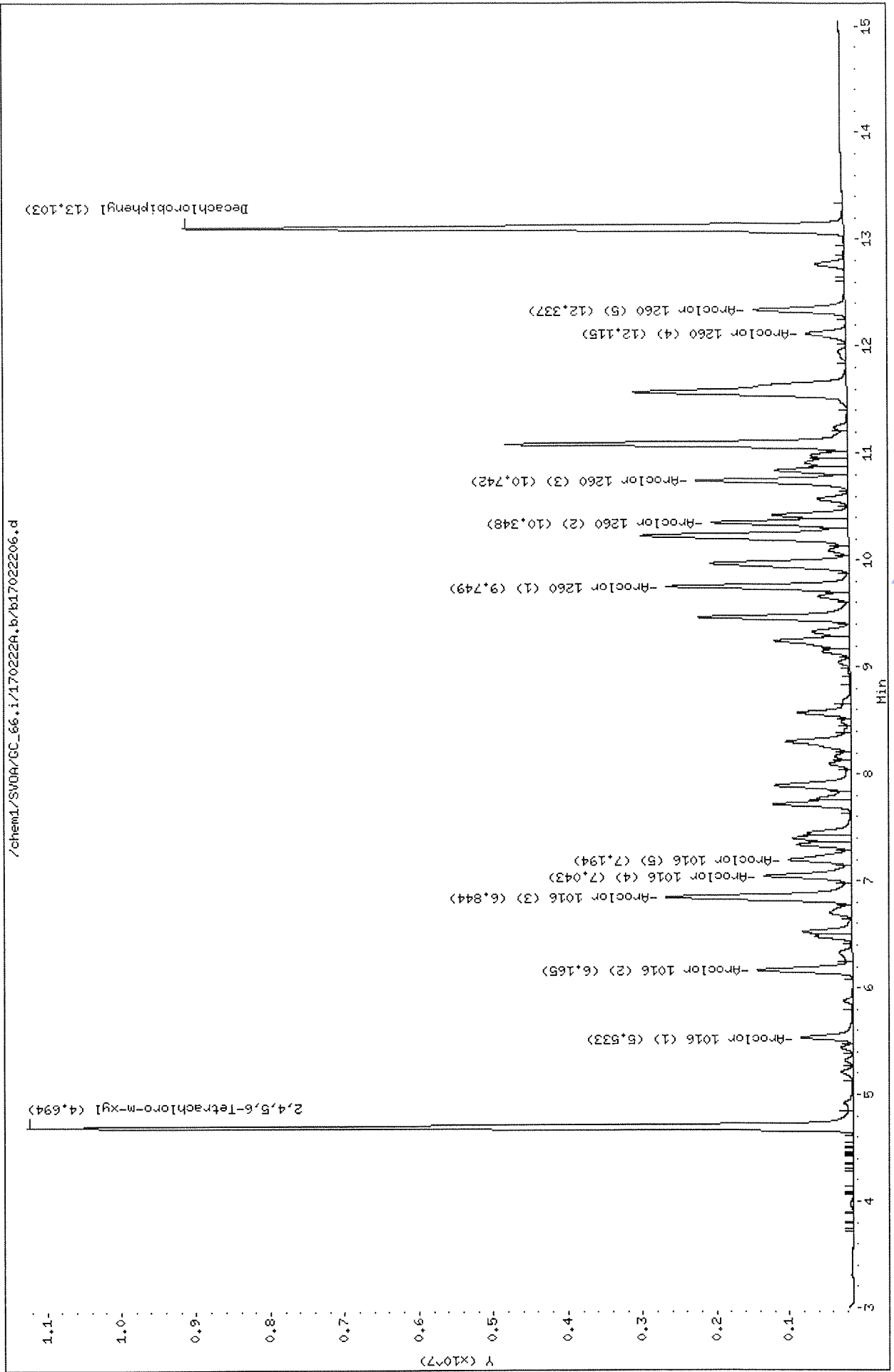
Sample Info: PCB ICV P021517H 500PPB

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022207.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 19:14
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB 1221/54 500PPB P120616I
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 27-JAN-2017 13:18 Cal File: b17012710.d
 Als bottle: 7 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1221_1254.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
M 14 Aroclor-1221				146604462	500.000	500
15 Aroclor 1221 (1)	3.949	3.949	0.000	21574609	500.000	500
16 Aroclor 1221 (2)	5.209	5.209	0.000	27788962	500.000	500
17 Aroclor 1221 (3)	5.436	5.436	0.000	17435037	500.000	500
18 Aroclor 1221 (4)	5.532	5.532	0.000	69267403	500.000	500
19 Aroclor 1221 (5)	6.166	6.166	0.000	10538451	500.000	500
M 38 Aroclor-1254				659429043	500.000	500
39 Aroclor 1254 (1)	8.300	8.300	0.000	132689273	500.000	500
40 Aroclor 1254 (2)	8.658	8.658	0.000	93074176	500.000	500
41 Arcolor 1254 (3)	9.199	9.199	0.000	175076127	500.000	500
42 Aroclor 1254 (4)	9.513	9.513	0.000	119868437	500.000	500
43 Aroclor 1254 (5)	9.976	9.976	0.000	138721030	500.000	500



Data File: /chem1/SV04/GC_66.i/170222A.b/b17022207.d

Date : 22-FEB-2017 19:14

Client ID:

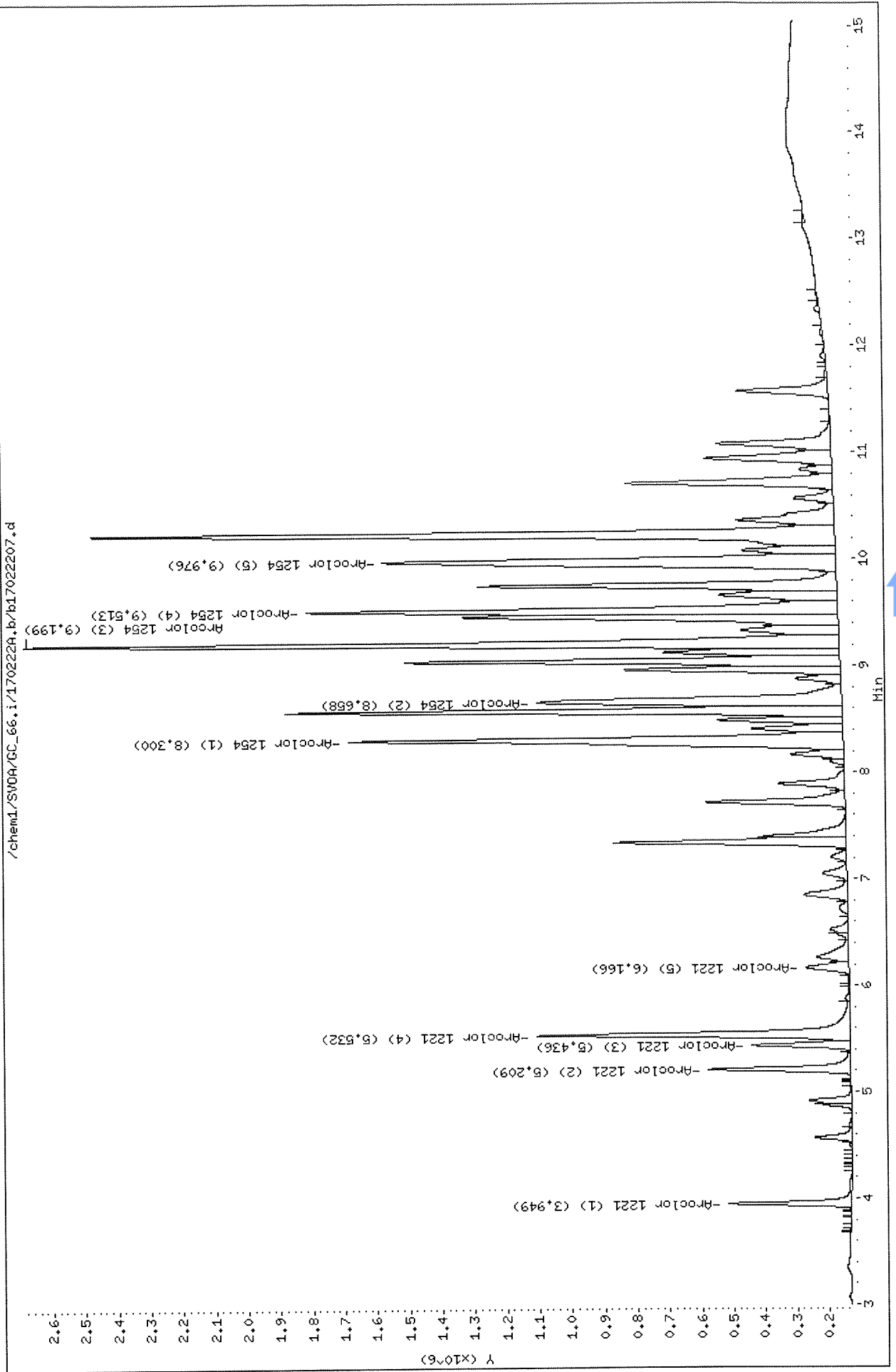
Sample Info: PCB 1221/54 500PPB P1206161

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022208.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 19:32
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB 1232/62 500PPB P120616J
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 27-JAN-2017 13:18 Cal File: b17012710.d
 Als bottle: 8 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1232_1262.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
M 20 Aroclor-1232				230600593	500.000	500
21 Aroclor 1232 (1)	5.533	5.533	0.000	46830292	500.000	500
22 Aroclor 1232 (2)	6.164	6.164	0.000	37990961	500.000	500
23 Aroclor 1232 (3)	6.846	6.846	0.000	80230398	500.000	500
24 Aroclor 1232 (4)	7.045	7.045	0.000	37042341	500.000	500
25 Aroclor 1232 (5)	7.198	7.198	0.000	28506601	500.000	500
M 44 Aroclor-1262				529581884	500.000	500
45 Aroclor 1262 (1)	9.749	9.749	0.000	99779332	500.000	500
46 Aroclor 1262 (2)	10.347	10.347	0.000	142884711	500.000	500
47 Aroclor 1262 (3)	10.742	10.742	0.000	132313903	500.000	500
48 Aroclor 1262 (4)	12.118	12.118	0.000	44885518	500.000	500
49 Aroclor 1262 (5)	12.337	12.337	0.000	109718420	500.000	500

Data File: /chem1/SV0A/GC_66.i/170222A.b/b17022208.d

Date : 22-FEB-2017 19:32

Client ID:

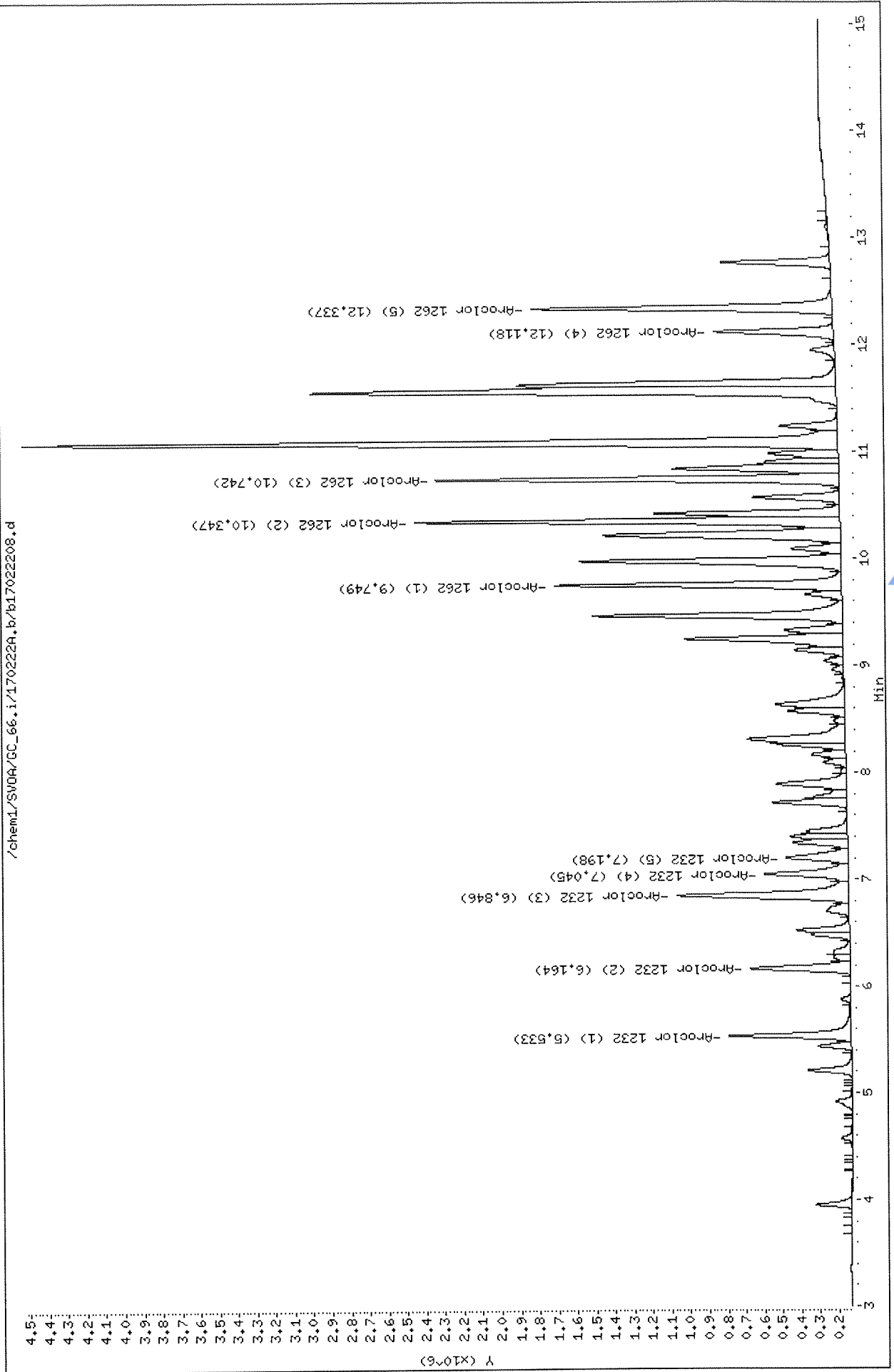
Sample Info: PCB 1232/62 500PPB F120616J

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022209.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 19:50
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB 1248/68 500PPB P120616K
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 27-JAN-2017 13:18 Cal File: b17012710.d
 Als bottle: 9 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1248_1268.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
M 32 Aroclor-1248				349356650	500.000	500
33 Aroclor 1248 (1)	6.845	6.164	0.681	92597659	500.000	500
34 Arcolor 1248 (2)	7.397	7.397	0.000	51572554	500.000	500
35 Aroclor 1248 (3)	7.445	7.445	0.000	38405884	500.000	500
36 Aroclor 1248 (4)	7.715	7.715	0.000	76011522	500.000	500
37 Aroclor 1248 (5)	7.888	7.888	0.000	90769029	500.000	500
M 50 Aroclor-1268				2132325871	500.000	500
51 Aroclor 1268 (1)	11.554	11.554	0.000	353673525	500.000	500
52 Aroclor 1268 (2)	11.615	11.615	0.000	354519915	500.000	500
53 Aroclor 1268 (3)	11.948	11.948	0.000	290210573	500.000	500
54 Aroclor 1268 (4)	12.334	12.334	0.000	127638418	500.000	500
55 Aroclor 1268 (5)	12.764	12.764	0.000	1006283437	500.000	500

Data File: /chem1/SV00A/GC_66.i/170222A.b/b17022209.d

Date : 22-FEB-2017 19:50

Client ID:

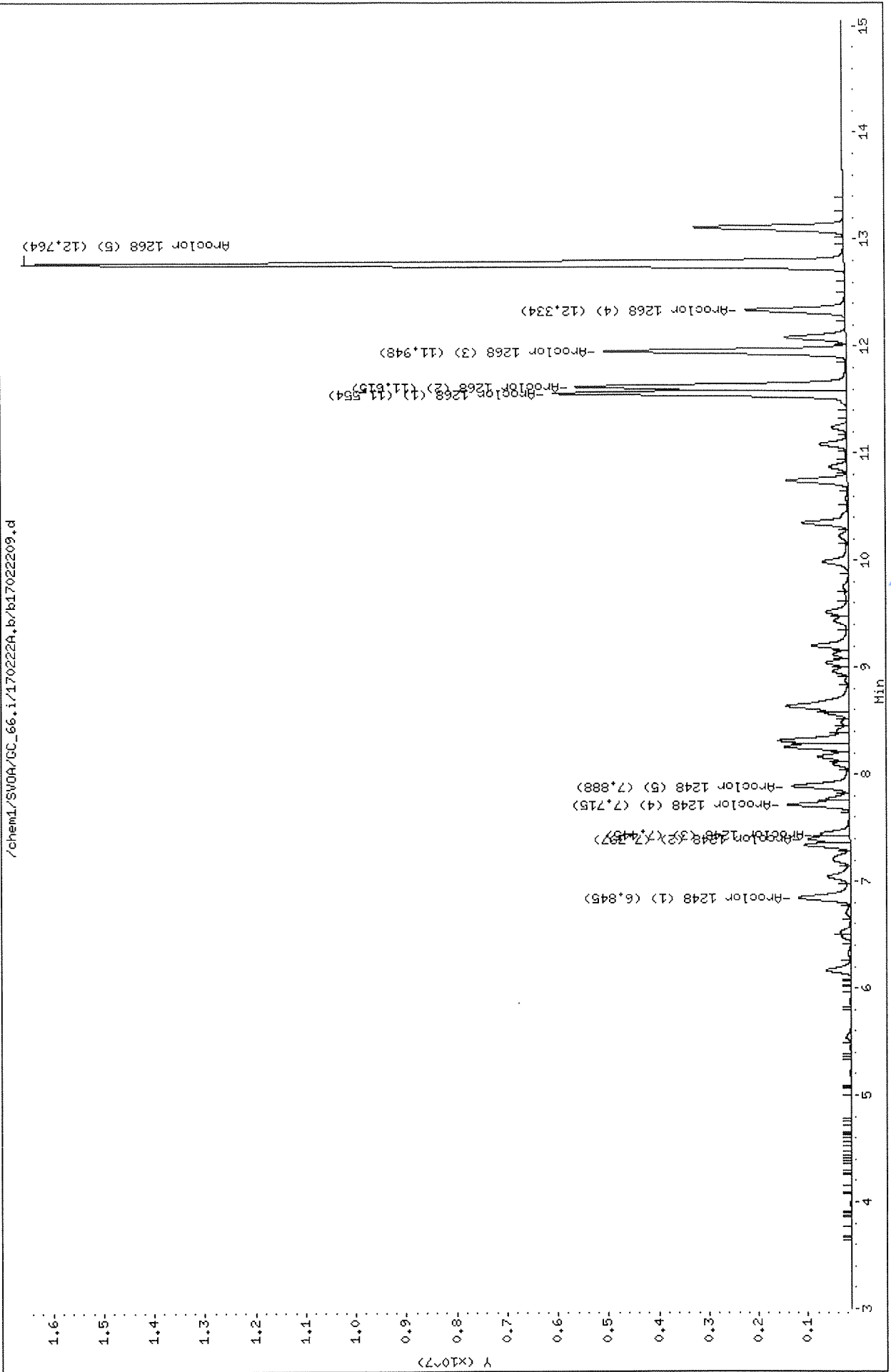
Sample Info: PCB 1248/68 500PPB P120616K

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170222A.b/b17022210.d
 Lab Smp Id:
 Inj Date : 22-FEB-2017 20:07
 Operator : 944 Inst ID: GC_66.i
 Smp Info : PCB 1242 500PPB P120616L
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170222A.b/b8082d-n2.m
 Meth Date : 23-Feb-2017 09:41 uj3k Quant Type: ESTD
 Cal Date : 27-JAN-2017 13:18 Cal File: b17012710.d
 Als bottle: 10 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1242-ical.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
M 26 Aroclor-1242				360695998	500.000	500
27 Aroclor 1242 (1)	5.534	5.534	0.000	35639904	500.000	500
28 Aroclor 1242 (2)	6.164	6.164	0.000	64753688	500.000	500
29 Aroclor 1242 (3)	6.844	6.844	0.000	143503184	500.000	500
30 Aroclor 1242 (4)	7.044	7.044	0.000	65451422	500.000	500
31 Aroclor 1242 (5)	7.195	7.195	0.000	51347800	500.000	500



Data File: /chem1/SWDH/GC_66.i/170222A.b/b17022210.d

Date : 22-FEB-2017 20:07

Client ID:

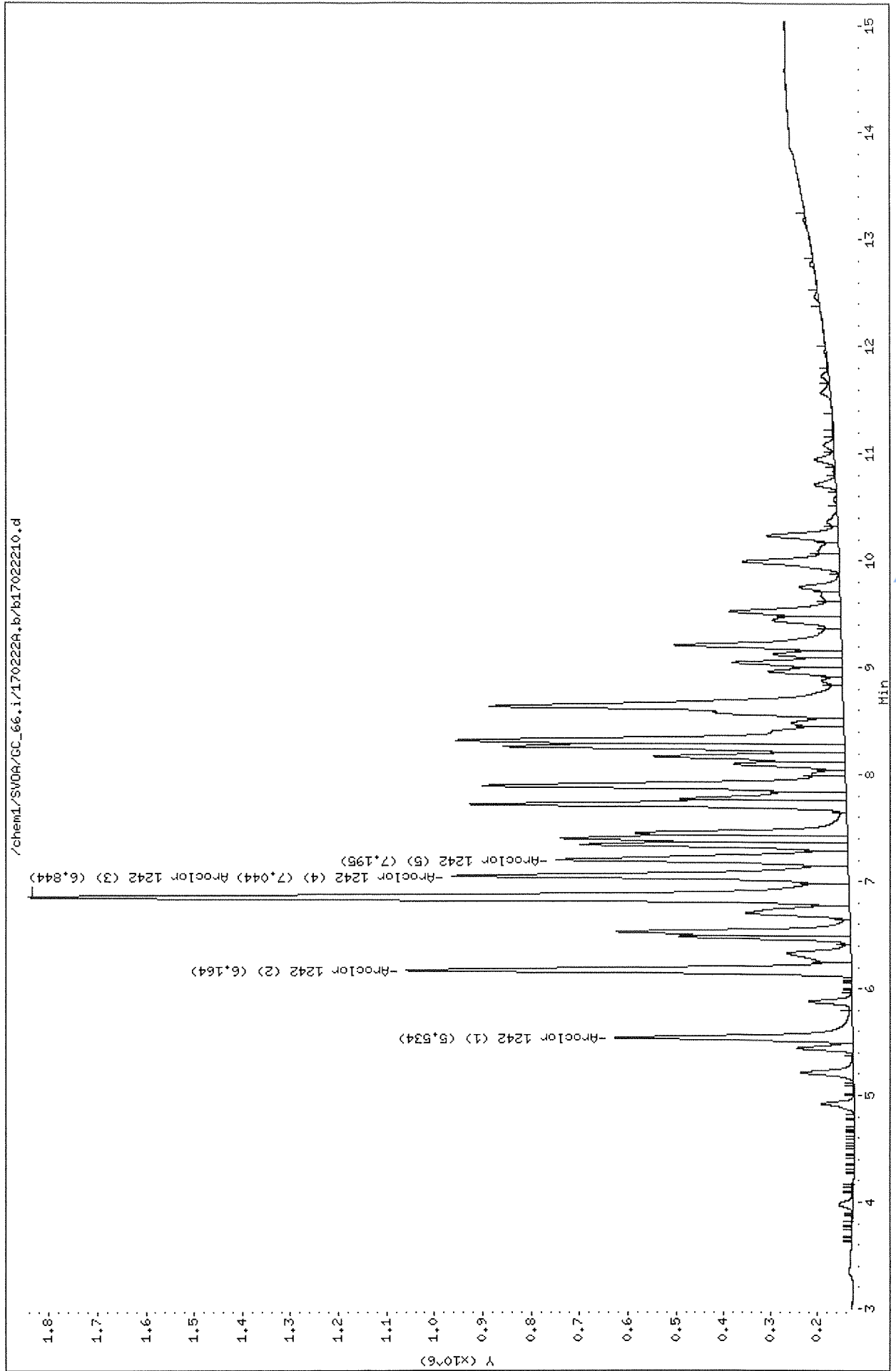
Sample Info: PCB 1242 500PPB P120616L

Instrument: GC_66.i

Operator: 944

Column diameter: 2.00

Column phase:



EPA METHOD 8082 PCB

Sample Data

RAW DATA SHEET
FOR METHOD: EPA 8082

WORK ORDER: 17-03-1557
INSTRUMENT: GC 66
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED: 2017-03-28 03:41
REVIEWED BY:
D/T REVIEWED: *M*

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282917032829

28 CLIENT SAMPLE NUMBER: B-DU1-ISM1-8

LCS/MB BATCH: 170327L05 SAMPLE VOLUME / WEIGHT: DEFAULT: 20.00 g / ACTUAL: 20.00 g
MS/MSD BATCH: 170327S05 FINAL VOLUME / WEIGHT: DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: ug/kg ADJUSTMENT RATIO TO PF: 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Aroclor-1016	0.000	1.00	ND	50	
Aroclor-1221	0.000	1.00	ND	50	
Aroclor-1232	0.000	1.00	ND	50	
Aroclor-1242	0.000	1.00	ND	50	
Aroclor-1248	0.000	1.00	ND	50	
Aroclor-1254	0.000	1.00	ND	50	
Aroclor-1260	0.000	1.00	ND	50	
Aroclor-1262	0.000	1.00	ND	50	
Aroclor-1268	0.000	1.00	ND	50	

Return to Contents

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032829.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032829.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 03:41
 Operator : 669 Inst ID: GC_66.i
 Smp Info : 17-03-1557-28
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 29
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000		384884629	64.8393	64.8
M 2 Aroclor-1016					Compound Not Detected.		
3 Aroclor 1016 (1)					Compound Not Detected.		
4 Aroclor 1016 (2)					Compound Not Detected.		
5 Aroclor 1016 (3)					Compound Not Detected.		
6 Aroclor 1016 (4)					Compound Not Detected.		
7 Aroclor 1016 (5)					Compound Not Detected.		
M 8 Aroclor-1260					Compound Not Detected.		
9 Aroclor 1260 (1)					Compound Not Detected.		
10 Aroclor 1260 (2)					Compound Not Detected.		
11 Aroclor 1260 (3)					Compound Not Detected.		
12 Aroclor 1260 (4)					Compound Not Detected.		
13 Aroclor 1260 (5)					Compound Not Detected.		
M 14 Aroclor-1221					Compound Not Detected.		
15 Aroclor 1221 (1)					Compound Not Detected.		
16 Aroclor 1221 (2)					Compound Not Detected.		
17 Aroclor 1221 (3)					Compound Not Detected.		
18 Aroclor 1221 (4)					Compound Not Detected.		
19 Aroclor 1221 (5)					Compound Not Detected.		
M 20 Aroclor-1232					Compound Not Detected.		
21 Aroclor 1232 (1)					Compound Not Detected.		

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032829.d
 Report Date: 28-Mar-2017 09:42

Page 2

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ppb)	FINAL (ug/Kg)
=====	==	=====	=====	=====	=====	=====	=====	=====
22 Aroclor 1232 (2)						Compound Not Detected.		
23 Aroclor 1232 (3)						Compound Not Detected.		
24 Aroclor 1232 (4)						Compound Not Detected.		
25 Aroclor 1232 (5)						Compound Not Detected.		
M 26 Aroclor-1242						Compound Not Detected.		
27 Aroclor 1242 (1)						Compound Not Detected.		
28 Aroclor 1242 (2)						Compound Not Detected.		
29 Aroclor 1242 (3)						Compound Not Detected.		
30 Aroclor 1242 (4)						Compound Not Detected.		
31 Aroclor 1242 (5)						Compound Not Detected.		
M 32 Aroclor-1248						Compound Not Detected.		
33 Aroclor 1248 (1)						Compound Not Detected.		
34 Aroclor 1248 (2)						Compound Not Detected.		
35 Aroclor 1248 (3)						Compound Not Detected.		
36 Aroclor 1248 (4)						Compound Not Detected.		
37 Aroclor 1248 (5)						Compound Not Detected.		
M 38 Aroclor-1254						Compound Not Detected.		
39 Aroclor 1254 (1)						Compound Not Detected.		
40 Aroclor 1254 (2)						Compound Not Detected.		
41 Aroclor 1254 (3)						Compound Not Detected.		
42 Aroclor 1254 (4)						Compound Not Detected.		
43 Aroclor 1254 (5)						Compound Not Detected.		
M 44 Aroclor-1262						Compound Not Detected.		
45 Aroclor 1262 (1)						Compound Not Detected.		
46 Aroclor 1262 (2)						Compound Not Detected.		
47 Aroclor 1262 (3)						Compound Not Detected.		
48 Aroclor 1262 (4)						Compound Not Detected.		
49 Aroclor 1262 (5)						Compound Not Detected.		
M 50 Aroclor-1268						Compound Not Detected.		
51 Aroclor 1268 (1)						Compound Not Detected.		
52 Aroclor 1268 (2)						Compound Not Detected.		
53 Aroclor 1268 (3)						Compound Not Detected.		
54 Aroclor 1268 (4)						Compound Not Detected.		
55 Aroclor 1268 (5)						Compound Not Detected.		
§ 56 Decachlorobiphenyl	13.098	13.098	0.000		422410796		76.7050	76.7

Data File: /chem1/SV04/GC_66.i/170328.b/b17032829.d

Date : 28-MAR-2017 03:41

Client ID:

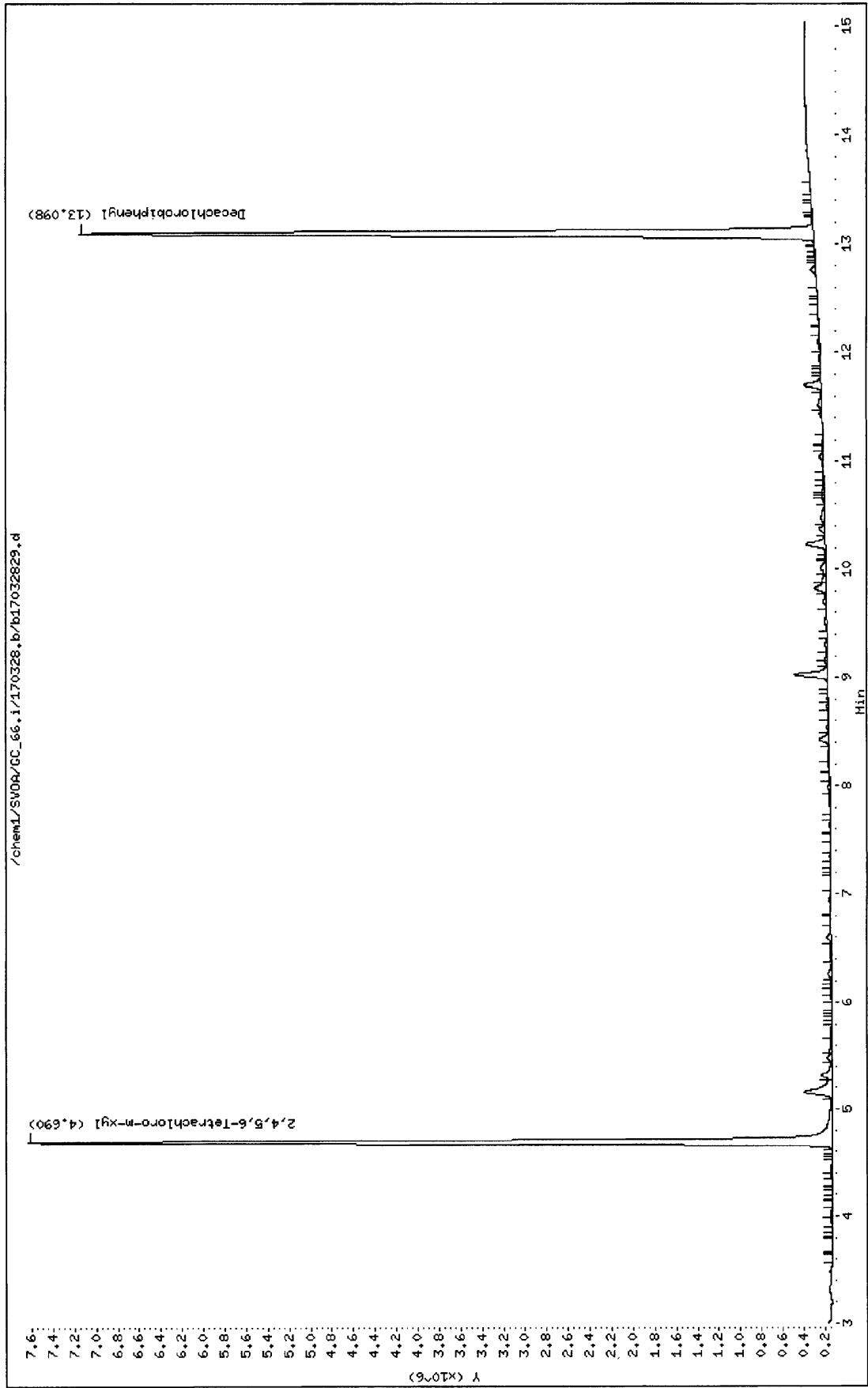
Sample Info: 17-03-1557-28

Instrument: GC_66.i

Operator: 669


Column diameter: 2.00

Column phase:



RAW DATA SHEET
FOR METHOD: EPA 8082

WORK ORDER: 17-03-1557
INSTRUMENT: GC 66
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED: 2017-03-28 03:59
REVIEWED BY: 
D/T REVIEWED:

DATA FILE: /chem1/SVOA/GC_66/170328/b1703283017032830

29 CLIENT SAMPLE NUMBER: B-DU1-ISM2-8

LCS/MB BATCH: 170327L05 SAMPLE VOLUME / WEIGHT: DEFAULT: 20.00 g / ACTUAL: 20.20 g
MS/MSD BATCH: 170327S05 FINAL VOLUME / WEIGHT: DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: ug/kg ADJUSTMENT RATIO TO PF: 0.99

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Aroclor-1016	0.000	1.00	ND	50	
Aroclor-1221	0.000	1.00	ND	50	
Aroclor-1232	0.000	1.00	ND	50	
Aroclor-1242	0.000	1.00	ND	50	
Aroclor-1248	0.000	1.00	ND	50	
Aroclor-1254	0.000	1.00	ND	50	
Aroclor-1260	0.000	1.00	ND	50	
Aroclor-1262	0.000	1.00	ND	50	
Aroclor-1268	0.000	1.00	ND	50	


Return to Contents

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032830.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032830.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 03:59
 Operator : 669 Inst ID: GC_66.i
 Smp Info : 17-03-1557-29
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 30
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.691	4.690	0.001	362632594	61.0906	61.1
M 2 Aroclor-1016				Compound Not Detected.		
3 Aroclor 1016 (1)				Compound Not Detected.		
4 Aroclor 1016 (2)				Compound Not Detected.		
5 Aroclor 1016 (3)				Compound Not Detected.		
6 Aroclor 1016 (4)				Compound Not Detected.		
7 Aroclor 1016 (5)				Compound Not Detected.		
M 8 Aroclor-1260				Compound Not Detected.		
9 Aroclor 1260 (1)				Compound Not Detected.		
10 Aroclor 1260 (2)				Compound Not Detected.		
11 Aroclor 1260 (3)				Compound Not Detected.		
12 Aroclor 1260 (4)				Compound Not Detected.		
13 Aroclor 1260 (5)				Compound Not Detected.		
M 14 Aroclor-1221				Compound Not Detected.		
15 Aroclor 1221 (1)				Compound Not Detected.		
16 Aroclor 1221 (2)				Compound Not Detected.		
17 Aroclor 1221 (3)				Compound Not Detected.		
18 Aroclor 1221 (4)				Compound Not Detected.		
19 Aroclor 1221 (5)				Compound Not Detected.		
M 20 Aroclor-1232				Compound Not Detected.		
21 Aroclor 1232 (1)				Compound Not Detected.		

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032830.d
 Report Date: 28-Mar-2017 09:42

Page 2

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ppb)	FINAL (ug/Kg)
=====	==	=====	=====	=====	=====	=====	=====	=====
22 Aroclor 1232 (2)						Compound Not Detected.		
23 Aroclor 1232 (3)						Compound Not Detected.		
24 Aroclor 1232 (4)						Compound Not Detected.		
25 Aroclor 1232 (5)						Compound Not Detected.		
M 26 Aroclor-1242						Compound Not Detected.		
27 Aroclor 1242 (1)						Compound Not Detected.		
28 Aroclor 1242 (2)						Compound Not Detected.		
29 Aroclor 1242 (3)						Compound Not Detected.		
30 Aroclor 1242 (4)						Compound Not Detected.		
31 Aroclor 1242 (5)						Compound Not Detected.		
M 32 Aroclor-1248						Compound Not Detected.		
33 Aroclor 1248 (1)						Compound Not Detected.		
34 Aroclor 1248 (2)						Compound Not Detected.		
35 Aroclor 1248 (3)						Compound Not Detected.		
36 Aroclor 1248 (4)						Compound Not Detected.		
37 Aroclor 1248 (5)						Compound Not Detected.		
M 38 Aroclor-1254						Compound Not Detected.		
39 Aroclor 1254 (1)						Compound Not Detected.		
40 Aroclor 1254 (2)						Compound Not Detected.		
41 Aroclor 1254 (3)						Compound Not Detected.		
42 Aroclor 1254 (4)						Compound Not Detected.		
43 Aroclor 1254 (5)						Compound Not Detected.		
M 44 Aroclor-1262						Compound Not Detected.		
45 Aroclor 1262 (1)						Compound Not Detected.		
46 Aroclor 1262 (2)						Compound Not Detected.		
47 Aroclor 1262 (3)						Compound Not Detected.		
48 Aroclor 1262 (4)						Compound Not Detected.		
49 Aroclor 1262 (5)						Compound Not Detected.		
M 50 Aroclor-1268						Compound Not Detected.		
51 Aroclor 1268 (1)						Compound Not Detected.		
52 Aroclor 1268 (2)						Compound Not Detected.		
53 Aroclor 1268 (3)						Compound Not Detected.		
54 Aroclor 1268 (4)						Compound Not Detected.		
55 Aroclor 1268 (5)						Compound Not Detected.		
§ 56 Decachlorobiphenyl	13.097	13.098	-0.001		403526302	73.2758	73.3	

Data File: /chem1/SV04/CC_66.i/170328.b/b17032830.d

Date: 28-MAR-2017 03:59

Client ID:

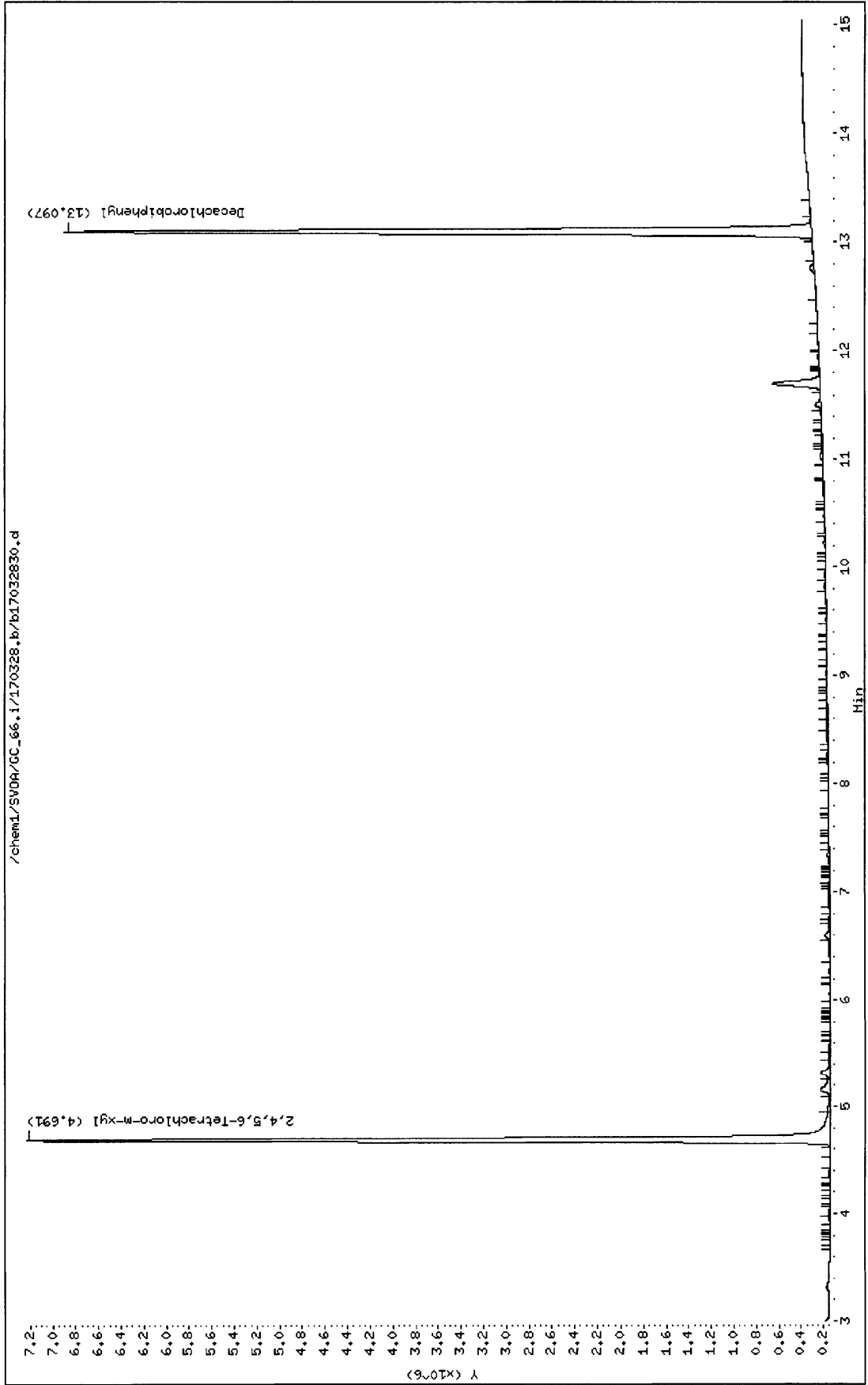
Sample Info: 17-03-1557-29

Instrument: GC_66.i

Operator: 669


Column diameter: 2.00

Column phase:



**RAW DATA SHEET
FOR METHOD: EPA 8082**

WORK ORDER: 17-03-1557
INSTRUMENT: GC 66
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED: 2017-03-28 04:35
REVIEWED BY: 
D/T REVIEWED:

DATA FILE: /chem1/SVOA/GC_66/170328/b1703283217032832

30 **CLIENT SAMPLE NUMBER: B-DU1-ISM3-8**

LCS/MB BATCH: 170327L05 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 20.00 g / ACTUAL: 20.10 g
MS/MSD BATCH: 170327S05 **FINAL VOLUME / WEIGHT:** DEFAULT: 10.00 ml / ACTUAL: 10.00 ml
UNITS: ug/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Aroclor-1016	0.000	1.00	ND	50	
Aroclor-1221	0.000	1.00	ND	50	
Aroclor-1232	0.000	1.00	ND	50	
Aroclor-1242	0.000	1.00	ND	50	
Aroclor-1248	0.000	1.00	ND	50	
Aroclor-1254	0.000	1.00	ND	50	
Aroclor-1260	0.000	1.00	ND	50	
Aroclor-1262	0.000	1.00	ND	50	
Aroclor-1268	0.000	1.00	ND	50	


Return to Contents

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032832.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032832.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 04:35
 Operator : 669 Inst ID: GC_66.i
 Smp Info : 17-03-1557-30
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 32
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000	414146348	69.7689	69.8
M 2 Aroclor-1016				Compound Not Detected.		
3 Aroclor 1016 (1)				Compound Not Detected.		
4 Aroclor 1016 (2)				Compound Not Detected.		
5 Aroclor 1016 (3)				Compound Not Detected.		
6 Aroclor 1016 (4)				Compound Not Detected.		
7 Aroclor 1016 (5)				Compound Not Detected.		
M 8 Aroclor-1260				Compound Not Detected.		
9 Aroclor 1260 (1)				Compound Not Detected.		
10 Aroclor 1260 (2)				Compound Not Detected.		
11 Aroclor 1260 (3)				Compound Not Detected.		
12 Aroclor 1260 (4)				Compound Not Detected.		
13 Aroclor 1260 (5)				Compound Not Detected.		
M 14 Aroclor-1221				Compound Not Detected.		
15 Aroclor 1221 (1)				Compound Not Detected.		
16 Aroclor 1221 (2)				Compound Not Detected.		
17 Aroclor 1221 (3)				Compound Not Detected.		
18 Aroclor 1221 (4)				Compound Not Detected.		
19 Aroclor 1221 (5)				Compound Not Detected.		
M 20 Aroclor-1232				Compound Not Detected.		
21 Aroclor 1232 (1)				Compound Not Detected.		

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032832.d
 Report Date: 28-Mar-2017 09:42

Page 2

Compounds	RT	EXP RT	DLT RT	RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ppb)	FINAL (ug/Kg)
=====	==	=====	=====	=====	=====	=====	=====
22 Aroclor 1232 (2)					Compound Not Detected.		
23 Aroclor 1232 (3)					Compound Not Detected.		
24 Aroclor 1232 (4)					Compound Not Detected.		
25 Aroclor 1232 (5)					Compound Not Detected.		
M 26 Aroclor-1242					Compound Not Detected.		
27 Aroclor 1242 (1)					Compound Not Detected.		
28 Aroclor 1242 (2)					Compound Not Detected.		
29 Aroclor 1242 (3)					Compound Not Detected.		
30 Aroclor 1242 (4)					Compound Not Detected.		
31 Aroclor 1242 (5)					Compound Not Detected.		
M 32 Aroclor-1248					Compound Not Detected.		
33 Aroclor 1248 (1)					Compound Not Detected.		
34 Aroclor 1248 (2)					Compound Not Detected.		
35 Aroclor 1248 (3)					Compound Not Detected.		
36 Aroclor 1248 (4)					Compound Not Detected.		
37 Aroclor 1248 (5)					Compound Not Detected.		
M 38 Aroclor-1254					Compound Not Detected.		
39 Aroclor 1254 (1)					Compound Not Detected.		
40 Aroclor 1254 (2)					Compound Not Detected.		
41 Aroclor 1254 (3)					Compound Not Detected.		
42 Aroclor 1254 (4)					Compound Not Detected.		
43 Aroclor 1254 (5)					Compound Not Detected.		
M 44 Aroclor-1262					Compound Not Detected.		
45 Aroclor 1262 (1)					Compound Not Detected.		
46 Aroclor 1262 (2)					Compound Not Detected.		
47 Aroclor 1262 (3)					Compound Not Detected.		
48 Aroclor 1262 (4)					Compound Not Detected.		
49 Aroclor 1262 (5)					Compound Not Detected.		
M 50 Aroclor-1268					Compound Not Detected.		
51 Aroclor 1268 (1)					Compound Not Detected.		
52 Aroclor 1268 (2)					Compound Not Detected.		
53 Aroclor 1268 (3)					Compound Not Detected.		
54 Aroclor 1268 (4)					Compound Not Detected.		
55 Aroclor 1268 (5)					Compound Not Detected.		
\$ 56 Decachlorobiphenyl	13.097	13.097	0.000		419388900	76.1563	76.2

Data File: /chem1/SV04/GC_66.i/170328.b/b17032832.d

Date: 28-Mar-2017 04:35

Client ID:

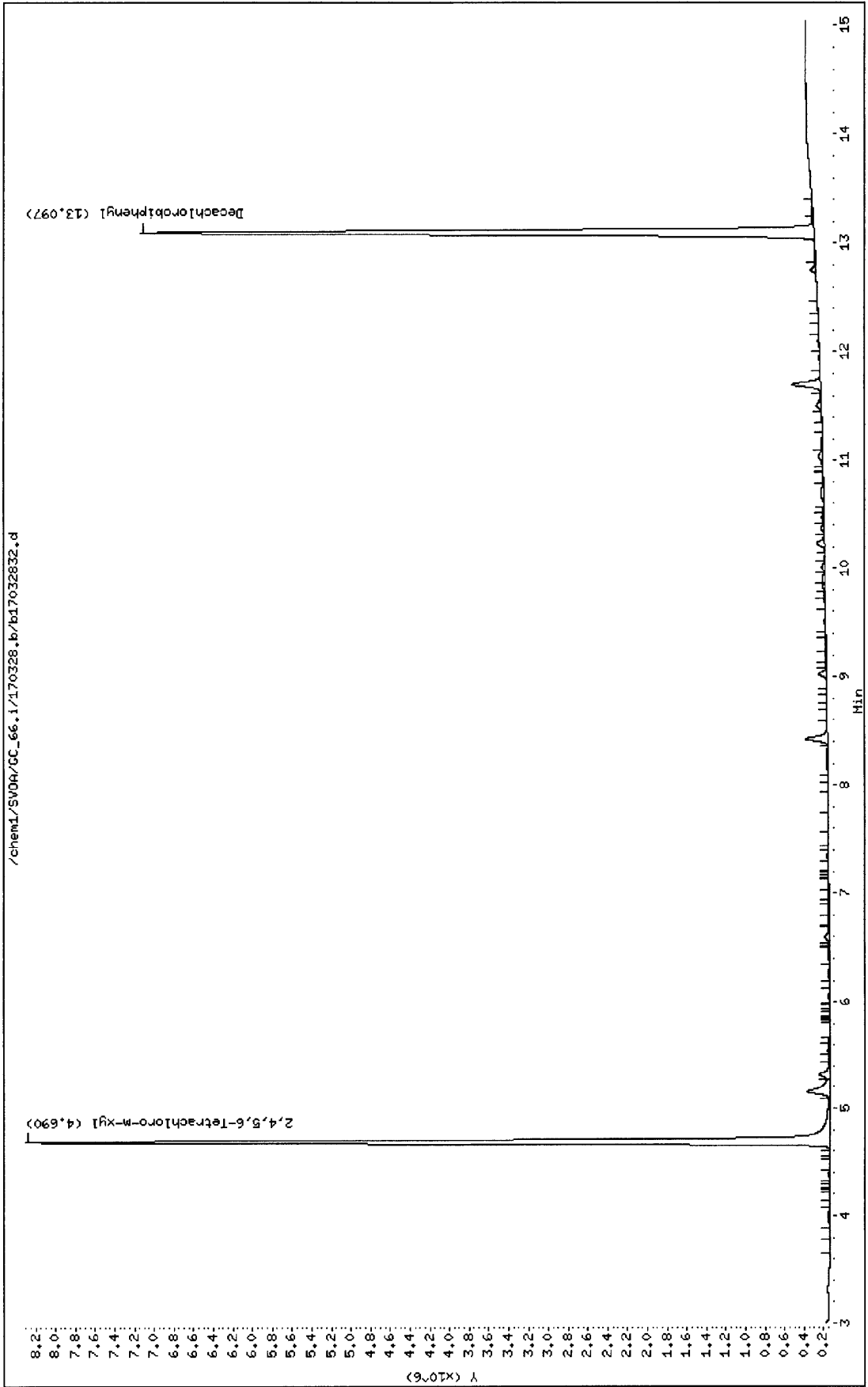
Sample Info: 17-03-1557-30

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phase:



EPA METHOD 8082 PCB

Quality Control

Method Blank LCS/LCSD MS/MSD

METHOD BLANK ASSOCIATION SUMMARY
FOR METHOD: EPA 8082

MB SAMPLE ID: 099-12-535-4123
MB BATCH ID: 170327L05
INSTRUMENT: GC 66
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED: 2017-03-28 02:30
REVIEWED BY: M
D/T REVIEWED:
MATRIX: Soil


DATA FILE: /chem1/SVOA/GC_66/170328/b1703282517032825

CLIENT WORK ORDER: 17-03-1557

<u>S#</u>	<u>RUN TYPE</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
28	B-DU1-ISM1-8		2017-03-28 03:41	/chem1/SVOA/GC_66/170328/b1703282917032829
29	B-DU1-ISM2-8		2017-03-28 03:59	/chem1/SVOA/GC_66/170328/b1703283017032830
30	B-DU1-ISM3-8		2017-03-28 04:35	/chem1/SVOA/GC_66/170328/b1703283217032832

RAW DATA SHEET FOR METHOD: EPA 8082

WORK ORDER: 099-12-535
INSTRUMENT: GC 66
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED: 2017-03-28 02:30
REVIEWED BY: 
D/T REVIEWED:

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282517032825

MB **CLIENT SAMPLE NUMBER: Method Blank**

LCS/MB BATCH: 170327L05 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 1.00 / ACTUAL: 1.00
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 1.00 ml / ACTUAL: 1.00 ml
UNITS: ug/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Aroclor-1016	0.000	1.00	ND	50	
Aroclor-1221	0.000	1.00	ND	50	
Aroclor-1232	0.000	1.00	ND	50	
Aroclor-1242	0.000	1.00	ND	50	
Aroclor-1248	0.000	1.00	ND	50	
Aroclor-1254	0.000	1.00	ND	50	
Aroclor-1260	0.000	1.00	ND	50	
Aroclor-1262	0.000	1.00	ND	50	
Aroclor-1268	0.000	1.00	ND	50	

↑
Return to Contents

LCS QUALITY CONTROL SHEET FOR METHOD: EPA 8082

LCS SAMPLE ID: 099-12-535-4123
LCS/MB BATCH ID: 170327L05
INSTRUMENT: GC 66

EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED: 2017-03-28 02:47
REVIEWED BY:
D/T REVIEWED:

~

DATA FILE: /chem1/SVOA/GC_66/170328/lb1703282617032826

<u>COMPOUND NAME</u>	<u>CONC ADDED</u>	<u>CONC REC</u>	<u>%RECOVERY</u>	<u>%REC CONTROL LIMIT</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Aroclor-1016	100.0	88.50	88	50-135	PASS	
Aroclor-1260	100.0	88.00	88	50-135	PASS	

**MATRIX SPIKE / MATRIX SPIKE DUPLICATE QUALITY CONTROL SHEET
FOR METHOD: EPA 8082**

SPIKED SAMPLE ID: 17-03-1557-28
MS/MSD BATCH: 170327S05
INSTRUMENTS:
 SAMPLE: GC 66
 MS: GC 66
 MSD: GC 66

EXTRACTION: EPA 3545
D/T EXTRACTED:
 SAMPLE: 2017-03-27 00:00
 MS: 2017-03-27 00:00
 MSD: 2017-03-27 00:00

ANALYZED BY: 1,028
D/T ANALYZED:
 SAMPLE: 2017-03-28 03:41
 MS: 2017-03-28 03:05
 MSD: 2017-03-28 03:23
REVIEWED BY: *W*
D/T REVIEWED:

COMMENT:

COMPOUND NAME	SAMPLE	INITIAL	FINAL	MS CONC	% MS.REC	MSD CONC	% MSD.REC	% REC CL	RPD	RPD CL	STATUS	QUALIFIERS
Aroclor-1016	ND	200.0	100.0	82.50	82	81.00	81	50-135	2	0-20	PASS	
Aroclor-1260	ND	200.0	100.0	85.00	85	86.00	86	50-135	1	0-20	PASS	

Data Files:

TYPE	DATA FILE	DATA FILE PATH
MS	17032827	/chem1/SVOA/GC_66/170328/b17032827
MSD	17032828	/chem1/SVOA/GC_66/170328/b17032828

SURROGATE RECOVERIES FOR METHOD: EPA 8082

WORK ORDER: 17-03-1557

BATCH ID:

LCS/MB: 170327L05**MS:** 170327S05

EXTRACTION : EPA 3545

REVIEWED BY: 

D/T REVIEWED:

28 **CLIENT SAMPLE NUMBER : B-DU1-ISM1-8**

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282917032829

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 03:41

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
Decachlorobiphenyl	77	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	65	25-145	PASS	

29 **CLIENT SAMPLE NUMBER : B-DU1-ISM2-8**

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703283017032830

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 03:59

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
Decachlorobiphenyl	73	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	61	25-145	PASS	

30 **CLIENT SAMPLE NUMBER : B-DU1-ISM3-8**

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703283217032832

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 04:35

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
Decachlorobiphenyl	76	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	70	25-145	PASS	

MB **CLIENT SAMPLE NUMBER : Method Blank**

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282517032825

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 02:30

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
Decachlorobiphenyl	70	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	69	25-145	PASS	

**SURROGATE RECOVERIES
FOR METHOD: EPA 8082**

WORK ORDER: 17-03-1557

BATCH ID:

LCS/MB: 170327L05

MS:

EXTRACTION : EPA 3545

REVIEWED BY: *21*

D/T REVIEWED:

LCS CLIENT SAMPLE NUMBER : Lab Control Sample

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282617032826

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 02:47

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Decachlorobiphenyl	71	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	71	25-145	PASS	

MS CLIENT SAMPLE NUMBER : Matrix Spike

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282717032827

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 03:05

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Decachlorobiphenyl	70	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	65	25-145	PASS	

MSD CLIENT SAMPLE NUMBER : Matrix Spike Duplicate

INSTRUMENT: GC 66

D/T EXTRACTED: 2017-03-27 00:00

DATA FILE: /chem1/SVOA/GC_66/170328/b1703282817032828

ANALYZED BY: 1,028

D/T ANALYZED 2017-03-28 03:23

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Decachlorobiphenyl	69	24-168	PASS	
2,4,5,6-Tetrachloro-m-Xylene	62	25-145	PASS	

Return to Contents

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032825.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032825.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 02:30
 Operator : 669 Inst ID: GC_66.i
 Smp Info : MB 170327L05
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 25
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000	407822147	68.7035	68.7
M 2 Aroclor-1016				Compound Not Detected.		
3 Aroclor 1016 (1)				Compound Not Detected.		
4 Aroclor 1016 (2)				Compound Not Detected.		
5 Aroclor 1016 (3)				Compound Not Detected.		
6 Aroclor 1016 (4)				Compound Not Detected.		
7 Aroclor 1016 (5)				Compound Not Detected.		
M 8 Aroclor-1260				Compound Not Detected.		
9 Aroclor 1260 (1)				Compound Not Detected.		
10 Aroclor 1260 (2)				Compound Not Detected.		
11 Aroclor 1260 (3)				Compound Not Detected.		
12 Aroclor 1260 (4)				Compound Not Detected.		
13 Aroclor 1260 (5)				Compound Not Detected.		
M 14 Aroclor-1221				Compound Not Detected.		
15 Aroclor 1221 (1)				Compound Not Detected.		
16 Aroclor 1221 (2)				Compound Not Detected.		
17 Aroclor 1221 (3)				Compound Not Detected.		
18 Aroclor 1221 (4)				Compound Not Detected.		
19 Aroclor 1221 (5)				Compound Not Detected.		
M 20 Aroclor-1232				Compound Not Detected.		
21 Aroclor 1232 (1)				Compound Not Detected.		

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032825.d
 Report Date: 28-Mar-2017 09:42

Page 2

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
22 Aroclor 1232 (2)				Compound Not Detected.		
23 Aroclor 1232 (3)				Compound Not Detected.		
24 Aroclor 1232 (4)				Compound Not Detected.		
25 Aroclor 1232 (5)				Compound Not Detected.		
M 26 Aroclor-1242				Compound Not Detected.		
27 Aroclor 1242 (1)				Compound Not Detected.		
28 Aroclor 1242 (2)				Compound Not Detected.		
29 Aroclor 1242 (3)				Compound Not Detected.		
30 Aroclor 1242 (4)				Compound Not Detected.		
31 Aroclor 1242 (5)				Compound Not Detected.		
M 32 Aroclor-1248				Compound Not Detected.		
33 Aroclor 1248 (1)				Compound Not Detected.		
34 Aroclor 1248 (2)				Compound Not Detected.		
35 Aroclor 1248 (3)				Compound Not Detected.		
36 Aroclor 1248 (4)				Compound Not Detected.		
37 Aroclor 1248 (5)				Compound Not Detected.		
M 38 Aroclor-1254				Compound Not Detected.		
39 Aroclor 1254 (1)				Compound Not Detected.		
40 Aroclor 1254 (2)				Compound Not Detected.		
41 Aroclor 1254 (3)				Compound Not Detected.		
42 Aroclor 1254 (4)				Compound Not Detected.		
43 Aroclor 1254 (5)				Compound Not Detected.		
M 44 Aroclor-1262				Compound Not Detected.		
45 Aroclor 1262 (1)				Compound Not Detected.		
46 Aroclor 1262 (2)				Compound Not Detected.		
47 Aroclor 1262 (3)				Compound Not Detected.		
48 Aroclor 1262 (4)				Compound Not Detected.		
49 Aroclor 1262 (5)				Compound Not Detected.		
M 50 Aroclor-1268				Compound Not Detected.		
51 Aroclor 1268 (1)				Compound Not Detected.		
52 Aroclor 1268 (2)				Compound Not Detected.		
53 Aroclor 1268 (3)				Compound Not Detected.		
54 Aroclor 1268 (4)				Compound Not Detected.		
55 Aroclor 1268 (5)				Compound Not Detected.		
\$ 56 Decachlorobiphenyl	13.097	13.098	-0.001	387709171	70.4036	70.4

Data File: /chem1/SV04/GC_66.i/170328.b/b17032825.d

Date : 28-MAR-2017 02:30

Client ID:

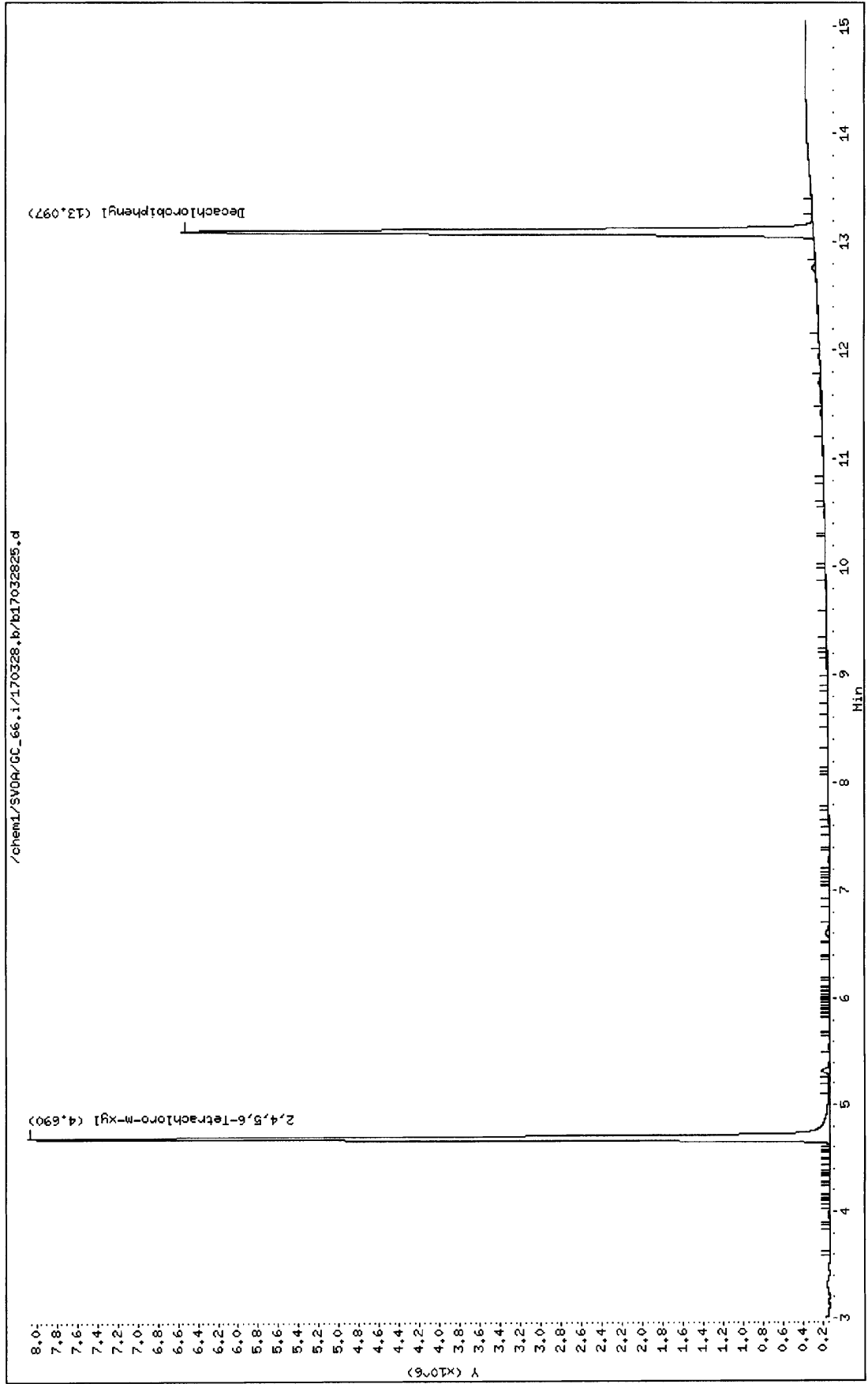
Sample Info: MB 170327L05

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_66.i/170328.b/b17032826.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032826.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 02:47
 Operator : 669 Inst ID: GC_66.i
 Smp Info : LCS 170327L05
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 26
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.689	4.690	-0.001	418868429	70.5644	70.6
M 2 Aroclor-1016				162212301	176.844	177
3 Aroclor 1016 (1)	5.531	5.529	0.002	17121353	187.422	187
4 Aroclor 1016 (2)	6.161	6.160	0.001	29111574	178.532	178
5 Aroclor 1016 (3)	6.843	6.840	0.003	62137562	170.426	170
6 Aroclor 1016 (4)	7.041	7.039	0.002	30054337	180.755	181
7 Aroclor 1016 (5)	7.194	7.190	0.004	23787475	180.240	180
M 8 Aroclor-1260				161879821	176.125	176
9 Aroclor 1260 (1)	9.744	9.743	0.001	50983712	192.483	192
10 Aroclor 1260 (2)	10.343	10.342	0.001	35963760	169.828	170
11 Aroclor 1260 (3)	10.736	10.736	0.000	39495834	170.016	170
12 Aroclor 1260 (4)	12.110	12.112	-0.002	11241783	171.972	172
13 Aroclor 1260 (5)	12.329	12.330	-0.001	24194732	167.089	167
\$ 56 Decachlorobiphenyl	13.097	13.098	-0.001	390988159	70.9990	71.0

Data File: /chem1/SV00A/GC_66.i/170328.b/b17032826.d

Date : 28-MAR-2017 02:47

Client ID:

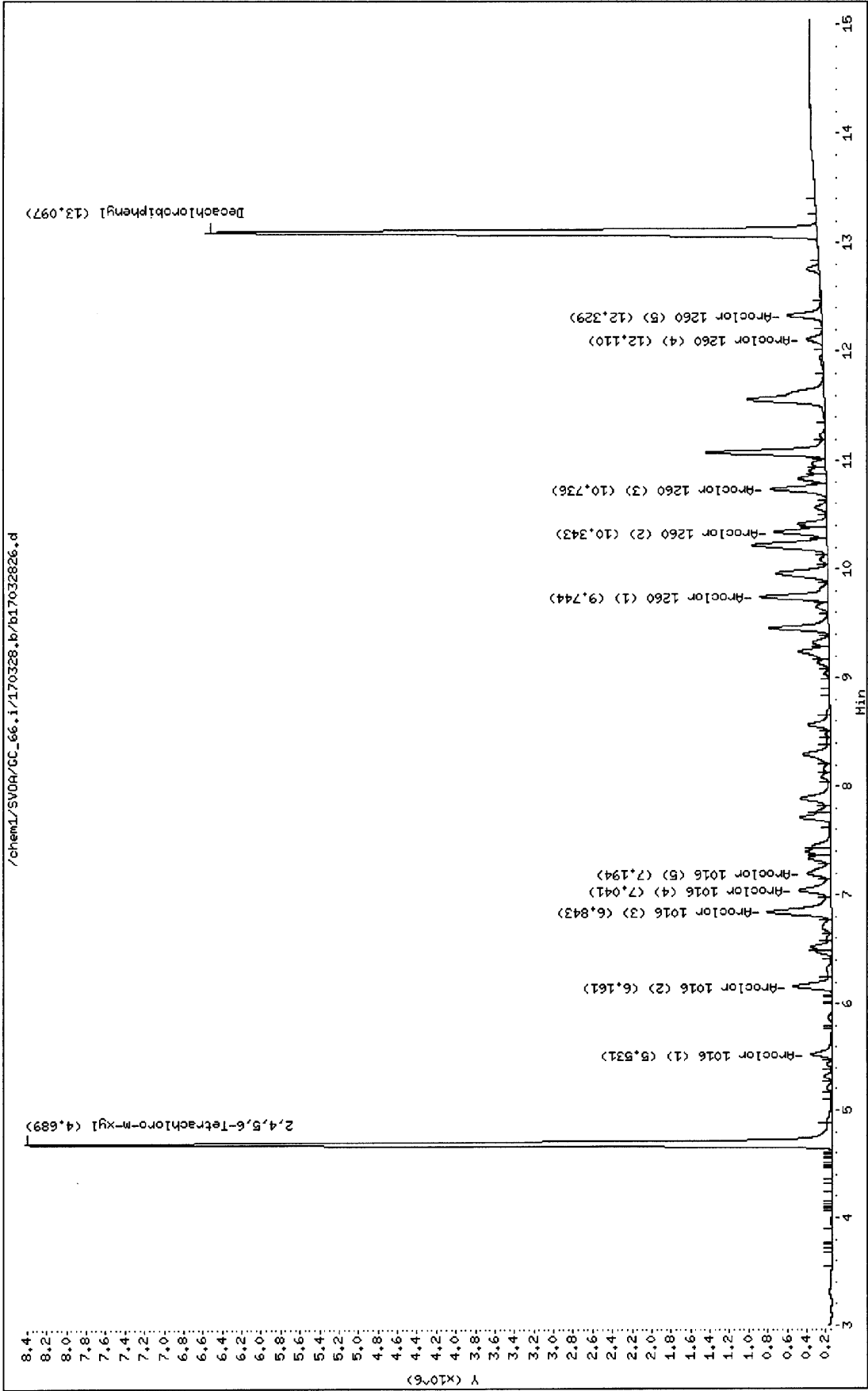
Sample Info: LCS 170327L05

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_66.i/170328.b/b17032827.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032827.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 03:05
 Operator : 669 Inst ID: GC_66.i
 Smp Info : MS 17-03-1557-28
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhnn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 27
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000	386619322	65.1315	65.1
M 2 Aroclor-1016				151406137	165.063	165
3 Aroclor 1016 (1)	5.531	5.529	0.002	16378707	179.293	179
4 Aroclor 1016 (2)	6.161	6.160	0.001	27202631	166.825	167
5 Aroclor 1016 (3)	6.842	6.840	0.002	58441294	160.288	160
6 Aroclor 1016 (4)	7.041	7.039	0.002	27913886	167.882	168
7 Aroclor 1016 (5)	7.192	7.190	0.002	21469619	162.677	163
M 8 Aroclor-1260				156034216	169.765	170
9 Aroclor 1260 (1)	9.744	9.743	0.001	47228372	178.305	178
10 Aroclor 1260 (2)	10.343	10.342	0.001	35058709	165.554	166
11 Aroclor 1260 (3)	10.737	10.736	0.001	38828487	167.143	167
12 Aroclor 1260 (4)	12.110	12.112	-0.002	10807824	165.333	165
13 Aroclor 1260 (5)	12.329	12.330	-0.001	24110824	166.510	166
\$ 56 Decachlorobiphenyl	13.097	13.098	-0.001	386749822	70.2294	70.2

Data File: /chem1/SV00A/GC_66.1/170328.b/17032827.d

Date : 28-HAR-2017 03:05

Client ID:

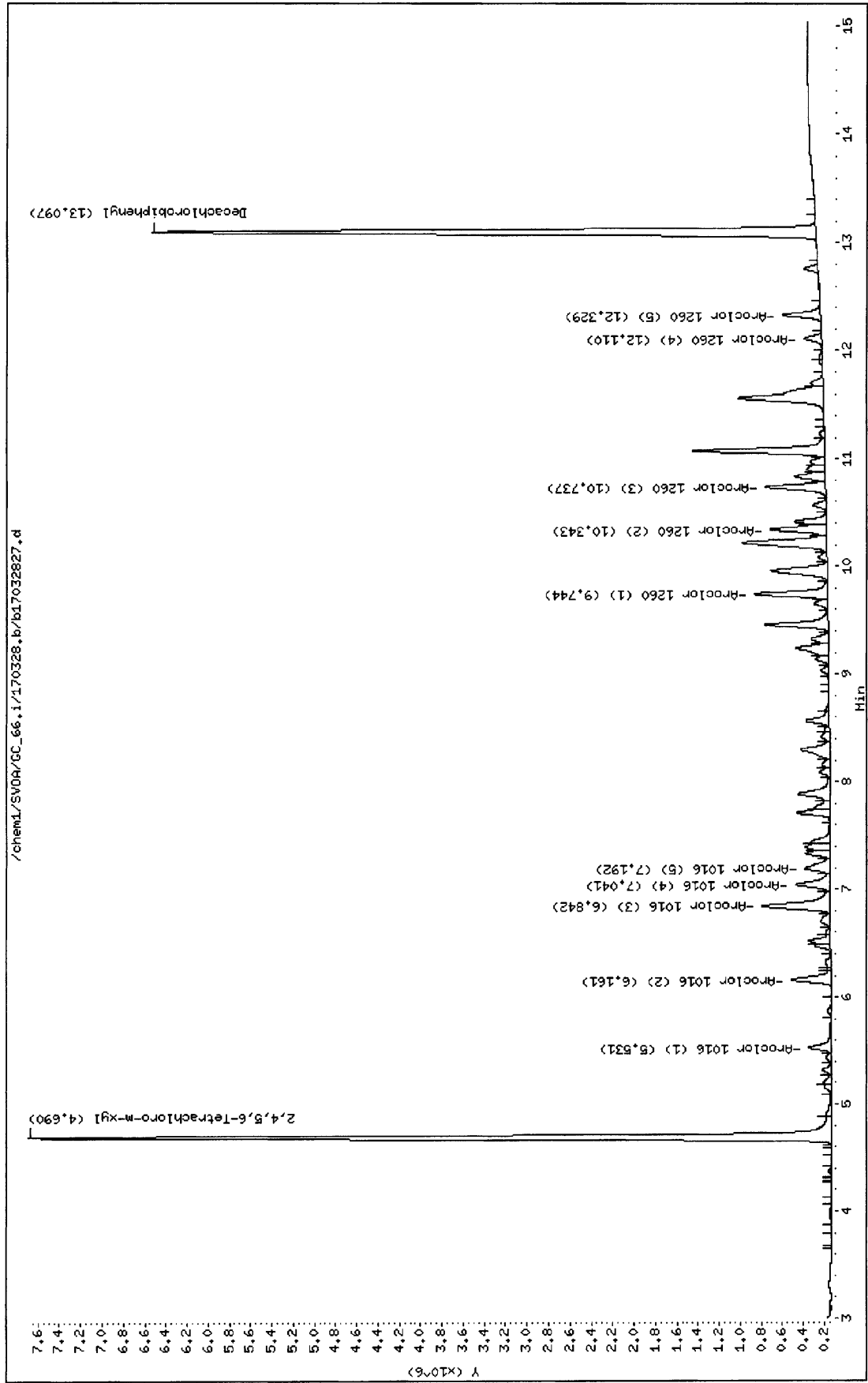
Sample Info: MS 17-03-1557-28

Instrument: GC_66.1

Operator: 669

Column diameter: 2.00

Column phase:



Data File: /chem1/SVOA/GC_66.i/170328.b/b17032828.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032828.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 03:23
 Operator : 669 Inst ID: GC_66.i
 Smp Info : MSD 17-03-1557-28
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 28
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ppb)	FINAL (ug/Kg)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000	365205260	61.5240	61.5
M 2 Aroclor-1016				148838746	162.264	162
3 Aroclor 1016 (1)	5.529	5.529	0.000	16185328	177.176	177
4 Aroclor 1016 (2)	6.160	6.160	0.000	26847354	164.646	165
5 Aroclor 1016 (3)	6.842	6.840	0.002	57517092	157.753	158
6 Aroclor 1016 (4)	7.041	7.039	0.002	27283345	164.089	164
7 Aroclor 1016 (5)	7.191	7.190	0.001	21005627	159.161	159
M 8 Aroclor-1260				157989844	171.893	172
9 Aroclor 1260 (1)	9.744	9.743	0.001	47702826	180.097	180
10 Aroclor 1260 (2)	10.343	10.342	0.001	35145838	165.966	166
11 Aroclor 1260 (3)	10.736	10.736	0.000	39512192	170.086	170
12 Aroclor 1260 (4)	12.110	12.112	-0.002	11137671	170.379	170
13 Aroclor 1260 (5)	12.329	12.330	-0.001	24491317	169.138	169
\$ 56 Decachlorobiphenyl	13.096	13.098	-0.002	378230462	68.6824	68.7

Data File: /chem1/SV00A/GC_66.i/170328.b/b17032828.d

Date : 28-MAR-2017 03:23

Client ID:

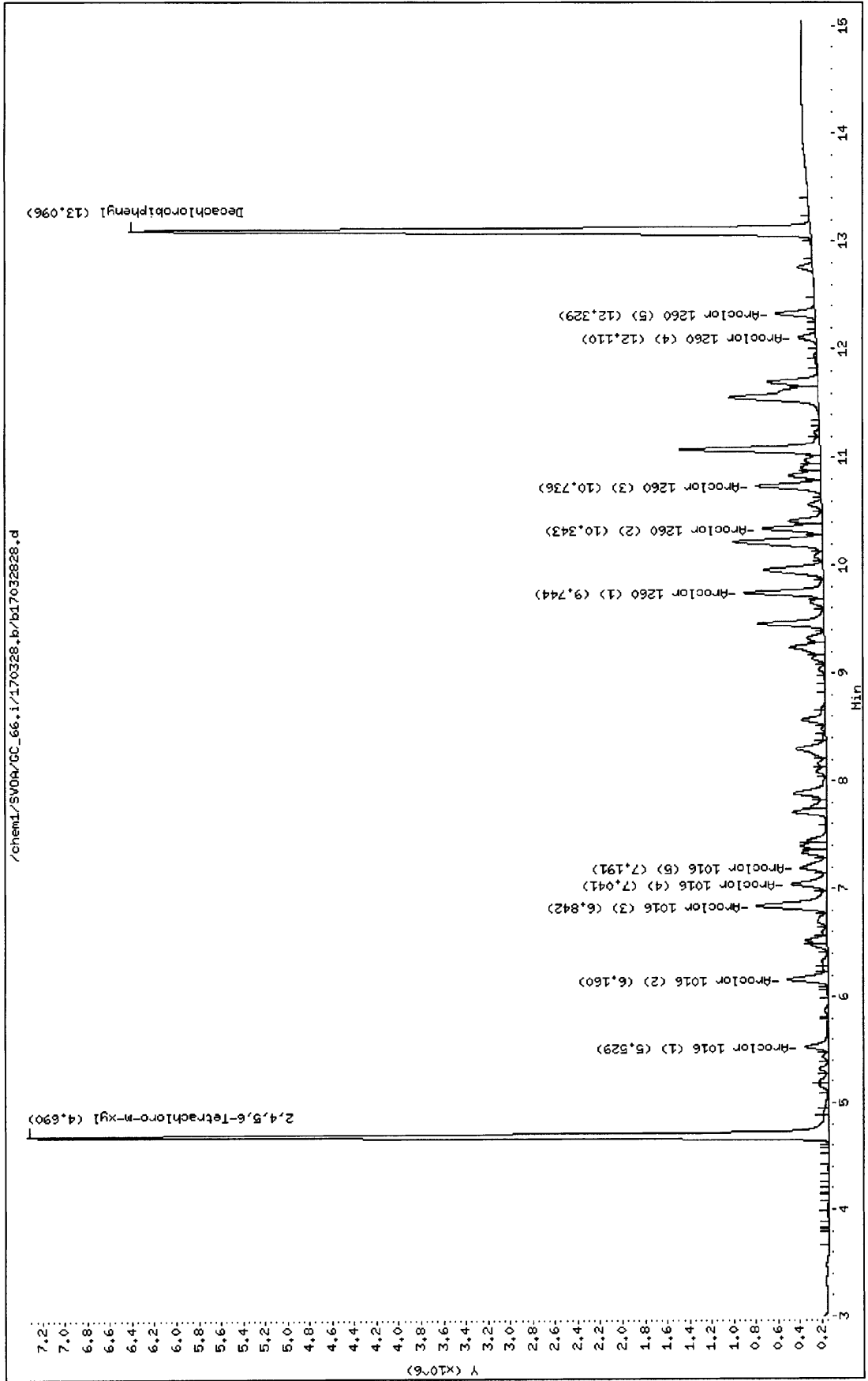
Sample Info: MSD 17-03-1557-28

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phase:



EPA METHOD 8082 PCB

Continuing Calibration

CCV ASSOCIATION SUMMARY
FOR METHOD: EPA 8082

BATCH ID: 170328A021
INSTRUMENT: GC 66

ANALYZED BY: 669

WORK ORDER: 099-12-532
MATRIX: Water

REVIEWED BY: *u*
D/T REVIEWED:

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
9514	Daily Calibration	2017-03-27 22:20	/chem1/SVOA/GC_66/170328/b1703281117032811

WORK ORDER: 17-03-1557
MATRIX: Soil

REVIEWED BY:
D/T REVIEWED:

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
28	B-DU1-ISM1-8	2017-03-28 03:05	/chem1/SVOA/GC_66/170328/b1703282717032827
28	B-DU1-ISM1-8	2017-03-28 03:23	/chem1/SVOA/GC_66/170328/b1703282817032828
28	B-DU1-ISM1-8	2017-03-28 03:41	/chem1/SVOA/GC_66/170328/b1703282917032829
29	B-DU1-ISM2-8	2017-03-28 03:59	/chem1/SVOA/GC_66/170328/b1703283017032830

CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8082

CCV WORK ORDER: 099-12-532-9514-5154

BATCH ID:

170222I003
170328A021
GC 66

ANALYZED BY: 669

D/T ANALYZED:

INITIAL:
CCV:
2017-02-22 17:27
2017-03-27 22:20

INSTRUMENT:

REVIEWED BY:
D/T REVIEWED:

M

DATA FILE: /chem1/SVOA/GC_66/170328/b170328117032811

<u>COMPOUND NAME</u>	<u>TYPE</u>	<u>CALIB MODEL</u>	<u>MIN RF</u>	<u>AVG RF</u>	<u>CCV RF</u>	<u>AMOUNT</u>	<u>CCV CONC</u>	<u>CCV %D</u>	<u>CCV %D CL</u>	<u>STATUS</u>
Aroclor-1016	C	Avg Resp	0.00	893304.438	847497.946			5	0-15	PASS
Aroclor-1260	C	Avg Resp	0.00	947781.003	853241.020			10	0-15	PASS

MIN RF: Method Specified Minimum Response Factor

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032811.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032811.d
 Lab Smp Id:
 Inj Date : 27-MAR-2017 22:20
 Operator : 669 Inst ID: GC_66.i
 Smp Info : PCB CCV 500PPB P021517I
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhhn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 11 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000	557400338	100.000	93.9
M 2 Aroclor-1016				423748973	500.000	462
3 Aroclor 1016 (1)	5.529	5.529	0.000	42854439	500.000	469
4 Aroclor 1016 (2)	6.160	6.160	0.000	75367011	500.000	462
5 Aroclor 1016 (3)	6.840	6.840	0.000	168894186	500.000	463
6 Aroclor 1016 (4)	7.039	7.039	0.000	76176510	500.000	458
7 Aroclor 1016 (5)	7.190	7.190	0.000	60456827	500.000	458
M 8 Aroclor-1260				426620510	500.000	464
9 Aroclor 1260 (1)	9.743	9.743	0.000	124792482	500.000	471
10 Aroclor 1260 (2)	10.342	10.342	0.000	97186700	500.000	459
11 Aroclor 1260 (3)	10.736	10.736	0.000	106686109	500.000	459
12 Aroclor 1260 (4)	12.112	12.112	0.000	30339923	500.000	464
13 Aroclor 1260 (5)	12.330	12.330	0.000	67615296	500.000	467
\$ 56 Decachlorobiphenyl	13.098	13.098	0.000	514389618	100.000	93.4

Data File: /chem1/SV04/GC_66.i/170328.b/b17032811.d

Date : 27-MAR-2017 22:20

Client ID:

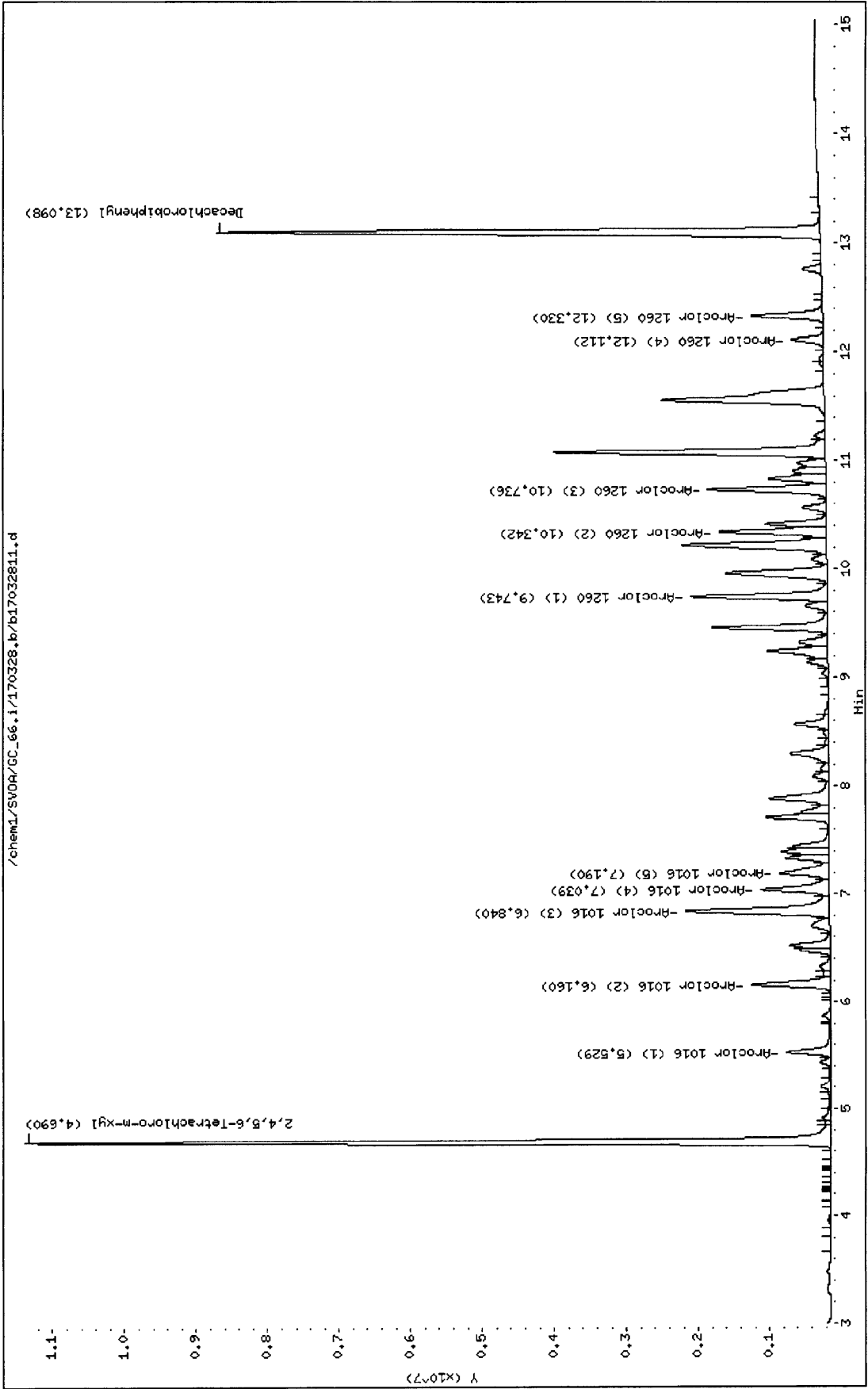
Sample Info: PCB CCV 500PPB F0215171

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phase:




CCV ASSOCIATION SUMMARY
FOR METHOD: EPA 8082

BATCH ID: 170328A022
INSTRUMENT: GC 66

ANALYZED BY: 669

WORK ORDER: 099-12-532
MATRIX: Water

REVIEWED BY: 
D/T REVIEWED:

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
9515	Daily Calibration	2017-03-28 04:17	/chem1/SVOA/GC_66/170328/b1703283117032831

WORK ORDER: 17-03-1557
MATRIX: Soil

REVIEWED BY:
D/T REVIEWED:

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
30	B-DU1-ISM3-8	2017-03-28 04:35	/chem1/SVOA/GC_66/170328/b1703283217032832

CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8082

CCV WORK ORDER: 099-12-532-9515-5154

BATCH ID:

170222I003
170328A022
GC 66

ANALYZED BY: 669

D/T ANALYZED:

2017-02-22 17:27
2017-03-28 04:17

INITIAL:
CCV:

REVIEWED BY:

D/T REVIEWED:

M

DATA FILE: /chem1/SVOA/GC_66/170328/b1703283117032831

<u>COMPOUND NAME</u>	<u>TYPE</u>	<u>CALIB MODEL</u>	<u>MIN RF</u>	<u>AVG RF</u>	<u>CCV RF</u>	<u>AMOUNT</u>	<u>CCV CONC</u>	<u>CCV %D</u>	<u>CCV %D CL</u>	<u>STATUS</u>
Atroclor-1016	C	Avg Resp	0.00	893304.438	825648.918			8	0-15	PASS
Atroclor-1260	C	Avg Resp	0.00	947781.003	836538.960			12	0-15	PASS

MIN RF: Method Specified Minimum Response Factor

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032831.d
 Report Date: 03/28/2017 10:34

Eurofins CalScience
 Calibration Verification Report

Instrument ID: GC_66.i Injection Date and Time: 28-MAR-2017 04:17
 Sample Name: PCB CCV 500PPB P021517I Initial Calibration Date(s): 22-FEB-2017 22-FEB-2017
 Sublist used: p1016_1260.sub Initial Calibration Time(s): 17:27 20:07
 Method used: /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m

Target Compounds	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift	Max%D / Drift	Curve Type
Aroclor 1260 (1)	270284.532	244053.526	0.01	10	15	Averaged
Aroclor 1260 (2)	213956.488	187468.718	0.01	12	15	Averaged
Aroclor 1260 (3)	233064.808	208078.902	0.01	11	15	Averaged
Aroclor 1260 (4)	83122.167	61148.526	0.01	26	15	Averaged
Aroclor-1260	947781.003	836538.960	0.01	12	15	Averaged
Aroclor-1016	893304.438	825648.918	0.01	8	15	Averaged
Aroclor 1016 (1)	90035.625	84114.300	0.01	7	15	Averaged
Aroclor 1016 (2)	159465.395	147869.816	0.01	7	15	Averaged
Aroclor 1016 (3)	354923.864	334754.466	0.01	6	15	Averaged
Aroclor 1016 (4)	162098.269	143114.172	0.01	12	15	Averaged
Aroclor 1016 (5)	126781.283	115796.164	0.01	9	15	Averaged
Aroclor 1260 (5)	147353.008	135789.288	0.01	8	15	Averaged
Surrogate Standards	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift	Max%D / Drift	Curve Type
2,4,5,6-Tetrachloro-m-xylene	5741865.174	5475699.490	0.01	5	15	Averaged
Decachlorobiphenyl	5459627.036	5076834.410	0.01	7	15	Averaged

<-Failed

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032831.d
 Report Date: 28-Mar-2017 09:42

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032831.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 04:17
 Operator : 669 Inst ID: GC_66.i
 Smp Info : PCB CCV 500PPB P021517I
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 09:42 uhnn Quant Type: ESTD
 Cal Date : 27-MAR-2017 20:32 Cal File: b17032805.d
 Als bottle: 31 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.690	4.690	0.000	547569949	100.000	92.2
M 2 Aroclor-1016				412824459	500.000	450
3 Aroclor 1016 (1)	5.528	5.528	0.000	42057150	500.000	460
4 Aroclor 1016 (2)	6.160	6.160	0.000	73934908	500.000	453
5 Aroclor 1016 (3)	6.838	6.838	0.000	167377233	500.000	459
6 Aroclor 1016 (4)	7.038	7.038	0.000	71557086	500.000	430
7 Aroclor 1016 (5)	7.188	7.188	0.000	57898082	500.000	439
M 8 Aroclor-1260				418269480	500.000	455
9 Aroclor 1260 (1)	9.742	9.742	0.000	122026763	500.000	461
10 Aroclor 1260 (2)	10.341	10.341	0.000	93734359	500.000	443
11 Aroclor 1260 (3)	10.735	10.735	0.000	104039451	500.000	448
12 Aroclor 1260 (4)	12.110	12.110	0.000	30574263	500.000	468
13 Aroclor 1260 (5)	12.329	12.329	0.000	67894644	500.000	469
\$ 56 Decachlorobiphenyl	13.097	13.097	0.000	507683441	100.000	92.2

Data File: /chem1/SV09/GC_66.i/170328.b/17032831.d

Date : 28-MAR-2017 04:17

Client ID:

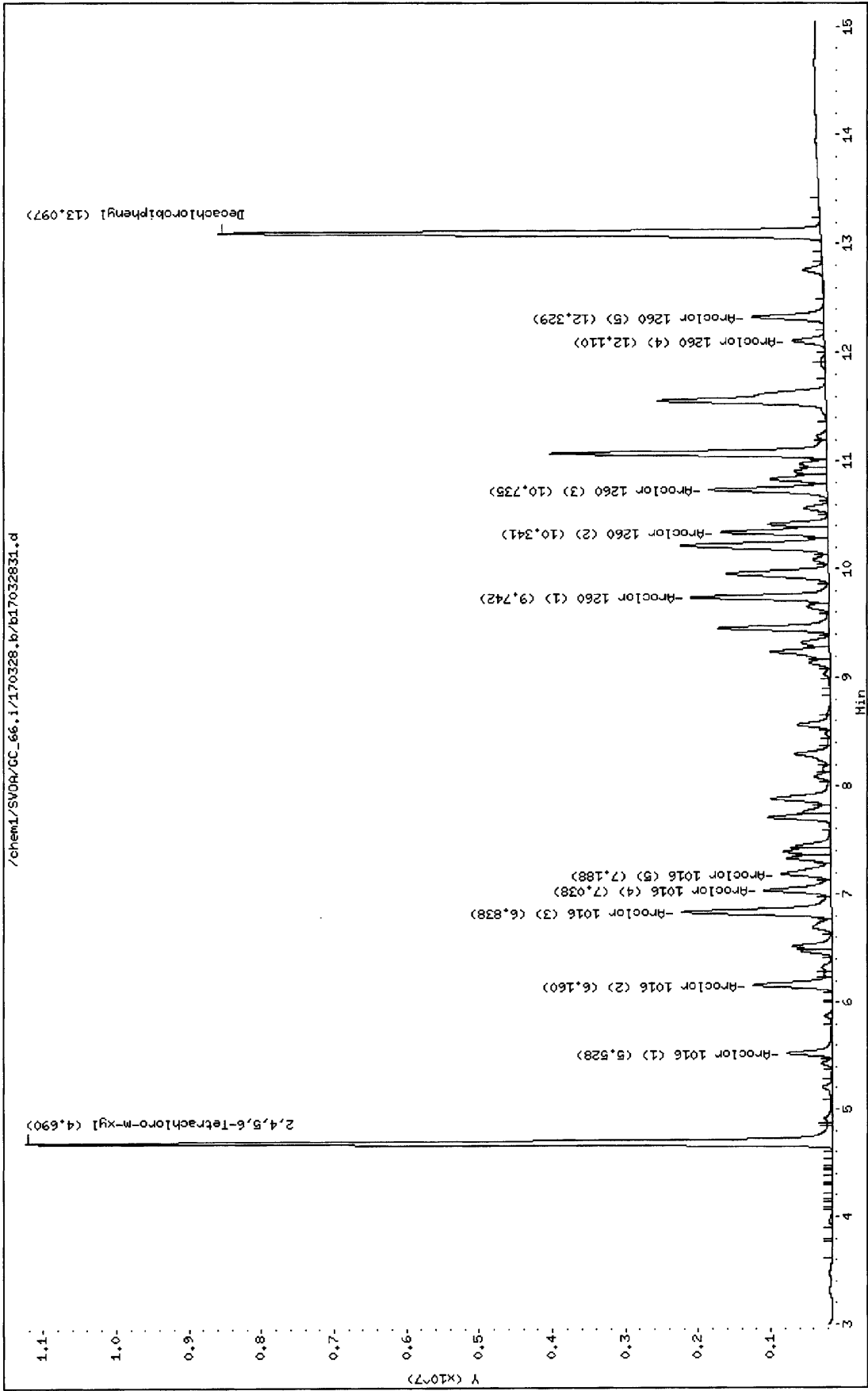
Sample Info: PCB CCV 500PPB P0215171

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phase:



CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8082

CCV WORK ORDER: 099-12-532-9516-5154

BATCH ID:

170222I003
170328A023
GC 66

ANALYZED BY: 669

D/T ANALYZED:

2017-02-22 17:27
2017-03-28 11:21

INITIAL:

CCV:

REVIEWED BY:

D/T REVIEWED:

~

DATA FILE: /chem1/SVOA/GC_66/170328/b1703285217032852

COMPOUND NAME	TYPE	CALIB MODEL	MIN RF	AVG RF	CCV RF	AMOUNT	CCV CONC	CCV %D	CCV %D CL	STATUS
Atroclor-1016	C	Avg Resp	0.00	893304.438	977934.848			-9	0-15	PASS
Atroclor-1260	C	Avg Resp	0.00	947781.003	979021.840			-3	0-15	PASS

MIN RF: Method Specified Minimum Response Factor

Data File: /chem1/SVOA/GC_66.i/170328.b/b17032852.d
 Report Date: 28-Mar-2017 11:49

Page 1

Eurofins Calscience

EPA 8082/A PCB analysis

Data file : /chem1/SVOA/GC_66.i/170328.b/b17032852.d
 Lab Smp Id:
 Inj Date : 28-MAR-2017 11:21
 Operator : 669 Inst ID: GC_66.i
 Smp Info : PCB CCV 500PPB P021517I
 Misc Info :
 Comment : Rtx-CLPesticide II
 Method : /chem1/SVOA/GC_66.i/170328.b/b8082d-n2.m
 Meth Date : 28-Mar-2017 11:49 uhhn Quant Type: ESTD
 Cal Date : 22-FEB-2017 20:07 Cal File: b17022210.d
 Als bottle: 52 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP Genie Compound Sublist: p1016_1260.sub
 Target Version: 3.50
 Processing Host: US26TAR4

Concentration Formula: Amt * DF * CpndVariable

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ppb)	ON-COL (ppb)
\$ 1 2,4,5,6-Tetrachloro-m-xylene	4.693	4.693	0.000	660349578	100.000	115
M 2 Aroclor-1016				488967424	500.000	547
3 Aroclor 1016 (1)	5.529	5.529	0.000	48777547	500.000	542
4 Aroclor 1016 (2)	6.160	6.160	0.000	87980479	500.000	552
5 Aroclor 1016 (3)	6.838	6.838	0.000	200141466	500.000	564
6 Aroclor 1016 (4)	7.038	7.038	0.000	83646827	500.000	516
7 Aroclor 1016 (5)	7.188	7.188	0.000	68421105	500.000	540
M 8 Aroclor-1260				489510920	500.000	516
9 Aroclor 1260 (1)	9.740	9.740	0.000	143191792	500.000	530
10 Aroclor 1260 (2)	10.341	10.341	0.000	110084698	500.000	514
11 Aroclor 1260 (3)	10.734	10.734	0.000	123038381	500.000	528
12 Aroclor 1260 (4)	12.108	12.108	0.000	35578667	500.000	428
13 Aroclor 1260 (5)	12.328	12.328	0.000	77617382	500.000	527
\$ 56 Decachlorobiphenyl	13.096	13.096	0.000	596309084	100.000	109

Data File: /chem1/SV04/GC_66.i/170328.b/b17032852.d

Date : 28-MAR-2017 11:21

Client ID:

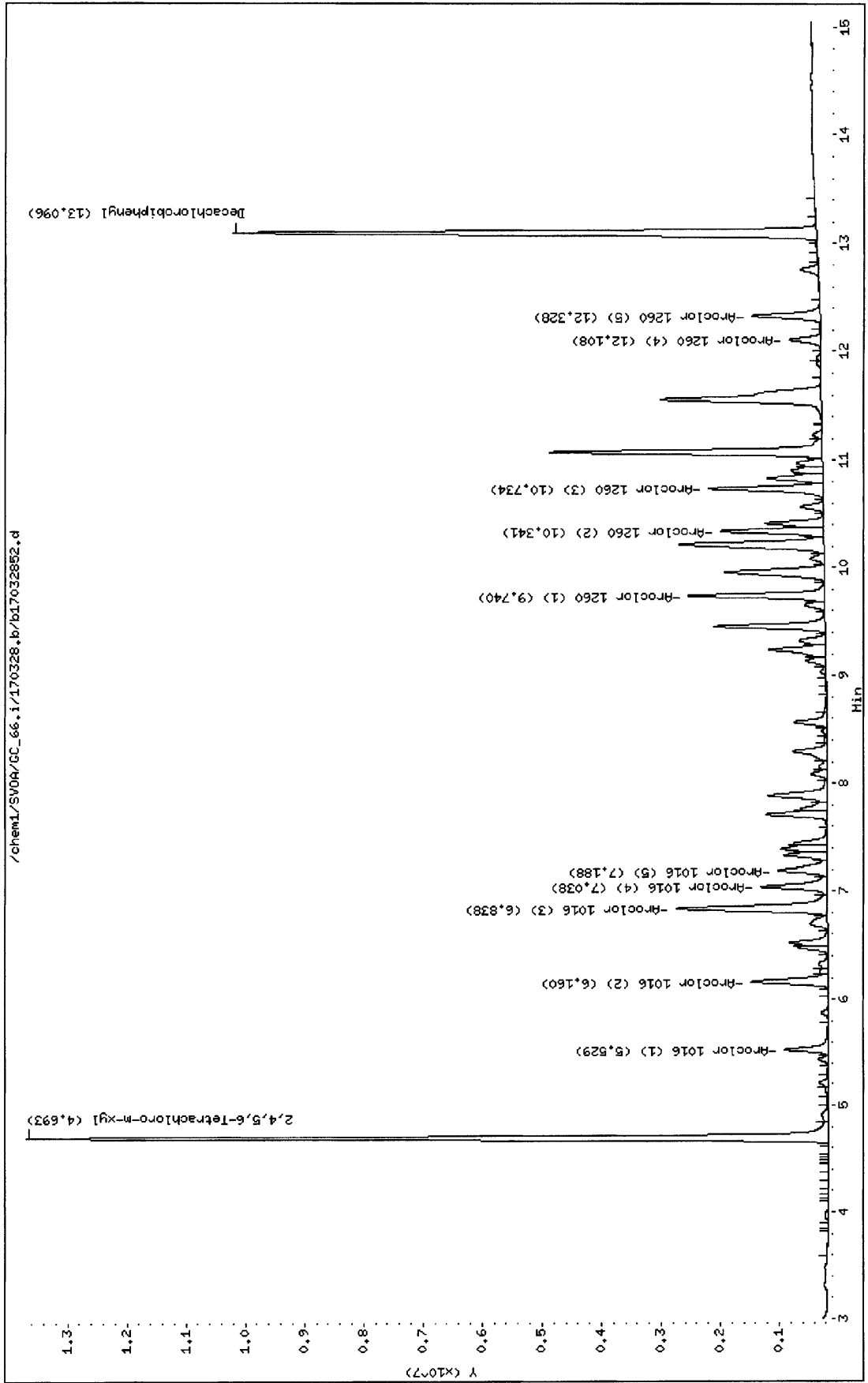
Sample Info: PCB CCV 500PPB P0215171

Instrument: GC_66.i

Operator: 669

Column diameter: 2.00

Column phaset



EPA METHOD 8082 PCB

Run Logs

Line	Vial	File	Name	Method	InjVolume	Acquired
1	100	17022200	IB S007-044-07	8082D-N2		22-Feb-17, 17:09:45
2	1	17022201	PCB ICAL1 P021517F 100PPB	8082D-N2		22-Feb-17, 17:27:32
3	2	17022202	PCB ICAL2 P021517E 250PPB	8082D-N2		22-Feb-17, 17:45:20
4	3	17022203	PCB ICAL3 P021517D 500PPB	8082D-N2		22-Feb-17, 18:03:10
5	4	17022204	PCB ICAL4 P021517C 750PPB	8082D-N2		22-Feb-17, 18:21:00
6	5	17022205	PCB ICAL5 P021517B 2000PPB	8082D-N2	1003	22-Feb-17, 18:38:49
7	6	17022206	PCB ICV P021517H 500PPB	8082D-N2		22-Feb-17, 18:56:37
8	7	17022207	PCB 1221/54 500PPB P120616I	8082D-N2		22-Feb-17, 19:14:26
9	8	17022208	PCB 1232/62 500PPB P120616J	8082D-N2		22-Feb-17, 19:32:18
10	9	17022209	PCB 1248/68 500PPB P120616K	8082D-N2		22-Feb-17, 19:50:07
11	10	17022210	PCB 1242 500PPB P120616L	8082D-N2		22-Feb-17, 20:07:55



Line	Vial	File	Name	Method	InjVolume	Acquired
1	150	17032800	S007-044-07	8082D-N2		27-Mar-17, 19:03:33
2	1	17032801	PCB CCV P021517H 500PPB	8082D-N2		27-Mar-17, 19:21:22
3	2	17032802	LCS 170323L16	8082D-N2		27-Mar-17, 19:39:12
4	3	17032803	MB 170323L16	8082D-N2		27-Mar-17, 19:57:03
5	4	17032804	MS 17-03-1551-1 170323L16	8082D-N2		27-Mar-17, 20:14:57
6	5	17032805	MSD 17-03-1551-1 170323L16	8082D-N2		27-Mar-17, 20:32:49
7	6	17032806	17-03-1551-1	8082D-N2		27-Mar-17, 20:50:42
8	7	17032807	PCB1221/54 P032617A	8082D-N2		27-Mar-17, 21:08:32
9	8	17032808	PCB1232/62 P032617B	8082D-N2		27-Mar-17, 21:26:27
10	9	17032809	PCB1248/68 P032617C	8082D-N2		27-Mar-17, 21:44:17
11	10	17032810	PCB1242 P032617D	8082D-N2		27-Mar-17, 22:02:17
12	11	17032811	PCB CCV 500PPB P021517I	8082D-N2		27-Mar-17, 22:20:07
13	12	17032812	MB 170324L03	8082D-N2		27-Mar-17, 22:38:00
14	13	17032813	LCS 170324L03	8082D-N2		27-Mar-17, 22:55:50
15	14	17032814	MS 17-03-1417-29 170324S03	8082D-N2		27-Mar-17, 23:13:39
16	15	17032815	MSD 17-03-1417-29 170324S03	8082D-N2		27-Mar-17, 23:31:29
17	16	17032816	17-03-1417-29	8082D-N2		27-Mar-17, 23:49:22
18	17	17032817	17-03-1417-28	8082D-N2		28-Mar-17, 00:07:14
19	18	17032818	17-03-1417-30	8082D-N2		28-Mar-17, 00:25:05
20	19	17032819	17-03-1418-28	8082D-N2		28-Mar-17, 00:42:56
21	20	17032820	17-03-1418-29	8082D-N2		28-Mar-17, 01:00:49
22	21	17032821	17-03-1418-30	8082D-N2		28-Mar-17, 01:18:39
23	22	17032822	17-03-1419-28	8082D-N2		28-Mar-17, 01:36:30
24	23	17032823	17-03-1419-29	8082D-N2		28-Mar-17, 01:54:21
25	24	17032824	17-03-1419-30	8082D-N2		28-Mar-17, 02:12:15
26	25	17032825	MB 170327L05	8082D-N2		28-Mar-17, 02:30:08
27	26	17032826	LCS 170327L05	8082D-N2		28-Mar-17, 02:47:56
28	27	17032827	MS 17-03-1557-28	8082D-N2		28-Mar-17, 03:05:48
29	28	17032828	MSD 17-03-1557-28	8082D-N2		28-Mar-17, 03:23:47
30	29	17032829	17-03-1557-28	8082D-N2		28-Mar-17, 03:41:39
31	30	17032830	17-03-1557-29	8082D-N2		28-Mar-17, 03:59:29
32	31	17032831	PCB CCV 500PPB P021517I	8082D-N2		28-Mar-17, 04:17:18
33	32	17032832	17-03-1557-30	8082D-N2		28-Mar-17, 04:35:12
34	33	17032833	17-03-1556-28	8082D-N2		28-Mar-17, 04:53:02
35	34	17032834	17-03-1556-29	8082D-N2		28-Mar-17, 05:10:54
36	35	17032835	17-03-1556-30	8082D-N2		28-Mar-17, 05:28:42
37	36	17032836	MB 170323L21	8082D-N2		28-Mar-17, 05:46:36
38	37	17032837	LCS 170323L21	8082D-N2		28-Mar-17, 06:04:24
39	38	17032838	LCSD 170323L21	8082D-N2		28-Mar-17, 06:22:13
40	39	17032839	MS 16-10-1328-47 10X	8082D-N2		28-Mar-17, 06:40:03
41	40	17032840	MSD 16-10-1328-47 10X	8082D-N2		28-Mar-17, 06:57:58
42	41	17032841	16-10-1328-47 10X	8082D-N2		28-Mar-17, 07:15:48
43	42	17032842	16-10-1328-48 10X	8082D-N2		28-Mar-17, 07:33:41
44	43	17032843	16-10-1328-49 10X	8082D-N2		28-Mar-17, 07:51:29
45	44	17032844	16-10-1328-50 10X	8082D-N2		28-Mar-17, 08:09:25
46	45	17032845	16-10-1328-51 10X	8082D-N2		28-Mar-17, 08:27:15
47	46	17032846	16-10-1328-52 10X	8082D-N2		28-Mar-17, 08:45:04
48	47	17032847	16-10-1328-53 10X	8082D-N2		28-Mar-17, 09:02:55
49	48	17032848	17-03-1950-1	8082D-N2		28-Mar-17, 09:20:51
50	49	17032849	17-03-1005-1	8082D-N2		28-Mar-17, 10:02:40
51	50	17032850	17-03-1005-2	8082D-N2		28-Mar-17, 10:33:04
52	51	17032851	16-10-1328-48 10X Hg	8082D-N2		28-Mar-17, 10:50:55
53	52	17032852	PCB CCV 500PPB P021517I	8082D-N2		28-Mar-17, 11:21:49
54	53	17032853	17-03-1005-3	8082D-N2		
55	54	17032854	17-03-1005-4	8082D-N2		
56	55	17032855	17-03-1005-6	8082D-N2		
57	56	17032856	17-03-1005-7	8082D-N2		
58	57	17032857	17-03-1005-8	8082D-N2		
59	58	17032858	17-03-1005-10	8082D-N2		
60	59	17032859	17-03-1005-11	8082D-N2		

EPA METHOD 8082 PCB

Preparation Logs

Analysis Method (EPA Method): 608 8081 8082 8141 8310 TO-13 TO-4
 8270 (Soil Soil SIM SUPER PAH SIM PAH SIM Pest SIM PCB cong. SIM FL)

Extraction Method (EPA Method): 3510 3520 3540 3541 3545 3550 3580

Analyst ID#: Measuring Sample- 610 Start Extraction- 610/787 Blow Down- 610 Clean Up- 47

Matrix: Soil Aqueous Oil Wipe Filter Tissue Air

Balance ID#: 70 Filter ID#: 507-17-17 ASE ID#: 7 Soxtherm ID#: Orbit Shaker ID#: Sonicator ID#:

Ext. Start Date/Time: 03/27/17 9:00 Ext. End Date/Time: 03/27/17 13:00

Sand or Wipe ID#: 507-19-19 Drying Agent: Na₂SO₄ Diatomaceous Earth
Drying Agent(s) ID#: 507-44-18 / 507-22-03

Surrogate Std ID# & Volume Added (mL): S101316A 0.5

Spike Std ID# & Volume Added (mL): S020317A 0.5 Spike Added to: LCS LCSD MS MSD

Extraction Solvent: MeCl₂ 1:1 Hexane-Acetone 1:1 MeCl₂-Acetone 9:1 Hexane-Diethyl-ether Acetonitrile

Extraction Solvent ID#: 507-44-07/507-44-08 Exchange Solvent (Hexane Acetonitrile) ID#: 507-44-07

Clean Up Start Date & Time: 3/27/17 16:30 Clean Up End Date & Time: 3/27/17 17:00

Clean Up: 3620 Florisil 3630 SGC 3660 Sulfur 3665 Acid Other Cartridge ID#: 4-127

Clean Up Reagent ID#: 507-52-12 Cartridge Conditioning Column Pre-Elution Reagent ID#:

MB/LCS/MS Batch #: 170327205	Sample W(g)/V (mL)		Clean Up Performed	Comments
	Initial	Final		
Cel ID#:				
MB	20.0	10	<input checked="" type="checkbox"/>	
LCS	20.0	10	<input checked="" type="checkbox"/>	
LCSD MA	-	-	<input type="checkbox"/>	
MS 17-03-1557-28B	20.0	10	<input checked="" type="checkbox"/>	
MSD -28B	20.0	10	<input checked="" type="checkbox"/>	
17-03-1556-28 B	20.0	10	<input checked="" type="checkbox"/>	
17-03-1556-29 B	20.1	10	<input checked="" type="checkbox"/>	
17-03-1556-30 B	20.1	10	<input checked="" type="checkbox"/>	
17-03-1557-28 B	20.0	10	<input checked="" type="checkbox"/>	
17-03-1557-29 B	20.2	10	<input checked="" type="checkbox"/>	
17-03-1557-30 B	20.1	10	<input checked="" type="checkbox"/>	
17-03-1558-28 B	20.0	10	<input checked="" type="checkbox"/>	Hold ↓ dirty sample
17-03-1558-29 B	20.1	10	<input checked="" type="checkbox"/>	
17-03-1558-30 B	20.1	10	<input checked="" type="checkbox"/>	
17-03-1950-1A	1.01	10	<input checked="" type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	

Peer Reviewed by: 471

Peer Reviewed Date: 3/27/17

Revision Date: 10/20/16

EPA METHOD 8270C PAHSIM

RAW DATA

EPA METHOD 8270C PAHSIM

Initial Calibration

INITIAL CALIBRATION QUALITY CONTROL SUMMARY FOR METHOD: EPA 8270C SIM PAHS

ICAL WORK ORDER: 099-06-009-4979-4741
ICAL BATCH ID: 1703131001
INSTRUMENT: GC/MS EEE

ANALYZED BY: 907
ICAL D/T ANALYZED: 2017-03-13 14:13
REVIEWED BY: 262
D/T REVIEWED: 2017-03-14 12:54

COMPOUND	COMP. TYPE	CALIB. MODEL	1	2	3	4	5	6	7	8	9	AVG. RF	MIN. RF	%RSD CL	%RSD CL	R of R ² CL	R of R ² CL	STATUS
Naphthalene		Avg RF	1.116	1.006	1.100	1.129	1.071					1.085	0.00	4	0-15	4.487		PASS
2-Methylnaphthalene		Avg RF	0.652	0.598	0.675	0.704	0.679					0.661	0.00	6	0-15	6.043		PASS
1-Methylnaphthalene		Avg RF	0.707	0.604	0.670	0.683	0.650					0.663	0.00	6	0-15	5.855		PASS
Acenaphthylene		Avg RF	2.585	2.328	2.575	2.640	2.513					2.528	0.00	5	0-15	4.766		PASS
Acenaphthene	C	Avg RF	1.539	1.413	1.552	1.581	1.501					1.517	0.00	4	0-15	4.263		PASS
Fluorene		Avg RF	1.678	1.586	1.751	1.767	1.664					1.689	0.00	4	0-15	4.314		PASS
Phenanthrene		Avg RF	1.080	0.981	1.087	1.105	1.053					1.061	0.00	5	0-15	4.587		PASS
Anthracene		Avg RF	1.036	0.984	0.969	1.116	1.083					1.037	0.00	6	0-15	6.047		PASS
Fluoranthene	C	Avg RF	1.376	1.262	1.391	1.425	1.371					1.365	0.00	4	0-15	4.487		PASS
Pyrene		Avg RF	1.387	1.261	1.393	1.422	1.344					1.362	0.00	5	0-15	4.620		PASS
Benzo (a) Anthracene		Avg RF	1.432	1.208	1.306	1.343	1.298					1.317	0.00	6	0-15	6.146		PASS
Chrysene		Avg RF	1.217	1.108	1.220	1.253	1.186					1.197	0.00	5	0-15	4.604		PASS
Benzo (k) Fluoranthene		Avg RF	1.363	1.261	1.425	1.424	1.358					1.366	0.00	5	0-15	4.922		PASS
Benzo (b) Fluoranthene		Avg RF	1.221	1.183	1.395	1.380	1.336					1.303	0.00	7	0-15	7.367		PASS
Benzo (a) Pyrene	C	Avg RF	1.253	1.166	1.268	1.329	1.265					1.256	0.00	5	0-15	4.648		PASS
Indeno (1,2,3-c,d) Pyrene		Avg RF	1.472	1.337	1.583	1.587	1.530					1.502	0.00	7	0-15	6.876		PASS
Dibenz (a,h) Anthracene		Avg RF	1.022	1.057	1.229	1.228	1.159					1.139	0.00	8	0-15	8.436		PASS
Benzo (g,h,i) Perylene		Avg RF	1.197	1.089	1.261	1.263	1.221					1.207	0.00	6	0-15	5.902		PASS

Data Files:

LR - E: Linear Regression (Equal Weight) LR - IC: Linear Regression (Inverse Concentration Weight) LR - ISC: Linear Regression (Inverse Square Concentration Weight)
Avg RF: Average Response Factor QR - E: Quadratic Regression (Equal Weight)



INITIAL CALIBRATION QUALITY CONTROL SUMMARY FOR METHOD: EPA 8270C SIM PAHS

ICAL WORK ORDER: 099-06-009-4979-4741
ICAL BATCH ID: 1703131001
INSTRUMENT: GC/MS EEE

ANALYZED BY: 907
ICAL D/T ANALYZED: 2017-03-13 14:13
REVIEWED BY: 262
D/T REVIEWED: 2017-03-14 12:54

LEVEL #	D/T ANALYZED	DATA FILE
1	2017-03-13 14:13	Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar010.d\13mar010.ir
2	2017-03-13 13:53	Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar009.d\13mar009.ir
3	2017-03-13 13:33	Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar008.d\13mar008.ir
4	2017-03-13 13:13	Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar007.d\13mar007.ir
5	2017-03-13 12:52	Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar006.d\13mar006.ir

LR - E: Linear Regression (Equal Weight)
Avg RF: Average Response Factor

LR - IC: Linear Regression (Inverse Concentration Weight)
QR - E: Quadratic Regression (Equal Weight)

LR - ISC: Linear Regression (Inverse Square Concentration Weight)



Return to Contents

INITIAL CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8270C SIM PAHS

ICV WORK ORDER: 099-06-009-4979-4741
 INITIAL BATCH ID: 1703131001
 INSTRUMENT: GC/MS EEE

ANALYZED BY: 907
 D/T ANALYZED: 2017-03-13 14:13
 INITIAL: 2017-03-13 14:34
 ICV: 262
 REVIEWED BY: 262
 D/T REVIEWED: 2017-03-14 12:54

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar011.d\13mar011.r

COMPOUND	COMP TYPE	CALIB MODEL	MIN RF	AVG RF	ICV RF	AMOUNT	ICV	ICV %D	ICV %D CL	STATUS
Naphthalene	Avg Resp		0.00	1.085	1.135			-5	0-20	PASS
2-Methylnaphthalene	Avg Resp		0.00	0.661	0.783			-18	0-20	PASS
1-Methylnaphthalene	Avg Resp		0.00	0.663	0.698			-5	0-20	PASS
Acenaphthylene	Avg Resp		0.00	2.528	2.579			-2	0-20	PASS
Acenaphthene	C Avg Resp		0.00	1.517	1.642			-8	0-20	PASS
Fluorene	Avg Resp		0.00	1.689	1.833			-9	0-20	PASS
Phenanthrene	Avg Resp		0.00	1.061	1.203			-13	0-20	PASS
Anthracene	Avg Resp		0.00	1.037	1.123			-8	0-20	PASS
Fluoranthene	C Avg Resp		0.00	1.365	1.484			-9	0-20	PASS
Pyrene	Avg Resp		0.00	1.362	1.480			-9	0-20	PASS
Benzo (a) Anthracene	Avg Resp		0.00	1.317	1.360			-3	0-20	PASS
Chrysene	Avg Resp		0.00	1.197	1.249			-4	0-20	PASS
Benzo (k) Fluoranthene	Avg Resp		0.00	1.366	1.424			-4	0-20	PASS
Benzo (b) Fluoranthene	Avg Resp		0.00	1.303	1.418			-9	0-20	PASS
Benzo (a) Pyrene	C Avg Resp		0.00	1.256	1.318			-5	0-20	PASS
Indeno (1,2,3-c,d) Pyrene	Avg Resp		0.00	1.502	1.570			-5	0-20	PASS
Dibenz (a,h) Anthracene	Avg Resp		0.00	1.139	1.243			-9	0-20	PASS
Benzo (g,h,i) Perylene	Avg Resp		0.00	1.207	1.334			-11	0-20	PASS

MIN RF: Method Specified Minimum Response Factor

Report Date : 13-Mar-2017 15:04

Page 1

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 03-JAN-2017 11:59
 End Cal Date : 13-MAR-2017 14:13
 Quant Method : ISTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP RTE
 Method file : /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m
 Cal Date : 13-Mar-2017 14:59 ev7p
 Curve Type : Average

Calibration File Names:

Level 1: /chem/SVOA/GCMS_EEE.i/170313.b/13mar010.d
 Level 2: /chem/SVOA/GCMS_EEE.i/170313.b/13mar009.d
 Level 3: /chem/SVOA/GCMS_EEE.i/170313.b/13mar008.d
 Level 4: /chem/SVOA/GCMS_EEE.i/170313.b/13mar007.d
 Level 5: /chem/SVOA/GCMS_EEE.i/170313.b/13mar006.d

Compound	0.10000	0.50000	1.000	2.000	5.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5		
4 Naphthalene	1.11581	1.00634	1.10029	1.12879	1.07133	1.08451	4
5 2-Methylnaphthalene	0.65197	0.59811	0.67457	0.70415	0.67863	0.66149	6
6 1-Methylnaphthalene	0.70747	0.60440	0.67039	0.68304	0.65028	0.66312	6
8 Biphenyl	2.35823	2.11929	2.30021	2.29717	2.12271	2.23952	5
9 2,6-Dimethylnaphthalene	1.61703	1.47875	1.60729	1.63394	1.52299	1.57200	4
10 Acenaphthylene	2.58512	2.32846	2.57549	2.63987	2.51288	2.52836	5
12 Acenaphthene	1.53912	1.41348	1.55163	1.58080	1.50094	1.51719	4
13 Dibenzofuran	2.16158	2.04374	2.26236	2.30541	2.19584	2.19379	5
14 1,6,7-Trimethylnaphthalene	1.63669	1.46201	1.59794	1.62883	1.52745	1.57058	5
15 Fluorene	1.67753	1.58594	1.75075	1.76676	1.66409	1.68901	4
16 Dibenzothiophene	4.72855	4.27529	4.71886	4.80185	4.58747	4.62241	5
18 Phenanthrene	1.07962	0.98082	1.08700	1.10520	1.05266	1.06106	5
19 Anthracene	1.03555	0.98375	0.96886	1.11560	1.08257	1.03727	6
20 1-Methylphenanthrene	0.97116	0.92998	1.02072	1.05271	1.00870	0.99665	5
21 Fluoranthene	1.37602	1.26225	1.39052	1.42534	1.37085	1.36500	4
22 Pyrene	1.38699	1.26059	1.39347	1.42216	1.34443	1.36153	5
24 Benzo (a) Anthracene	1.43152	1.20758	1.30623	1.34262	1.29815	1.31722	6
26 Chrysene	1.21702	1.10798	1.22017	1.25314	1.18561	1.19678	5
27 Benzo (b) Fluoranthene	1.22069	1.18251	1.39522	1.38038	1.33563	1.30289	7
28 Benzo (k) Fluoranthene	1.36330	1.26057	1.42543	1.42402	1.35775	1.36621	5
29 Benzo (e) pyrene	1.59986	1.40940	1.63930	1.60515	1.53257	1.55726	6
30 Benzo (a) Pyrene	1.25298	1.16587	1.26774	1.32867	1.26501	1.25606	5

Report Date : 13-Mar-2017 15:04

Page 2

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 03-JAN-2017 11:59
 End Cal Date : 13-MAR-2017 14:13
 Quant Method : ISTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP RTE
 Method file : /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m
 Cal Date : 13-Mar-2017 14:59 ev7p
 Curve Type : Average

Compound	0.10000	0.50000	1.000	2.000	5.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5		
32 Perylene	1.53641	1.41079	1.50876	1.54399	1.46685	1.49336	4
33 Indeno (1,2,3-c,d) Pyrene	1.47205	1.33679	1.58268	1.58673	1.52950	1.50155	7
34 Dibenz (a,h) Anthracene	1.02180	1.05709	1.22915	1.22831	1.15856	1.13898	8
35 Benzo (g,h,i) Perylene	1.19737	1.08929	1.26132	1.26340	1.22141	1.20656	6
\$ 2 Nitrobenzene-d5	0.30130	0.29496	0.33923	0.35910	0.36136	0.33119	10
\$ 7 2-Fluorobiphenyl	1.88477	1.69943	1.88233	1.85964	1.69885	1.80500	5
\$ 23 p-Terphenyl-d14	0.78803	0.71467	0.79649	0.80490	0.77019	0.77486	5

Report Date : 13-Mar-2017 15:04

Page 3

Eurofins Calscience

INITIAL CALIBRATION DATA

Start Cal Date : 03-JAN-2017 11:59
End Cal Date : 13-MAR-2017 14:13
Quant Method : ISTD
Origin : Disabled
Target Version : 3.50
Integrator : HP RTE
Method file : /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m
Cal Date : 13-Mar-2017 14:59 ev7p
Curve Type : Average

Average %RSD Results.	
=====	
Calculated Average %RSD =	5.35528
Maximum Average %RSD =	15.00000
* Passed Average %RSD Test.	

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar011.d
 Report Date: 03/13/2017 15:04

Eurofins CalScience
 Calibration Verification Report

Instrument ID: GCMS_EEE.i Injection Date and Time: 13-MAR-2017 14:34
 Sample Name: ICV S010317I 1PPM Initial Calibration Date(s): 03-JAN-2017 13-MAR-2017
 Sublist used: all.sub Initial Calibration Time(s): 11:59 14:13
 Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m

Target Compounds	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift	Max%D / Drift	Curve Type
Naphthalene	1.085	1.135	0.00	-5	20	Averaged
2-Methylnaphthalene	0.661	0.783	0.00	-18	20	Averaged
1-Methylnaphthalene	0.663	0.698	0.00	-5	20	Averaged
Acenaphthylene	2.528	2.579	0.00	-2	20	Averaged
Acenaphthene	1.517	1.642	0.00	-8	20	Averaged
Fluorene	1.689	1.833	0.00	-9	20	Averaged
Phenanthrene	1.061	1.203	0.00	-13	20	Averaged
Anthracene	1.037	1.123	0.00	-8	20	Averaged
Fluoranthene	1.365	1.484	0.00	-9	20	Averaged
Pyrene	1.362	1.480	0.00	-9	20	Averaged
Benzo (a) Anthracene	1.317	1.360	0.00	-3	20	Averaged
Chrysene	1.197	1.249	0.00	-4	20	Averaged
Benzo (b) Fluoranthene	1.303	1.418	0.00	-9	20	Averaged
Benzo (k) Fluoranthene	1.366	1.424	0.00	-4	20	Averaged
Benzo (a) Pyrene	1.256	1.318	0.00	-5	20	Averaged
Indeno (1,2,3-c,d) Pyrene	1.502	1.570	0.00	-5	20	Averaged
Dibenz (a,h) Anthracene	1.139	1.243	0.00	-9	20	Averaged
Benzo (g,h,i) Perylene	1.207	1.334	0.00	-11	20	Averaged
Biphenyl	2.240	2.047	0.00	9	20	Averaged
2,6-Dimethylnaphthalene	1.572	1.418	0.00	10	20	Averaged
1,6,7-Trimethylnaphthalene	1.571	1.354	0.00	14	20	Averaged
Dibenzothiophene	4.622	4.193	0.00	9	20	Averaged
1-Methylphenanthrene	0.997	0.912	0.00	9	20	Averaged
Benzo (e) pyrene	1.557	1.395	0.00	10	20	Averaged
Perylene	1.493	1.264	0.00	15	20	Averaged
Dibenzofuran	2.194	2.418	0.00	-10	20	Averaged
Surrogate Standards	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / %Drift	Max%D / Drift	Curve Type
Nitrobenzene-d5	0.331	0.397	0.00	-20	20	Averaged
2-Fluorobiphenyl	1.805	2.154	0.00	-19	20	Averaged
p-Terphenyl-d14	0.775	0.898	0.00	-16	20	Averaged

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar006.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 12:52 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICAL-1 S010317D 5PPM Misc Info: 170313I001
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.415	152	39443	5.000	0.00
3)*Naphthalene-d8	(2)	4.657	136	113690	5.000	0.00
11)*Acenaphthene-d10	(3)	6.542	164	52246	5.000	0.00
17)*Phenanthrene-d10	(4)	8.138	188	176924	5.000	0.00
31)*Perylene-d12	(6)	12.826	264	181619	5.000	0.00
25)*Chrysene-d12	(5)	11.060	240	185442	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.966	82	41083	5.455	0.00
SpikedAmount		5.000		Recovery =	0.000	
7)\$2-Fluorobiphenyl	(3)	5.833	172	88758	4.706	0.00
SpikedAmount		5.000		Recovery =	0.000	
23)\$p-Terphenyl-d14	(5)	9.934	244	142825	4.970	0.00
SpikedAmount		5.000		Recovery =	0.000	
Target Compounds						
4) Naphthalene	(2)	4.677	128	121800	4.939	99
5) 2-Methylnaphthalene	(2)	5.405	142	77154	5.130	98
6) 1-Methylnaphthalene	(2)	5.518	142	73930	4.903	100
10) Acenaphthylene	(3)	6.373	152	131288	4.969	99
12) Acenaphthene	(3)	6.575	153	78418	4.946	99
15) Fluorene	(3)	7.125	166	86942	4.926	100
18) Phenanthrene	(4)	8.163	178	186241	4.960	100
19) Anthracene	(4)	8.211	178	191533	5.218	100
21) Fluoranthene	(4)	9.471	202	242537	5.021	99
22) Pyrene	(5)	9.711	202	249314	4.937	99
24) Benzo (a) Anthracene	(5)	11.042	228	240732	4.928	100
26) Chrysene	(5)	11.088	228	219861	4.953	99
27) Benzo (b) Fluoranthene	(6)	12.366	252	242576	5.126	99
28) Benzo (k) Fluoranthene	(6)	12.396	252	246593	4.969	99
30) Benzo (a) Pyrene	(6)	12.757	252	229750	5.036	99
33) Indeno (1,2,3-c,d) Pyrene	(6)	14.070	276	277787	5.093	92
34) Dibenz (a,h) Anthracene	(6)	14.094	278	210417	5.086	98
35) Benzo (g,h,i) Perylene	(6)	14.348	276	221831	5.062	99
8) Biphenyl	(3)	5.922	154	110903	4.739	99
9) 2,6-Dimethylnaphthalene	(3)	6.084	156	79570	4.844	99
14) 1,6,7-Trimethylnaphthalene	(3)	7.006	170	79803	4.863	99
16) Dibenzothiophene	(3)	8.027	184	239677	4.962	100

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar006.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 12:52 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

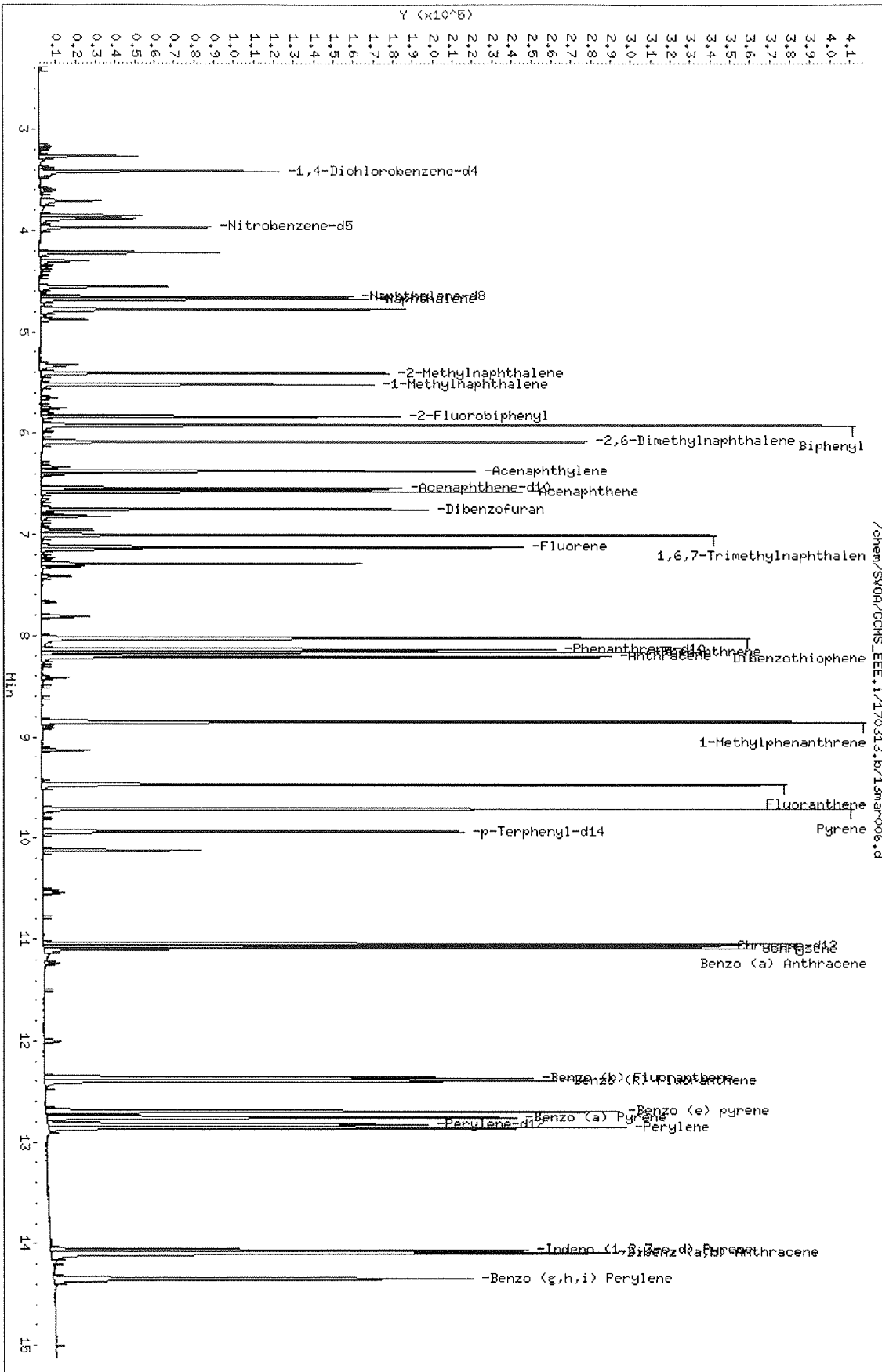
Sample Name: ICAL-1 S010317D 5PPM Misc Info: 170313I001
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.852	192	187056	5.060	98
29) Benzo (e) pyrene	(6)	12.697	252	278344	4.921	97
32) Perylene	(6)	12.859	252	266407	4.911	98
13) Dibenzofuran	(3)	6.749	168	114724	5.005	99

page 2 of 2

Data File: /chem/SW08/GCHS_EEE.i/170313.b/13mar006.d
Date: 13-MAR-2017 12:52
Client ID:
Sample Info: ICAL-1 SOL0317D SPPM
Column phase: J&M DB-SMS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



Return to Contents

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar007.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 13:13 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICAL-2 S010317E 2PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
=====						
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.415	152	43537	5.000	0.00
3)*Naphthalene-d8	(2)	4.656	136	128027	5.000	0.00
11)*Acenaphthene-d10	(3)	6.541	164	58252	5.000	0.00
17)*Phenanthrene-d10	(4)	8.138	188	195969	5.000	0.00
31)*Perylene-d12	(6)	12.826	264	200691	5.000	0.00
25)*Chrysene-d12	(5)	11.058	240	204739	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.965	82	18390	2.169	0.00
SpikedAmount 2.000				Recovery = 204.796		
7)\$2-Fluorobiphenyl	(3)	5.832	172	43331	2.061	0.00
SpikedAmount 2.000				Recovery = 201.236		
23)\$p-Terphenyl-d14	(5)	9.931	244	65918	2.078	0.00
SpikedAmount 2.000				Recovery = 204.185		
Target Compounds						
4) Naphthalene	(2)	4.676	128	57806	2.082	99
5) 2-Methylnaphthalene	(2)	5.404	142	36060	2.129	97
6) 1-Methylnaphthalene	(2)	5.518	142	34979	2.060	99
10) Acenaphthylene	(3)	6.371	152	61511	2.088	99
12) Acenaphthene	(3)	6.574	153	36834	2.084	99
15) Fluorene	(3)	7.123	166	41167	2.092	100
18) Phenanthrene	(4)	8.161	178	86634	2.083	100
19) Anthracene	(4)	8.209	178	87449	2.151	99
21) Fluoranthene	(4)	9.469	202	111729	2.088	99
22) Pyrene	(5)	9.708	202	116469	2.089	100
24) Benzo (a) Anthracene	(5)	11.039	228	109955	2.039	100
26) Chrysene	(5)	11.085	228	102627	2.094	99
27) Benzo (b) Fluoranthene	(6)	12.363	252	110812	2.119	100
28) Benzo (k) Fluoranthene	(6)	12.391	252	114315	2.085	99
30) Benzo (a) Pyrene	(6)	12.752	252	106661	2.116	100
33) Indeno (1,2,3-c,d) Pyrene	(6)	14.066	276	127377	2.113	91
34) Dibenz (a,h) Anthracene	(6)	14.093	278	98604	2.157	98
35) Benzo (g,h,i) Perylene	(6)	14.343	276	101421	2.094	100
8) Biphenyl	(3)	5.921	154	53526	2.051	99
9) 2,6-Dimethylnaphthalene	(3)	6.081	156	38072	2.079	99
14) 1,6,7-Trimethylnaphthalene	(3)	7.004	170	37953	2.074	99
16) Dibenzothiophene	(3)	8.024	184	111887	2.078	100

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar007.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 13:13 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

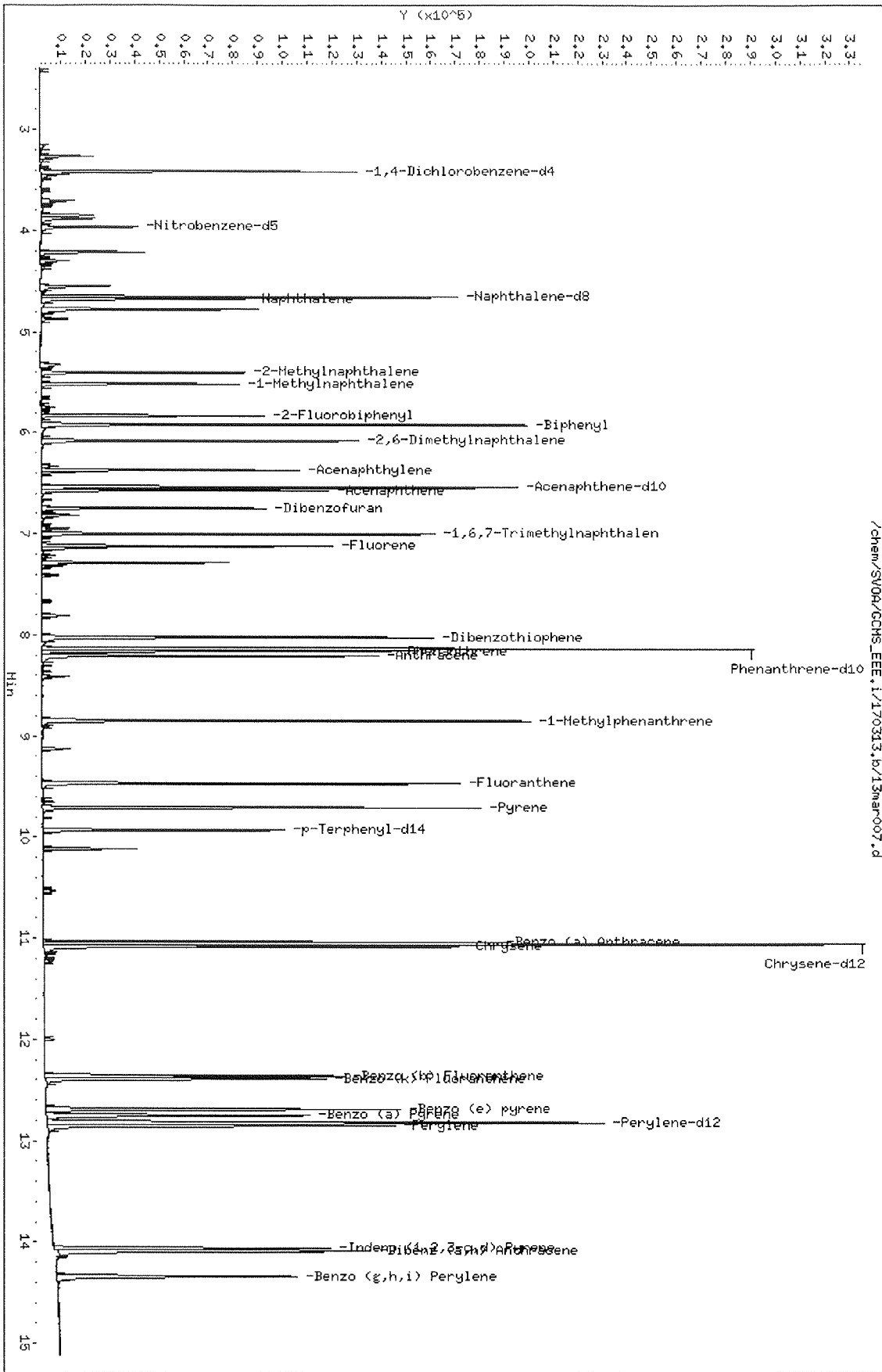
Sample Name: ICAL-2 S010317E 2PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.850	192	86212	2.112	99
29) Benzo (e) pyrene	(6)	12.694	252	128856	2.062	98
32) Perylene	(6)	12.855	252	123946	2.068	99
13) Dibenzofuran	(3)	6.749	168	53718	2.102	99

page 2 of 2

Data File: /chem/SV04/GCHS_EEE.i/170313.b/13mar007.d
Date: 13-MAR-2017 13:13
Client ID:
Sample Info: ICAI-2 SOL0317E 2PPM
Column phase: J&M DB-SMS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



[Return to Contents](#)

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar008.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 13:33 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICAL-3 S010317F 1PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.414	152	46489	5.000	0.01
3)*Naphthalene-d8	(2)	4.656	136	140022	5.000	0.00
11)*Acenaphthene-d10	(3)	6.541	164	63153	5.000	0.00
17)*Phenanthrene-d10	(4)	8.136	188	210133	5.000	0.00
31)*Perylene-d12	(6)	12.823	264	208071	5.000	0.00
25)*Chrysene-d12	(5)	11.056	240	220711	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.966	82	9500	1.024	0.00
SpikedAmount 1.000				Recovery = 97.033		
7)\$2-Fluorobiphenyl	(3)	5.833	172	23775	1.043	0.00
SpikedAmount 1.000				Recovery = 100.830		
23)\$p-Terphenyl-d14	(5)	9.932	244	35159	1.028	0.00
SpikedAmount 1.000				Recovery = 101.689		
Target Compounds						
4) Naphthalene	(2)	4.676	128	30813	1.015	100
5) 2-Methylnaphthalene	(2)	5.404	142	18891	1.020	100
6) 1-Methylnaphthalene	(2)	5.517	142	18774	1.011	100
10) Acenaphthylene	(3)	6.371	152	32530	1.019	100
12) Acenaphthene	(3)	6.574	153	19598	1.023	100
15) Fluorene	(3)	7.123	166	22113	1.037	100
18) Phenanthrene	(4)	8.159	178	45683	1.024	100
19) Anthracene	(4)	8.211	178	40718	0.934	100
21) Fluoranthene	(4)	9.470	202	58439	1.019	100
22) Pyrene	(5)	9.707	202	61511	1.023	100
24) Benzo (a) Anthracene	(5)	11.040	228	57660	0.992	100
26) Chrysene	(5)	11.084	228	53861	1.020	100
27) Benzo (b) Fluoranthene	(6)	12.362	252	58061	1.071	100
28) Benzo (k) Fluoranthene	(6)	12.391	252	59318	1.043	100
30) Benzo (a) Pyrene	(6)	12.750	252	52756	1.009	100
33) Indeno (1,2,3-c,d) Pyrene	(6)	14.064	276	65862	1.054	100
34) Dibenz (a,h) Anthracene	(6)	14.094	278	51150	1.079	100
35) Benzo (g,h,i) Perylene	(6)	14.342	276	52489	1.045	100
8) Biphenyl	(3)	5.922	154	29053	1.027	100
9) 2,6-Dimethylnaphthalene	(3)	6.081	156	20301	1.022	100
14) 1,6,7-Trimethylnaphthalene	(3)	7.003	170	20183	1.017	100
16) Dibenzothiophene	(3)	8.024	184	59602	1.021	100

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar008.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 13:33 Analyst ID: 907

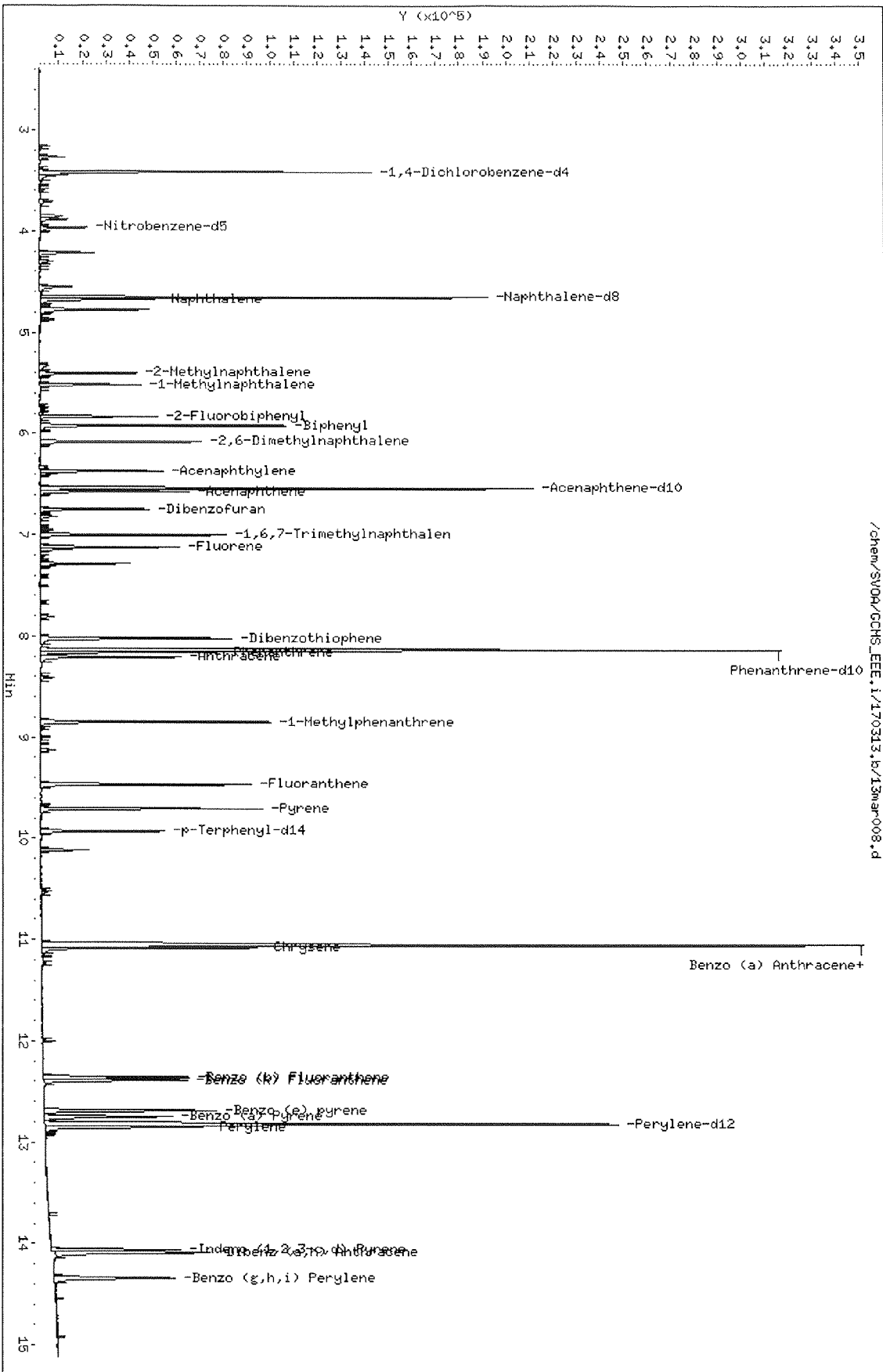
Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICAL-3 S010317F 1PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.850	192	45057	1.024	100
29) Benzo (e) pyrene	(6)	12.690	252	68218	1.053	100
32) Perylene	(6)	12.854	252	62786	1.010	100
13) Dibenzofuran	(3)	6.749	168	28575	1.031	100

Data File: /chem/SV09/GCHS_EEE.i/170313.b/13mar008.d
Date: 13-MAR-2017 13:33
Client ID:
Sample Info: ICAL-3 5010317F 1PPH
Column phase: J&W DB-SMS

Instrument: GCHS_EEE.1
Operator: 907
Column diameter: 0.18



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar009.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 13:53 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICAL-4 S010317G 0.5PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.414	152	46816	5.000	0.01
3)*Naphthalene-d8	(2)	4.654	136	141612	5.000	0.00
11)*Acenaphthene-d10	(3)	6.540	164	65116	5.000	0.00
17)*Phenanthrene-d10	(4)	8.136	188	215147	5.000	0.00
31)*Perylene-d12	(6)	12.825	264	224201	5.000	0.00
25)*Chrysene-d12	(5)	11.057	240	226005	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.967	82	4177	0.445	0.00
SpikedAmount 0.500			Recovery = 43.133			
7)\$2-Fluorobiphenyl	(3)	5.831	172	11066	0.471	0.00
SpikedAmount 0.500			Recovery = 45.627			
23)\$p-Terphenyl-d14	(5)	9.931	244	16152	0.461	0.00
SpikedAmount 0.500			Recovery = 45.890			
Target Compounds						
4) Naphthalene	(2)	4.676	128	14251	0.464	98
5) 2-Methylnaphthalene	(2)	5.404	142	8470	0.452	96
6) 1-Methylnaphthalene	(2)	5.516	142	8559	0.456	99
10) Acenaphthylene	(3)	6.371	152	15162	0.460	99
12) Acenaphthene	(3)	6.574	153	9204	0.466	99
15) Fluorene	(3)	7.123	166	10327	0.469	99
18) Phenanthrene	(4)	8.160	178	21102	0.462	99
19) Anthracene	(4)	8.209	178	21165	0.474	99
21) Fluoranthene	(4)	9.469	202	27157	0.462	99
22) Pyrene	(5)	9.707	202	28490	0.463	99
24) Benzo (a) Anthracene	(5)	11.040	228	27292	0.458	99
26) Chrysene	(5)	11.085	228	25041	0.463	99
27) Benzo (b) Fluoranthene	(6)	12.361	252	26512	0.454	99
28) Benzo (k) Fluoranthene	(6)	12.393	252	28262	0.461	98
30) Benzo (a) Pyrene	(6)	12.752	252	26139	0.464	99
33) Indeno (1,2,3-c,d) Pyrene	(6)	14.065	276	29971	0.445	92
34) Dibenz (a,h) Anthracene	(6)	14.093	278	23700	0.464	99
35) Benzo (g,h,i) Perylene	(6)	14.344	276	24422	0.451	100
8) Biphenyl	(3)	5.921	154	13800	0.473	100
9) 2,6-Dimethylnaphthalene	(3)	6.081	156	9629	0.470	100
14) 1,6,7-Trimethylnaphthalene	(3)	7.001	170	9520	0.465	99
16) Dibenzothiophene	(3)	8.024	184	27839	0.462	99

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar009.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 13:53 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

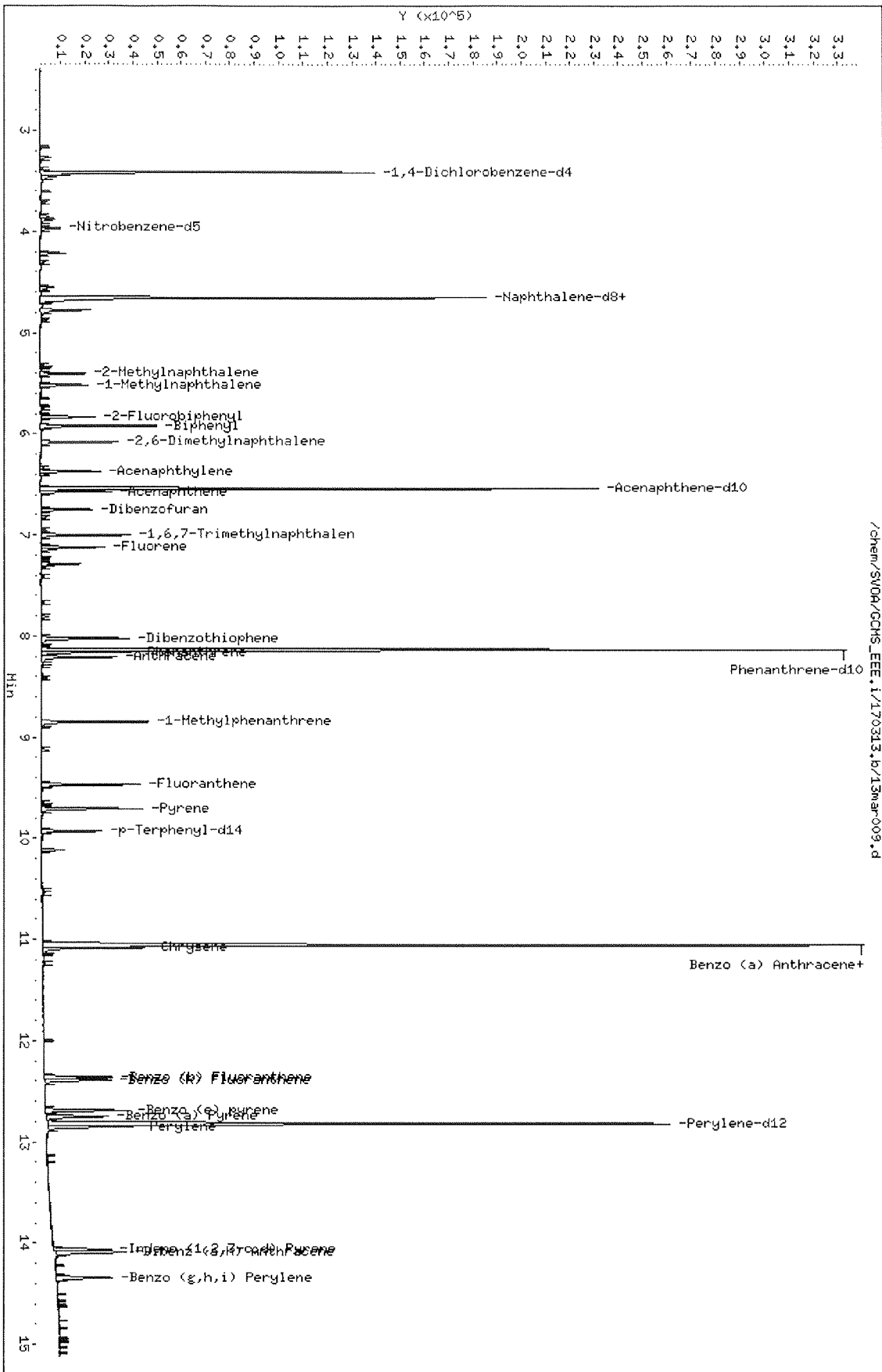
Sample Name: ICAL-4 S010317G 0.5PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.850	192	21018	0.467	99
29) Benzo (e) pyrene	(6)	12.691	252	31599	0.453	99
32) Perylene	(6)	12.853	252	31630	0.472	100
13) Dibenzofuran	(3)	6.749	168	13308	0.466	100

page 2 of 2

Data File: /chem/SW04/GCHS_EEE.i/170313.b/13mar009.d
 Date : 13-MAR-2017 13:53
 Client ID:
 Sample Info: ICAL-4 5010317C 0.5PPM
 Column phase: J&W DB-SMS

Instrument: GCHS_EEE.1
 Operator: 907
 Column diameter: 0.18



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar010.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 14:13 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICAL-5 S010317H 0.1PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.414	152	46155	5.000	0.01
3)*Naphthalene-d8	(2)	4.655	136	138733	5.000	0.00
11)*Acenaphthene-d10	(3)	6.539	164	66109	5.000	0.00
17)*Phenanthrene-d10	(4)	8.135	188	215586	5.000	0.00
31)*Perylene-d12	(6)	12.823	264	222988	5.000	0.00
25)*Chrysene-d12	(5)	11.057	240	227975	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.971	82	836	0.091	-0.01
SpikedAmount 0.100			Recovery = 8.992			
7)\$2-Fluorobiphenyl	(3)	5.833	172	2492	0.104	0.00
SpikedAmount 0.100			Recovery = 10.221			
23)\$p-Terphenyl-d14	(5)	9.933	244	3593	0.102	0.00
SpikedAmount 0.100			Recovery = 10.183			
Target Compounds						
4) Naphthalene	(2)	4.676	128	3096	0.103	100
5) 2-Methylnaphthalene	(2)	5.404	142	1809	0.099	97
6) 1-Methylnaphthalene	(2)	5.517	142	1963	0.107	96
10) Acenaphthylene	(3)	6.371	152	3418	0.102	99
12) Acenaphthene	(3)	6.571	153	2035	0.101	97
15) Fluorene	(3)	7.124	166	2218	0.099	99
18) Phenanthrene	(4)	8.162	178	4655	0.102	98
19) Anthracene	(4)	8.211	178	4465	0.100	100
21) Fluoranthene	(4)	9.471	202	5933	0.101	100
22) Pyrene	(5)	9.708	202	6324	0.102	99
24) Benzo (a) Anthracene	(5)	11.039	228	6527	0.109	97
26) Chrysene	(5)	11.084	228	5549	0.102	99
27) Benzo (b) Fluoranthene	(6)	12.361	252	5444	0.094	93
28) Benzo (k) Fluoranthene	(6)	12.393	252	6080	0.100	91
30) Benzo (a) Pyrene	(6)	12.752	252	5588	0.100	93
33) Indeno (1,2,3-c,d) Pyrene	(6)	14.068	276	6565	0.098	81
34) Dibenz (a,h) Anthracene	(6)	14.096	278	4557	0.090	91
35) Benzo (g,h,i) Perylene	(6)	14.346	276	5340	0.099	97
8) Biphenyl	(3)	5.922	154	3118	0.105	90
9) 2,6-Dimethylnaphthalene	(3)	6.081	156	2138	0.103	99
14) 1,6,7-Trimethylnaphthalene	(3)	7.003	170	2164	0.104	100
16) Dibenzothiophene	(3)	8.024	184	6252	0.102	98

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar010.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 14:13 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

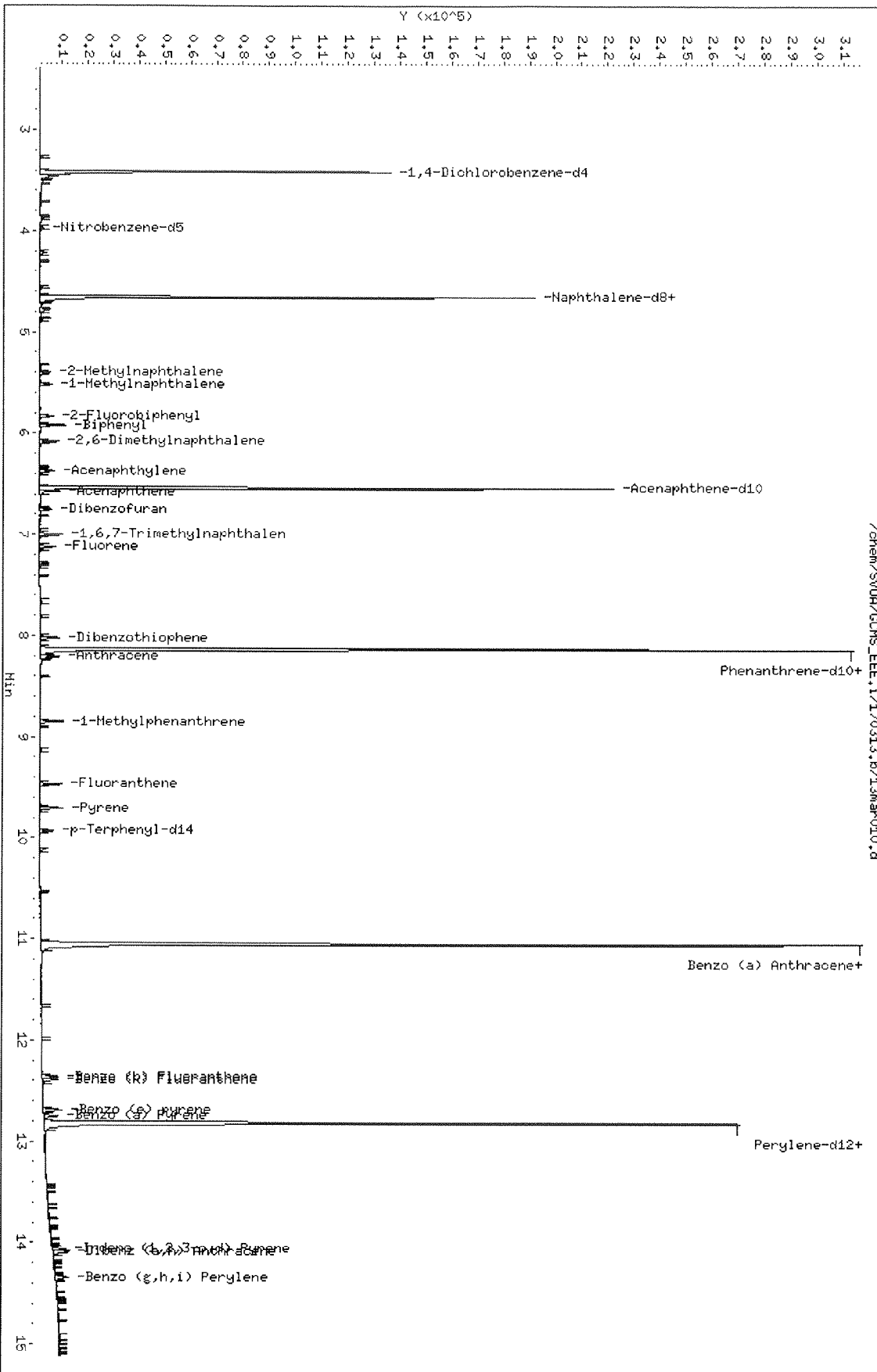
Sample Name: ICAL-5 S010317H 0.1PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.850	192	4428	0.097	98
29) Benzo (e) pyrene	(6)	12.691	252	7135	0.103	99
32) Perylene	(6)	12.853	252	6852	0.103	98
13) Dibenzofuran	(3)	6.748	168	2858	0.099	98

page 2 of 2

Date File: /chem/SV04/GCHS_EEE.i/170313.b/13mar010.d
Date : 13-MAR-2017 14:13
Client ID:
Sample Info: ICAL-5 S010317H 0.1PPM
Column phase: J&M DB-SMS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar011.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 14:34 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

Sample Name: ICV S010317I 1PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.419	152	44664	5.000	0.00
3)*Naphthalene-d8	(2)	4.653	136	134376	5.000	0.00
11)*Acenaphthene-d10	(3)	6.538	164	61082	5.000	0.00
17)*Phenanthrene-d10	(4)	8.136	188	200519	5.000	0.00
31)*Perylene-d12	(6)	12.823	264	200860	5.000	0.00
25)*Chrysene-d12	(5)	11.055	240	207148	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.965	82	10658	1.197	0.00
SpikedAmount 1.000			Recovery =	0.000		
7)\$2-Fluorobiphenyl	(3)	5.831	172	26320	1.194	0.00
SpikedAmount 1.000			Recovery =	0.000		
23)\$p-Terphenyl-d14	(5)	9.929	244	37191	1.159	0.00
SpikedAmount 1.000			Recovery =	0.000		
Target Compounds						
4) Naphthalene	(2)	4.673	128	30493	1.046	100
5) 2-Methylnaphthalene	(2)	5.402	142	21043	1.184	100
6) 1-Methylnaphthalene	(2)	5.516	142	18763	1.053	100
10) Acenaphthylene	(3)	6.370	152	31506	1.020	100
12) Acenaphthene	(3)	6.572	153	20057	1.082	100
15) Fluorene	(3)	7.122	166	22387	1.085	100
18) Phenanthrene	(4)	8.160	178	48240	1.134	100
19) Anthracene	(4)	8.209	178	45040	1.083	100
21) Fluoranthene	(4)	9.469	202	59504	1.087	100
22) Pyrene	(5)	9.707	202	61311	1.087	100
24) Benzo (a) Anthracene	(5)	11.038	228	56357	1.033	100
26) Chrysene	(5)	11.083	228	51734	1.043	100
27) Benzo (b) Fluoranthene	(6)	12.360	252	56960	1.088	100
28) Benzo (k) Fluoranthene	(6)	12.391	252	57224	1.043	100
30) Benzo (a) Pyrene	(6)	12.752	252	52949	1.049	100
33) Indeno (1,2,3-c,d) Pyrene	(6)	14.065	276	63053	1.045	100
34) Dibenz (a,h) Anthracene	(6)	14.091	278	49952	1.092	100
35) Benzo (g,h,i) Perylene	(6)	14.339	276	53577	1.105	100
8) Biphenyl	(3)	5.919	154	25012	0.914	100
9) 2,6-Dimethylnaphthalene	(3)	6.079	156	17327	0.902	100
14) 1,6,7-Trimethylnaphthalene	(3)	7.001	170	16542	0.862	100
16) Dibenzothiophene	(3)	8.024	184	51223	0.907	100

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170313.b/13mar011.d Instrument ID: GCMS_EEE.i
 Injection date and time: 13-MAR-2017 14:34 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170313.b/simpah-extra.m Sublist used: all
 Calibration date and time: 13-MAR-2017 14:59
 Date, time and analyst ID of latest file update: 13-Mar-2017 14:59 ev7p

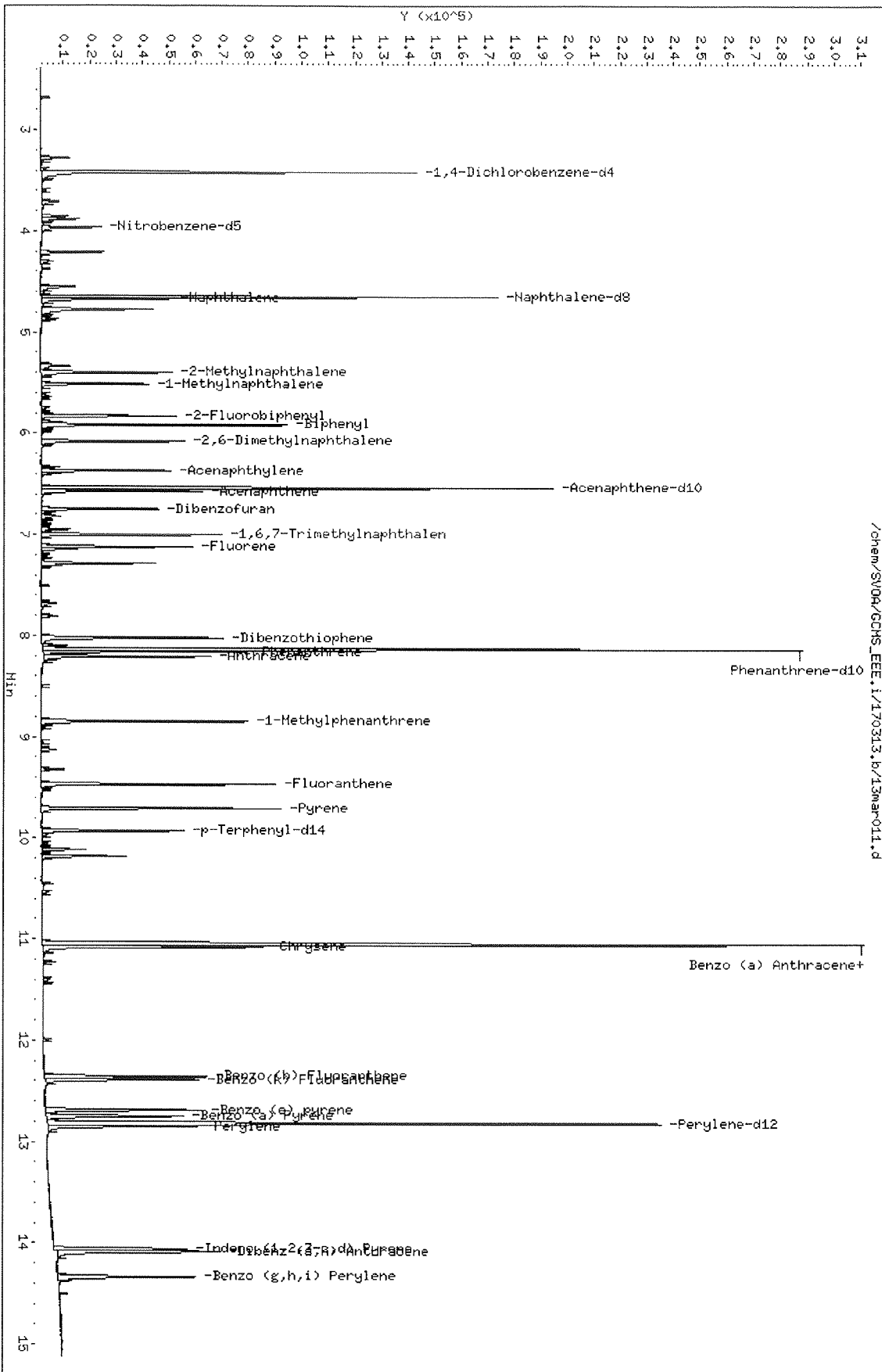
Sample Name: ICV S010317I 1PPM Misc Info:
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.848	192	37778	0.915	100
29) Benzo (e) pyrene	(6)	12.691	252	56048	0.896	100
32) Perylene	(6)	12.851	252	50780	0.846	100
13) Dibenzofuran	(3)	6.748	168	29537	1.102	100

page 2 of 2

Date File: /chem/SV04/GCHS_EEE.i/170313.b/13mar011.d
Date : 13-MAR-2017 14:34
Client ID:
Sample Info: ICV S0103171 1PPH
Column phase: J&M DB-SMS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



EPA METHOD 8270C PAHSIM

Sample Data

RAW DATA SHEET FOR METHOD: EPA 8270C SIM PAHs

WORK ORDER: 17-03-1557
INSTRUMENT: GC/MS EEE
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED: 2017-03-27 13:57
REVIEWED BY:
D/T REVIEWED:

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar011.d\27mar011.rr

25 **CLIENT SAMPLE NUMBER: B-DU1-ISM1-8**

LCS/MB BATCH: 170324L17 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 20.00 g / ACTUAL: 20.00 g
MS/MSD BATCH: 170324S17 **FINAL VOLUME / WEIGHT:** DEFAULT: 2.00 ml / ACTUAL: 2.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Naphthalene	0.000	1.00	ND	0.010	
2-Methylnaphthalene	0.000	1.00	ND	0.010	
1-Methylnaphthalene	0.000	1.00	ND	0.010	
Acenaphthylene	0.000	1.00	ND	0.010	
Acenaphthene	0.000	1.00	ND	0.010	
Fluorene	0.000	1.00	ND	0.010	
Phenanthrene	0.000	1.00	ND	0.010	
Anthracene	0.000	1.00	ND	0.010	
Fluoranthene	0.0162	1.00	ND	0.010	
Pyrene	0.0216	1.00	ND	0.010	
Benzo (a) Anthracene	0.0243	1.00	ND	0.010	
Chrysene	0.0271	1.00	ND	0.010	
Benzo (k) Fluoranthene	0.0226	1.00	ND	0.010	
Benzo (b) Fluoranthene	0.0236	1.00	ND	0.010	
Benzo (a) Pyrene	0.0230	1.00	ND	0.010	
Indeno (1,2,3-c,d) Pyrene	0.0308	1.00	ND	0.010	
Dibenz (a,h) Anthracene	0.000	1.00	ND	0.010	
Benzo (g,h,i) Perylene	0.0475	1.00	ND	0.010	

Return to Contents

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar011.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 13:57 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 14:38 ev7p

Sample Name: 17-03-1557-25 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV (Min)
=====						
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.302	152	59986	5.000	0.00
3)*Naphthalene-d8	(2)	4.535	136	169237	5.000	0.00
11)*Acenaphthene-d10	(3)	6.418	164	80719	5.000	0.00
17)*Phenanthrene-d10	(4)	8.017	188	267245	5.000	0.00
31)*Perylene-d12	(6)	12.679	264	254564	5.000	0.01
25)*Chrysene-d12	(5)	10.934	240	260860	5.000	0.01
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.854	82	8826	0.787	0.00
SpikedAmount 1.000				Recovery = 78.735		
7)\$2-Fluorobiphenyl	(3)	5.719	172	24177	0.830	0.00
SpikedAmount 1.000				Recovery = 82.970		
23)\$p-Terphenyl-d14	(5)	9.810	244	37848	0.936	0.01
SpikedAmount 1.000				Recovery = 93.625		
Target Compounds						
						QValue
4) Naphthalene	(2)	0.000		0		N.D.
5) 2-Methylnaphthalene	(2)	0.000		0		N.D.
6) 1-Methylnaphthalene	(2)	0.000		0		N.D.
10) Acenaphthylene	(3)	0.000		0		N.D.
12) Acenaphthene	(3)	0.000		0		N.D.
15) Fluorene	(3)	0.000		0		N.D.
18) Phenanthrene	(4)	0.000		0		N.D.
19) Anthracene	(4)	0.000		0		N.D.
21) Fluoranthene	(4)	9.353	202	1181	0.016	33
22) Pyrene	(5)	9.592	202	1533	0.022	82
24) Benzo (a) Anthracene	(5)	10.918	228	1669	0.024	82
26) Chrysene	(5)	10.961	228	1694	0.027	96
27) Benzo (b) Fluoranthene	(6)	12.229	252	1567	0.024	96
28) Benzo (k) Fluoranthene	(6)	12.240	252	1573A	0.023	86
30) Benzo (a) Pyrene	(6)	12.613	252	1472	0.023	84
33) Indeno (1,2,3-c,d) Pyrene	(6)	13.931	276	2351	0.031	68
34) Dibenz (a,h) Anthracene	(6)	0.000		0		N.D.
35) Benzo (g,h,i) Perylene	(6)	14.204	276	2919	0.048	83
8) Biphenyl	(3)	0.000		0		N.D.
9) 2,6-Dimethylnaphthalene	(3)	0.000		0		N.D.
14) 1,6,7-Trimethylnaphthalene	(3)	0.000		0		N.D.
16) Dibenzothiophene	(3)	0.000		0		N.D.

A = User selected an alternate hit.
 * = Compound is an internal standard.
 \$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar011.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 13:57 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 14:38 ev7p

Sample Name: 17-03-1557-25 Misc Info: S032317A 10UL
 Response via Initial Calibration

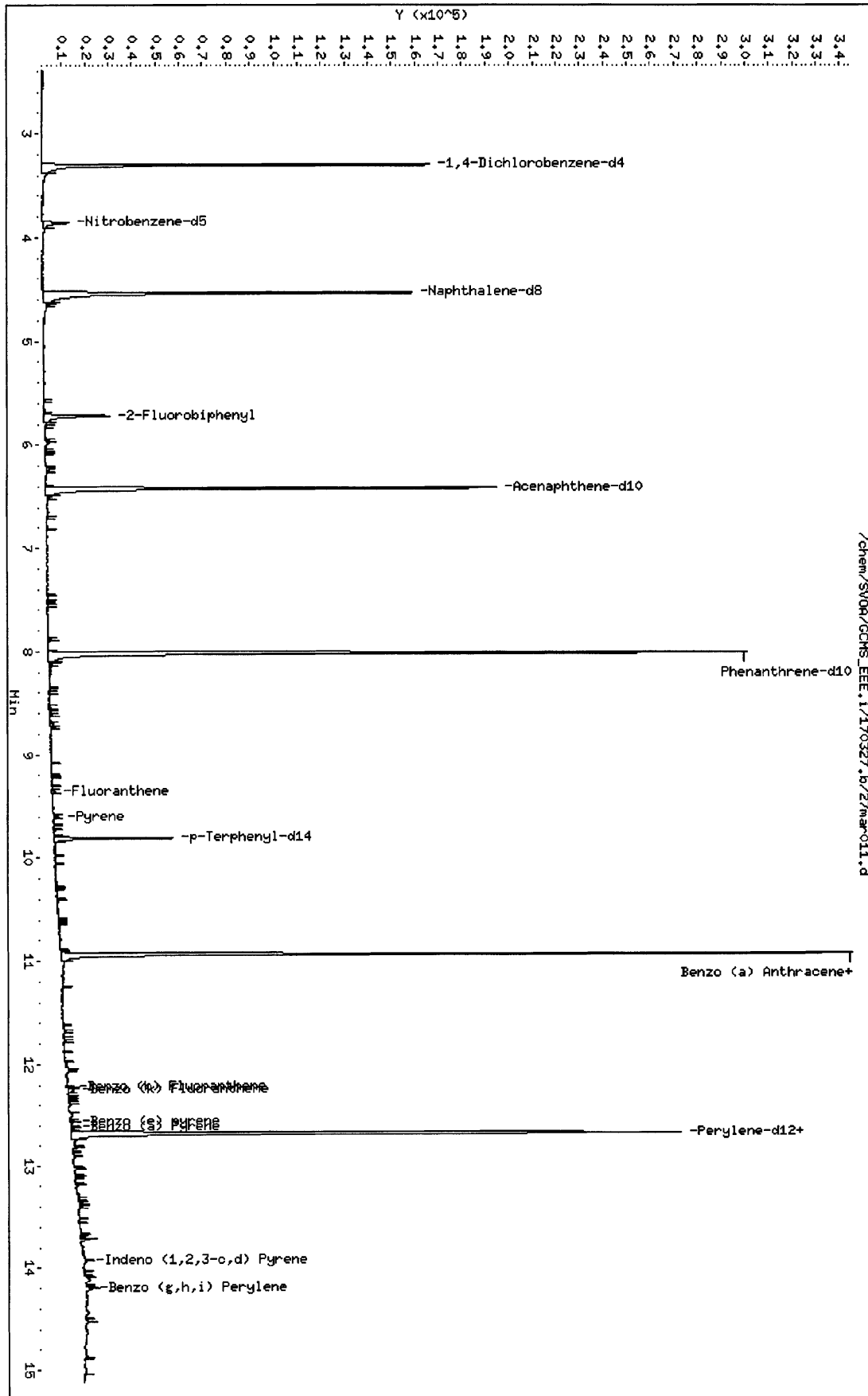
Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	0.000		0	N.D.	
29) Benzo (e) pyrene	(6)	12.548	252	1587	0.020	84
32) Perylene	(6)	12.709	252	1305A	0.017	95
13) Dibenzofuran	(3)	0.000		0	N.D.	

A = User selected an alternate hit.

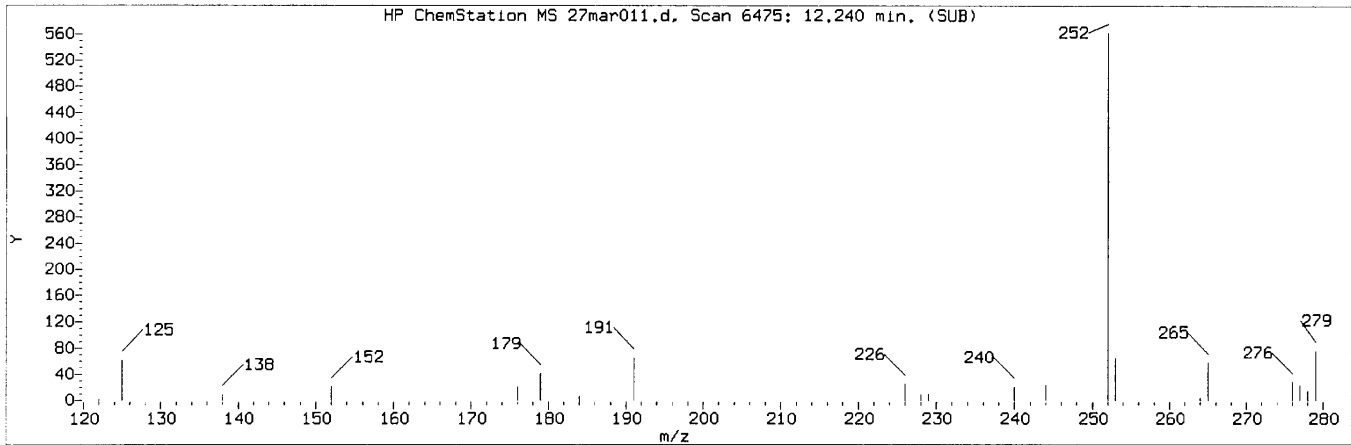
page 2 of 2

Data File: /chem/SVOR/GCHS_EEE.i/170327.b/27mar011.d
Date: 27-MAR-2017 13:57
Client ID:
Sample Info: 17-03-1557-25
Column phase: 3M DB-SMS

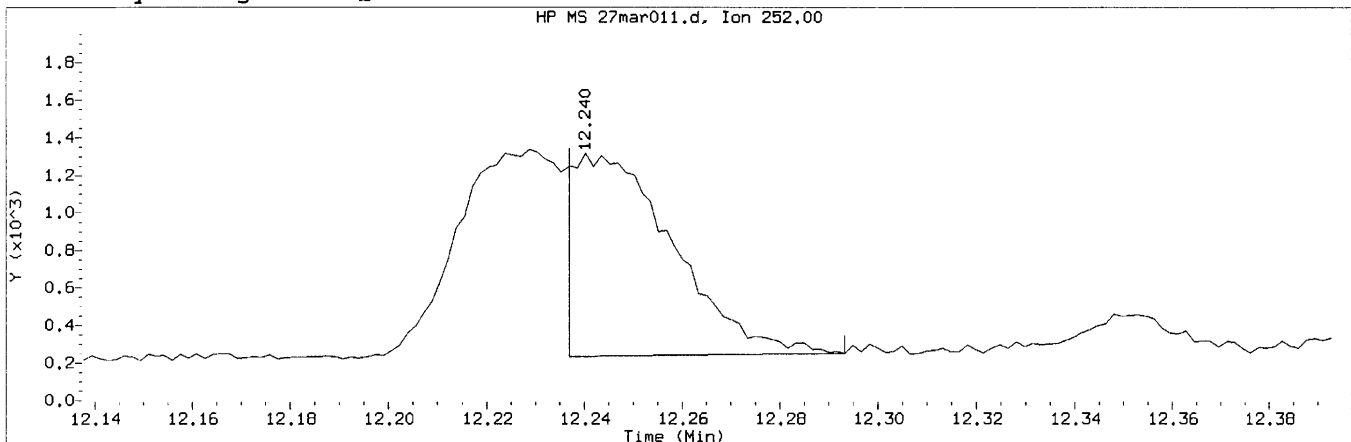
Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



Sample Spectrum (Background Subtracted)



Manually Integrated Quant Ion



Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar011.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 13:57 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 14:38 ev7p

Sample Name: 17-03-1557-25

Compound Number : 28
 Compound Name : Benzo (k) Fluoranthene
 Scan Number : 6475
 Retention Time (minutes): 12.240
 Quant Ion : 252.00
 Area (flag) : 1573A
 On-Column Amount (mg/L) : 0.0226
 Integration start scan : 6472 Integration stop scan: 6506
 Y at integration start : 234 Y at integration end: 250

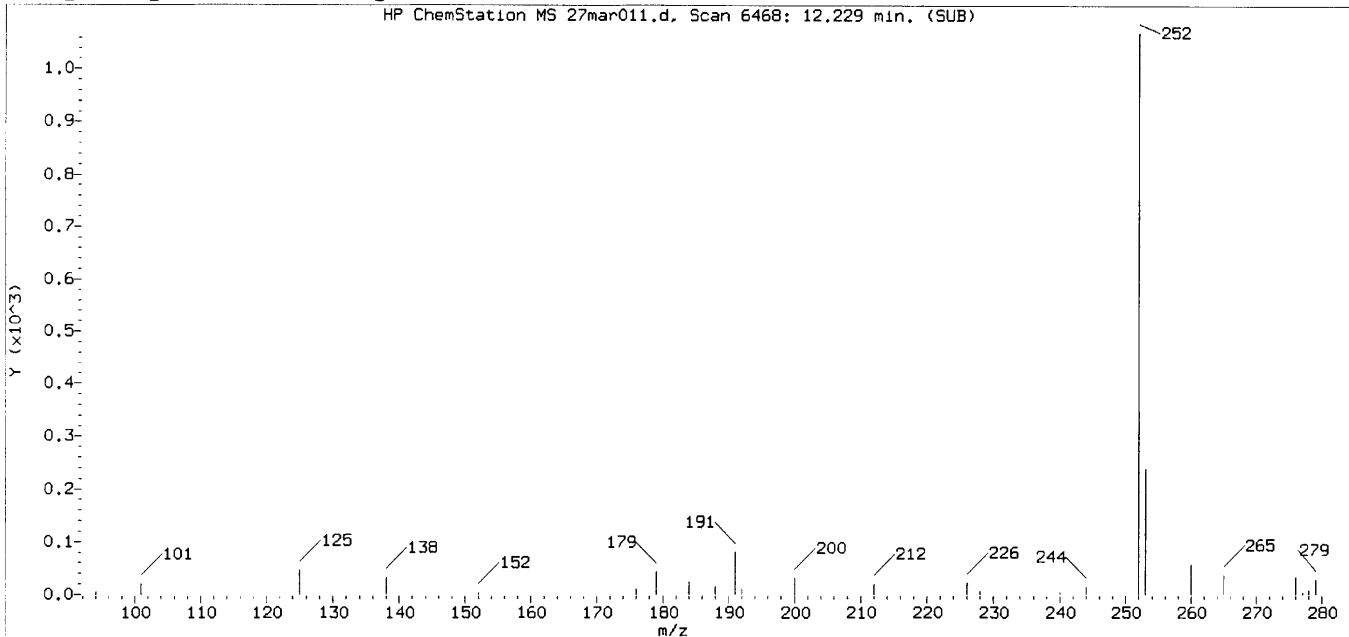
Reason for manual integration: improper integration

Digitally signed by Jeremy Huynh
 Analyst responsible for change: on 03/28/2017 at 09:12.
 Target 3.5 esignature user ID: ev7p

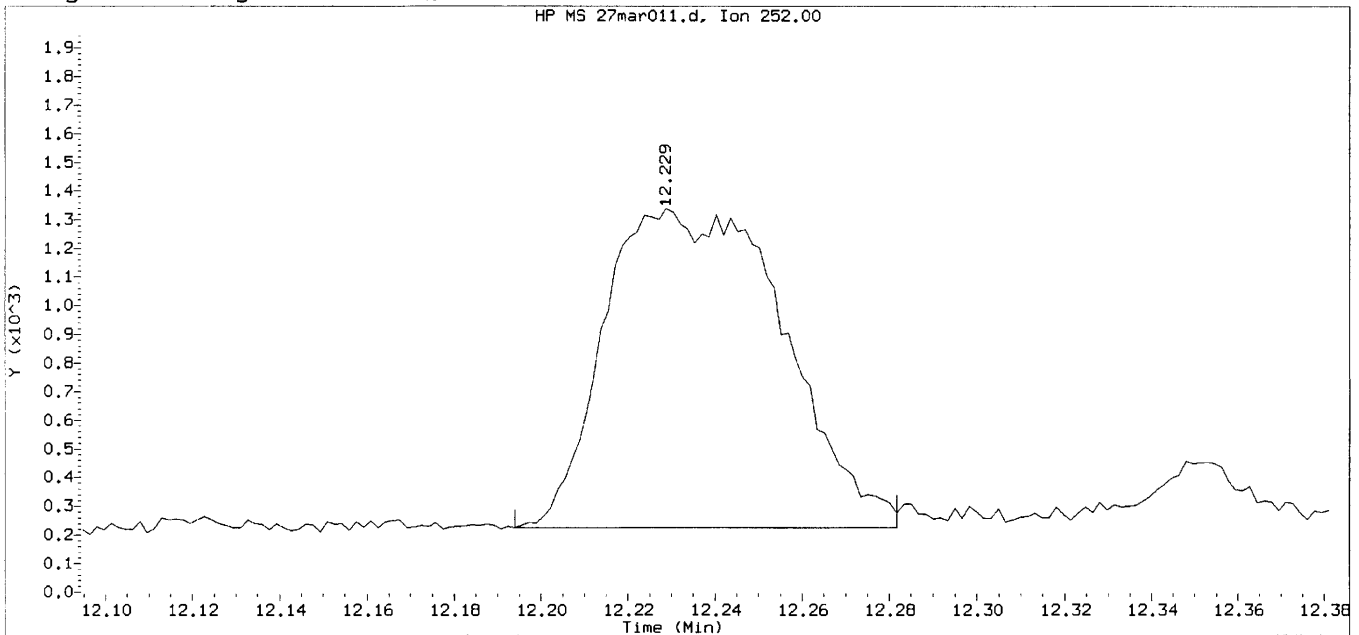
GC/MS audit/management approval: _____

Handwritten signature and date: 3/28/17

Sample Spectrum (Background Subtracted)



Original Integration of Quant Ion



Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar011.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 13:57 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 14:20 Unknown

Sample Name: 17-03-1557-25

Compound Number	: 28	
Compound Name	: Benzo (k) Fluoranthene	
Scan Number	: 6468	
Retention Time (minutes)	: 12.229	
Quant Ion	: 252.00	
Area	: 3182	
On-column Amount (mg/L)	: 0.0458	
Integration start scan	: 6446	Integration stop scan: 6499
Y at integration start	: 224	Y at integration end: 224

**RAW DATA SHEET
FOR METHOD: EPA 8270C SIM PAHs**

WORK ORDER: 17-03-1557
INSTRUMENT: GC/MS EEE
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED: 2017-03-27 14:17
REVIEWED BY:
D/T REVIEWED:

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar012.d\27mar012.rr

26 **CLIENT SAMPLE NUMBER: B-DU1-ISM2-8**

LCS/MB BATCH: 170324L17 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 20.00 g / ACTUAL: 20.20 g
MS/MSD BATCH: 170324S17 **FINAL VOLUME / WEIGHT:** DEFAULT: 2.00 ml / ACTUAL: 2.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 0.99

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Naphthalene	0.000	1.00	ND	0.0099	
2-Methylnaphthalene	0.000	1.00	ND	0.0099	
1-Methylnaphthalene	0.000	1.00	ND	0.0099	
Acenaphthylene	0.000	1.00	ND	0.0099	
Acenaphthene	0.000	1.00	ND	0.0099	
Fluorene	0.000	1.00	ND	0.0099	
Phenanthrene	0.000	1.00	ND	0.0099	
Anthracene	0.000	1.00	ND	0.0099	
Fluoranthene	0.0212	1.00	ND	0.0099	
Pyrene	0.0249	1.00	ND	0.0099	
Benzo (a) Anthracene	0.0285	1.00	ND	0.0099	
Chrysene	0.0231	1.00	ND	0.0099	
Benzo (k) Fluoranthene	0.000	1.00	ND	0.0099	
Benzo (b) Fluoranthene	0.000	1.00	ND	0.0099	
Benzo (a) Pyrene	0.000	1.00	ND	0.0099	
Indeno (1,2,3-c,d) Pyrene	0.000	1.00	ND	0.0099	
Dibenz (a,h) Anthracene	0.000	1.00	ND	0.0099	
Benzo (g,h,i) Perylene	0.000	1.00	ND	0.0099	

Return to Contents

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar012.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:17 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 16:17 ev7p

Sample Name: 17-03-1557-26 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S.			Area	On-Column	
	Ref.	RT	QIon		Amount (mg/L)	DEV (Min)
=====						
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.301	152	60589	5.000	0.00
3)*Naphthalene-d8	(2)	4.534	136	171804	5.000	0.00
11)*Acenaphthene-d10	(3)	6.418	164	81274	5.000	0.00
17)*Phenanthrene-d10	(4)	8.017	188	274687	5.000	0.00
31)*Perylene-d12	(6)	12.683	264	271055	5.000	0.00
25)*Chrysene-d12	(5)	10.936	240	279862	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.854	82	9817	0.863	0.00
SpikedAmount 1.000	Recovery = 86.270					
7)\$2-Fluorobiphenyl	(3)	5.716	172	26564	0.905	0.01
SpikedAmount 1.000	Recovery = 90.539					
23)\$p-Terphenyl-d14	(5)	9.812	244	42475	0.979	0.01
SpikedAmount 1.000	Recovery = 97.937					
Target Compounds						
						QValue
4) Naphthalene	(2)	0.000		0		N.D.
5) 2-Methylnaphthalene	(2)	0.000		0		N.D.
6) 1-Methylnaphthalene	(2)	0.000		0		N.D.
10) Acenaphthylene	(3)	0.000		0		N.D.
12) Acenaphthene	(3)	0.000		0		N.D.
15) Fluorene	(3)	0.000		0		N.D.
18) Phenanthrene	(4)	0.000		0		N.D.
19) Anthracene	(4)	0.000		0		N.D.
21) Fluoranthene	(4)	9.352	202	1590	0.021	94
22) Pyrene	(5)	9.590	202	1898	0.025	93
24) Benzo (a) Anthracene	(5)	10.923	228	2097	0.028	76
26) Chrysene	(5)	10.963	228	1549	0.023	86
27) Benzo (b) Fluoranthene	(6)	0.000		0		N.D.
28) Benzo (k) Fluoranthene	(6)	0.000		0		N.D.
30) Benzo (a) Pyrene	(6)	0.000		0		N.D.
33) Indeno (1,2,3-c,d) Pyrene	(6)	0.000		0		N.D.
34) Dibenz (a,h) Anthracene	(6)	0.000		0		N.D.
35) Benzo (g,h,i) Perylene	(6)	0.000		0		N.D.
8) Biphenyl	(3)	0.000		0		N.D.
9) 2,6-Dimethylnaphthalene	(3)	0.000		0		N.D.
14) 1,6,7-Trimethylnaphthalene	(3)	0.000		0		N.D.
16) Dibenzothiophene	(3)	0.000		0		N.D.

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar012.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:17 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 16:17 ev7p

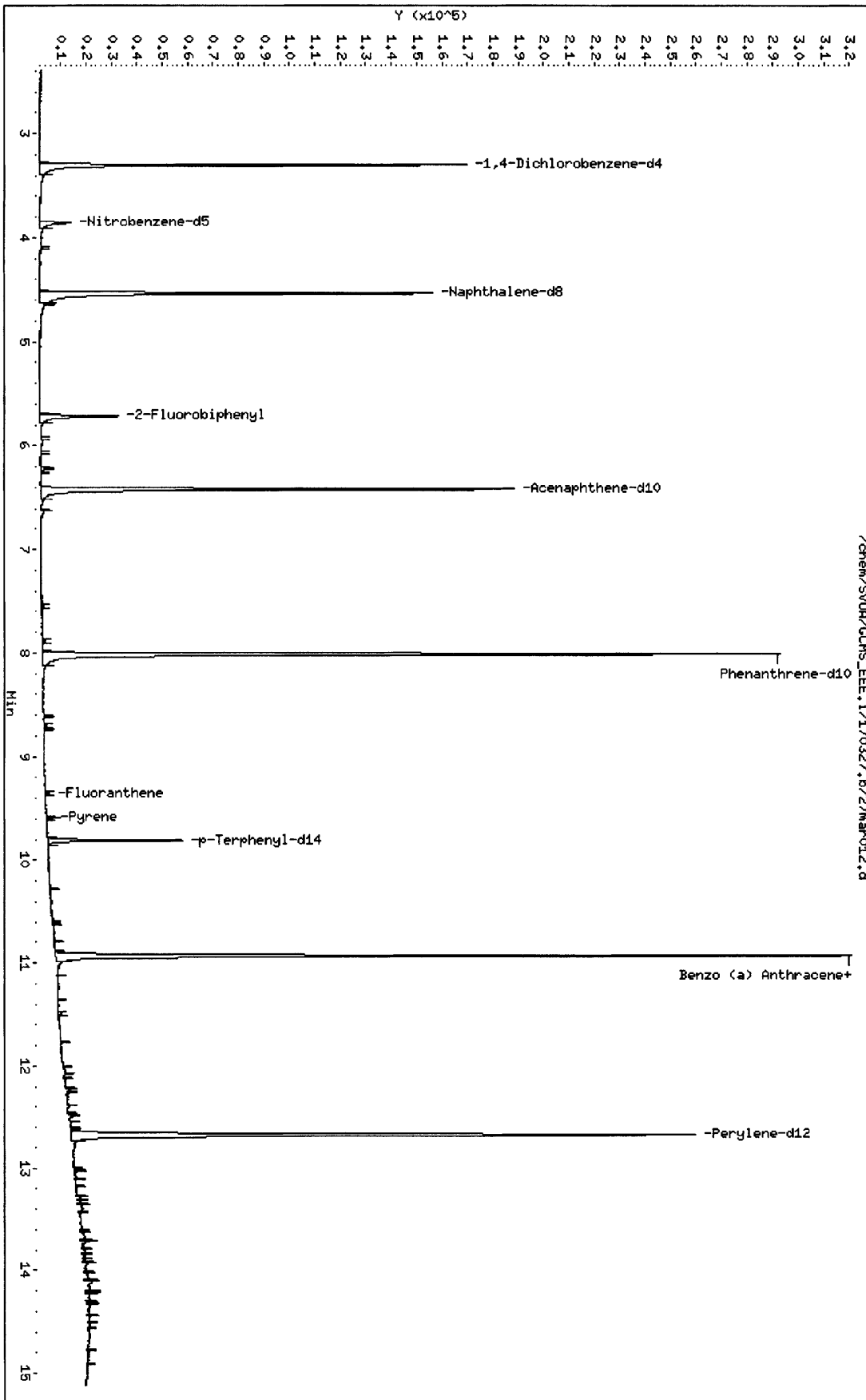
Sample Name: 17-03-1557-26 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	0.000		0	N.D.	
29) Benzo (e) pyrene	(6)	0.000		0	N.D.	
32) Perylene	(6)	0.000		0	N.D.	
13) Dibenzofuran	(3)	0.000		0	N.D.	

page 2 of 2

Data File: /chem/SV09/GCHS_EEE.i/170327.b/27mar012.d
Date : 27-MAR-2017 14:17
Client ID:
Sample Info: 17-03-1557-26
Column phase: J&M DB-SMS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



**RAW DATA SHEET
FOR METHOD: EPA 8270C SIM PAHs**

WORK ORDER: 17-03-1557
INSTRUMENT: GC/MS EEE
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED: 2017-03-27 14:38
REVIEWED BY:
D/T REVIEWED:

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar013.d\27mar013.rr

27 **CLIENT SAMPLE NUMBER: B-DU1-ISM3-8**

LCS/MB BATCH: 170324L17 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 20.00 g / ACTUAL: 20.20 g
MS/MSD BATCH: 170324S17 **FINAL VOLUME / WEIGHT:** DEFAULT: 2.00 ml / ACTUAL: 2.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 0.99

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Naphthalene	0.000	1.00	ND	0.0099	
2-Methylnaphthalene	0.000	1.00	ND	0.0099	
1-Methylnaphthalene	0.000	1.00	ND	0.0099	
Acenaphthylene	0.000	1.00	ND	0.0099	
Acenaphthene	0.000	1.00	ND	0.0099	
Fluorene	0.000	1.00	ND	0.0099	
Phenanthrene	0.0246	1.00	ND	0.0099	
Anthracene	0.000	1.00	ND	0.0099	
Fluoranthene	0.0549	1.00	ND	0.0099	
Pyrene	0.0626	1.00	ND	0.0099	
Benzo (a) Anthracene	0.0439	1.00	ND	0.0099	
Chrysene	0.0454	1.00	ND	0.0099	
Benzo (k) Fluoranthene	0.0371	1.00	ND	0.0099	
Benzo (b) Fluoranthene	0.0331	1.00	ND	0.0099	
Benzo (a) Pyrene	0.0406	1.00	ND	0.0099	
Indeno (1,2,3-c,d) Pyrene	0.0267	1.00	ND	0.0099	
Dibenz (a,h) Anthracene	0.000	1.00	ND	0.0099	
Benzo (g,h,i) Perylene	0.0396	1.00	ND	0.0099	

Return to Contents

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar013.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:38 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 15:03 ev7p

Sample Name: 17-03-1557-27 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.301	152	61080	5.000	0.00
3)*Naphthalene-d8	(2)	4.536	136	171171	5.000	0.00
11)*Acenaphthene-d10	(3)	6.420	164	82099	5.000	0.00
17)*Phenanthrene-d10	(4)	8.019	188	275585	5.000	0.00
31)*Perylene-d12	(6)	12.688	264	268126	5.000	0.00
25)*Chrysene-d12	(5)	10.940	240	277786	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.853	82	11423	1.008	0.00
SpikedAmount 1.000				Recovery = 100.755		
7)\$2-Fluorobiphenyl	(3)	5.718	172	29263	0.987	0.00
SpikedAmount 1.000				Recovery = 98.735		
23)\$p-Terphenyl-d14	(5)	9.812	244	38881	0.903	0.00
SpikedAmount 1.000				Recovery = 90.319		
Target Compounds						
4) Naphthalene	(2)	0.000		0	N.D.	
5) 2-Methylnaphthalene	(2)	0.000		0	N.D.	
6) 1-Methylnaphthalene	(2)	0.000		0	N.D.	
10) Acenaphthylene	(3)	0.000		0	N.D.	
12) Acenaphthene	(3)	0.000		0	N.D.	
15) Fluorene	(3)	0.000		0	N.D.	
18) Phenanthrene	(4)	8.044	178	1436	0.025	96
19) Anthracene	(4)	0.000		0D	N.D.	
21) Fluoranthene	(4)	9.355	202	4133	0.055	95
22) Pyrene	(5)	9.590	202	4735	0.063	98
24) Benzo (a) Anthracene	(5)	10.923	228	3210	0.044	87
26) Chrysene	(5)	10.964	228	3017	0.045	97
27) Benzo (b) Fluoranthene	(6)	12.229	252	2313	0.033	70
28) Benzo (k) Fluoranthene	(6)	12.249	252	2719M	0.037	50
30) Benzo (a) Pyrene	(6)	12.620	252	2733A	0.041	97
33) Indeno (1,2,3-c,d) Pyrene	(6)	13.938	276	2148	0.027	84
34) Dibenz (a,h) Anthracene	(6)	0.000		0	N.D.	
35) Benzo (g,h,i) Perylene	(6)	14.211	276	2560	0.040	89
8) Biphenyl	(3)	0.000		0	N.D.	
9) 2,6-Dimethylnaphthalene	(3)	0.000		0	N.D.	
14) 1,6,7-Trimethylnaphthalene	(3)	0.000		0	N.D.	
16) Dibenzothiophene	(3)	0.000		0	N.D.	

M = Compound was manually integrated.
 D = Compound was deleted.
 A = User selected an alternate hit.
 * = Compound is an internal standard.
 \$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar013.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:38 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 15:03 ev7p

Sample Name: 17-03-1557-27 Misc Info: S032317A 10UL
 Response via Initial Calibration

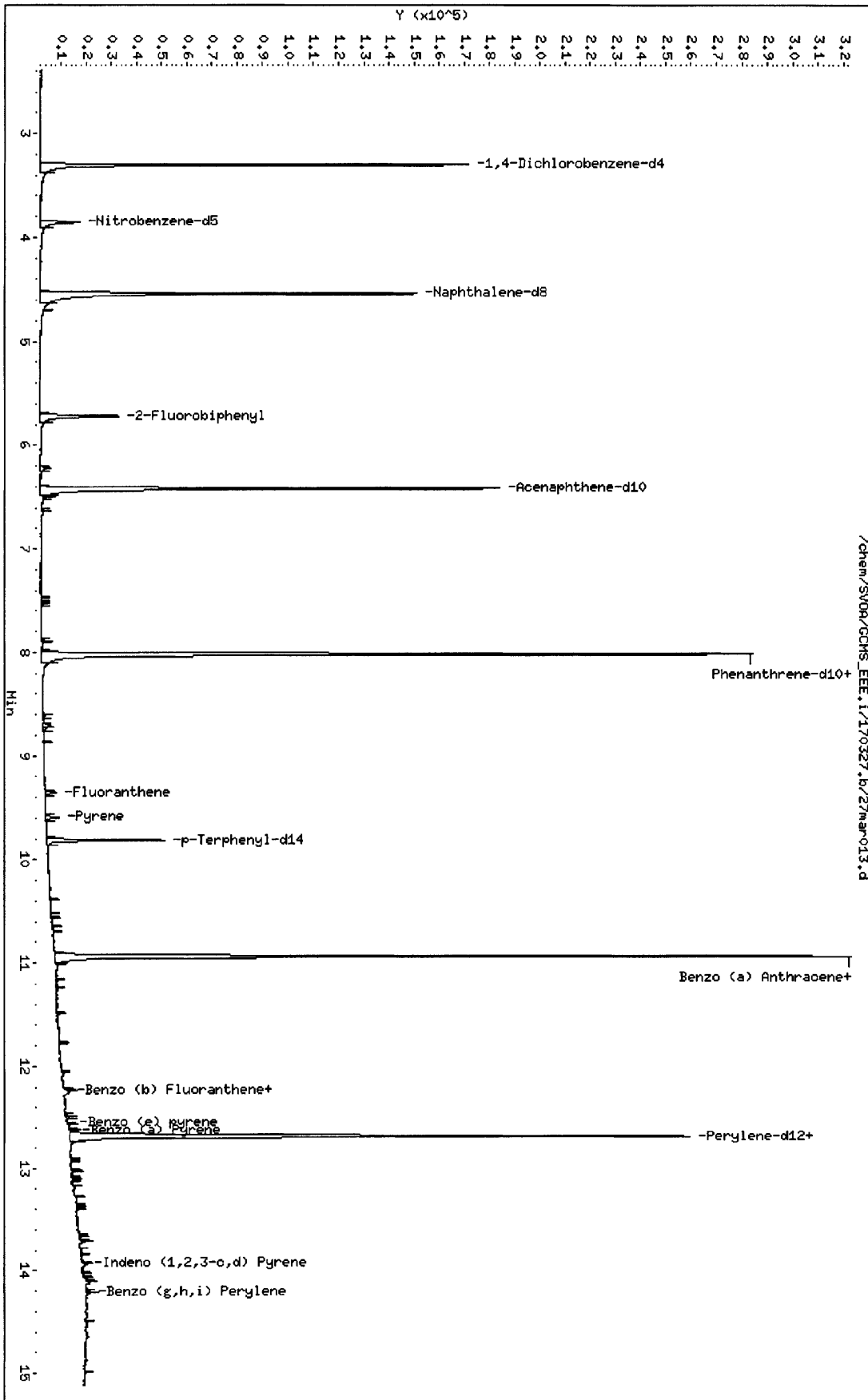
Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	0.000		0	N.D.	
29) Benzo (e) pyrene	(6)	12.555	252	1785	0.021	79
32) Perylene	(6)	12.714	252	1327A	0.017	72
13) Dibenzofuran	(3)	0.000		0	N.D.	

A = User selected an alternate hit.

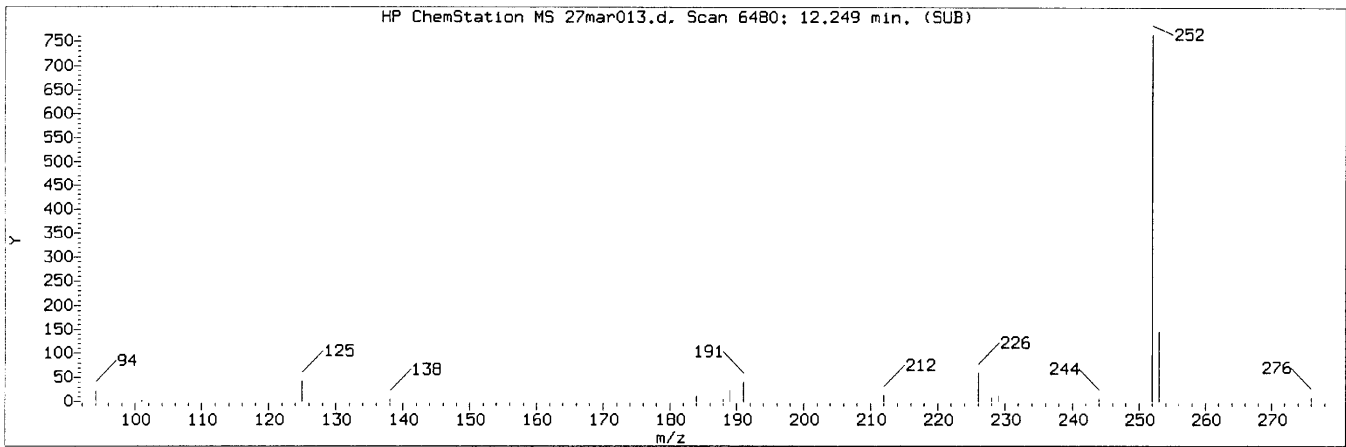
page 2 of 2

Data File: /chem/SV09/GCHS_EEE.i/170327.b/27mar013.d
Date: 27-MAR-2017 14:38
Client ID:
Sample Info: 17-03-1557-27
Column phase: J&W DB-5MS

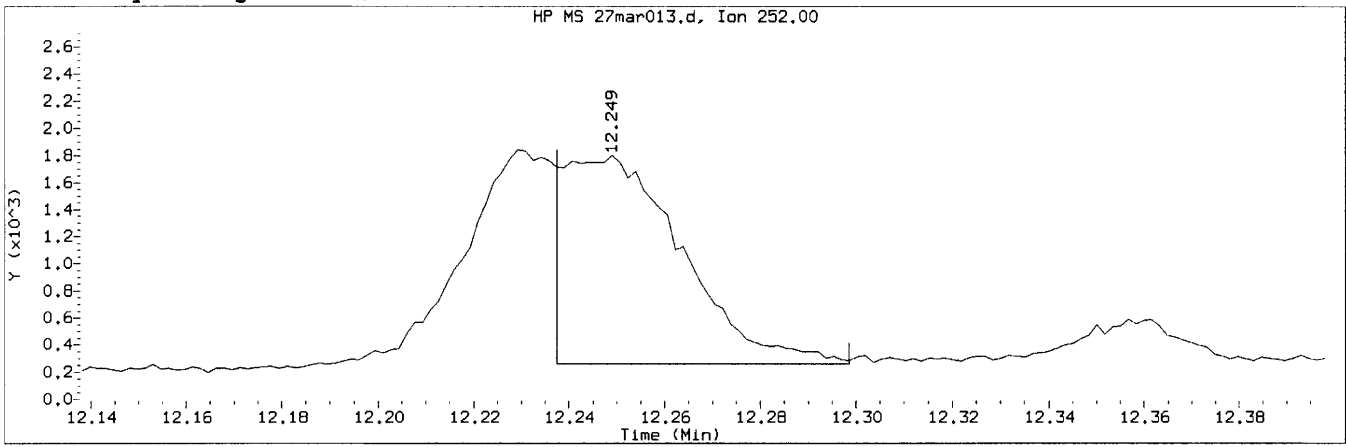
Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



Sample Spectrum (Background Subtracted)



Manually Integrated Quant Ion



Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar013.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:38 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 15:03 ev7p

Sample Name: 17-03-1557-27

Compound Number : 28
 Compound Name : Benzo (k) Fluoranthene
 Scan Number : 6480
 Retention Time (minutes): 12.249
 Quant Ion : 252.00
 Area (flag) : 2719M
 On-Column Amount (mg/L) : 0.0371
 Integration start scan : 6472 Integration stop scan: 6509
 Y at integration start : 259 Y at integration end: 259

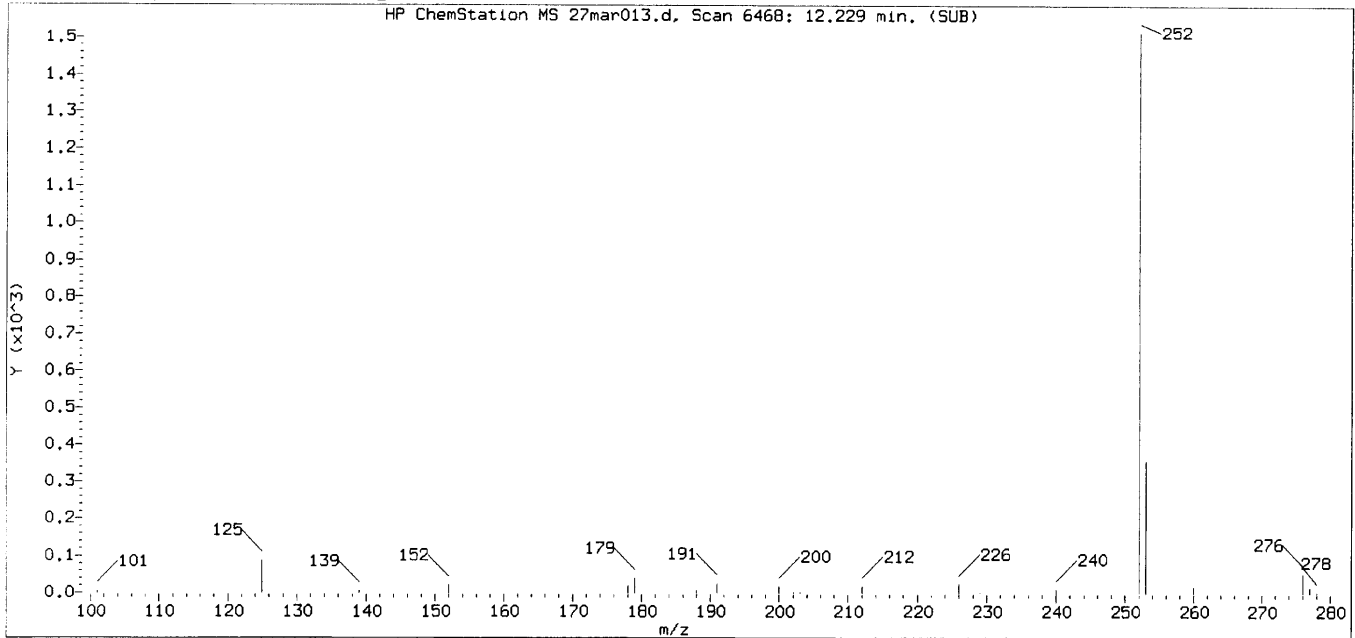
Reason for manual integration: improper integration

Analyst responsible for change: Digitally signed by Jeremy Huynh
 on 03/28/2017 at 09:12.
 Target 3.5 esignature user ID: ev7p

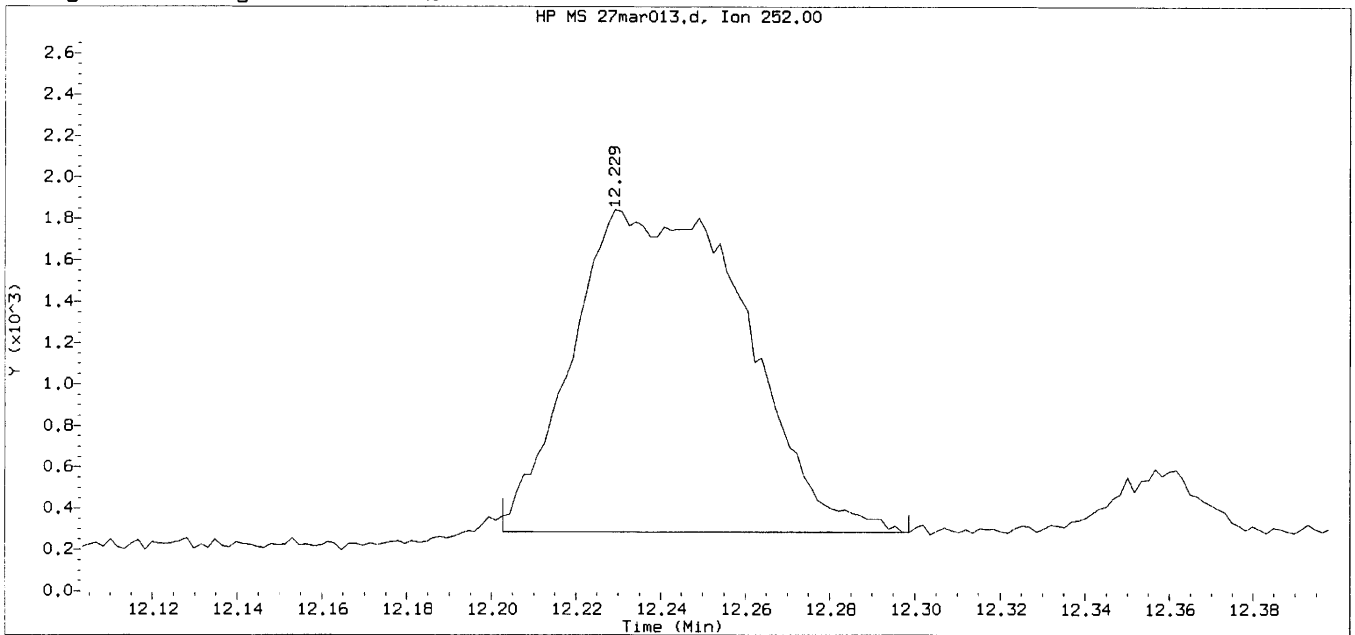
GC/MS audit/management approval: _____

Handwritten signature and date: ev 3/28/17

Sample Spectrum (Background Subtracted)



Original Integration of Quant Ion



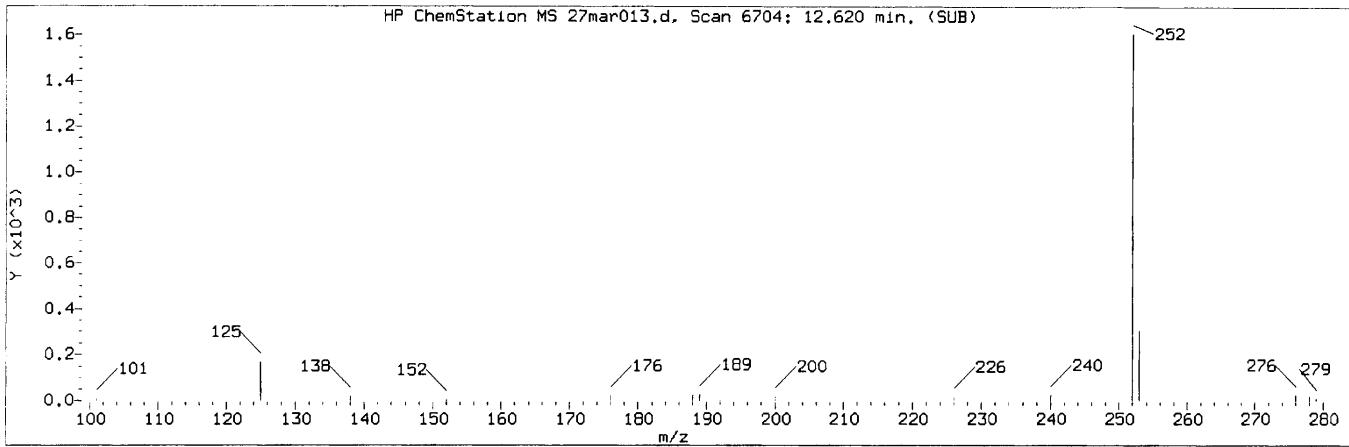
Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar013.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:38 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 15:00 Unknown

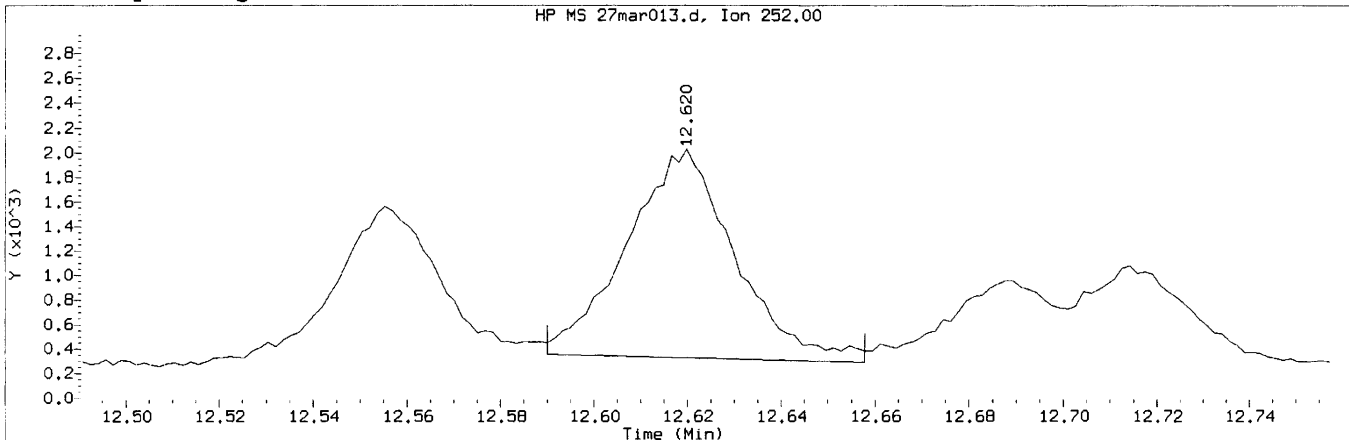
Sample Name: 17-03-1557-27

Compound Number	: 28	
Compound Name	: Benzo (k) Fluoranthene	
Scan Number	: 6468	
Retention Time (minutes)	: 12.229	
Quant Ion	: 252.00	
Area	: 4454	
On-column Amount (mg/L)	: 0.0608	
Integration start scan	: 6451	Integration stop scan: 6509
Y at integration start	: 286	Y at integration end: 286

Sample Spectrum (Background Subtracted)



Manually Integrated Quant Ion



Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar013.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:38 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 15:03 ev7p

Sample Name: 17-03-1557-27

Compound Number : 30
 Compound Name : Benzo (a) Pyrene
 Scan Number : 6704
 Retention Time (minutes): 12.620
 Quant Ion : 252.00
 Area (flag) : 2733A
 On-Column Amount (mg/L) : 0.0406
 Integration start scan : 6685 Integration stop scan: 6726
 Y at integration start : 359 Y at integration end: 288

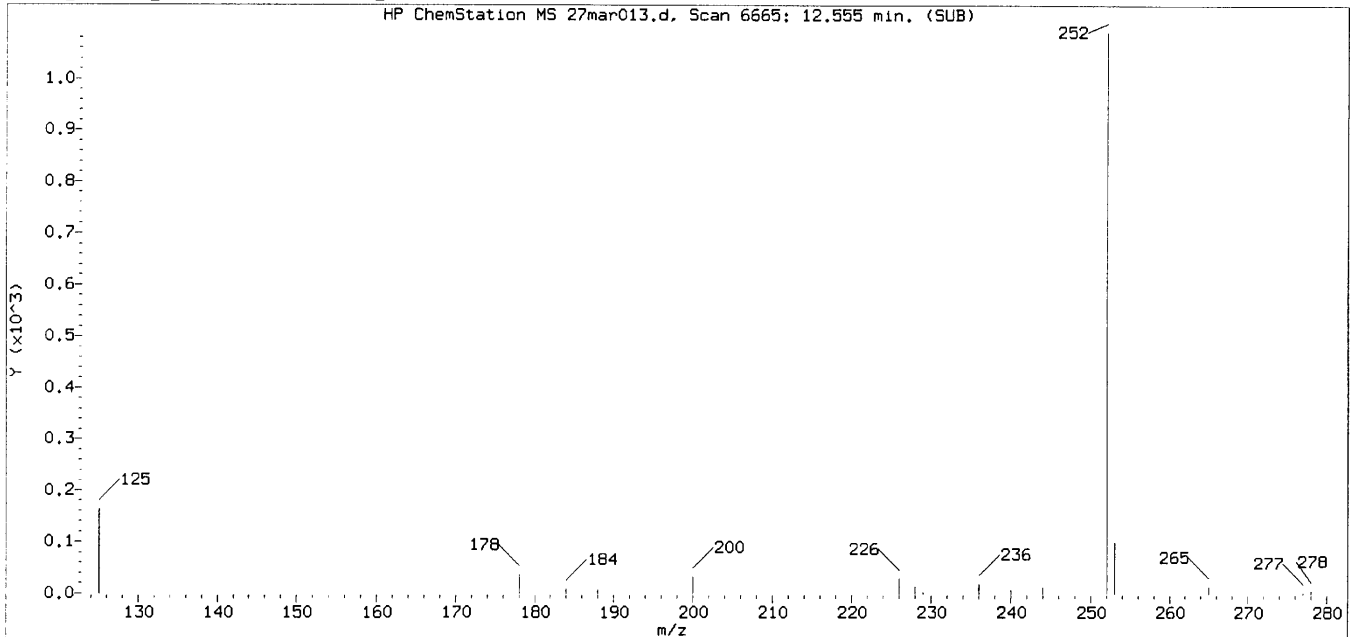
Reason for manual integration: improper integration

Analyst responsible for change: Digitally signed by Jeremy Huynh
 on 03/28/2017 at 09:12.
 Target 3.5 esignature user ID: ev7p

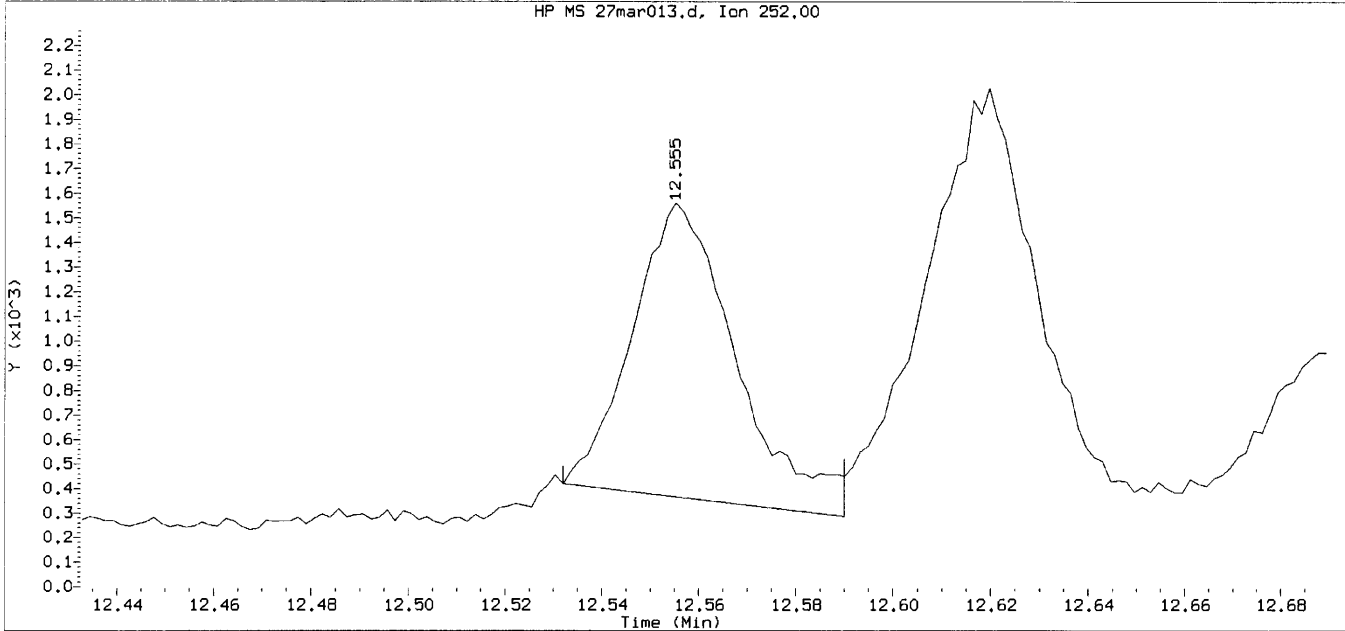
GC/MS audit/management approval: _____

Handwritten signature and date: 3/28/17

Sample Spectrum (Background Subtracted)



Original Integration of Quant Ion



Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar013.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 14:38 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 15:00 Unknown

Sample Name: 17-03-1557-27

Compound Number	: 30	
Compound Name	: Benzo (a) Pyrene	
Scan Number	: 6665	
Retention Time (minutes)	: 12.555	
Quant Ion	: 252.00	
Area	: 1785	
On-column Amount (mg/L)	: 0.0265	
Integration start scan	: 6650	Integration stop scan: 6685
Y at integration start	: 420	Y at integration end: 288

EPA METHOD 8270C PAHSIM

Quality Control

Method Blank LCS/LCSD MS/MSD

METHOD BLANK ASSOCIATION SUMMARY
FOR METHOD: EPA 8270C SIM PAHs

MB SAMPLE ID: 099-14-035-386
MB BATCH ID: 170324L17
INSTRUMENT: GC/MS EEE
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED: 2017-03-27 11:15
REVIEWED BY: 262
D/T REVIEWED: 2017-03-27 13:13
MATRIX: Soil

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar003.d\27mar003.rr

CLIENT WORK ORDER: 17-03-1557

<u>S#</u>	<u>RUN TYPE</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
25	B-DU1-ISM1-8		2017-03-27 13:57	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar011.d\27mar011.rr
26	B-DU1-ISM2-8		2017-03-27 14:17	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar012.d\27mar012.rr
27	B-DU1-ISM3-8		2017-03-27 14:38	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar013.d\27mar013.rr

RAW DATA SHEET FOR METHOD: EPA 8270C SIM PAHs

WORK ORDER: 099-14-035
INSTRUMENT: GC/MS EEE
EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED: 2017-03-27 11:15
REVIEWED BY: 262
D/T REVIEWED: 2017-03-27 13:13

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar003.d\27mar003.rr

MB **CLIENT SAMPLE NUMBER:** Method Blank

LCS/MB BATCH: 170324L17 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 20.00 g / ACTUAL: 20.00 g
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 2.00 ml / ACTUAL: 2.00 ml
UNITS: mg/kg **ADJUSTMENT RATIO TO PF:** 1.00

COMMENT:

<u>COMPOUND</u>	<u>ON COL CONC</u>	<u>DF</u>	<u>CONC</u>	<u>RL</u>	<u>QUAL</u>
Naphthalene	0.000	1.00	ND	0.010	
2-Methylnaphthalene	0.000	1.00	ND	0.010	
1-Methylnaphthalene	0.000	1.00	ND	0.010	
Acenaphthylene	0.000	1.00	ND	0.010	
Acenaphthene	0.000	1.00	ND	0.010	
Fluorene	0.000	1.00	ND	0.010	
Phenanthrene	0.000	1.00	ND	0.010	
Anthracene	0.000	1.00	ND	0.010	
Fluoranthene	0.000	1.00	ND	0.010	
Pyrene	0.000	1.00	ND	0.010	
Benzo (a) Anthracene	0.000	1.00	ND	0.010	
Chrysene	0.000	1.00	ND	0.010	
Benzo (k) Fluoranthene	0.000	1.00	ND	0.010	
Benzo (b) Fluoranthene	0.000	1.00	ND	0.010	
Benzo (a) Pyrene	0.000	1.00	ND	0.010	
Indeno (1,2,3-c,d) Pyrene	0.000	1.00	ND	0.010	
Dibenz (a,h) Anthracene	0.000	1.00	ND	0.010	
Benzo (g,h,i) Perylene	0.000	1.00	ND	0.010	

Return to Contents

LCS QUALITY CONTROL SHEET FOR METHOD: EPA 8270C SIM PAHS

LCS SAMPLE ID: 099-14-035-386
LCS/MB BATCH ID: 170324L17
INSTRUMENT: GC/MS EEE

EXTRACTION: EPA 3545
D/T EXTRACTED: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED: 2017-03-27 11:35
REVIEWED BY: 262
D/T REVIEWED: 2017-03-27 13:13

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar004.d\27mar004.tr

<u>COMPOUND</u>	<u>CONC.</u>	<u>CONC REC</u>	<u>%REC</u>	<u>%REC CL</u>	<u>ME CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
Naphthalene	0.1000	0.09899	99	51-129	38-142	PASS	
2-Methylnaphthalene	0.1000	0.1132	113	50-127	37-140	PASS	
1-Methylnaphthalene	0.1000	0.1008	101	54-132	41-145	PASS	
Acenaphthylene	0.1000	0.09919	99	50-123	38-135	PASS	
Acenaphthene	0.1000	0.1036	104	53-125	41-137	PASS	
Fluorene	0.1000	0.1041	104	55-127	43-139	PASS	
Phenanthrene	0.1000	0.1044	104	50-122	38-134	PASS	
Anthracene	0.1000	0.1059	106	50-132	36-146	PASS	
Fluoranthene	0.1000	0.1039	104	55-127	43-139	PASS	
Pyrene	0.1000	0.1021	102	50-134	36-148	PASS	
Benzo (a) Anthracene	0.1000	0.09779	98	50-133	36-147	PASS	
Chrysene	0.1000	0.1023	102	51-129	38-142	PASS	
Benzo (k) Fluoranthene	0.1000	0.09607	96	49-150	32-167	PASS	
Benzo (b) Fluoranthene	0.1000	0.1068	107	50-142	35-157	PASS	
Benzo (a) Pyrene	0.1000	0.09631	96	50-134	36-148	PASS	
Indeno (1,2,3-c,d) Pyrene	0.1000	0.09683	97	50-148	34-164	PASS	
Dibenz (a,h) Anthracene	0.1000	0.1004	100	50-133	36-147	PASS	
Benzo (g,h,i) Perylene	0.1000	0.1028	103	50-130	37-143	PASS	

Total number of LCS compounds: 18
 Total number of ME compounds: 0
 Total number of ME compounds allowed: 1
 LCS ME CL validation result: Pass

MATRIX SPIKE / MATRIX SPIKE DUPLICATE QUALITY CONTROL SHEET FOR METHOD: EPA 8270C SIM PAHS

SPIKED SAMPLE ID: 17-03-1753-6
MS/MSD BATCH: 170324S17

INSTRUMENTS:
SAMPLE: GC/MS EEE
MS: GC/MS EEE
MSD: GC/MS EEE

EXTRACTION: EPA 3545
D/T EXTRACTED:

SAMPLE: 2017-03-24 00:00
MS: 2017-03-24 00:00
MSD: 2017-03-24 00:00

ANALYZED BY: 907
D/T ANALYZED:

SAMPLE: 2017-03-27 12:36
MS: 2017-03-27 11:56
MSD: 2017-03-27 12:16
REVIEWED BY: 262
D/T REVIEWED: 2017-03-27 13:13

COMMENT:

COMPOUND NAME	SAMPLE	INITIAL	FINAL	MS CONC	% MS.REC	MSD CONC	% MSD.REC	% REC CL	RPD	RPD CL	STATUS	QUALIFIERS
Naphthalene	ND	1.000	0.1000	0.1123	112	0.1093	109	20-150	3	0-33	PASS	
2-Methylnaphthalene	ND	1.000	0.1000	0.1381	138	0.1341	134	29-137	3	0-31	FAIL	3F
1-Methylnaphthalene	0.01270	1.000	0.1000	0.1196	107	0.1189	106	34-136	1	0-29	PASS	
Acenaphthylene	ND	1.000	0.1000	0.1041	104	0.1049	105	29-131	1	0-32	PASS	
Acenaphthene	ND	1.000	0.1000	0.1076	108	0.1064	106	29-137	1	0-28	PASS	
Fluorene	ND	1.000	0.1000	0.1148	115	0.1155	115	36-132	1	0-27	PASS	
Phenanthrene	ND	1.000	0.1000	0.1057	106	0.1076	108	20-144	2	0-27	PASS	
Anthracene	ND	1.000	0.1000	0.1124	112	0.1092	109	26-134	3	0-27	PASS	
Fluoranthene	ND	1.000	0.1000	0.1066	107	0.1059	106	20-151	1	0-26	PASS	
Pyrene	ND	1.000	0.1000	0.1054	105	0.1084	108	20-150	3	0-32	PASS	
Benzo (a) Anthracene	ND	1.000	0.1000	0.1010	101	0.1040	104	24-150	3	0-24	PASS	
Chrysene	ND	1.000	0.1000	0.1012	101	0.09957	100	25-145	2	0-28	PASS	
Benzo (k) Fluoranthene	ND	1.000	0.1000	0.08966	90	0.08590	86	28-148	4	0-26	PASS	
Benzo (b) Fluoranthene	ND	1.000	0.1000	0.1132	113	0.1146	115	21-153	1	0-26	PASS	
Benzo (a) Pyrene	ND	1.000	0.1000	0.09968	100	0.09760	98	29-149	2	0-22	PASS	
Indeno (1,2,3-c,d) Pyrene	ND	1.000	0.1000	0.1004	100	0.09966	100	20-154	1	0-25	PASS	
Dibenz (a,h) Anthracene	ND	1.000	0.1000	0.1039	104	0.1048	105	20-132	1	0-26	PASS	
Benzo (g,h,i) Perylene	ND	1.000	0.1000	0.1071	107	0.1078	108	20-148	1	0-27	PASS	

Data Files:

TYPE	DATA FILE	DATA FILE PATH
MS	27mar005.rr	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar005.d\
MSD	27mar006.rr	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar006.d\

SURROGATE RECOVERIES FOR METHOD: EPA 8270C SIM PAHs

WORK ORDER: 17-03-1557

BATCH ID:

LCS/MB: 170324L17**MS:** 170324S17

EXTRACTION: EPA 3545

REVIEWED BY:

D/T REVIEWED:

25 **CLIENT SAMPLE NUMBER : B-DU1-ISM1-8**

INSTRUMENT: GC/MS EEE

D/T EXTRACTED: 2017-03-24 00:00

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar011.d\27mar011.rr

ANALYZED BY: 907

D/T ANALYZED 2017-03-27 13:57

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
2-Fluorobiphenyl	83	13-127	PASS	
Nitrobenzene-d5	79	17-137	PASS	
p-Terphenyl-d14	94	4-160	PASS	

26 **CLIENT SAMPLE NUMBER : B-DU1-ISM2-8**

INSTRUMENT: GC/MS EEE

D/T EXTRACTED: 2017-03-24 00:00

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar012.d\27mar012.rr

ANALYZED BY: 907

D/T ANALYZED 2017-03-27 14:17

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
p-Terphenyl-d14	98	4-160	PASS	
Nitrobenzene-d5	86	17-137	PASS	
2-Fluorobiphenyl	91	13-127	PASS	

27 **CLIENT SAMPLE NUMBER : B-DU1-ISM3-8**

INSTRUMENT: GC/MS EEE

D/T EXTRACTED: 2017-03-24 00:00

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar013.d\27mar013.rr

ANALYZED BY: 907

D/T ANALYZED 2017-03-27 14:38

COMMENT:

COMPOUND	% REC	% REC CL	STATUS	QUALIFIERS
2-Fluorobiphenyl	99	13-127	PASS	
Nitrobenzene-d5	101	17-137	PASS	
p-Terphenyl-d14	90	4-160	PASS	

**SURROGATE RECOVERIES
FOR METHOD: EPA 8270C SIM PAHs**

WORK ORDER: 17-03-1557

REVIEWED BY: 262

BATCH ID:

D/T REVIEWED: 2017-03-27 13:13

LCS/MB: 170324L17

MS:

EXTRACTION: EPA 3545

MB CLIENT SAMPLE NUMBER : Method Blank

INSTRUMENT: GC/MS EEE

ANALYZED BY: 907

D/T EXTRACTED: 2017-03-24 00:00

D/T ANALYZED 2017-03-27 11:15

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar003.d\27mar003.rr

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
2-Fluorobiphenyl	107	13-127	PASS	
Nitrobenzene-d5	109	17-137	PASS	
p-Terphenyl-d14	104	4-160	PASS	

LCS CLIENT SAMPLE NUMBER : Lab Control Sample

INSTRUMENT: GC/MS EEE

ANALYZED BY: 907

D/T EXTRACTED: 2017-03-24 00:00

D/T ANALYZED 2017-03-27 11:35

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar004.d\27mar004.rr

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
p-Terphenyl-d14	100	4-160	PASS	
Nitrobenzene-d5	111	17-137	PASS	
2-Fluorobiphenyl	97	13-127	PASS	

MS CLIENT SAMPLE NUMBER : Matrix Spike

INSTRUMENT: GC/MS EEE

ANALYZED BY: 907

D/T EXTRACTED: 2017-03-24 00:00

D/T ANALYZED 2017-03-27 11:56

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar005.d\27mar005.rr

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
p-Terphenyl-d14	99	4-160	PASS	
Nitrobenzene-d5	117	17-137	PASS	
2-Fluorobiphenyl	90	13-127	PASS	

MSD CLIENT SAMPLE NUMBER : Matrix Spike Duplicate

INSTRUMENT: GC/MS EEE

ANALYZED BY: 907

D/T EXTRACTED: 2017-03-24 00:00

D/T ANALYZED 2017-03-27 12:16

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar006.d\27mar006.rr

COMMENT:

<u>COMPOUND</u>	<u>% REC</u>	<u>% REC CL</u>	<u>STATUS</u>	<u>QUALIFIERS</u>
2-Fluorobiphenyl	93	13-127	PASS	
Nitrobenzene-d5	115	17-137	PASS	
p-Terphenyl-d14	102	4-160	PASS	

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar003.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 11:15 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 11:58 ev7p

Sample Name: MB 170324 L17 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S.		QIon	Area	On-Column	
	Ref.	RT			Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.297	152	53792	5.000	-0.01
3)*Naphthalene-d8	(2)	4.531	136	151636	5.000	0.00
11)*Acenaphthene-d10	(3)	6.411	164	71653	5.000	0.00
17)*Phenanthrene-d10	(4)	8.004	188	255973	5.000	0.00
31)*Perylene-d12	(6)	12.651	264	249943	5.000	0.00
25)*Chrysene-d12	(5)	10.919	240	264101	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.848	82	10990	1.094	0.00
SpikedAmount 1.000				Recovery = 109.418		
7)\$2-Fluorobiphenyl	(3)	5.710	172	27608	1.067	0.00
SpikedAmount 1.000				Recovery = 106.732		
23)\$p-Terphenyl-d14	(5)	9.800	244	42587	1.041	0.00
SpikedAmount 1.000				Recovery = 104.053		
Target Compounds						
						QValue
4) Naphthalene	(2)	0.000		0		N.D.
5) 2-Methylnaphthalene	(2)	0.000		0		N.D.
6) 1-Methylnaphthalene	(2)	0.000		0		N.D.
10) Acenaphthylene	(3)	0.000		0		N.D.
12) Acenaphthene	(3)	0.000		0		N.D.
15) Fluorene	(3)	0.000		0		N.D.
18) Phenanthrene	(4)	0.000		0		N.D.
19) Anthracene	(4)	0.000		0		N.D.
21) Fluoranthene	(4)	0.000		0		N.D.
22) Pyrene	(5)	0.000		0		N.D.
24) Benzo (a) Anthracene	(5)	0.000		0		N.D.
26) Chrysene	(5)	0.000		0D		N.D.
27) Benzo (b) Fluoranthene	(6)	0.000		0		N.D.
28) Benzo (k) Fluoranthene	(6)	0.000		0		N.D.
30) Benzo (a) Pyrene	(6)	0.000		0		N.D.
33) Indeno (1,2,3-c,d) Pyrene	(6)	0.000		0		N.D.
34) Dibenz (a,h) Anthracene	(6)	0.000		0		N.D.
35) Benzo (g,h,i) Perylene	(6)	0.000		0		N.D.
8) Biphenyl	(3)	0.000		0		N.D.
9) 2,6-Dimethylnaphthalene	(3)	0.000		0		N.D.
14) 1,6,7-Trimethylnaphthalene	(3)	0.000		0		N.D.
16) Dibenzothiophene	(3)	0.000		0		N.D.

D = Compound was deleted.

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar003.d Instrument ID: GCMS_EEE.i
Injection date and time: 27-MAR-2017 11:15 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
Calibration date and time: 27-MAR-2017 11:23
Date, time and analyst ID of latest file update: 27-Mar-2017 11:58 ev7p

Sample Name: MB 170324 L17 Misc Info: S032317A 10UL
Response via Initial Calibration

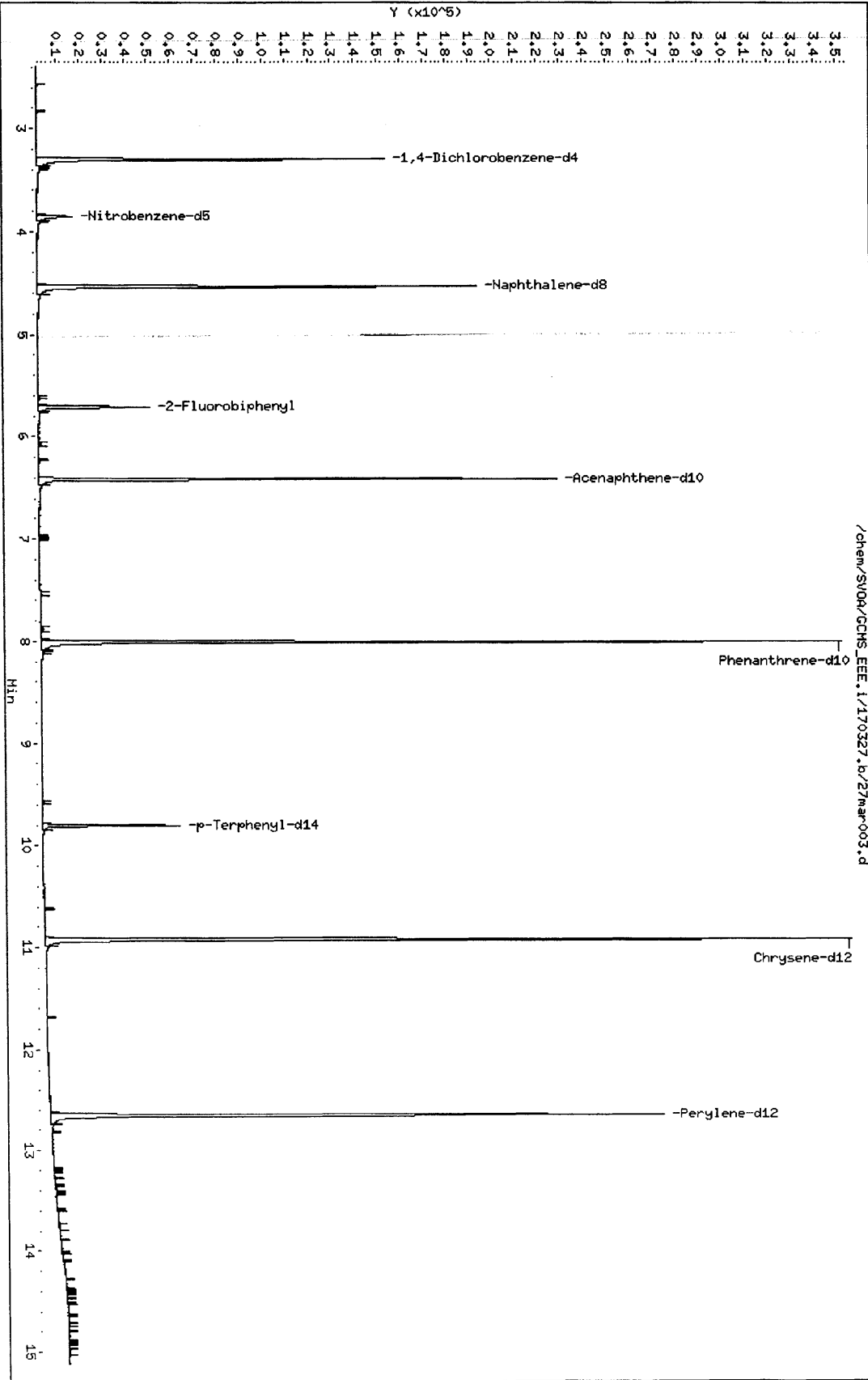
Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	0.000		0	N.D.	
29) Benzo (e) pyrene	(6)	0.000		0	N.D.	
32) Perylene	(6)	0.000		0	N.D.	
13) Dibenzofuran	(3)	0.000		0	N.D.	

page 2 of 2

Data File: /chem/SVDR/CCHS_EEE.i/170327.b/27mar003.d
Date: 27-MAR-2017 11:15
Client ID:
Sample Info: HB 170324 L17
Column phase: J&W DB-SMS

Instrument: CCHS_EEE.i
Operator: 907
Column diameter: 0.18

/chem/SVDR/CCHS_EEE.i/170327.b/27mar003.d



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar004.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 11:35 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 11:57 Unknown

Sample Name: LCS 170324 L17 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column	
					Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.296	152	59177	5.000	0.00
3)*Naphthalene-d8	(2)	4.530	136	161558	5.000	0.00
11)*Acenaphthene-d10	(3)	6.411	164	75915	5.000	0.00
17)*Phenanthrene-d10	(4)	8.006	188	258057	5.000	0.00
31)*Perylene-d12	(6)	12.651	264	264968	5.000	0.00
25)*Chrysene-d12	(5)	10.918	240	273070	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.846	82	11840	1.106	0.00
SpikedAmount	1.000			Recovery = 110.648		
7)\$2-Fluorobiphenyl	(3)	5.708	172	26532	0.968	0.00
SpikedAmount	1.000			Recovery = 96.815		
23)\$p-Terphenyl-d14	(5)	9.800	244	42412	1.002	0.00
SpikedAmount	1.000			Recovery = 100.223		
Target Compounds						
4) Naphthalene	(2)	4.551	128	34688	0.990	99
5) 2-Methylnaphthalene	(2)	5.279	142	24194	1.132	100
6) 1-Methylnaphthalene	(2)	5.391	142	21604	1.008	98
10) Acenaphthylene	(3)	6.242	152	38079	0.992	99
12) Acenaphthene	(3)	6.445	153	23876	1.036	100
15) Fluorene	(3)	6.996	166	26694	1.041	97
18) Phenanthrene	(4)	8.029	178	57171	1.044	99
19) Anthracene	(4)	8.079	178	56701	1.059	99
21) Fluoranthene	(4)	9.335	202	73166	1.039	99
22) Pyrene	(5)	9.572	202	75887	1.021	100
24) Benzo (a) Anthracene	(5)	10.901	228	70352	0.978	100
26) Chrysene	(5)	10.944	228	66849	1.023	100
27) Benzo (b) Fluoranthene	(6)	12.194	252	73767	1.068	95
28) Benzo (k) Fluoranthene	(6)	12.221	252	69552	0.961	95
30) Benzo (a) Pyrene	(6)	12.582	252	64106	0.963	99
33) Indeno (1,2,3-c,d) Pyrene	(6)	13.889	276	77051	0.968	99
34) Dibenz (a,h) Anthracene	(6)	13.912	278	60607	1.004	99
35) Benzo (g,h,i) Perylene	(6)	14.158	276	65733	1.028	99
8) Biphenyl	(3)	5.796	154	31974	0.940	99
9) 2,6-Dimethylnaphthalene	(3)	5.956	156	21642	0.907	100
14) 1,6,7-Trimethylnaphthalene	(3)	6.876	170	21756	0.912	100
16) Dibenzothiophene	(3)	7.893	184	70317	1.002	99

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar004.d Instrument ID: GCMS_EEE.i
Injection date and time: 27-MAR-2017 11:35 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
Calibration date and time: 27-MAR-2017 11:23

Date, time and analyst ID of latest file update: 27-Mar-2017 11:57 Unknown

Sample Name: LCS 170324 L17

Misc Info: S032317A 10UL

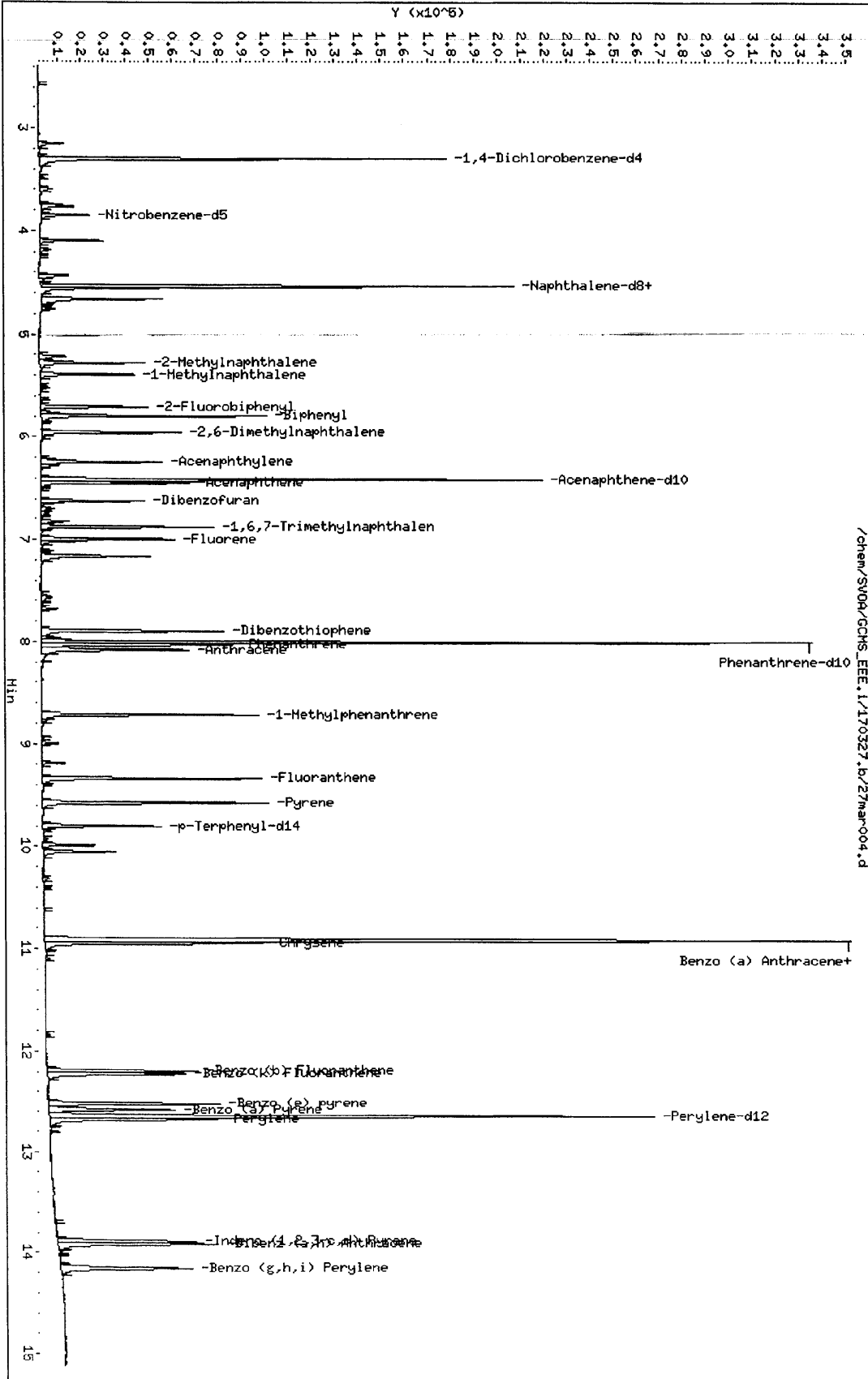
Response via Initial Calibration

Compounds	I.S.	RT	QIon	Area	On-Column	QValue
	Ref.				Amount (mg/L)	
20) 1-Methylphenanthrene	(5)	8.716	192	52870	0.971	100
29) Benzo (e) pyrene	(6)	12.519	252	78369	0.950	99
32) Perylene	(6)	12.679	252	73990	0.935	98
13) Dibenzofuran	(3)	6.623	168	34622	1.039	100

page 2 of 2

Data File: /chem/SV06/GCHS_EEE.i/170327.b/27mar004.d
Date: 27-Mar-2017 11:35
Client ID:
Sample Info: LCS 170324 L17
Column phase: J&W DB-SHS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar005.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 11:56 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 12:49 ev7p

Sample Name: 17-03-1753-6 MS Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S.		QIon	Area	On-Column	
	Ref.	RT			Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.299	152	57973	5.000	-0.01
3)*Naphthalene-d8	(2)	4.532	136	150857	5.000	0.00
11)*Acenaphthene-d10	(3)	6.416	164	71434	5.000	-0.01
17)*Phenanthrene-d10	(4)	8.009	188	252691	5.000	0.00
31)*Perylene-d12	(6)	12.661	264	256493	5.000	-0.01
25)*Chrysene-d12	(5)	10.922	240	265492	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.845	82	11711	1.172	0.00
SpikedAmount	1.000					Recovery = 117.202
7)\$2-Fluorobiphenyl	(3)	5.712	172	23148	0.898	0.00
SpikedAmount	1.000					Recovery = 89.764
23)\$p-Terphenyl-d14	(5)	9.803	244	40645	0.988	0.00
SpikedAmount	1.000					Recovery = 98.788
Target Compounds						
4) Naphthalene	(2)	4.552	128	36731	1.123	85
5) 2-Methylnaphthalene	(2)	5.283	142	27560	1.381	100
6) 1-Methylnaphthalene	(2)	5.395	142	23933	1.196	95
10) Acenaphthylene	(3)	6.248	152	37605	1.041	87
12) Acenaphthene	(3)	6.449	153	23317	1.076	98
15) Fluorene	(3)	6.998	166	27707	1.148	100
18) Phenanthrene	(4)	8.032	178	56669	1.057	99
19) Anthracene	(4)	8.082	178	58947	1.124	98
21) Fluoranthene	(4)	9.340	202	73523	1.066	98
22) Pyrene	(5)	9.577	202	76232	1.054	100
24) Benzo (a) Anthracene	(5)	10.908	228	70603	1.009	99
26) Chrysene	(5)	10.951	228	64286	1.012	99
27) Benzo (b) Fluoranthene	(6)	12.202	252	75658	1.132	97
28) Benzo (k) Fluoranthene	(6)	12.232	252	62835	0.897	95
30) Benzo (a) Pyrene	(6)	12.591	252	64226	0.997	98
33) Indeno (1,2,3-c,d) Pyrene	(6)	13.899	276	77376	1.005	97
34) Dibenz (a,h) Anthracene	(6)	13.924	278	60692	1.039	99
35) Benzo (g,h,i) Perylene	(6)	14.172	276	66307	1.071	96
8) Biphenyl	(3)	5.800	154	34323	1.073	97
9) 2,6-Dimethylnaphthalene	(3)	5.962	156	30916	1.377	99
14) 1,6,7-Trimethylnaphthalene	(3)	6.878	170	32486	1.448	99
16) Dibenzothiophene	(3)	7.895	184	70477	1.067	99

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar005.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 11:56 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 12:49 ev7p

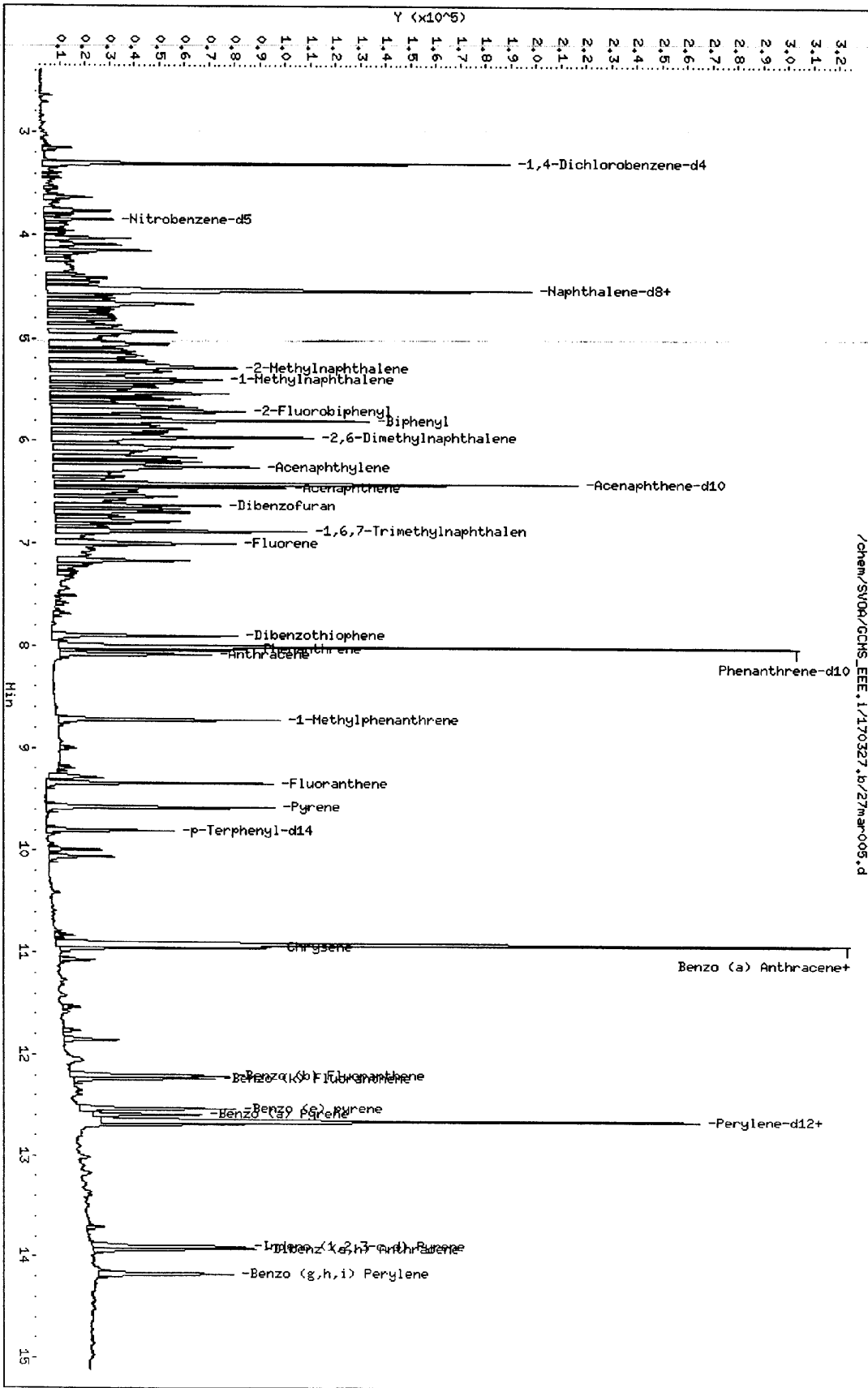
Sample Name: 17-03-1753-6 MS Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column	
					Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.721	192	53620	1.013	100
29) Benzo (e) pyrene	(6)	12.532	252	77418	0.969	97
32) Perylene	(6)	12.692	252	76413	0.997	98
13) Dibenzofuran	(3)	6.626	168	36035	1.150	96

page 2 of 2

Data File: /chem/SV08/GCHS_EEE.i/170327.b/27mar005.d
Date: 27-Mar-2017 11:56
Client ID:
Sample Info: 17-03-1753-6 HS
Column phase: J&W DB-5MS

Instrument: GCHS_EEE.i
Operator: 907
Column diameter: 0.18



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar006.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 12:16 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23

Date, time and analyst ID of latest file update: 27-Mar-2017 12:49 ev7p

Sample Name: 17-03-1753-6 MSD
 Response via Initial Calibration

Misc Info: S032317A 10UL

Compounds	I.S.		QIon	Area	On-Column Amount	
	Ref.	RT			(mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.300	152	58479	5.000	-0.01
3)*Naphthalene-d8	(2)	4.534	136	152866	5.000	0.00
11)*Acenaphthene-d10	(3)	6.418	164	72569	5.000	-0.01
17)*Phenanthrene-d10	(4)	8.011	188	255933	5.000	-0.01
31)*Perylene-d12	(6)	12.671	264	255038	5.000	-0.02
25)*Chrysene-d12	(5)	10.928	240	257728	5.000	-0.01
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.846	82	11596	1.145	0.00
SpikedAmount 1.000				Recovery = 114.523		
7)\$2-Fluorobiphenyl	(3)	5.714	172	24339	0.929	-0.01
SpikedAmount 1.000				Recovery = 92.907		
23)\$p-Terphenyl-d14	(5)	9.805	244	40891	1.024	0.00
SpikedAmount 1.000				Recovery = 102.381		
Target Compounds						
4) Naphthalene	(2)	4.554	128	36254	1.093	82
5) 2-Methylnaphthalene	(2)	5.285	142	27117	1.341	97
6) 1-Methylnaphthalene	(2)	5.397	142	24102	1.189	95
10) Acenaphthylene	(3)	6.249	152	38508	1.049	81
12) Acenaphthene	(3)	6.450	153	23432	1.064	97
15) Fluorene	(3)	7.000	166	28303	1.155	99
18) Phenanthrene	(4)	8.034	178	58419	1.076	100
19) Anthracene	(4)	8.084	178	57960	1.092	99
21) Fluoranthene	(4)	9.343	202	73969	1.059	99
22) Pyrene	(5)	9.580	202	76041	1.084	99
24) Benzo (a) Anthracene	(5)	10.909	228	70638	1.040	100
26) Chrysene	(5)	10.954	228	61424	0.996	98
27) Benzo (b) Fluoranthene	(6)	12.211	252	76123	1.145	94
28) Benzo (k) Fluoranthene	(6)	12.237	252	59864	0.859	96
30) Benzo (a) Pyrene	(6)	12.600	252	62531	0.976	96
33) Indeno (1,2,3-c,d) Pyrene	(6)	13.911	276	76328	0.997	96
34) Dibenz (a,h) Anthracene	(6)	13.936	278	60882	1.048	99
35) Benzo (g,h,i) Perylene	(6)	14.184	276	66337	1.078	98
8) Biphenyl	(3)	5.802	154	35389	1.089	96
9) 2,6-Dimethylnaphthalene	(3)	5.963	156	31736	1.391	95
14) 1,6,7-Trimethylnaphthalene	(3)	6.881	170	33039	1.449	99
16) Dibenzothiophene	(3)	7.900	184	70636	1.053	99

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar006.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 12:16 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 12:49 ev7p

Sample Name: 17-03-1753-6 MSD Misc Info: S032317A 10UL
 Response via Initial Calibration

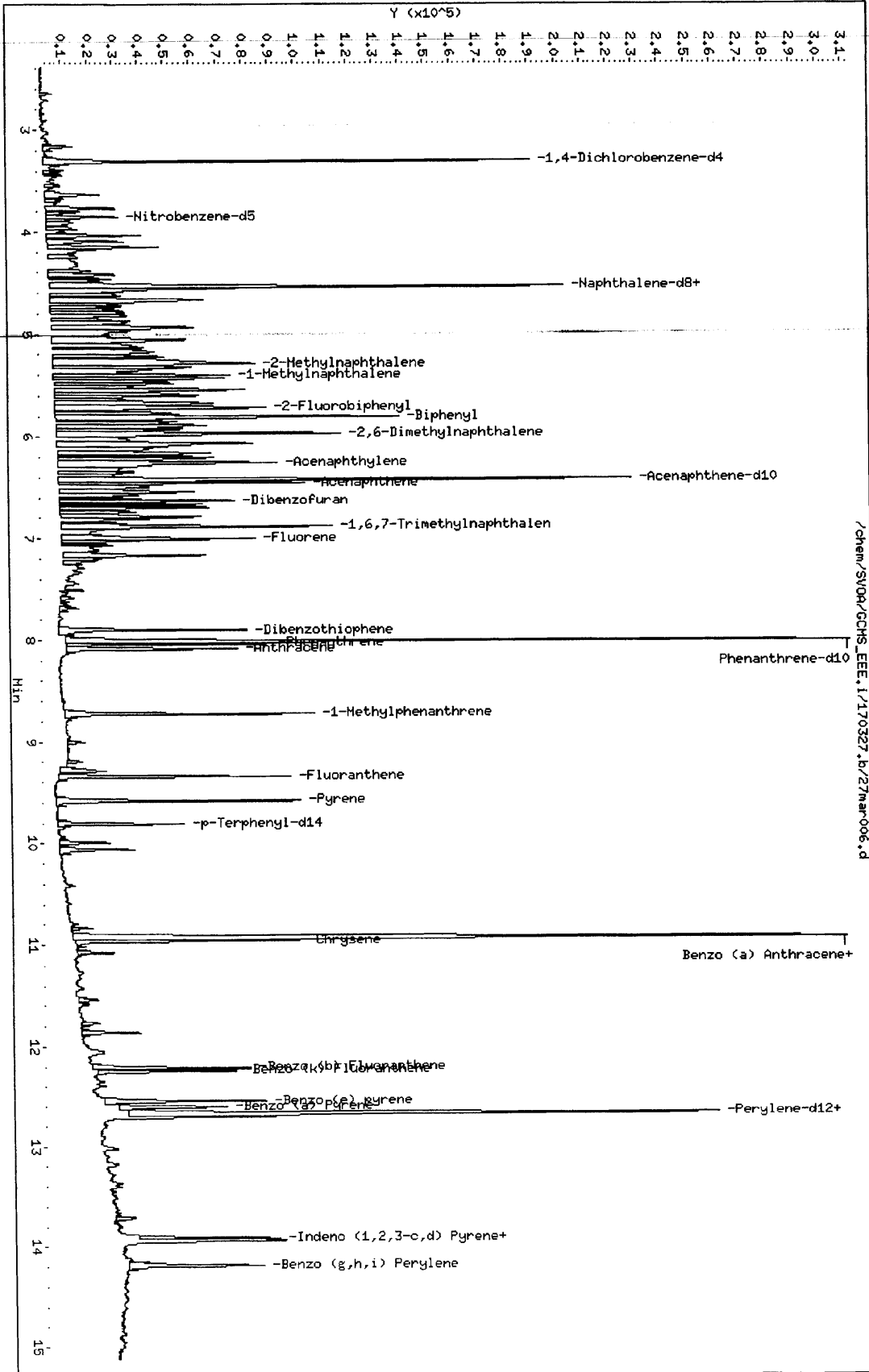
Compounds	I.S. Ref.	RT	QIon	Area	On-Column	QValue
					Amount (mg/L)	
20) 1-Methylphenanthrene	(5)	8.724	192	54048	1.052	99
29) Benzo (e) pyrene	(6)	12.540	252	78269	0.985	91
32) Perylene	(6)	12.701	252	75577	0.992	99
13) Dibenzofuran	(3)	6.626	168	37097	1.165	98

page 2 of 2

Data File: /chem/SV004/CCHS_EEE.1/170327.b/27mar006.d
Date: 27-MAR-2017 12:16
Client ID:
Sample Info: 17-03-1783-6_HSD
Column phase: J&M DB-SMS

Instrument: CCHS_EEE.1
Operator: 907
Column diameter: 0.18

/chem/SV004/CCHS_EEE.1/170327.b/27mar006.d



Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar007.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 12:36 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 13:02 ev7p

Sample Name: 17-03-1753-6 Misc Info: S032317A 10UL
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV (Min)	
Internal Standards							
1)*1,4-Dichlorobenzene-d4	(1)	3.301	152	59052	5.000	-0.01	
3)*Naphthalene-d8	(2)	4.536	136	155982	5.000	-0.01	
11)*Acenaphthene-d10	(3)	6.421	164	73544	5.000	-0.01	
17)*Phenanthrene-d10	(4)	8.013	188	259932	5.000	-0.01	
31)*Perylene-d12	(6)	12.676	264	264839	5.000	-0.02	
25)*Chrysene-d12	(5)	10.933	240	269341	5.000	-0.01	
System Monitoring Compounds							
2)\$Nitrobenzene-d5	(2)	3.850	82	9955	0.964	-0.01	
SpikedAmount	1.000			Recovery = 96.353			
7)\$2-Fluorobiphenyl	(3)	5.716	172	24362	0.918	-0.01	
SpikedAmount	1.000			Recovery = 91.762			
23)\$p-Terphenyl-d14	(5)	9.807	244	43951	1.053	-0.01	
SpikedAmount	1.000			Recovery = 105.298			
Target Compounds							
4) Naphthalene	(2)	4.570	128	2881	0.085	1	QValue
5) 2-Methylnaphthalene	(2)	5.287	142	1073	0.052	42	
6) 1-Methylnaphthalene	(2)	5.400	142	2640	0.128	74	
10) Acenaphthylene	(3)	0.000		0D	N.D.		
12) Acenaphthene	(3)	0.000		0D	N.D.		
15) Fluorene	(3)	7.005	166	2153	0.087	88	
18) Phenanthrene	(4)	8.036	178	920	0.017	94	
19) Anthracene	(4)	0.000		0D	N.D.		
21) Fluoranthene	(4)	0.000		0	N.D.		
22) Pyrene	(5)	9.582	202	1126	0.015	73	
24) Benzo (a) Anthracene	(5)	0.000		0D	N.D.		
26) Chrysene	(5)	0.000		0D	N.D.		
27) Benzo (b) Fluoranthene	(6)	0.000		0	N.D.		
28) Benzo (k) Fluoranthene	(6)	0.000		0	N.D.		
30) Benzo (a) Pyrene	(6)	0.000		0D	N.D.		
33) Indeno (1,2,3-c,d) Pyrene	(6)	0.000		0	N.D.		
34) Dibenz (a,h) Anthracene	(6)	0.000		0	N.D.		
35) Benzo (g,h,i) Perylene	(6)	0.000		0	N.D.		
8) Biphenyl	(3)	5.805	154	3337	0.101	75	
9) 2,6-Dimethylnaphthalene	(3)	5.968	156	9498	0.411	91	
14) 1,6,7-Trimethylnaphthalene	(3)	6.880	170	11105	0.481	98	
16) Dibenzothiophene	(3)	0.000		0	N.D.		

D = Compound was deleted.

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar007.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 12:36 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 13:02 ev7p

Sample Name: 17-03-1753-6 Misc Info: S032317A 10UL
 Response via Initial Calibration

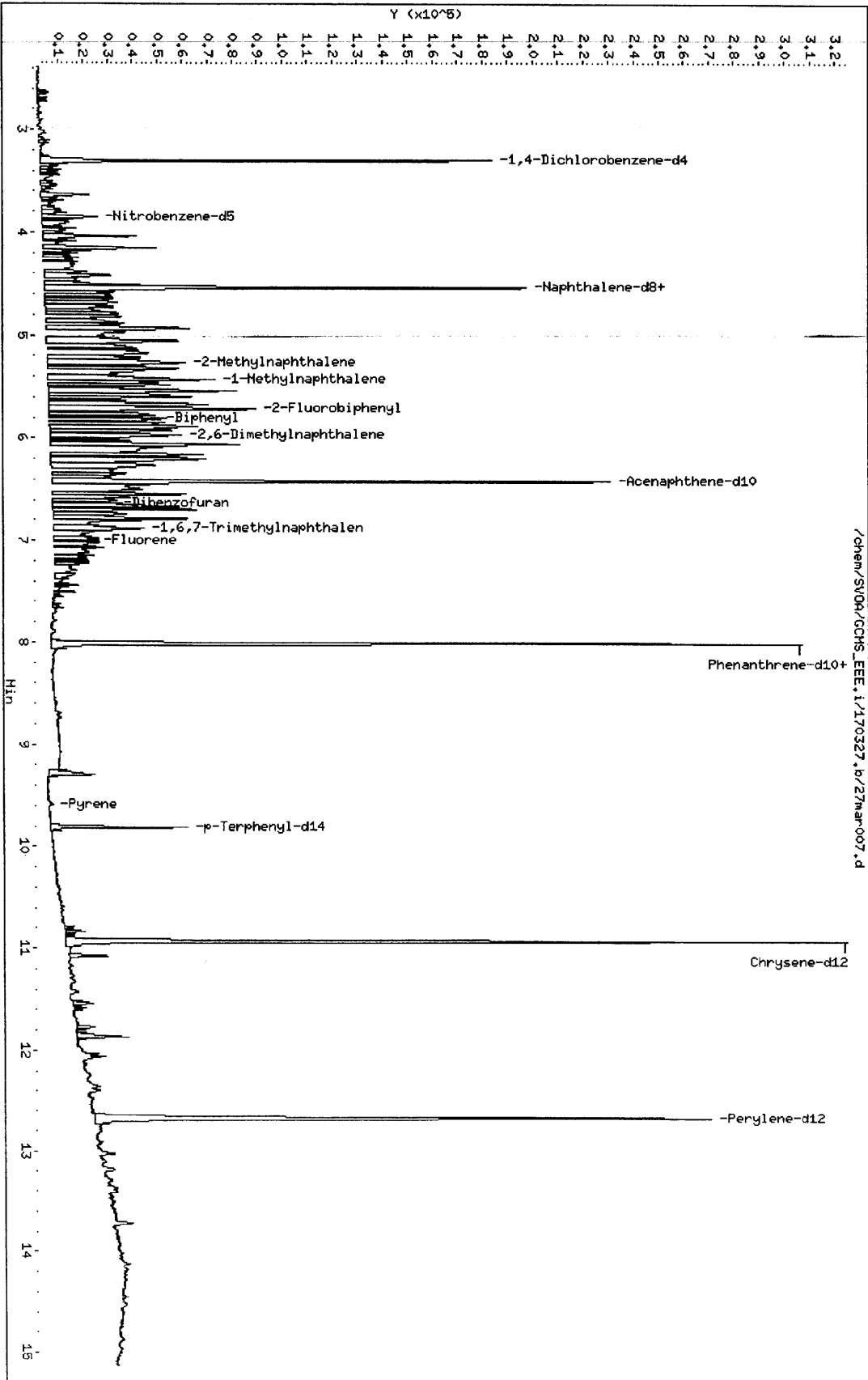
Compounds	I.S. Ref.	RT	QIon	Area	On-Column	QValue
					Amount (mg/L)	
20) 1-Methylphenanthrene	(5)	0.000		0	N.D.	
29) Benzo (e) pyrene	(6)	0.000		0	N.D.	
32) Perylene	(6)	0.000		0	N.D.	
13) Dibenzofuran	(3)	6.631	168	2167	0.067	10

page 2 of 2

Data File: /chem/SV09/CCHS_EEE.i/170327.b/27mar007.d
Date: 27-Mar-2017 12:36
Client ID:
Sample Info: 17-03-1753-6
Column phase: J&M DB-5MS

Instrument: CCHS_EEE.i
Operator: 907
Column diameter: 0.18

/chem/SV09/CCHS_EEE.i/170327.b/27mar007.d



EPA METHOD 8270C PAHSIM

Continuing Calibration

CCV ASSOCIATION SUMMARY
FOR METHOD: EPA 8270C SIM PAHs

BATCH ID: 170327A010
INSTRUMENT: GC/MS EEE

ANALYZED BY: 907

WORK ORDER: 099-06-009
MATRIX: Water

REVIEWED BY: 262
D/T REVIEWED: 2017-03-27 13:13

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
4997	Daily Calibration	2017-03-27 10:55	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar002.d\27mar002.r

WORK ORDER: 17-03-1557
MATRIX: Soil

REVIEWED BY:
D/T REVIEWED:

<u>CEL SAMPLE #</u>	<u>CLIENT SAMPLE ID</u>	<u>D/T ANALYZED</u>	<u>DATA FILE</u>
25	B-DU1-ISM1-8	2017-03-27 13:57	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar011.d\27mar011.r
26	B-DU1-ISM2-8	2017-03-27 14:17	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar012.d\27mar012.r
27	B-DU1-ISM3-8	2017-03-27 14:38	Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar013.d\27mar013.r

CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET FOR METHOD: EPA 8270C SIM PAHS

CCV WORK ORDER: 099-06-009-4997-4741
INSTRUMENT: GC/MS EEE
BATCH ID: 1703131001
INITIAL: 170327A010
CCV:

ANALYZED BY: 907
D/T ANALYZED: 2017-03-13 14:13
INITIAL: 2017-03-27 10:55
CCV: 262
REVIEWED BY: 2017-03-27 13:13
D/T REVIEWED:

Data File: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar002.d\27mar002.r

COMPOUND	TYPE	CALIB MODEL	MIN RF	AVG RF	CCV RF	AMOUNT	CCV CONC	CCV %D	CCV %D CL	STATUS
Naphthalene	Avg Resp		0.00	1.085	1.196			-10	0-20	PASS
2-Methylnaphthalene	Avg Resp		0.00	0.661	0.759			-15	0-20	PASS
1-Methylnaphthalene	Avg Resp		0.00	0.663	0.699			-5	0-20	PASS
Acenaphthylene	Avg Resp		0.00	2.528	2.700			-7	0-20	PASS
Acenaphthene	Avg Resp	C	0.00	1.517	1.630			-7	0-20	PASS
Fluorene	Avg Resp		0.00	1.689	1.784			-6	0-20	PASS
Phenanthrene	Avg Resp		0.00	1.061	1.111			-5	0-20	PASS
Anthracene	Avg Resp		0.00	1.037	1.121			-8	0-20	PASS
Fluoranthene	Avg Resp	C	0.00	1.365	1.448			-6	0-20	PASS
Pyrene	Avg Resp		0.00	1.362	1.469			-8	0-20	PASS
Benzo (a) Anthracene	Avg Resp		0.00	1.317	1.349			-2	0-20	PASS
Chrysene	Avg Resp		0.00	1.197	1.323			-11	0-20	PASS
Benzo (k) Fluoranthene	Avg Resp		0.00	1.366	1.436			-5	0-20	PASS
Benzo (b) Fluoranthene	Avg Resp		0.00	1.303	1.437			-10	0-20	PASS
Benzo (a) Pyrene	Avg Resp	C	0.00	1.256	1.339			-7	0-20	PASS
Indeno (1,2,3-c,d) Pyrene	Avg Resp		0.00	1.502	1.558			-4	0-20	PASS
Dibenz (a,h) Anthracene	Avg Resp		0.00	1.139	1.237			-9	0-20	PASS
Benzo (g,h,i) Perylene	Avg Resp		0.00	1.207	1.261			-4	0-20	PASS

MIN RF: Method Specified Minimum Response Factor



INTERNAL STANDARD COMPOUNDS AREA REPORT FOR METHOD: EPA 8270C SIM PAHs

ICAL BATCH ID: 170313I001

CCV BATCH ID: 170327A010

ICAL MIDPOINT

SAMPLE ID: 099-06-009-4979

D/T ANALYZED: 2017-03-13 13:33

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar008.d\13mar008.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>RETENTION TIME</u>
1,4-Dichlorobenzene-d4	46489	3.41
Naphthalene-d8	140022	4.66
Acenaphthene-d10	63153	6.54
Phenanthrene-d10	210133	8.14
Chrysene-d12	220711	11.06
Perylene-d12	208071	12.82

ICV

SAMPLE ID 099-06-009-4979

D/T ANALYZED: 2017-03-13 14:34

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170313\13mar011.d\13mar011.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	44664	23244	92978	3.42	PASS
Naphthalene-d8	134376	70011	280044	4.65	PASS
Acenaphthene-d10	61082	31576	126306	6.54	PASS
Phenanthrene-d10	200519	105066	420266	8.14	PASS
Chrysene-d12	207148	110356	441422	11.06	PASS
Perylene-d12	200860	104036	416142	12.82	PASS

CCV

SAMPLE ID 099-06-009-4997

D/T ANALYZED: 2017-03-27 10:55

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar002.d\27mar002.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	55771	23244	92978	3.29	PASS
Naphthalene-d8	153732	70011	280044	4.53	PASS
Acenaphthene-d10	70535	31576	126306	6.41	PASS
Phenanthrene-d10	237225	105066	420266	8.00	PASS
Chrysene-d12	244953	110356	441422	10.92	PASS
Perylene-d12	230206	104036	416142	12.65	PASS

MB

SAMPLE ID 099-14-035-386

D/T ANALYZED: 2017-03-27 11:15

DATA FILE: Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar003.d\27mar003.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	53792	27886	111542	3.30	PASS
Naphthalene-d8	151636	76866	307464	4.53	PASS
Acenaphthene-d10	71653	35268	141070	6.41	PASS
Phenanthrene-d10	255973	118612	474450	8.00	PASS
Chrysene-d12	264101	122476	489906	10.92	PASS
Perylene-d12	249943	115103	460412	12.65	PASS

LCS**SAMPLE ID** 099-14-035-386**D/T ANALYZED:** 2017-03-27 11:35**DATA FILE:** Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar004.d\27mar004.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	59177	27886	111542	3.30	PASS
Naphthalene-d8	161558	76866	307464	4.53	PASS
Acenaphthene-d10	75915	35268	141070	6.41	PASS
Phenanthrene-d10	258057	118612	474450	8.01	PASS
Chrysene-d12	273070	122476	489906	10.92	PASS
Perylene-d12	264968	115103	460412	12.65	PASS

MS**SAMPLE ID** 17-03-1753-6**D/T ANALYZED:** 2017-03-27 11:56**DATA FILE:** Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar005.d\27mar005.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	57973	27886	111542	3.30	PASS
Naphthalene-d8	150857	76866	307464	4.53	PASS
Acenaphthene-d10	71434	35268	141070	6.42	PASS
Phenanthrene-d10	252691	118612	474450	8.01	PASS
Chrysene-d12	265492	122476	489906	10.92	PASS
Perylene-d12	256493	115103	460412	12.66	PASS

MSD**SAMPLE ID** 17-03-1753-6**D/T ANALYZED:** 2017-03-27 12:16**DATA FILE:** Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar006.d\27mar006.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	58479	27886	111542	3.30	PASS
Naphthalene-d8	152866	76866	307464	4.53	PASS
Acenaphthene-d10	72569	35268	141070	6.42	PASS
Phenanthrene-d10	255933	118612	474450	8.01	PASS
Chrysene-d12	257728	122476	489906	10.93	PASS
Perylene-d12	255038	115103	460412	12.67	PASS

CS**SAMPLE ID** 17-03-1557-25**D/T ANALYZED:** 2017-03-27 13:57**DATA FILE:** Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar011.d\27mar011.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	59986	27886	111542	3.30	PASS
Naphthalene-d8	169237	76866	307464	4.54	PASS
Acenaphthene-d10	80719	35268	141070	6.42	PASS
Phenanthrene-d10	267245	118612	474450	8.02	PASS
Chrysene-d12	260860	122476	489906	10.93	PASS
Perylene-d12	254564	115103	460412	12.68	PASS

CS

SAMPLE ID 17-03-1557-26**D/T ANALYZED:** 2017-03-27 14:17**DATA FILE:** Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar012.d\27mar012.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	60589	27886	111542	3.30	PASS
Naphthalene-d8	171804	76866	307464	4.53	PASS
Acenaphthene-d10	81274	35268	141070	6.42	PASS
Phenanthrene-d10	274687	118612	474450	8.02	PASS
Chrysene-d12	279862	122476	489906	10.94	PASS
Perylene-d12	271055	115103	460412	12.68	PASS

CS

SAMPLE ID 17-03-1557-27**D/T ANALYZED:** 2017-03-27 14:38**DATA FILE:** Z:\GCMS_EEE\GCMS_EEE_data\2017\170327\27mar013.d\27mar013.rr

<u>COMPOUND</u>	<u>AREA</u>	<u>LOWER AREA LIMIT</u>	<u>UPPER AREA LIMIT</u>	<u>RETENTION TIME</u>	<u>STATUS</u>
1,4-Dichlorobenzene-d4	61080	27886	111542	3.30	PASS
Naphthalene-d8	171171	76866	307464	4.54	PASS
Acenaphthene-d10	82099	35268	141070	6.42	PASS
Phenanthrene-d10	275585	118612	474450	8.02	PASS
Chrysene-d12	277786	122476	489906	10.94	PASS
Perylene-d12	268126	115103	460412	12.69	PASS

Notes:

For all samples including QC, all internal standard area responses must be within 50% to 200% of the mean area response in the initial calibration.

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar002.d
 Report Date: 03/27/2017 13:02

Eurofins CalScience
 Calibration Verification Report

Instrument ID: GCMS_EEE.i Injection Date and Time: 27-MAR-2017 10:55
 Sample Name: CCV S010317F 1PPM Initial Calibration Date(s): 03-JAN-2017 13-MAR-2017
 Sublist used: all.sub Initial Calibration Time(s): 11:59 14:13
 Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m

Target Compounds	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / Drift /Drift	Max%D	Curve Type
Naphthalene	1.085	1.196	0.00	-10	20	Averaged
2-Methylnaphthalene	0.661	0.759	0.00	-15	20	Averaged
1-Methylnaphthalene	0.663	0.699	0.00	-5	20	Averaged
Acenaphthylene	2.528	2.700	0.00	-7	20	Averaged
Acenaphthene	1.517	1.630	0.00	-7	20	Averaged
Fluorene	1.689	1.784	0.00	-6	20	Averaged
Phenanthrene	1.061	1.111	0.00	-5	20	Averaged
Anthracene	1.037	1.121	0.00	-8	20	Averaged
Fluoranthene	1.365	1.448	0.00	-6	20	Averaged
Pyrene	1.362	1.469	0.00	-8	20	Averaged
Benzo (a) Anthracene	1.317	1.349	0.00	-2	20	Averaged
Chrysene	1.197	1.323	0.00	-11	20	Averaged
Benzo (b) Fluoranthene	1.303	1.437	0.00	-10	20	Averaged
Benzo (k) Fluoranthene	1.366	1.436	0.00	-5	20	Averaged
Benzo (a) Pyrene	1.256	1.339	0.00	-7	20	Averaged
Indeno (1,2,3-c,d) Pyrene	1.502	1.558	0.00	-4	20	Averaged
Dibenz (a,h) Anthracene	1.139	1.237	0.00	-9	20	Averaged
Benzo (g,h,i) Perylene	1.207	1.261	0.00	-4	20	Averaged
Biphenyl	2.240	2.370	0.00	-6	20	Averaged
2,6-Dimethylnaphthalene	1.572	1.656	0.00	-5	20	Averaged
1,6,7-Trimethylnaphthalene	1.571	1.650	0.00	-5	20	Averaged
Dibenzothiophene	4.622	4.970	0.00	-8	20	Averaged
1-Methylphenanthrene	0.997	1.065	0.00	-7	20	Averaged
Benzo (e) pyrene	1.557	1.644	0.00	-6	20	Averaged
Perylene	1.493	1.583	0.00	-6	20	Averaged
Dibenzofuran	2.194	2.370	0.00	-8	20	Averaged
Surrogate Standards	ICAL RRF or Amount	ICV RRF	Min. RRF	%D / Drift /Drift	Max%D	Curve Type
Nitrobenzene-d5	0.331	0.379	0.00	-15	20	Averaged
2-Fluorobiphenyl	1.805	1.918	0.00	-6	20	Averaged
p-Terphenyl-d14	0.775	0.813	0.00	-5	20	Averaged

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar002.d Instrument ID: GCMS_EEE.i
 Injection date and time: 27-MAR-2017 10:55 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
 Calibration date and time: 27-MAR-2017 11:23
 Date, time and analyst ID of latest file update: 27-Mar-2017 11:23 ev7p

Sample Name: CCV S010317F 1PPM Misc Info: 170327A010
 Response via Initial Calibration

Compounds	I.S. Ref.	RT	QIon	Area	On-Column Amount (mg/L)	DEV(Min)
Internal Standards						
1)*1,4-Dichlorobenzene-d4	(1)	3.292	152	55771	5.000	0.00
3)*Naphthalene-d8	(2)	4.530	136	153732	5.000	0.00
11)*Acenaphthene-d10	(3)	6.411	164	70535	5.000	0.00
17)*Phenanthrene-d10	(4)	8.004	188	237225	5.000	0.00
31)*Perylene-d12	(6)	12.655	264	230206	5.000	0.00
25)*Chrysene-d12	(5)	10.920	240	244953	5.000	0.00
System Monitoring Compounds						
2)\$Nitrobenzene-d5	(2)	3.845	82	11664	1.145	0.00
SpikedAmount 1.000			Recovery =	0.000		
7)\$2-Fluorobiphenyl	(3)	5.709	172	27056	1.063	0.00
SpikedAmount 1.000			Recovery =	0.000		
23)\$p-Terphenyl-d14	(5)	9.801	244	39849	1.050	0.00
SpikedAmount 1.000			Recovery =	0.000		
Target Compounds						
4) Naphthalene	(2)	4.551	128	36766	1.103	100
5) 2-Methylnaphthalene	(2)	5.279	142	23323	1.147	100
6) 1-Methylnaphthalene	(2)	5.391	142	21477	1.053	100
10) Acenaphthylene	(3)	6.243	152	38084	1.068	100
12) Acenaphthene	(3)	6.444	153	22989	1.074	100
15) Fluorene	(3)	6.995	166	25168	1.056	100
18) Phenanthrene	(4)	8.029	178	52735	1.048	100
19) Anthracene	(4)	8.079	178	53189	1.081	100
21) Fluoranthene	(4)	9.337	202	68699	1.061	100
22) Pyrene	(5)	9.574	202	71943	1.079	100
24) Benzo (a) Anthracene	(5)	10.903	228	66098	1.024	100
26) Chrysene	(5)	10.946	228	64808	1.105	100
27) Benzo (b) Fluoranthene	(6)	12.196	252	66156	1.103	100
28) Benzo (k) Fluoranthene	(6)	12.224	252	66131	1.051	100
30) Benzo (a) Pyrene	(6)	12.583	252	61661	1.066	100
33) Indeno (1,2,3-c,d) Pyrene	(6)	13.888	276	71745	1.038	100
34) Dibenz (a,h) Anthracene	(6)	13.918	278	56952	1.086	100
35) Benzo (g,h,i) Perylene	(6)	14.159	276	58066	1.045	100
8) Biphenyl	(3)	5.796	154	33437	1.058	100
9) 2,6-Dimethylnaphthalene	(3)	5.955	156	23361	1.053	100
14) 1,6,7-Trimethylnaphthalene	(3)	6.876	170	23270	1.050	100
16) Dibenzothiophene	(3)	7.893	184	70117	1.075	100

* = Compound is an internal standard.

\$ = Compound is a surrogate standard.

Quant Report

Target Revision 3.5

Data File: /chem/SVOA/GCMS_EEE.i/170327.b/27mar002.d Instrument ID: GCMS_EEE.i
Injection date and time: 27-MAR-2017 10:55 Analyst ID: 907

Method used: /chem/SVOA/GCMS_EEE.i/170327.b/simpah-extra.m Sublist used: all
Calibration date and time: 27-MAR-2017 11:23
Date, time and analyst ID of latest file update: 27-Mar-2017 11:23 ev7p

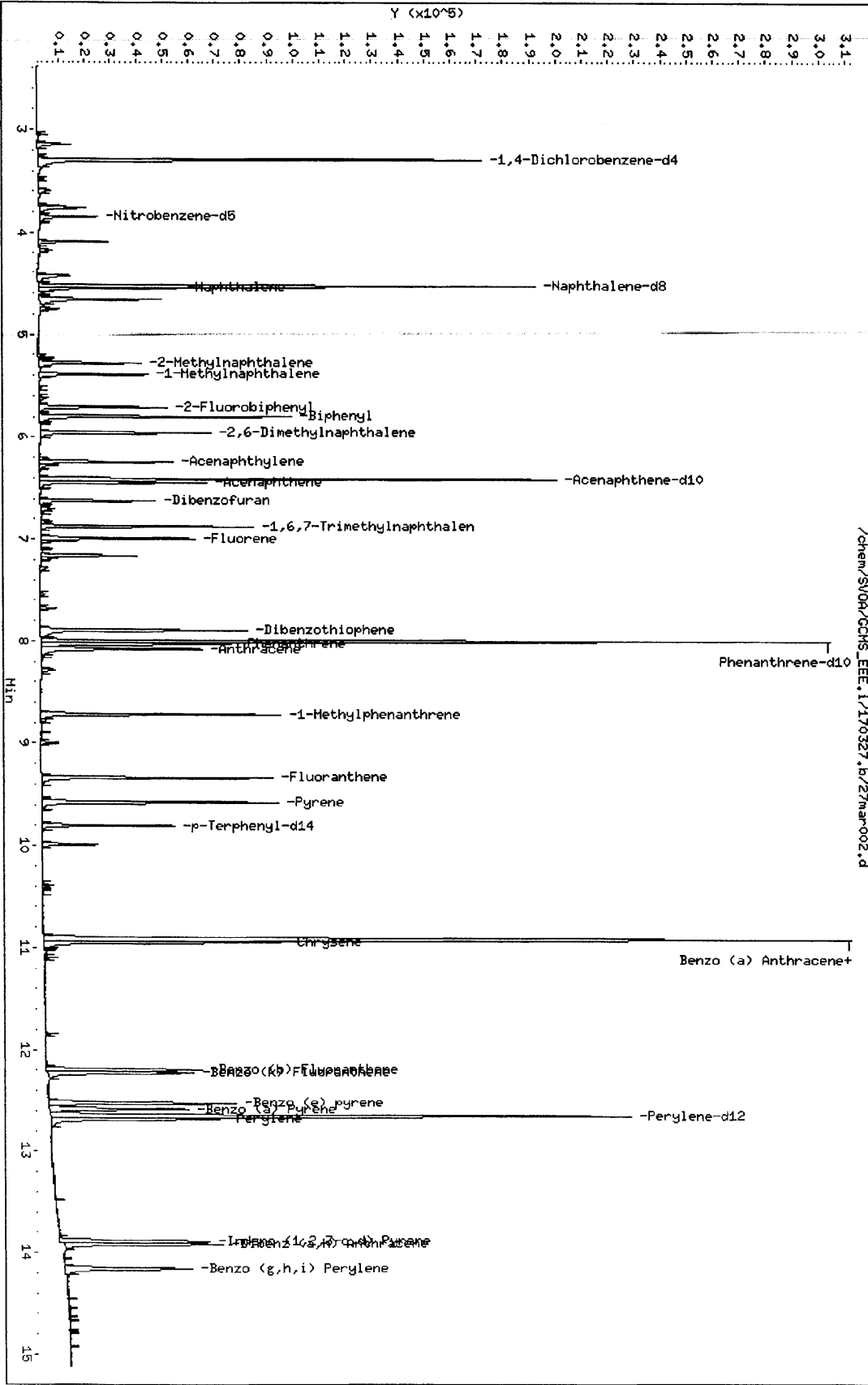
Sample Name: CCV S010317F 1PPM Misc Info: 170327A010
Response via Initial Calibration

Compounds	I.S.		QIon	Area	On-Column	
	Ref.	RT			Amount (mg/L)	QValue
20) 1-Methylphenanthrene	(5)	8.718	192	52168	1.068	100
29) Benzo (e) pyrene	(6)	12.522	252	75681	1.056	100
32) Perylene	(6)	12.683	252	72865	1.060	100
13) Dibenzofuran	(3)	6.621	168	33429	1.080	100

page 2 of 2

Data File: /chem/SV09/CCHS_EEE.1/170327.b/27mar002.d
Date: 27-MAR-2017 10:55
Client ID:
Sample Info: CCV S010317F 1PPM
Column phase: J&W DB-5MS

Instrument: CCHS_EEE.i
Operator: 907
Column diameter: 0.18



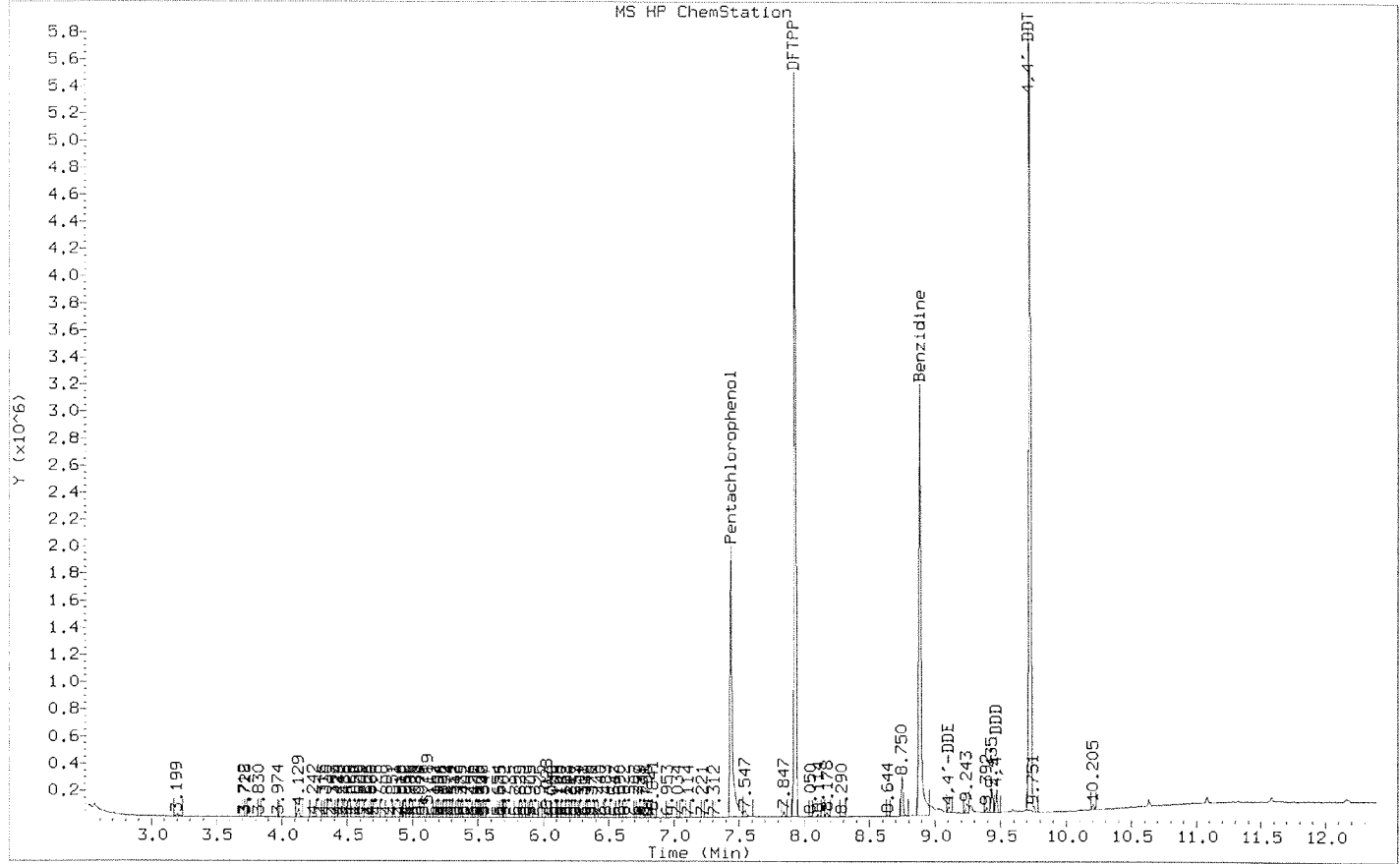
EPA METHOD 8270C PAHSIM

Tuning Reports

DFTPP TUNE/TAILING FACTOR/DEGRADATION SAMPLE AND GRAPHIC REPORT

Report Generated Time Mon Mar 13 15:07:34 2017

Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 1
 Acq on : 13-MAR-2017 11:49 Operator : 907
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Last Update : 07-MAR-2017 10:28



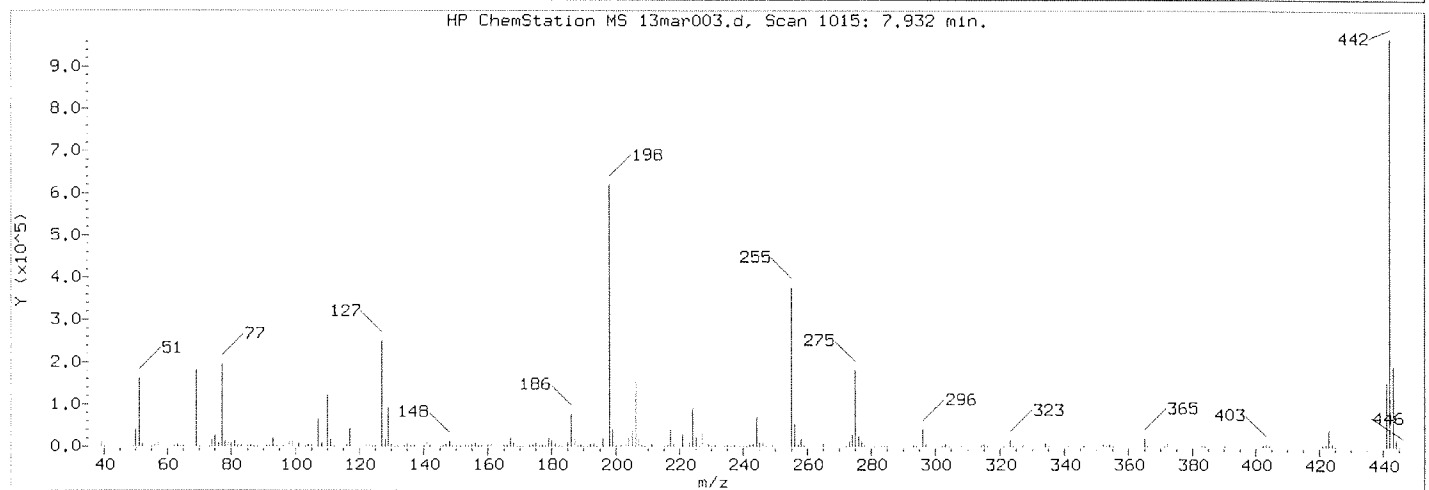
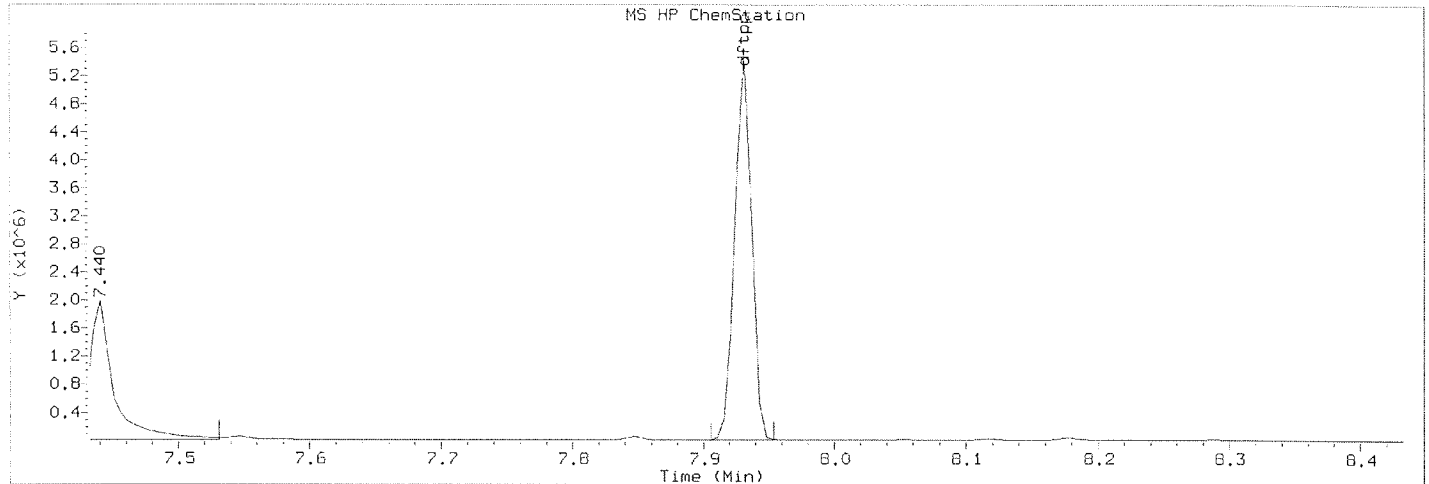
Tune *** PASSED ***
 Pentachlorophenol Tailing *** PASSED ***
 Benzidine Tailing *** PASSED ***
 DDT degradation *** PASSED ***

Tuning Sample, /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d , *** PASSED ***



Report Generated Time Mon Mar 13 15:07:34 2017

Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 1
 Acq on : 13-MAR-2017 11:49 Operator : 907
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170313.b/dftpptune.m
 Last Update : 07-MAR-2017 10:28



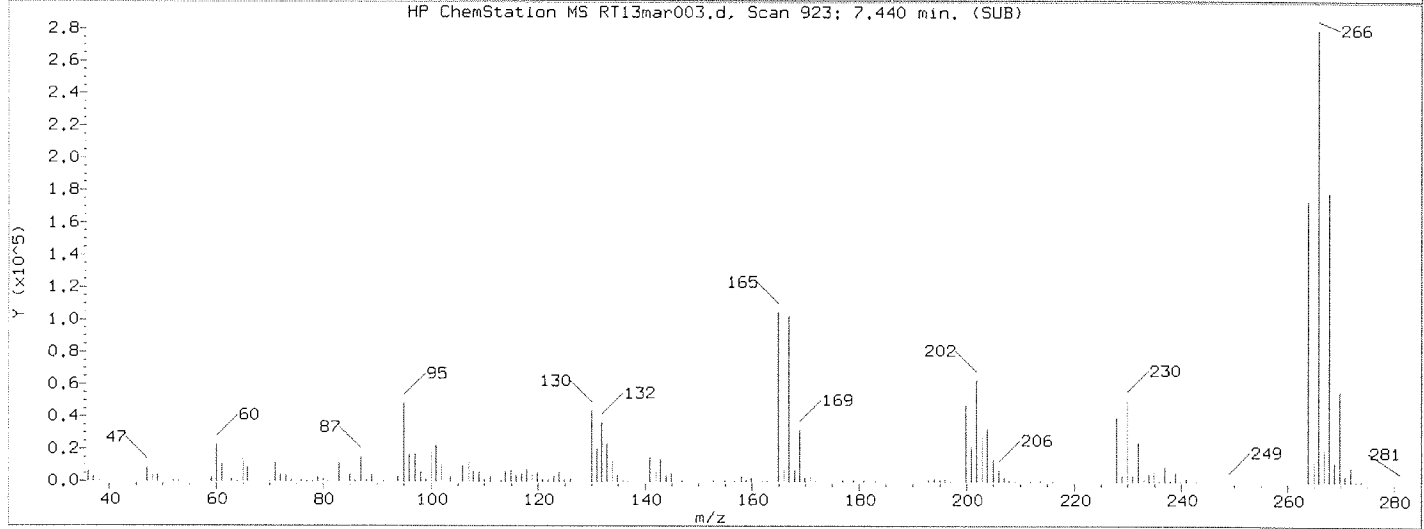
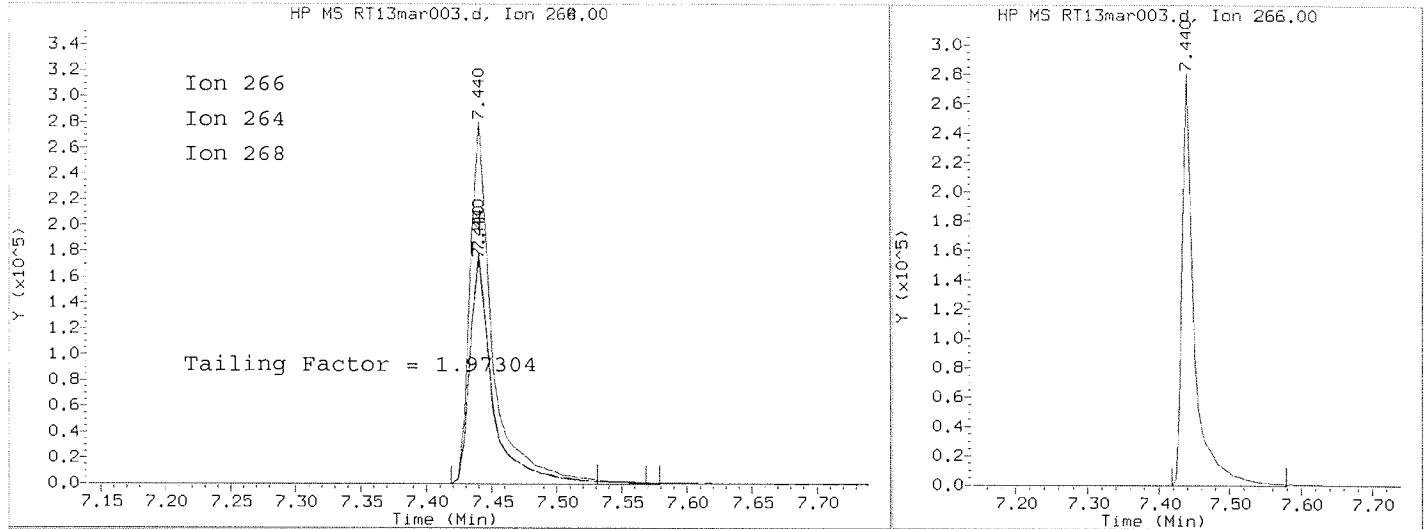
Spectrum: Avg. Scans 1014-1016 (7.93), Background Scan 1009

DFTPP Ion Abundance/Ratio Criteria Chart

Ion	Abundance Criteria	Base Peak	Response	Test
198	Base Peak, 100% relative abundance	100.00	461312	PASS
51	30 - 60% of mass 198	33.56	154816	PASS
68	Less than 2% of mass 69	0.00	0	PASS
69	Less than mass 198	36.32	167552	PASS
70	Less than 2% of mass 69	0.62	1045	PASS
127	40 - 60% of mass 198	44.77	206528	PASS
197	0 - 1% of mass 198	0.00	0	PASS
199	5 - 9% of mass 198	6.99	32256	PASS
275	10 - 30% of mass 198	25.66	118368	PASS
365	1 - 100% of mass 198	2.66	12269	PASS
441	Present, but less than mass 443	74.01	76992	PASS
442	40 - 200% of mass 198	109.96	507264	PASS
443	17 - 23% of mass 442	20.51	104024	PASS

Report Generated Time Mon Mar 13 15:07:34 2017

Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 3
 Acq on : 13-MAR-2017 11:49 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d/resolut.m
 Last Update : 13-MAR-2017 12:08



Pentachlorophenol

=====
 Exp. RT = 7.467
 Found RT = 7.440

Mass	Area	Ratio
266	343885	100.00
264	216440	62.94
268	218874	63.65

Peak baseline front width (sec) : 0.816
 Peak baseline tail width (sec) : 1.610
 Tail Factor = 1.610 / 0.816

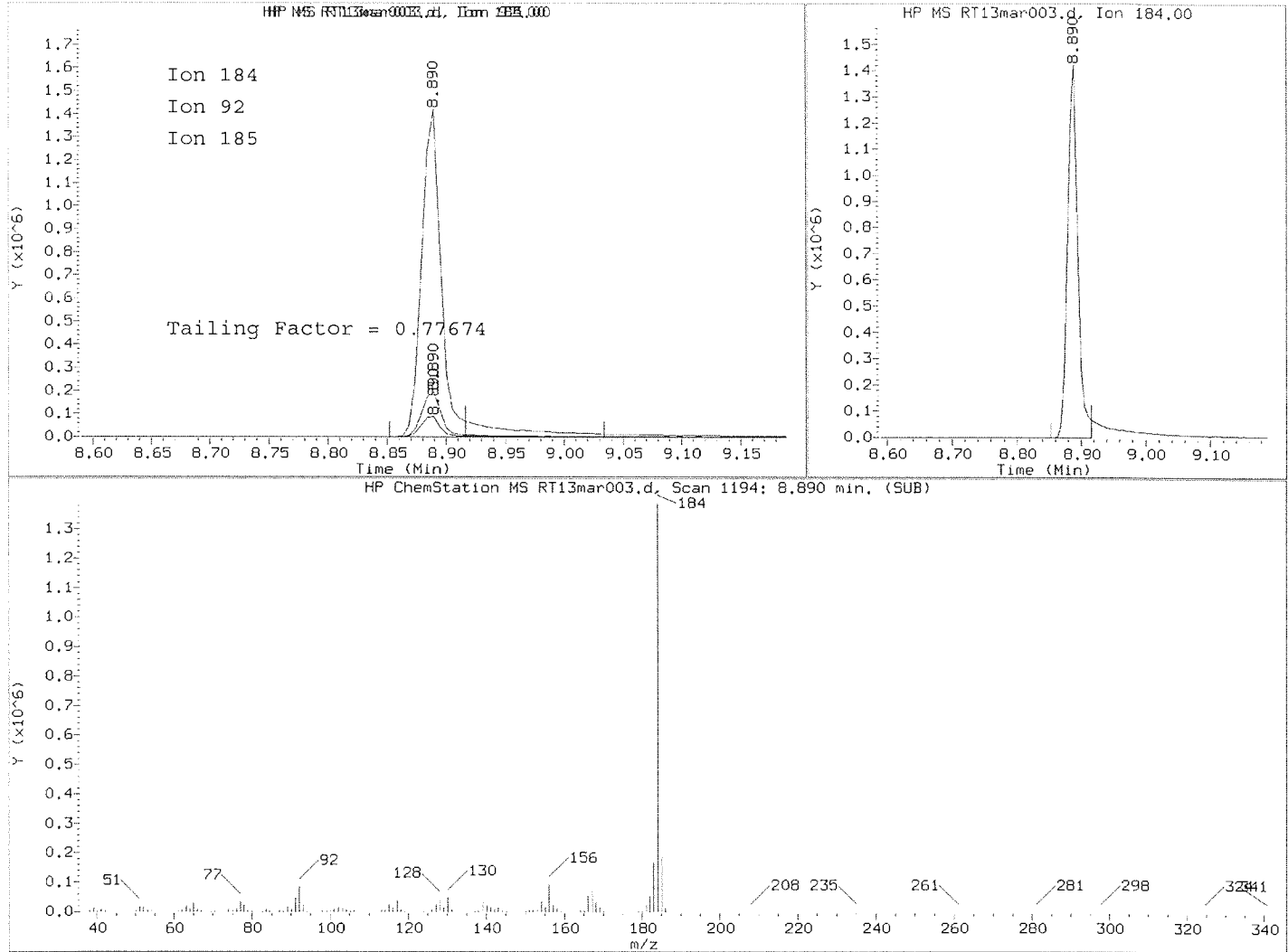
Tailing factor for Pentachlorophenol OK

Tail Factor = 1.973 Maximum Allowed = 3.0



Report Generated Time Mon Mar 13 15:07:34 2017

Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 3
 Acq on : 13-MAR-2017 11:49 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d/resolut.m
 Last Update : 13-MAR-2017 12:08



Benzidine

=====
 Exp. RT = 8.911
 Found RT = 8.890

Mass	Area	Ratio
184	1603178	100.00
92	113552	7.08
185	222697	13.89

Peak baseline front width (sec) : 1.066
 Peak baseline tail width (sec) : 0.828
 Tail Factor = 0.828/ 1.066

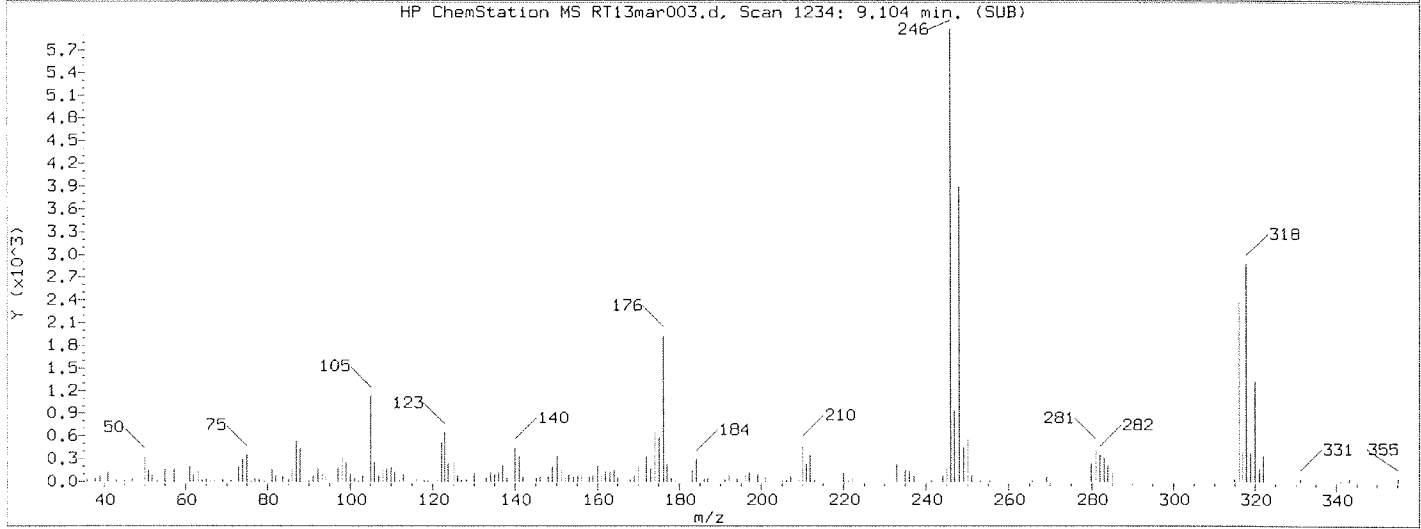
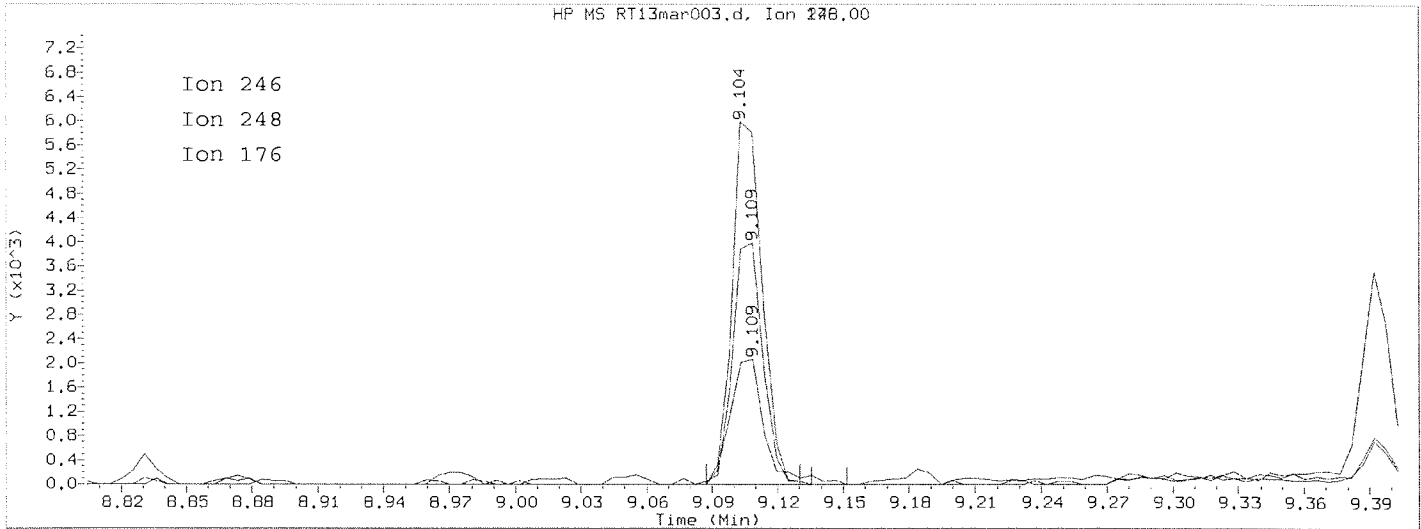
Tailing factor for Benzidine OK

Tail Factor = 0.777 Maximum Allowed = 3.0



Report Generated Time Mon Mar 13 15:07:34 2017

Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 3
 Acq on : 13-MAR-2017 11:49 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d/resolut.m
 Last Update : 13-MAR-2017 12:08



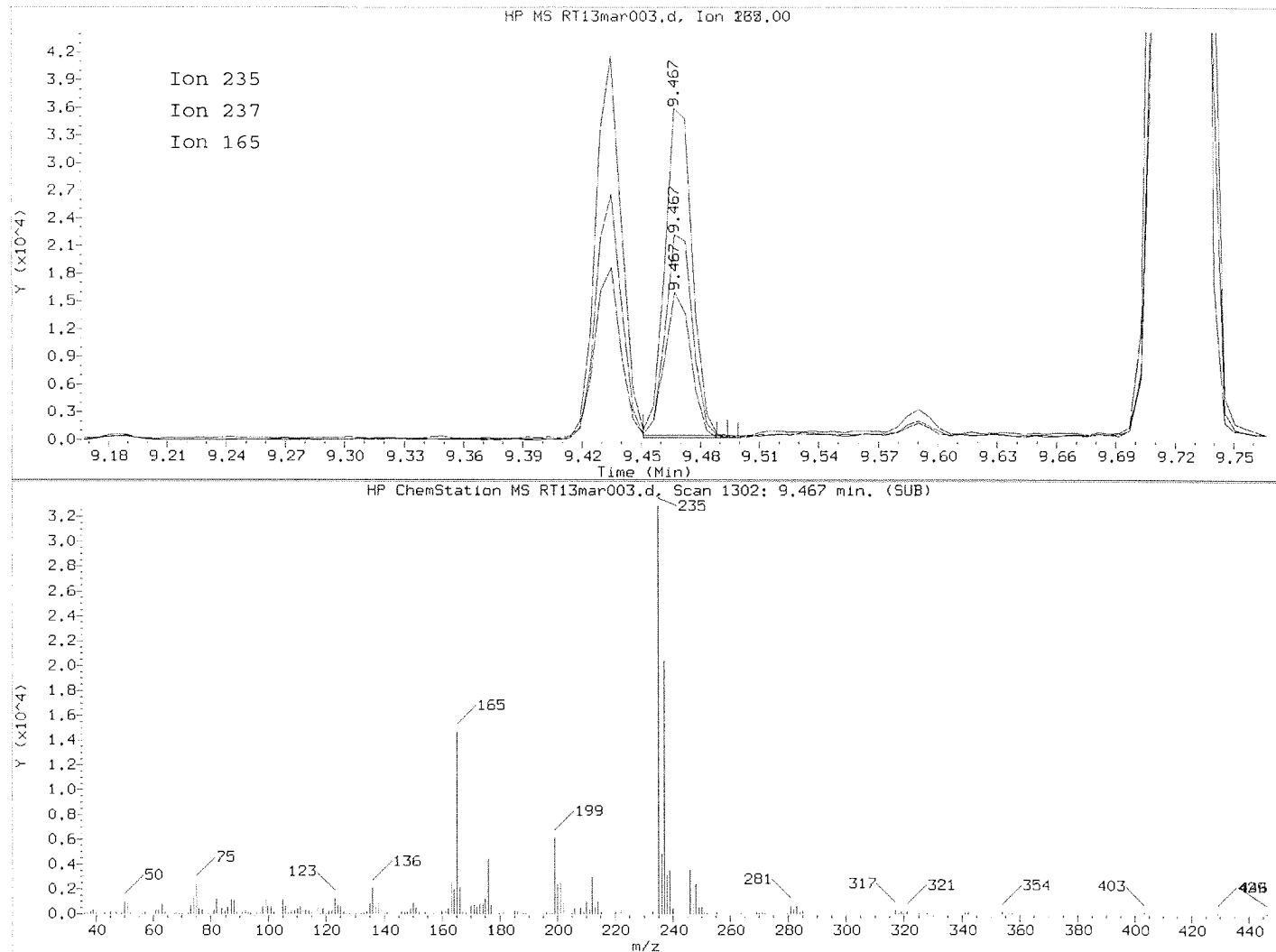
4,4'-DDE
 =====
 Exp. RT = 9.344
 Found RT = 9.104

Mass	Area	Ratio
246	5722	100.00
248	3838	67.08
176	2263	39.55

Return to Contents

Report Generated Time Mon Mar 13 15:07:34 2017

Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 3
 Acq on : 13-MAR-2017 11:49 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d/resolut.m
 Last Update : 13-MAR-2017 12:08

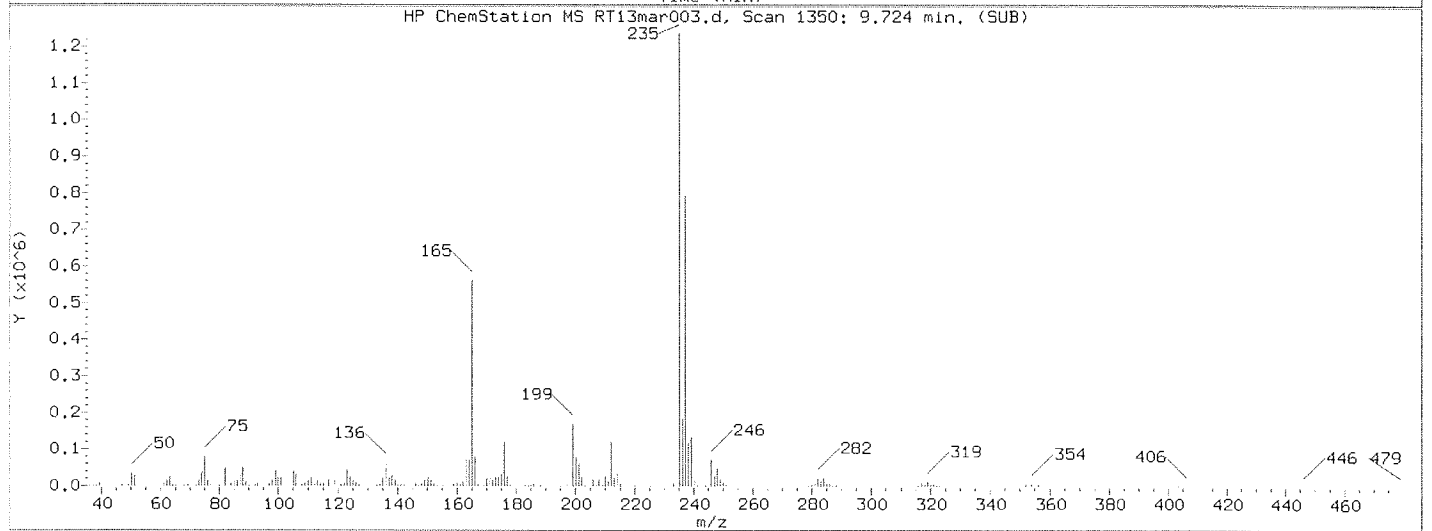
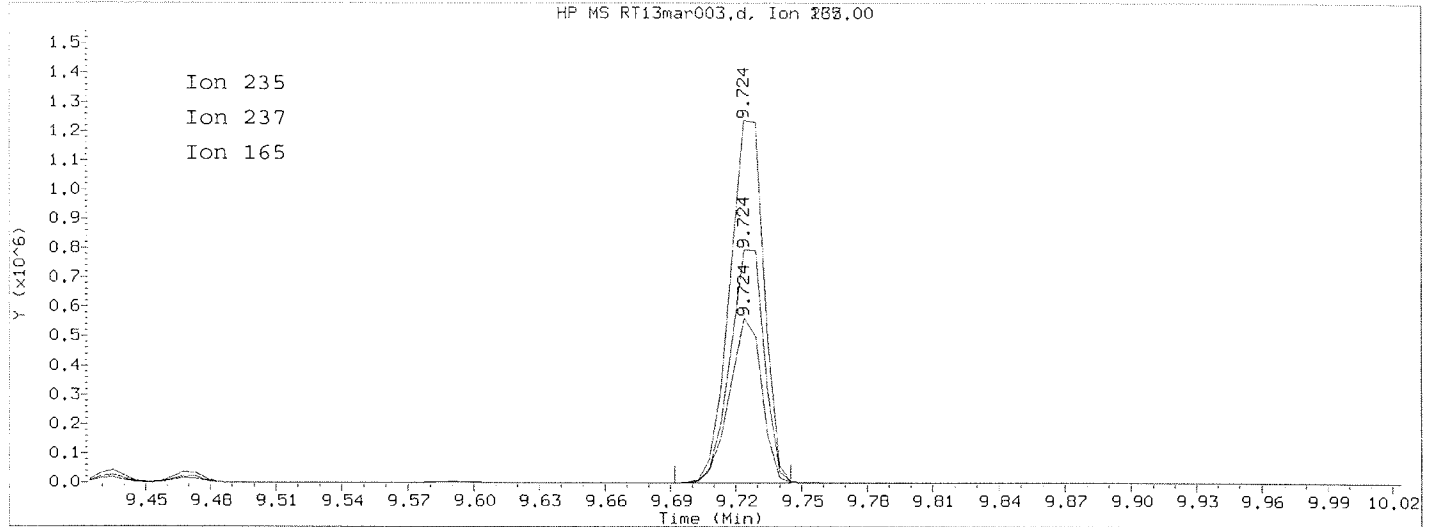


=====
 Exp. RT = 9.499
 Found RT = 9.467

Mass	Area	Ratio
235	34099	100.00
237	21311	62.50
165	14557	42.69

Report Generated Time Mon Mar 13 15:07:34 2017

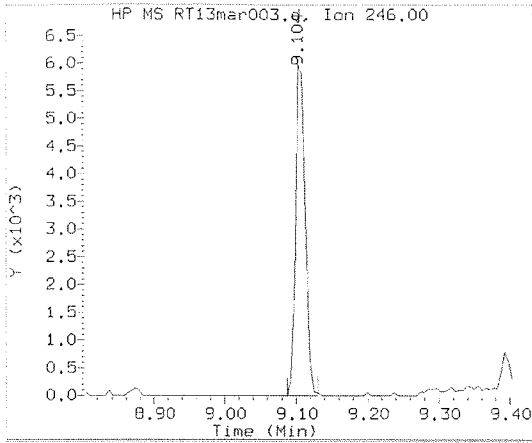
Data File : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d
 ALS Vial : 3
 Acq on : 13-MAR-2017 11:49 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170313.b/13mar003.d/resolut.m
 Last Update : 13-MAR-2017 12:08



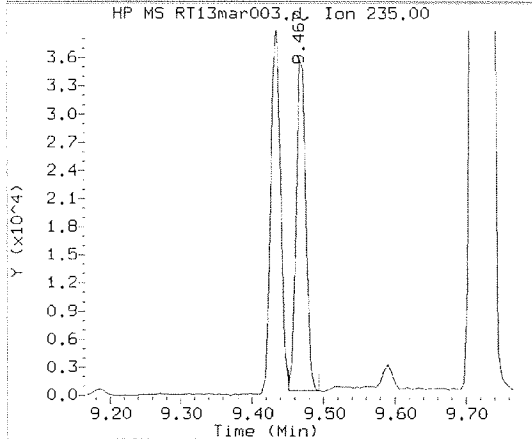
4,4'-DDT

=====
 Exp. RT = 9.756
 Found RT = 9.724

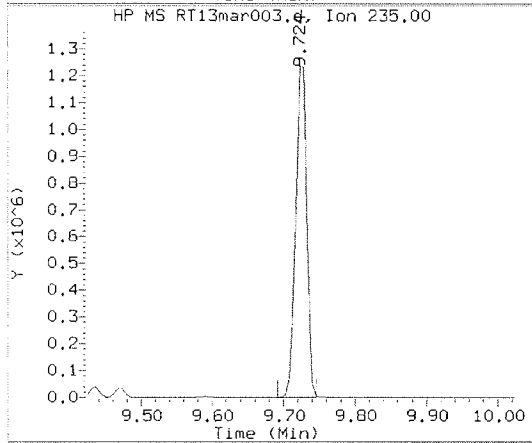
Mass	Area	Ratio
235	1352361	100.00
237	866214	64.05
165	581673	43.01



Compound: 4,4'-DDE
 Quant Mass: 246
 RT: 9.104
 Area: 5722



Compound: 4,4'-DDD
 Quant Mass: 235
 RT: 9.467
 Area: 34099



Compound: 4,4'-DDT
 Quant Mass: 235
 RT: 9.724
 Area: 1352361

DDT DEGRADATION BREAKDOWN ANALYSIS SUMMARY

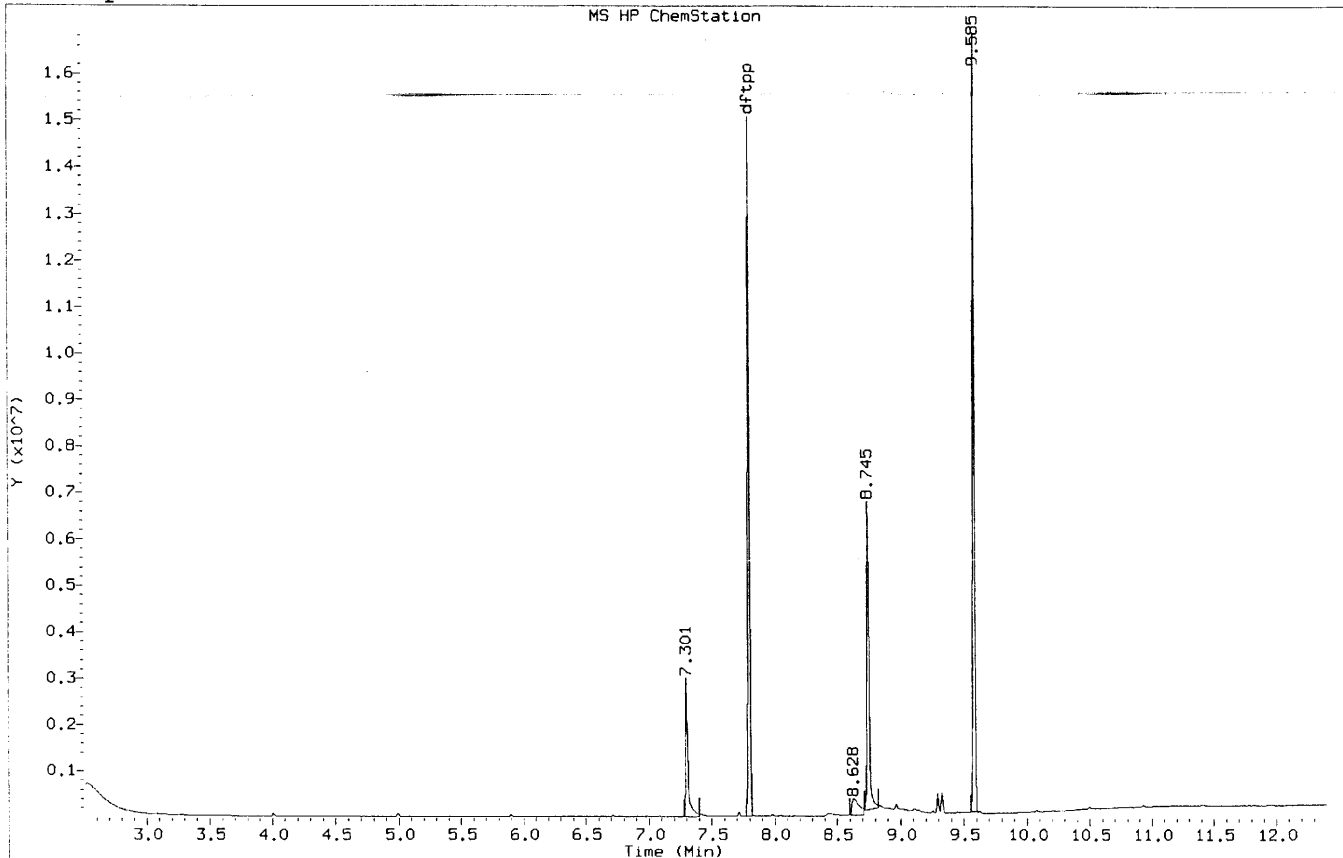
Compound	Response	%Breakdown	Max Allowed	Test
4,4-DDT	1352361			N/A
4,4-DDE	5722	0.42	20.0	PASS
4,4-DDD	34099	2.46	20.0	PASS
4,4-DDD + DDE	39821	2.9	20.0	PASS

TUNE SAMPLE *****
 *** PASSED *** DDT BREAKDOWN TEST

DFTPP TUNE/TAILING FACTOR/DEGRADATION SAMPLE AND GRAPHIC REPORT

Report Generated Time Mon Mar 27 10:55:20 2017

Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 1
 Acq on : 27-MAR-2017 10:33 Operator : 907
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Last Update : 07-MAR-2017 10:28

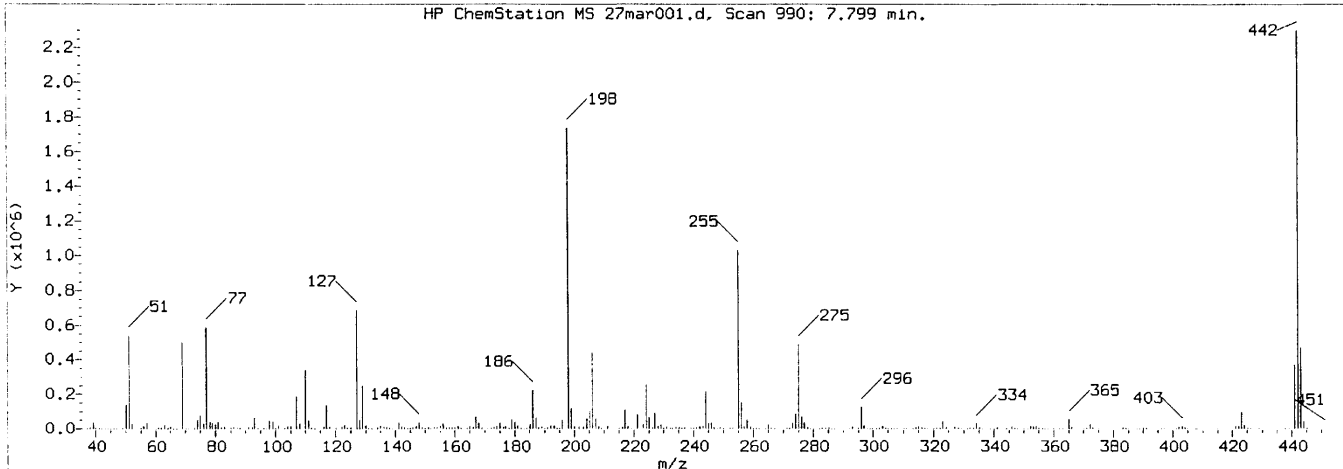
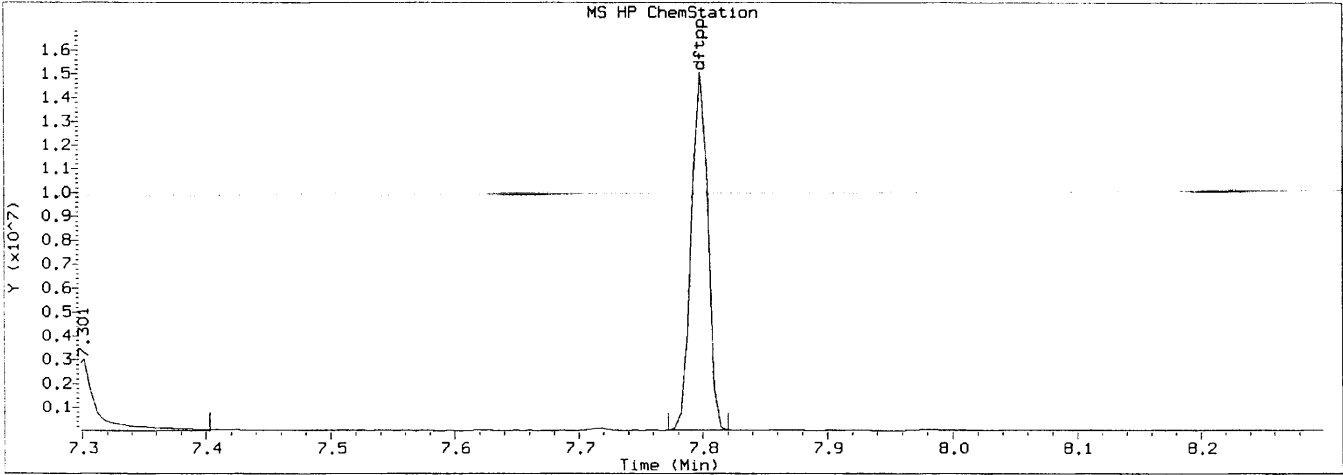


Tune *** PASSED ***
 Pentachlorophenol Tailing *** PASSED ***
 Benzidine Tailing *** PASSED ***
 DDT degradation *** PASSED ***

Tuning Sample, /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d ,*** PASSED ***

Report Generated Time Mon Mar 27 10:55:20 2017

Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 1
 Acq on : 27-MAR-2017 10:33 Operator : 907
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170327.b/dftpptune.m
 Last Update : 07-MAR-2017 10:28



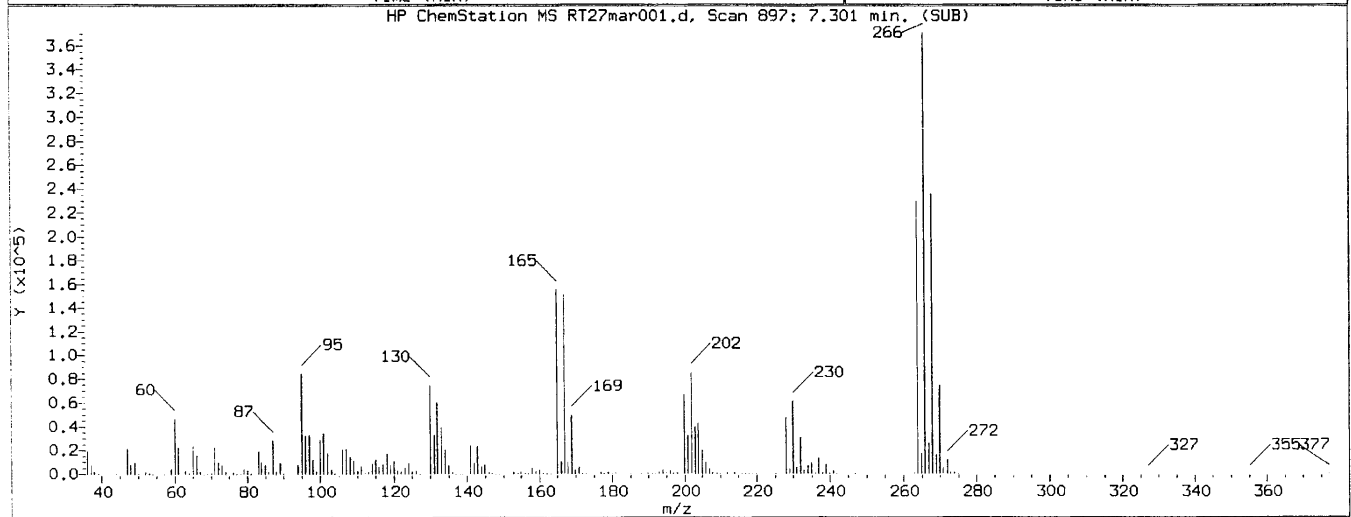
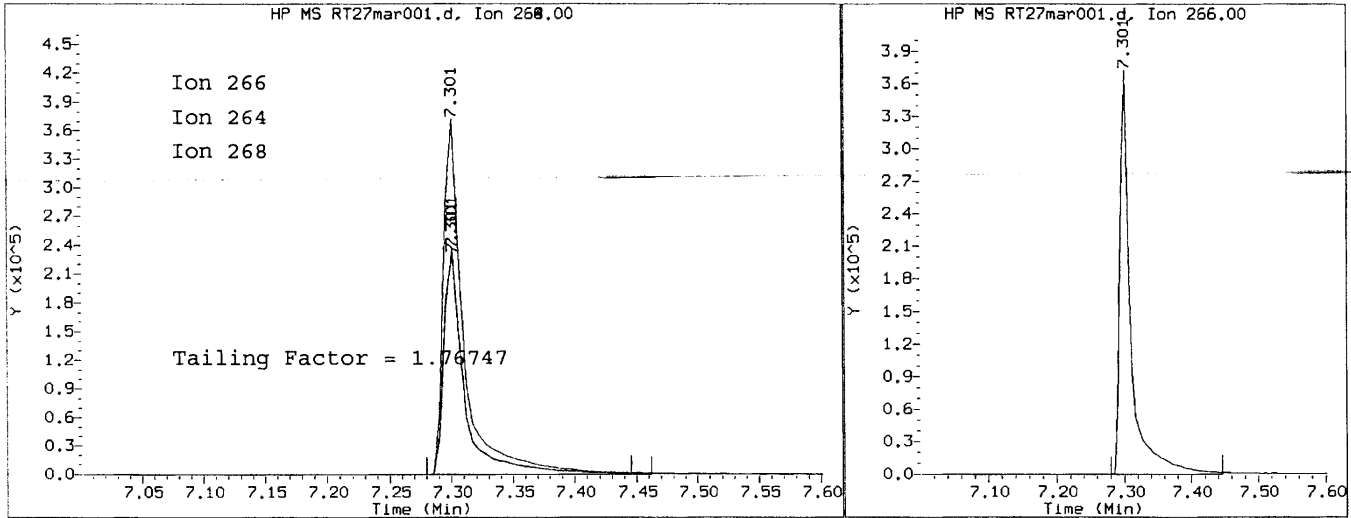
Spectrum: Avg. Scans 989-991 (7.80), Background Scan 984

DFTPP Ion Abundance/Ratio Criteria Chart

Ion	Abundance Criteria	Base Peak	Response	Test
198	Base Peak, 100% relative abundance	100.00	1337856	PASS
51	30 - 60% of mass 198	38.86	519936	PASS
68	Less than 2% of mass 69	0.00	0	PASS
69	Less than mass 198	34.78	465344	PASS
70	Less than 2% of mass 69	0.58	2710	PASS
127	40 - 60% of mass 198	43.81	586176	PASS
197	0 - 1% of mass 198	0.00	0	PASS
199	5 - 9% of mass 198	7.03	94072	PASS
275	10 - 30% of mass 198	23.51	314560	PASS
365	1 - 100% of mass 198	2.44	32632	PASS
441	Present, but less than mass 443	77.72	156736	PASS
442	40 - 200% of mass 198	74.63	998400	PASS
443	17 - 23% of mass 442	20.20	201664	PASS

Report Generated Time Mon Mar 27 10:55:20 2017

Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 3
 Acq on : 27-MAR-2017 10:33 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d/resolut.m
 Last Update : 27-MAR-2017 10:53



Pentachlorophenol

=====
 Exp. RT = 7.467
 Found RT = 7.301

Mass	Area	Ratio
266	435189	100.00
264	274310	63.03
268	276302	63.49

Peak baseline front width (sec) : 0.787
 Peak baseline tail width (sec) : 1.391
 Tail Factor = 1.391/ 0.787

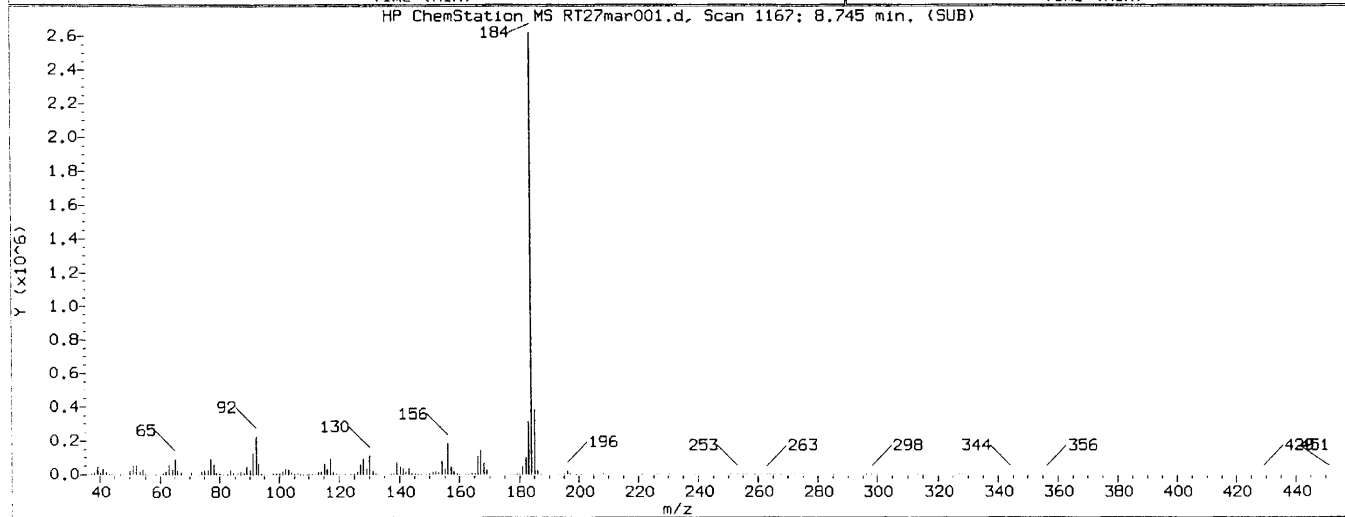
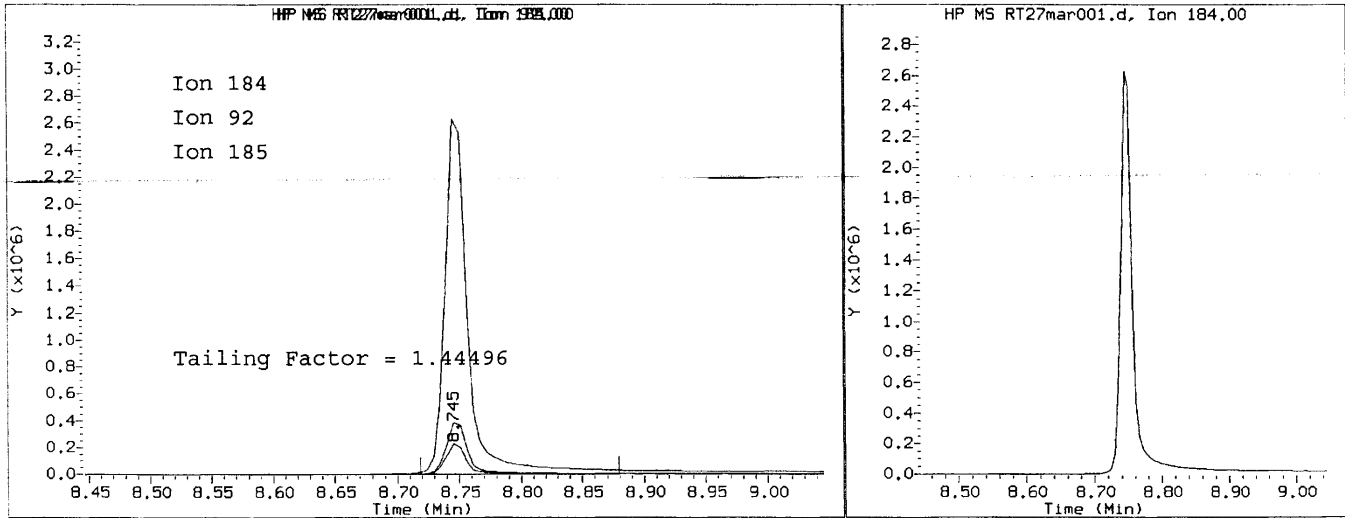
Tailing factor for Pentachlorophenol OK

Tail Factor = 1.767 Maximum Allowed = 3.0

Return to Contents

Report Generated Time Mon Mar 27 10:55:20 2017

Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 3
 Acq on : 27-MAR-2017 10:33 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d/resolut.m
 Last Update : 27-MAR-2017 10:53



Benzidine

=====
 Exp. RT = 8.911
 Found RT = 8.745

Mass	Area	Ratio
184	3677779	100.00
92	265404	7.22
185	484808	13.18

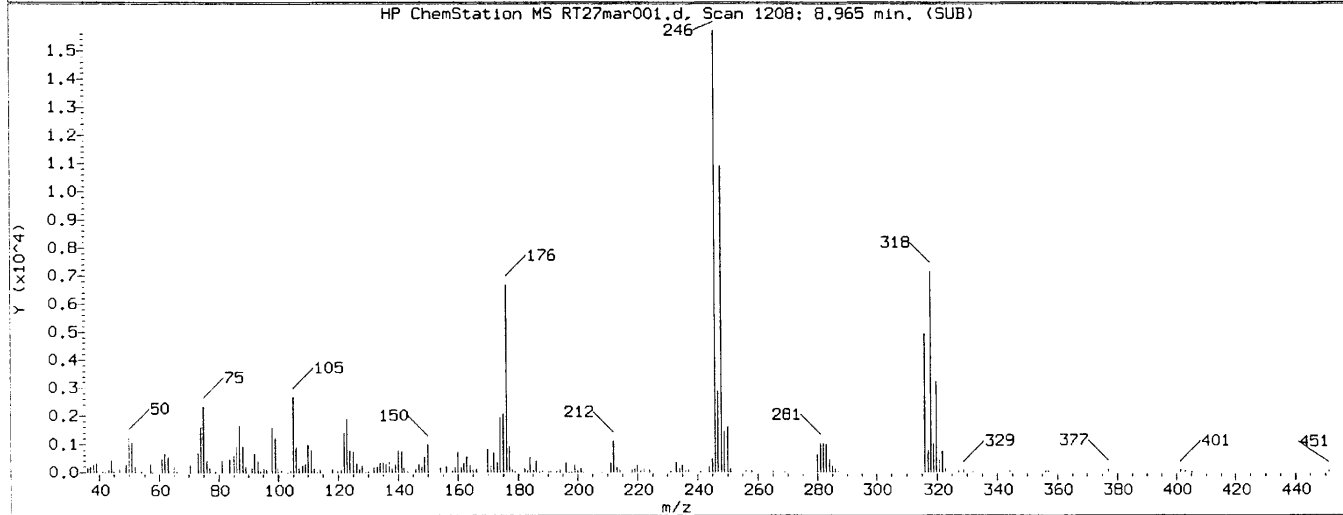
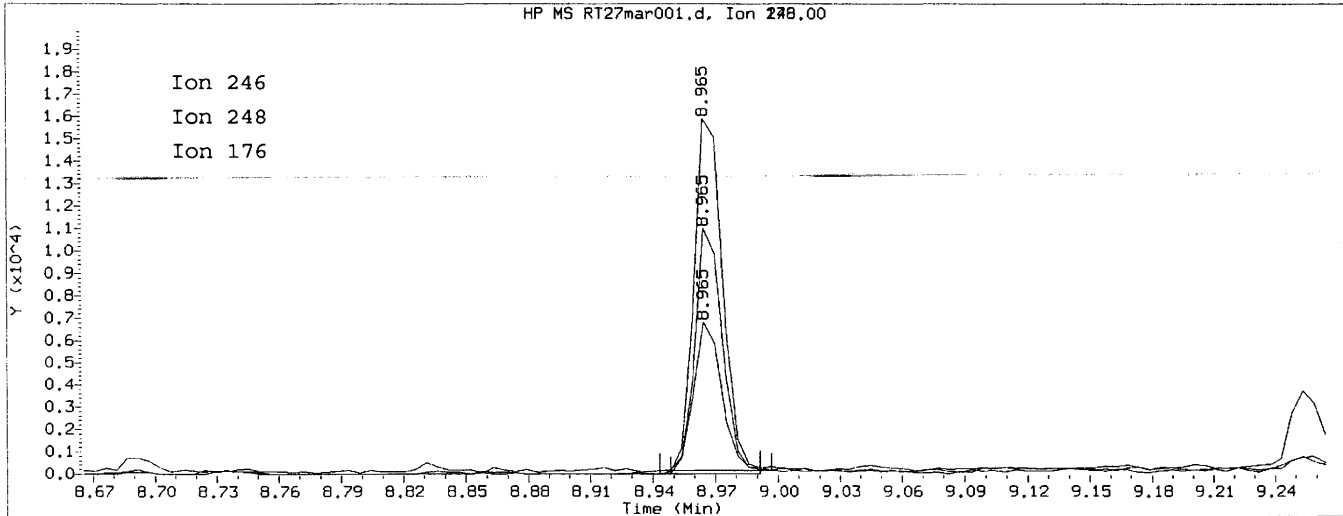
Peak baseline front width (sec) : 0.854
 Peak baseline tail width (sec) : 1.234
 Tail Factor = 1.234/ 0.854

Tailing factor for Benzidine OK

Tail Factor = 1.445 Maximum Allowed = 3.0

Report Generated Time Mon Mar 27 10:55:20 2017

Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 3
 Acq on : 27-MAR-2017 10:33 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d/resolut.m
 Last Update : 27-MAR-2017 10:53



4,4'-DDE

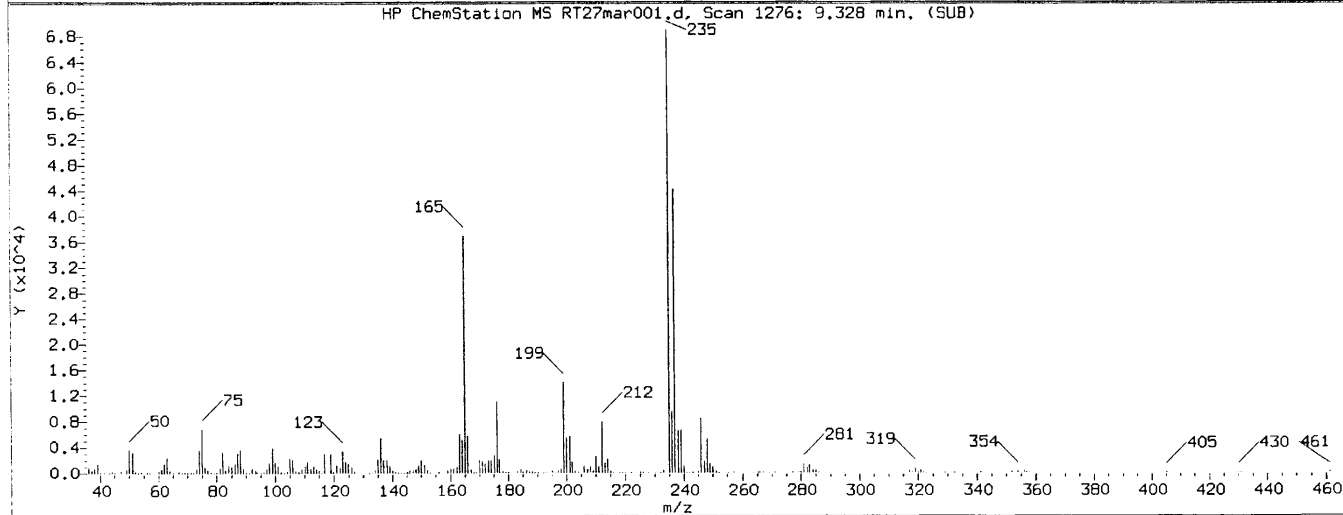
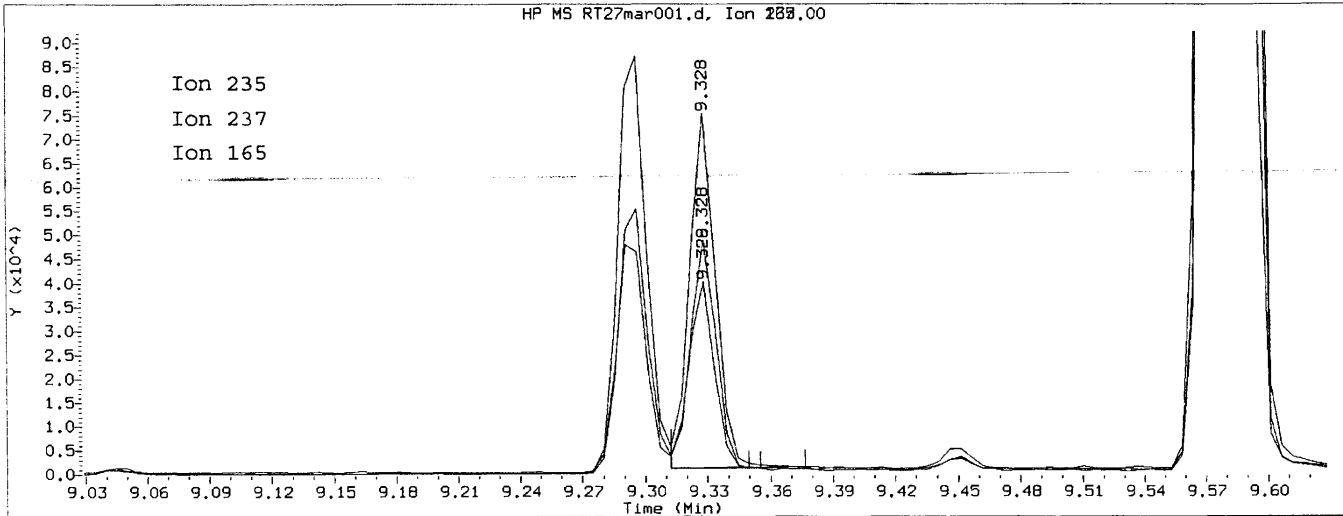
=====
 Exp. RT = 9.029
 Found RT = 8.965

Mass	Area	Ratio
246	15374	100.00
248	10139	65.95
176	6269	40.78

Return to Contents

Report Generated Time Mon Mar 27 10:55:20 2017

Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 3
 Acq on : 27-MAR-2017 10:33 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d/resolut.m
 Last Update : 27-MAR-2017 10:53



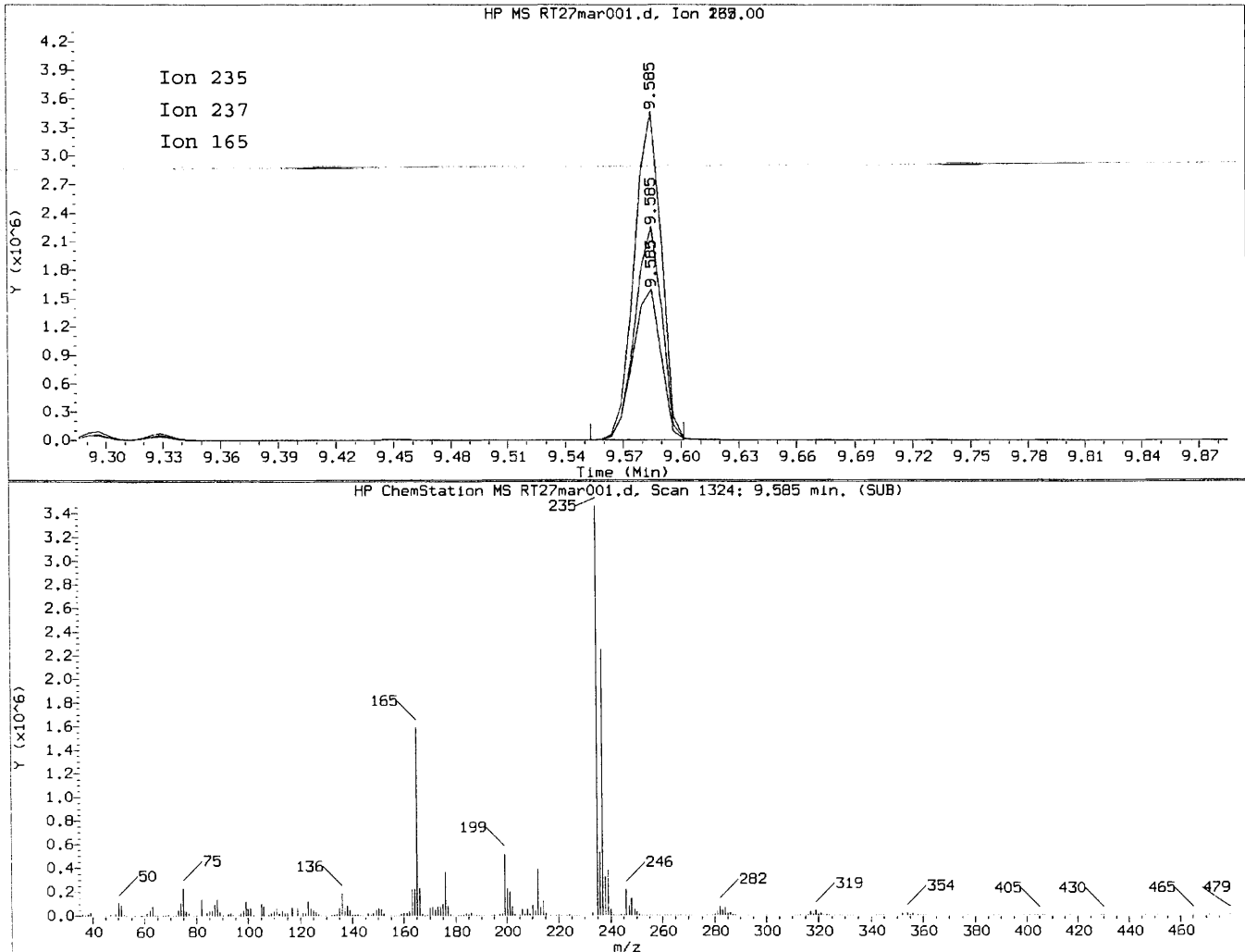
4,4'-DDD

=====
 Exp. RT = 9.380
 Found RT = 9.328

Mass	Area	Ratio
235	64320	100.00
237	40933	63.64
165	33791	52.54

Report Generated Time Mon Mar 27 10:55:20 2017

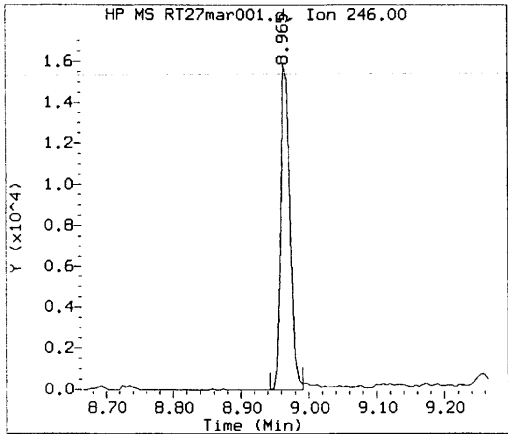
Data File : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d
 ALS Vial : 3
 Acq on : 27-MAR-2017 10:33 Operator : Tim Matthews
 Sample : TUNE S101716A DFTPP Inst : GCMS_EEE
 Misc : Multiplier : 1
 Integrator Type : HP RTE
 Method : /chem/SVOA/GCMS_EEE.i/170327.b/27mar001.d/resolut.m
 Last Update : 27-MAR-2017 10:53



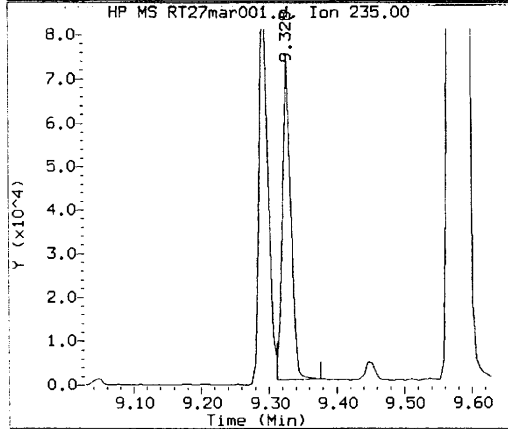
4,4'-DDT

=====
 Exp. RT = 9.756
 Found RT = 9.585

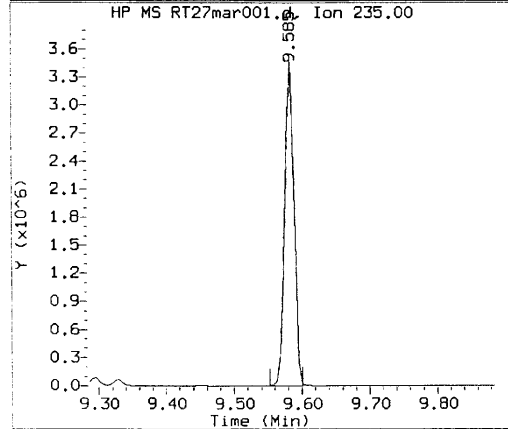
Mass	Area	Ratio
235	3360569	100.00
237	2171871	64.63
165	1603931	47.73



Compound: 4,4'-DDE
 Quant Mass: 246
 RT: 8.965
 Area: 15374



Compound: 4,4'-DDD
 Quant Mass: 235
 RT: 9.328
 Area: 64320



Compound: 4,4'-DDT
 Quant Mass: 235
 RT: 9.585
 Area: 3360569

DDT DEGRADATION BREAKDOWN ANALYSIS SUMMARY

Compound	Response	%Breakdown	Max Allowed	Test
4,4-DDT	3360569			N/A
4,4-DDE	15374	0.46	20.0	PASS
4,4-DDD	64320	1.88	20.0	PASS
4,4-DDD + DDE	79694	2.3	20.0	PASS

 TUNE SAMPLE *** PASSED *** DDT BREAKDOWN TEST

EPA METHOD 8270C PAHSIM

Run Logs

Injection Log

Directory: W:\GCMS_EEE\GCMS_EEE_DATA\2017\170313

Page 395 of 398

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	13mar001.d	1.	TUNE S101716A DFTPP		13 Mar 2017 09:34
2	1	13mar002.d	1.	TUNE S101716A DFTPP		13 Mar 2017 10:04
3	1	13mar003.d	1.	TUNE S101716A DFTPP		13 Mar 2017 11:49
4	2	13mar004.d	1.	CCV S010317F 1PPM	170313A001	13 Mar 2017 12:10
5	3	13mar005.d	1.	BLANK	S101716B 10UL	13 Mar 2017 12:32
6	28	13mar006.d	1.	ICAL-1 S010317D 5PPM	170313I001	13 Mar 2017 12:52
7	29	13mar007.d	1.	ICAL-2 S010317E 2PPM		13 Mar 2017 13:13
8	30	13mar008.d	1.	ICAL-3 S010317F 1PPM		13 Mar 2017 13:33
9	31	13mar009.d	1.	ICAL-4 S010317G 0.5PPM		13 Mar 2017 13:53
10	32	13mar010.d	1.	ICAL-5 S010317H 0.1PPM		13 Mar 2017 14:13
11	33	13mar011.d	1.	ICV S010317I 1PPM		13 Mar 2017 14:34
12	1	13mar012.d	1.	TUNE S101716A DFTPP		13 Mar 2017 14:59
13	2	13mar013.d	1.	CCV S010317F 1PPM	170313A001	13 Mar 2017 15:19
14	24	13mar014.d	1.	17-03-0323-3 10X	S101716B 10UL	13 Mar 2017 15:40
15		13mar015.d	1.	No MS or GC data present		

*Leaking
change new
column
dev > 30%*



Injection Log

Directory: W:\GCMS_EEE\GCMS_EEE_DATA\2017\170327

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	27mar001.d	1.	TUNE S101716A DFTPP		27 Mar 2017 10:33
2	2	27mar002.d	1.	CCV S010317F 1PPM	170327A010	27 Mar 2017 10:55
3	3	27mar003.d	1.	MB 170324 L17	S032317A 10UL	27 Mar 2017 11:15
4	4	27mar004.d	1.	LCS 170324 L17	S032317A 10UL	27 Mar 2017 11:35
5	5	27mar005.d	1.	17-03-1753-6 MS	S032317A 10UL	27 Mar 2017 11:56
6	6	27mar006.d	1.	17-03-1753-6 MSD	S032317A 10UL	27 Mar 2017 12:16
7	7	27mar007.d	1.	17-03-1753-6	S032317A 10UL	27 Mar 2017 12:36
8	8	27mar008.d	1.	17-03-1556-25	S032317A 10UL	27 Mar 2017 12:56
9	9	27mar009.d	1.	17-03-1556-26	S032317A 10UL	27 Mar 2017 13:16
10	10	27mar010.d	1.	17-03-1556-27	S032317A 10UL	27 Mar 2017 13:37
11	11	27mar011.d	1.	17-03-1557-25	S032317A 10UL	27 Mar 2017 13:57
12	12	27mar012.d	1.	17-03-1557-26	S032317A 10UL	27 Mar 2017 14:17
13	13	27mar013.d	1.	17-03-1557-27	S032317A 10UL	27 Mar 2017 14:38
14	14	27mar014.d	1.	MB 170325 L01	S032317A 10UL	27 Mar 2017 14:58
15	15	27mar015.d	1.	LCS 170325 L01	S032317A 10UL	27 Mar 2017 15:18
16	16	27mar016.d	1.	LCS 170325 L01	S032317A 10UL	27 Mar 2017 15:38
17	17	27mar017.d	1.	17-03-1795-2 - 10x, 100x	S032317A 10UL	27 Mar 2017 15:58
18	18	27mar018.d	1.	17-03-1795-3	S032317A 10UL	27 Mar 2017 16:19
19	19	27mar019.d	1.	17-03-1795-4 } 10x, 100x	S032317A 10UL	27 Mar 2017 16:39
20	20	27mar020.d	1.	17-03-1795-5 } 10x, 100x	S032317A 10UL	27 Mar 2017 16:59
21	21	27mar021.d	1.	17-03-1795-6	S032317A 10UL	27 Mar 2017 17:19
22	22	27mar022.d	1.	17-03-1795-7 } 10x	S032317A 10UL	27 Mar 2017 17:40
23	23	27mar023.d	1.	17-03-1795-8 } 10x	S032317A 10UL	27 Mar 2017 18:00
24	24	27mar024.d	1.	17-03-1795-9	S032317A 10UL	27 Mar 2017 18:20
25	25	27mar025.d	1.	17-03-1795-11	S032317A 10UL	27 Mar 2017 18:41
26	26	27mar026.d	1.	17-03-1795-12	S032317A 10UL	27 Mar 2017 18:41
27	27	27mar027.d	1.	17-03-1795-13 - 10x	S032317A 10UL	27 Mar 2017 19:01
28	28	27mar028.d	1.	17-03-1795-14	S032317A 10UL	27 Mar 2017 19:21
29	29	27mar029.d	1.	17-03-1795-15	S032317A 10UL	27 Mar 2017 19:41
30	30	27mar030.d	1.	17-03-1795-16	S032317A 10UL	27 Mar 2017 20:01
31	31	27mar031.d	1.	17-03-1795-17 } 10x	S032317A 10UL	27 Mar 2017 20:21
32	32	27mar032.d	1.	17-03-1795-18 } 10x	S032317A 10UL	27 Mar 2017 20:42
33	33	27mar033.d	1.	MB 170327 L12	S032317A 10UL	27 Mar 2017 21:02
34	34	27mar034.d	1.	LCS 170327 L12	S032317A 10UL	27 Mar 2017 21:22
35	35	27mar035.d	1.	17-03-1930-1 MS	S032317A 10UL	27 Mar 2017 21:42
36	36	27mar036.d	1.	17-03-1930-1 MSD	S032317A 10UL	27 Mar 2017 22:02
37	37	27mar037.d	1.	17-03-1930-1 - over 12 hour	S032317A 10UL	27 Mar 2017 22:23
						27 Mar 2017 22:43

EPA METHOD 8270C PAHSIM

Sample Preparation Logs

Analysis Method (EPA Method): 608 8081 8082 8141 8310 TO-13 TO-4
 8270 (Soil Soil SIM SUPER PAH SIM PAH SIM Pest SIM PCB cong. SIM FL)

Extraction Method (EPA Method): 3510 3520 3540 3541 3545 3550 3580

Analyst ID#: Measuring Sample- 1107 Start Extraction- 2/7/17 Blow Down- 2/7/17 Clean Up-

Matrix: Soil Aqueous Oil Wipe Filter Tissue Air

Balance ID#: 70 Filter ID#: 807-17-17 ASE ID#: 9 Soxtherm ID#: Orbit Shaker ID#: Sonicator ID#:

Ext. Start Date/Time: 03/24/17 - 14:00/20:00 Ext. End Date/Time: 03/24/17 - 16:00/22:30

Sand or Wipe ID#: 807-19-19 Drying Agent: Na₂SO₄ Diatomaceous Earth
 Drying Agent(s) ID#: 807-44-18 / 807-22-03

Surrogate Std ID# & Volume Added (mL): 8020117c - 0.5 mL

Spike Std ID# & Volume Added (mL): 8021017A - 0.5 mL Spike Added to: LCS LCSD MS MSD

Extraction Solvent: MeCl₂ 1:1 Hexane-Acetone 1:1 MeCl₂-Acetone 9:1 Hexane-Diethyl-ether Acetonitrile

Extraction Solvent ID#: 807-55-01 / 807-44-08 Exchange Solvent (Hexane Acetonitrile) ID#:

Clean Up Start Date & Time: Clean Up End Date & Time:

Clean Up: 3620 Florisil 3630 SGC 3660 Sulfur 3665 Acid Other Cartridge ID#:

Clean Up Reagent ID#: Cartridge Conditioning Column Pre-Elution Reagent ID#:

MB/LCS/MS Batch #: 170324 L17	Sample W(g) / V (mL)		Clean Up Performed	Comments
	Initial	Final		
Cel ID#:				
MB	20.0	2	<input type="checkbox"/>	
LCS	20.1	2	<input type="checkbox"/>	
LCSD NA	✓	✓	<input type="checkbox"/>	
MS 17-03-1753-6A	20.0	2	<input type="checkbox"/>	
MSD 1 - 1753-6A	20.0	2	<input type="checkbox"/>	
17-03-1753-6A	20.1	2	<input type="checkbox"/>	
17-03-1556-25D	20.1	2	<input type="checkbox"/>	
↓ ↓ 26D	20.2	2	<input type="checkbox"/>	
↓ ↓ 27D	20.1	2	<input type="checkbox"/>	
17-03-1557-25D	20.0	2	<input type="checkbox"/>	
↓ ↓ 26D	20.2	2	<input type="checkbox"/>	
↓ ↓ 27D	20.2	2	<input type="checkbox"/>	
17-03-1558-25D	20.1	2	<input type="checkbox"/>	Extract & Hold
↓ ↓ 26D	20.0	2	<input type="checkbox"/>	↓
↓ ↓ 27D	20.0	2	<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	
			<input type="checkbox"/>	

Peer Reviewed by: 2/7

Peer Reviewed Date: 03/25/17

Revision Date: 10/20/16

Return to Contents