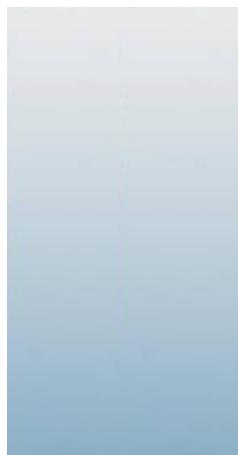




10/07/13



Technical Report for

XTO Energy

FRU 197-31A

1111-02A Cut 1 Contents

Accutest Job Number: D51123

Sampling Date: 09/30/13

Report to:

KRW Consulting, Inc.
8000 West 14th Avenue
Lakewood, CO 80214
dknudson@krwconsulting.com; jhess@krwconsulting.com;
crachak@krwconsulting.com; rrasic@krwconsulting.com;
ATTN: Dwayne Knudson

Total number of pages in report: 144



Test results contained within this data package meet the requirements
of the National Environmental Laboratory Accreditation Conference
and/or state specific certification programs as applicable.

A handwritten signature in black ink that appears to read "Scott Heideman".

Scott Heideman
Laboratory Director

Client Service contact: Renea Jackson 303-425-6021

Certifications: CO (CO00049), ID, NE (CO00049), ND (R-027), NJ (CO 0007), OK (D9942), UT (NELAP CO00049),
TX (T104704511)

This report shall not be reproduced, except in its entirety, without the written approval of Accutest Laboratories.
Test results relate only to samples analyzed.

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Sample Summary

XTO Energy

Job No: D51123

FRU 197-31A

Project No: 1111-02A Cut 1 Contents

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D51123-1	09/30/13	12:25 DS	10/01/13	SO	Soil	CUT 1 CONTENTS
D51123-1A	09/30/13	12:25 DS	10/01/13	SO	Soil	CUT 1 CONTENTS

Soil samples reported on a dry weight basis unless otherwise indicated on result page.



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: XTO Energy

Job No D51123

Site: FRU 197-31A

Report Date 10/7/2013 3:13:55 PM

On 10/01/2013, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 3.3 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D51123 was assigned to the project. The lab sample ID, client sample ID, and date of sample collection are detailed in the report's Results Summary.

Specified quality control criteria were achieved for this job except as noted below. For more information, please refer to the analytical results and QC summary pages.

Volatiles by GCMS By Method SW846 8260B

Matrix: SO

Batch ID: V5V1763

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51039-1MS, D51039-1MSD were used as the QC samples indicated.

Extractables by GCMS By Method SW846 8270C BY SIM

Matrix: SO

Batch ID: OP8670

- All samples were extracted within the recommended method holding time.
- All samples were analyzed and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.

Volatiles by GC By Method SW846 8015B

Matrix: SO

Batch ID: GGB1230

- All samples were analyzed within the recommended method holding time.
- Sample(s) D51008-1MS, D51008-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

Extractables by GC By Method SW846-8015B

Matrix: SO

Batch ID: OP8666

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D51127-1MS, D51127-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

Metals By Method SW846 6010C

Matrix: AQ

Batch ID: MP11305

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51224-6AMS, D51224-6AMSD, D51224-6ASDL were used as the QC samples for the metals analysis.

Matrix: SO

Batch ID: MP11267

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51122-1MS, D51122-1MSD, D51122-1SDL were used as the QC samples for the metals analysis.
- The matrix spike duplicate (MSD) recovery(s) of Zinc are outside control limits. Probable cause due to matrix interference.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The serial dilution RPD(s) for Cadmium, Selenium, Silver are outside control limits for sample MP11267-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP11267-SD1 for Nickel, Zinc: Serial dilution indicates possible matrix interference.

Metals By Method SW846 6020A

Matrix: SO

Batch ID: MP11268

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51122-1MS, D51122-1MSD, D51122-1SDL were used as the QC samples for the metals analysis.

Metals By Method SW846 7471B

Matrix: SO

Batch ID: MP11269

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51122-1MS, D51122-1MSD were used as the QC samples for the metals analysis.

Wet Chemistry By Method ASTM D1498-76M

Matrix: SO

Batch ID: GN22168

- Sample(s) D51122-1DUP were used as the QC samples for the Redox Potential Vs H₂ analysis.

Wet Chemistry By Method SM2540B-2011 M

Matrix: SO

Batch ID: GN22110

- The data for SM2540B-2011 M meets quality control requirements.

Wet Chemistry By Method SW846 3060A/7196A

Matrix: SO

Batch ID: GP11063

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51041-1MS, D51041-1MSD, D51041-1DUP were used as the QC samples for the Chromium, Hexavalent analysis.
- The duplicate RPD(s) for Chromium, Hexavalent are outside control limits for sample GP11063-D1. RPD acceptable due to low duplicate and sample concentrations.

Wet Chemistry By Method SW846 3060A/7196A M

Matrix: SO

Batch ID: R18903

- The data for SW846 3060A/7196A M meets quality control requirements.
- D51123-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Wet Chemistry By Method SW846 9045D

Matrix: SO

Batch ID: GN22154

- The following samples were run outside of holding time for method SW846 9045D: D51123-1

Wet Chemistry By Method USDA HANDBOOK 60

Matrix: SO

Batch ID: MP11305

- D51123-1A for Sodium Adsorption Ratio: Calculated as: $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})] / 2}$

AMS certifies that data reported for samples received, listed on the associated custody chain or analytical task order, were produced to specifications meeting AMS's Quality System precision, accuracy and completeness objectives except as noted.

Estimated non-standard method measurement uncertainty data is available on request, based on quality control bias and implicit for standard methods. Acceptable uncertainty requires tested parameter quality control data to meet method criteria.

AMS is not responsible for data quality assumptions if partial reports are used and recommends that this report be used in its entirety. This report is authorized by AMS indicated via signature on the report cover.

Summary of Hits

Page 1 of 1

Job Number: D51123
Account: XTO Energy
Project: FRU 197-31A
Collected: 09/30/13

3

Lab Sample ID	Client Sample ID	Result/ Analyte	Qual	RL	MDL	Units	Method
---------------	------------------	--------------------	------	----	-----	-------	--------

D51123-1 CUT 1 CONTENTS

Toluene	0.0993 J	0.16	0.078	mg/kg	SW846 8260B
Ethylbenzene	0.117 J	0.16	0.030	mg/kg	SW846 8260B
Chrysene	0.0365	0.011	0.0055	mg/kg	SW846 8270C BY SIM
Fluoranthene	0.0148	0.011	0.0055	mg/kg	SW846 8270C BY SIM
Naphthalene	0.266	0.015	0.013	mg/kg	SW846 8270C BY SIM
Pyrene	0.0238	0.011	0.0055	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	29.3	16	7.8	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	863	8.5	6.4	mg/kg	SW846-8015B
Arsenic	13.1	0.13		mg/kg	SW846 6020A
Barium	4070	6.4		mg/kg	SW846 6010C
Chromium	20.0	1.3		mg/kg	SW846 6010C
Copper	33.9	1.3		mg/kg	SW846 6010C
Lead	24.2	6.4		mg/kg	SW846 6010C
Nickel	16.6	3.9		mg/kg	SW846 6010C
Zinc	57.0	3.9		mg/kg	SW846 6010C
Specific Conductivity	17400	1.0		umhos/cm	SM 2510B-2011 MOD
Chromium, Trivalent ^a	19.9	2.3		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	233			mv	ASTM D1498-76M
pH	9.92			su	SW846 9045D

D51123-1A CUT 1 CONTENTS

Calcium	259	2.0	mg/l	SW846 6010C
Magnesium	10.4	1.0	mg/l	SW846 6010C
Sodium	3840	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	63.6		ratio	USDA HANDBOOK 60

(a) Calculated as: (Chromium) - (Chromium, Hexavalent)

(b) Calculated as: (Na meq/L) / sqrt [(Ca meq/L)+(Mg meq/L)/2]



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Sample Results

Report of Analysis

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 1 CONTENTS**Lab Sample ID:** D51123-1**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** FRU 197-31A**Date Sampled:** 09/30/13**Date Received:** 10/01/13**Percent Solids:** 77.9

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	5V29344.D	1	10/01/13	BD	n/a	n/a	V5V1763
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.04 g	5.0 ml	100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	0.078	0.039	mg/kg	
108-88-3	Toluene	0.0993	0.16	0.078	mg/kg	J
100-41-4	Ethylbenzene	0.117	0.16	0.030	mg/kg	
1330-20-7	Xylene (total)	ND	0.31	0.16	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	99%		64-130%
460-00-4	4-Bromofluorobenzene	98%		62-131%
17060-07-0	1,2-Dichloroethane-D4	100%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 1 CONTENTS	Date Sampled:	09/30/13
Lab Sample ID:	D51123-1	Date Received:	10/01/13
Matrix:	SO - Soil	Percent Solids:	77.9
Method:	SW846 8270C BY SIM	SW846 3546	
Project:	FRU 197-31A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G16541.D	1	10/03/13	DC	10/03/13	OP8670	E3G817
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	1.0 ml
Run #2		

COGCC Table 910-1 PAH List

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	0.011	0.0055	mg/kg	
120-12-7	Anthracene	ND	0.011	0.0055	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.011	0.0055	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.011	0.0055	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.011	0.0055	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.011	0.0055	mg/kg	
218-01-9	Chrysene	0.0365	0.011	0.0055	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.011	0.0055	mg/kg	
206-44-0	Fluoranthene	0.0148	0.011	0.0055	mg/kg	
86-73-7	Fluorene	ND	0.011	0.0064	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.011	0.0055	mg/kg	
91-20-3	Naphthalene	0.266	0.015	0.013	mg/kg	
129-00-0	Pyrene	0.0238	0.011	0.0055	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	69%		10-175%
321-60-8	2-Fluorobiphenyl	75%		25-130%
1718-51-0	Terphenyl-d14	113%		41-133%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

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Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 1 CONTENTS**Lab Sample ID:** D51123-1**Matrix:** SO - Soil**Method:** SW846 8015B**Project:** FRU 197-31A**Date Sampled:** 09/30/13**Date Received:** 10/01/13**Percent Solids:** 77.9

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB22383.D	1	10/02/13	EV	n/a	n/a	GGB1230
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.0 g	5.0 ml	100 ul
Run #2			

CAS No.	Compound	Result	RL	MDL	Units	Q
----------------	-----------------	---------------	-----------	------------	--------------	----------

TPH-GRO (C6-C10)	29.3	16	7.8	mg/kg
------------------	------	----	-----	-------

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
----------------	-----------------------------	---------------	---------------	---------------

120-82-1	1,2,4-Trichlorobenzene	76%		60-140%
----------	------------------------	-----	--	---------

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

4

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 1 CONTENTS
Lab Sample ID: D51123-1
Matrix: SO - Soil
Method: SW846-8015B SW846 3546
Project: FRU 197-31A

Date Sampled: 09/30/13
Date Received: 10/01/13
Percent Solids: 77.9

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FI09508.D	1	10/03/13	TU	10/02/13	OP8666	GFI637
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	863	8.5	6.4	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	83%		20-130%		

ND = Not detected MDL - Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 1 CONTENTS	Date Sampled:	09/30/13
Lab Sample ID:	D51123-1	Date Received:	10/01/13
Matrix:	SO - Soil	Percent Solids:	77.9
Project:	FRU 197-31A		

Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	13.1	0.13	mg/kg	5	10/02/13	10/04/13	JB	SW846 6020A ³
Barium	4070	6.4	mg/kg	5	10/02/13	10/03/13	JM	SW846 6010C ¹
Cadmium	< 1.3	1.3	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Chromium	20.0	1.3	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Copper	33.9	1.3	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Lead	24.2	6.4	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Mercury	< 0.11	0.11	mg/kg	1	10/04/13	10/04/13	JB	SW846 7471B ²
Nickel	16.6	3.9	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Selenium	< 6.4	6.4	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Silver	< 3.9	3.9	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹
Zinc	57.0	3.9	mg/kg	1	10/02/13	10/02/13	JM	SW846 6010C ¹

- (1) Instrument QC Batch: MA4027
- (2) Instrument QC Batch: MA4035
- (3) Instrument QC Batch: MA4036
- (4) Prep QC Batch: MP11267
- (5) Prep QC Batch: MP11268
- (6) Prep QC Batch: MP11269

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 1 CONTENTS**Lab Sample ID:** D51123-1**Matrix:** SO - Soil**Date Sampled:** 09/30/13**Date Received:** 10/01/13**Percent Solids:** 77.9**Project:** FRU 197-31A**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	17400	1.0	umhos/cm	1	10/03/13	JD	SM 2510B-2011 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	10/02/13	JD	SW846 3060A/7196A
Chromium, Trivalent ^a	19.9	2.3	mg/kg	1	10/02/13 16:28	JM	SW846 3060A/7196A M
Redox Potential Vs H2	233		mv	1	10/04/13	AK	ASTM D1498-76M
Solids, Percent	77.9		%	1	10/01/13	SWT	SM2540B-2011 M
pH	9.92		su	1	10/03/13 12:30	AK	SW846 9045D

(a) Calculated as: (Chromium) - (Chromium, Hexavalent)

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 1 CONTENTS	Date Sampled:	09/30/13
Lab Sample ID:	D51123-1A	Date Received:	10/01/13
Matrix:	SO - Soil	Percent Solids:	77.9
Project:	FRU 197-31A		

SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	259	2.0	mg/l	1	10/04/13	10/04/13 JM	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	10.4	1.0	mg/l	1	10/04/13	10/04/13 JM	SW846 6010C ¹	SW846 3010A/M ²
Sodium	3840	2.0	mg/l	1	10/04/13	10/04/13 JM	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA4038

(2) Prep QC Batch: MP11305

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 1 CONTENTS**Lab Sample ID:** D51123-1A**Matrix:** SO - Soil**Project:** FRU 197-31A**Date Sampled:** 09/30/13**Date Received:** 10/01/13**Percent Solids:** 77.9**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	63.6		ratio	1	10/04/13 20:38	JM	USDA HANDBOOK 60

(a) Calculated as: (Na meq/L) / sqrt [(Ca meq/L)+ (Mg meq/L)/2]

RL = Reporting Limit



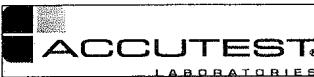
Misc. Forms

5

Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody



CHAIN OF CUSTODY

PAGE 1 OF 1

4036 Youngfield Street, Wheat Ridge, CO 80033
TEL. 303-425-6021 FAX: 303-425-6854
www.accutest.com

FED-EX Tracking #	Bottle Order Control #
	D51123

Client / Reporting Information		Project Information						Requested Analysis (see TEST CODE sheet)						Matrix Codes				
Company Name KRW Consulting	Project Name: XTO FRV 197-31A	Street Address 8000 West 14th Street; Suite 200	Street	Billing Information (If different from Report to)						Company Name XTO Energy	City Rifle, CO 81650	State CO	City Rifle, CO 81650	State CO	Attention: Jessica Dooling	Number of preserved bottles	DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank	
Project Contact: Dwayne Knudson	Project # 111-02A	Phone # 970-488-1088	Client Purchase Order #	Project Manager Joe Hess	Address 21469 CR 5	City Rifle, CO 81650	State CO	Attention: Jessica Dooling	Number of preserved bottles	HCl NaOH HN03 HN04 NONE DI Water MEOH ENCORE Bottles								
Acutest Sample #	Field ID / Point of Collection Cut 1 Contents	MEOH/DI Vial #	Date 9/30/13	Time 1225	Sampled by	# of bottles 5									LAB USE ONLY			
<i>Table 910</i>														01				
9/30/13																		
Turnaround Time (Business days)														Data Deliverable Information	Comments / Special Instructions			
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 5 Business Days (By contract only) <input type="checkbox"/> 3 Day Emergency <input type="checkbox"/> 2 Day Emergency <input type="checkbox"/> 1 Day Emergency														<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> COMM BN <input type="checkbox"/> COMM BN+	<input type="checkbox"/> State Forms Required <input type="checkbox"/> Send Forms to State <input type="checkbox"/> Report by Fax <input checked="" type="checkbox"/> Report by PDF ONLY <input type="checkbox"/> EDD Format	Please email to: KRW Piceance Team		
Emergency & Rush T/A data available VIA Lablink														Commercial "A" = Results Only Commercial "B" = Results + QC Summary Commercial BN = Results/QC narrative (+ = chromatograms)				
Sample Custody must be documented below each time samples change possession, including courier delivery.																		
Relinquished by Sampler: 1	Date Time: 9/30/13 1700	Received By: 1	Relinquished By: 2	Date Time:	Received By:													
Relinquished by Sampler: 3	Date Time:	Received By: 3	Relinquished By: 4	Date Time:	Received By: 4													
Relinquished by: 5	Date Time:	Received By: 5	Custody Seal # CO	<input type="checkbox"/> Intact <input type="checkbox"/> Not intact	Preserved where applicable <input type="checkbox"/>	On Ice <input type="checkbox"/>	Cooler Temp. 343											

D51123: Chain of Custody
Page 1 of 2



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D51123

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 10/1/2013 11:50:00 AM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO

Airbill #'s: CO

Cooler Security Y or N

- | | | | | | |
|---------------------------|-------------------------------------|--------------------------|-----------------------|-------------------------------------|--------------------------|
| 1. Custody Seals Present: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 3. COC Present: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Custody Seals Intact: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 4. Smpl Dates/Time OK | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

Cooler Temperature Y or N

- | | | |
|------------------------------|-------------------------------------|--------------------------|
| 1. Temp criteria achieved: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Cooler temp verification: | Infrared gun | |
| 3. Cooler media: | Ice (bag) | |

Quality Control Preservation Y or N N/A

- | | | |
|---------------------------------|-------------------------------------|--------------------------|
| 1. Trip Blank present / cooler: | <input type="checkbox"/> | <input type="checkbox"/> |
| 2. Trip Blank listed on COC: | <input type="checkbox"/> | <input type="checkbox"/> |
| 3. Samples preserved properly: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 4. VOCs headspace free: | <input type="checkbox"/> | <input type="checkbox"/> |

Sample Integrity - Documentation

- | | | |
|--|-------------------------------------|--------------------------|
| 1. Sample labels present on bottles: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Container labeling complete: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Sample container label / COC agree: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

Sample Integrity - Condition

- | | | |
|----------------------------------|-------------------------------------|--------------------------|
| 1. Sample rcvd within HT: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. All containers accounted for: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Condition of sample: | Intact | |

Sample Integrity - Instructions

- | | | |
|---|-------------------------------------|-------------------------------------|
| 1. Analysis requested is clear: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Bottles received for unspecified tests | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| 3. Sufficient volume rec'd for analysis: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 4. Compositing instructions clear: | <input type="checkbox"/> | <input type="checkbox"/> |
| 5. Filtering instructions clear: | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

Comments

Accutest Laboratories
V:(303) 425-6021

4036 Youngfield Street
F: (303) 425-6854

Wheat Ridge, CO
www.accutest.com

5.1

5

D51123: Chain of Custody

Page 2 of 2



GC/MS Volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries



Method Blank Summary

Page 1 of 1

Job Number: D51123
Account: XTOKRWR XTO Energy
Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1763-MB	5V29325.D	1	10/01/13	BD	n/a	n/a	V5V1763

The QC reported here applies to the following samples:

Method: SW846 8260B

D51123-1

6.11

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	50	25	ug/kg	
100-41-4	Ethylbenzene	ND	100	19	ug/kg	
108-88-3	Toluene	ND	100	50	ug/kg	
1330-20-7	Xylene (total)	ND	200	100	ug/kg	

CAS No.	Surrogate Recoveries	Limits
2037-26-5	Toluene-D8	98%
460-00-4	4-Bromofluorobenzene	87%
17060-07-0	1,2-Dichloroethane-D4	105%

Blank Spike Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1763-BS	5V29326.D	1	10/01/13	BD	n/a	n/a	V5V1763

The QC reported here applies to the following samples:

Method: SW846 8260B

D51123-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	2500	2650	106	70-130
100-41-4	Ethylbenzene	2500	2840	114	70-130
108-88-3	Toluene	2500	2740	110	70-130
1330-20-7	Xylene (total)	7500	8900	119	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	103%	64-130%
460-00-4	4-Bromofluorobenzene	101%	62-131%
17060-07-0	1,2-Dichloroethane-D4	93%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D51039-1MS	5V29328.D	1	10/01/13	BD	n/a	n/a	V5V1763
D51039-1MSD	5V29329.D	1	10/01/13	BD	n/a	n/a	V5V1763
D51039-1	5V29327.D	1	10/01/13	BD	n/a	n/a	V5V1763

The QC reported here applies to the following samples:

Method: SW846 8260B

D51123-1

CAS No.	Compound	D51039-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	ND		3620	3780	104	3940	109	4	64-139/30
100-41-4	Ethylbenzene	ND		3620	3880	107	3890	107	0	68-136/30
108-88-3	Toluene	ND		3620	3640	101	3610	100	1	60-130/30
1330-20-7	Xylene (total)	ND		10900	12300	113	12400	114	1	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D51039-1	Limits
2037-26-5	Toluene-D8	95%	93%	97%	64-130%
460-00-4	4-Bromofluorobenzene	108%	107%	96%	62-131%
17060-07-0	1,2-Dichloroethane-D4	94%	94%	99%	70-130%

* = Outside of Control Limits.

6.3.1
6



GC/MS Volatiles

Raw Data

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5100113.S\
 Data File : 5V29344.D
 Acq On : 1 Oct 2013 10:23 pm
 Operator : BRETD
 Sample : D51123-1
 Misc : MS6474,V5V1763,5.035,,100,5,1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Oct 02 09:24:01 2013
 Quant Method : C:\msdchem\1\METHODS\V5AP1728TVH1728.M
 Quant Title : 8260
 QLast Update : Tue Aug 20 09:59:22 2013
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.613	168	159687	50.00	ug/l	0.00
37) 1,4-Difluorobenzene	12.412	114	216574	50.00	ug/l	0.00
56) Chlorobenzene-d5	15.061	117	215985	50.00	ug/l	0.00
77) 1,4-Dichlorobenzene-d4	17.036	152	163895	50.00	ug/l	0.00

System Monitoring Compounds						
35) 1,2-Dichloroethane-d4	12.013	102	16153	49.78	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	99.56%
64) Toluene-d8	13.805	98	242951	49.65	ug/l	-0.01
Spiked Amount	50.000	Range	70 - 130	Recovery	=	99.30%
72) 4-Bromofluorobenzene	16.009	95	111306	48.80	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	97.60%

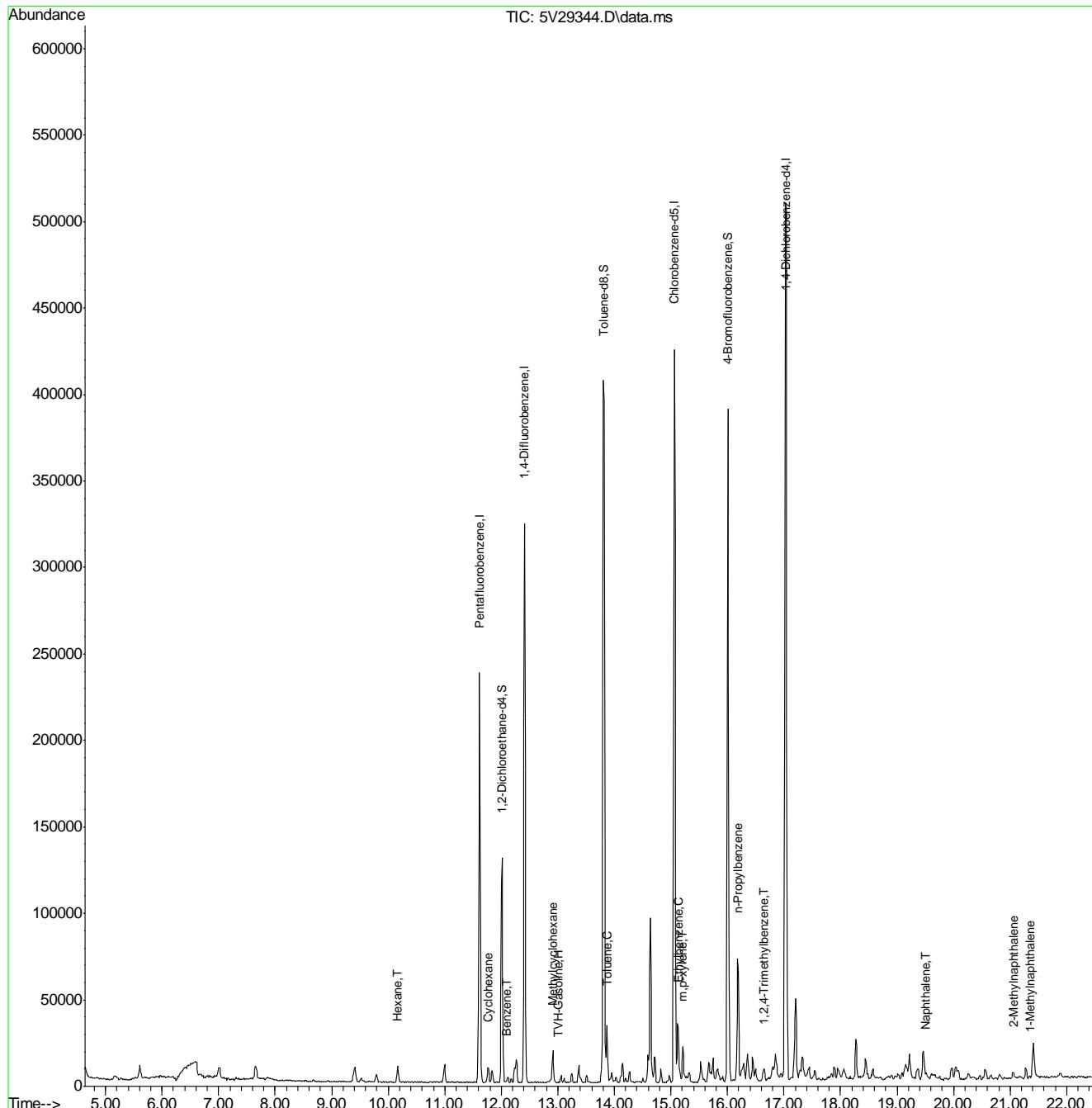
Target Compounds					Qvalue
1) TVH-Gasoline	13.006	TIC	856963m	120.14	ug/l
33) Cyclohexane	11.762	56	4727	15.81	ug/l 89
43) Hexane	10.163	57	3958	2.18	ug/l 100
47) Methylcyclohexane	12.915	83	4836	2.92	ug/l # 85
53) Benzene	12.093	78	763	0.16	ug/l 100
65) Toluene	13.874	92	5041	1.27	ug/l 98
69) Ethylbenzene	15.130	91	7995	1.50	ug/l 99
75) m,p-xylene	15.209	106	4953	1.97	ug/l 88
80) n-Propylbenzene	16.180	91	3408	0.47	ug/l 83
85) 1,2,4-Trimethylbenzene	16.648	105	1684	0.78	ug/l 91
94) Naphthalene	19.502	128	2884	1.20	ug/l 100
97) 2-Methylnaphthalene	21.055	142	1793	1.81	ug/l 89
98) 1-Methylnaphthalene	21.352	142	884	1.56	ug/l 95

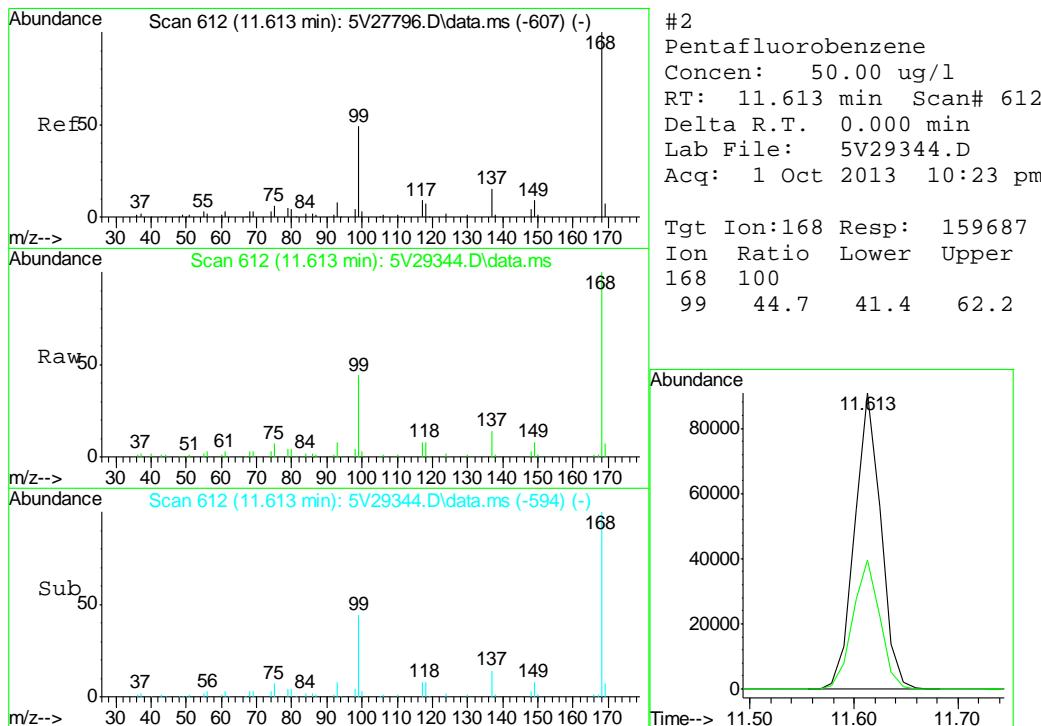
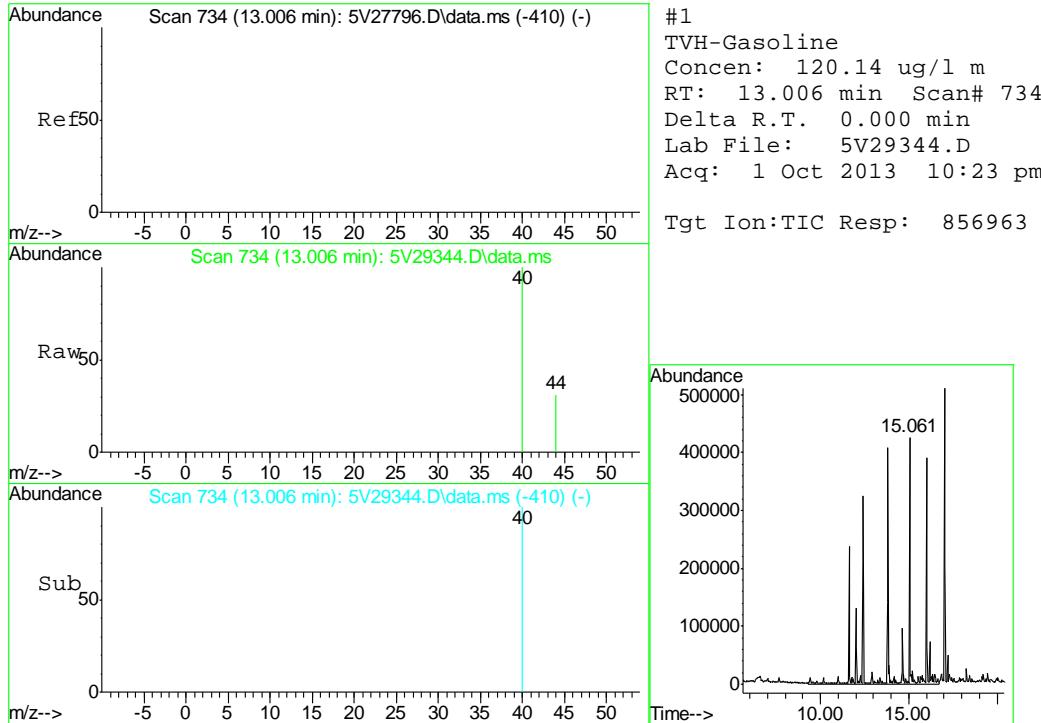
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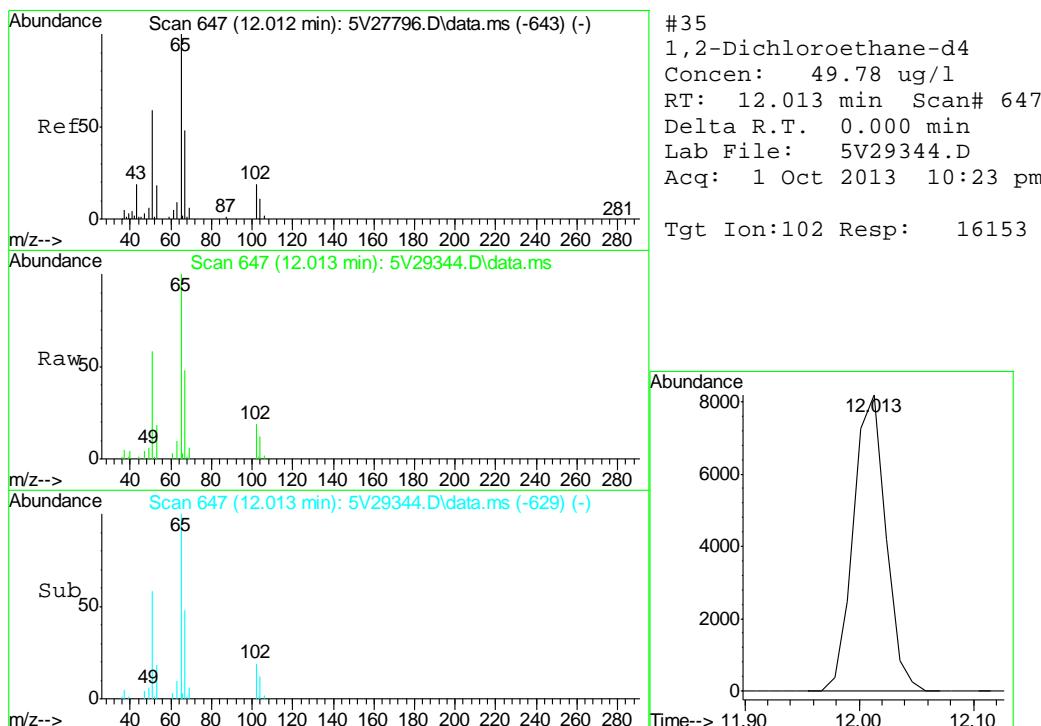
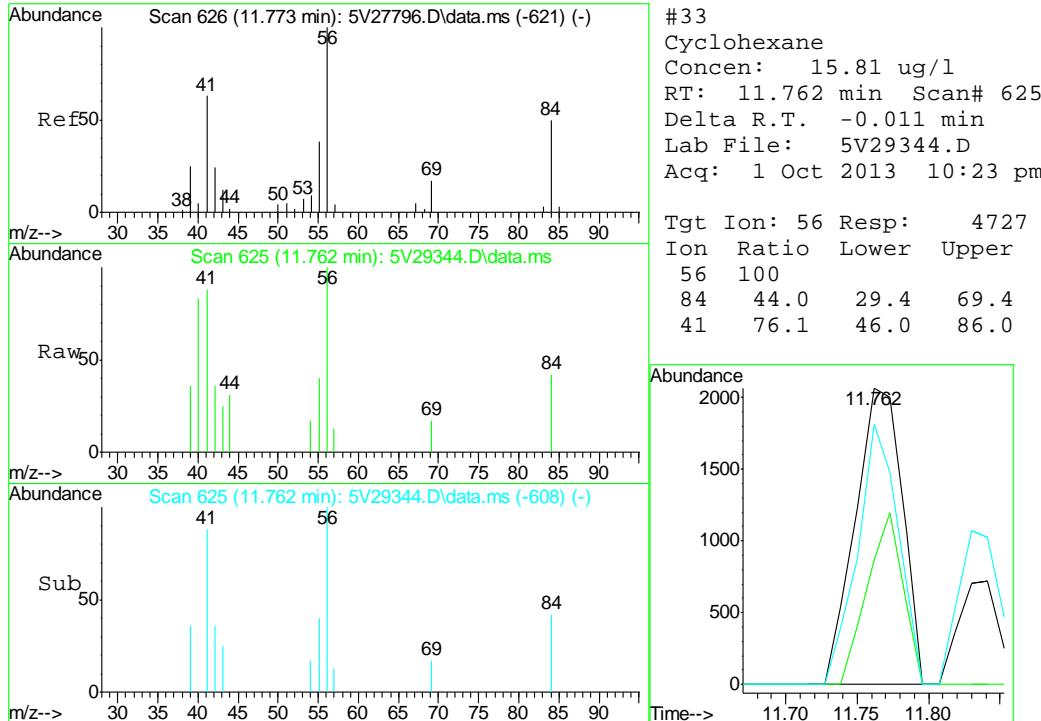
Quantitation Report (QT Reviewed)

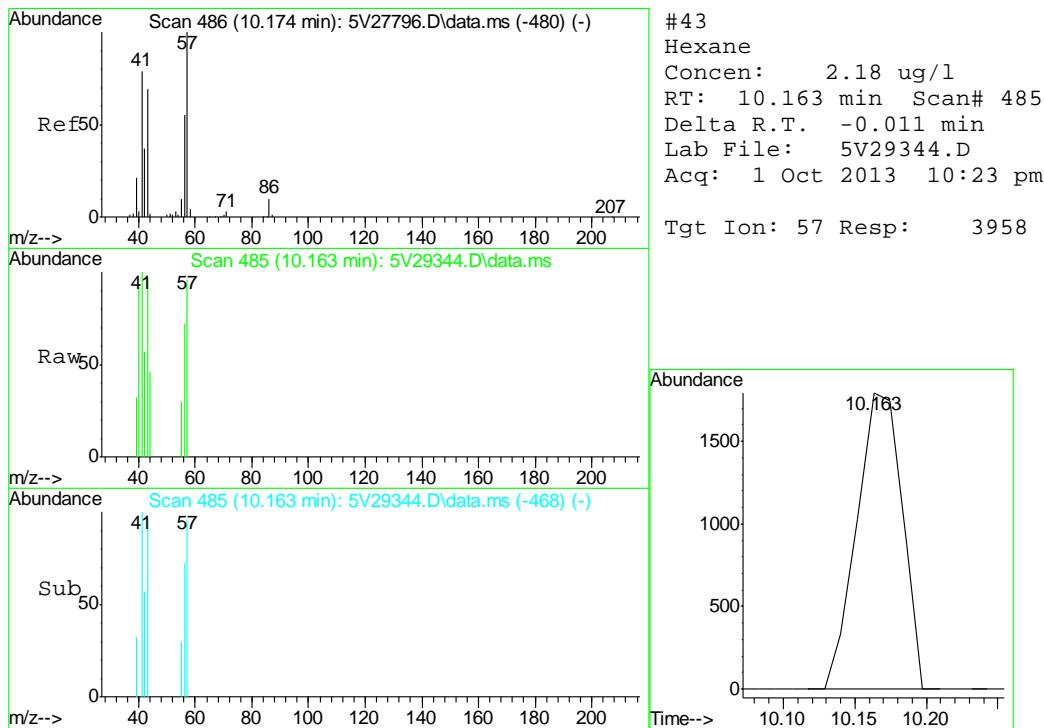
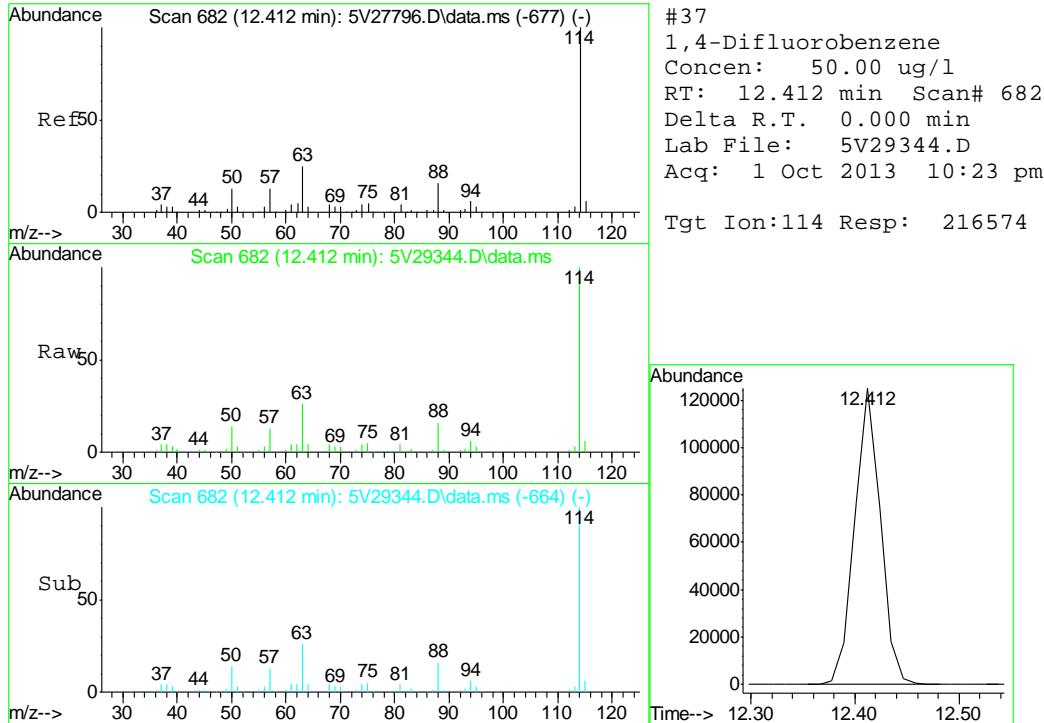
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 Operator : BRETD
 Sample : D51123-1
 Misc : MS6474,V5V1763,,5.035,,100,,5,1
 ALS Vial : 22 Sample Multiplier: 1

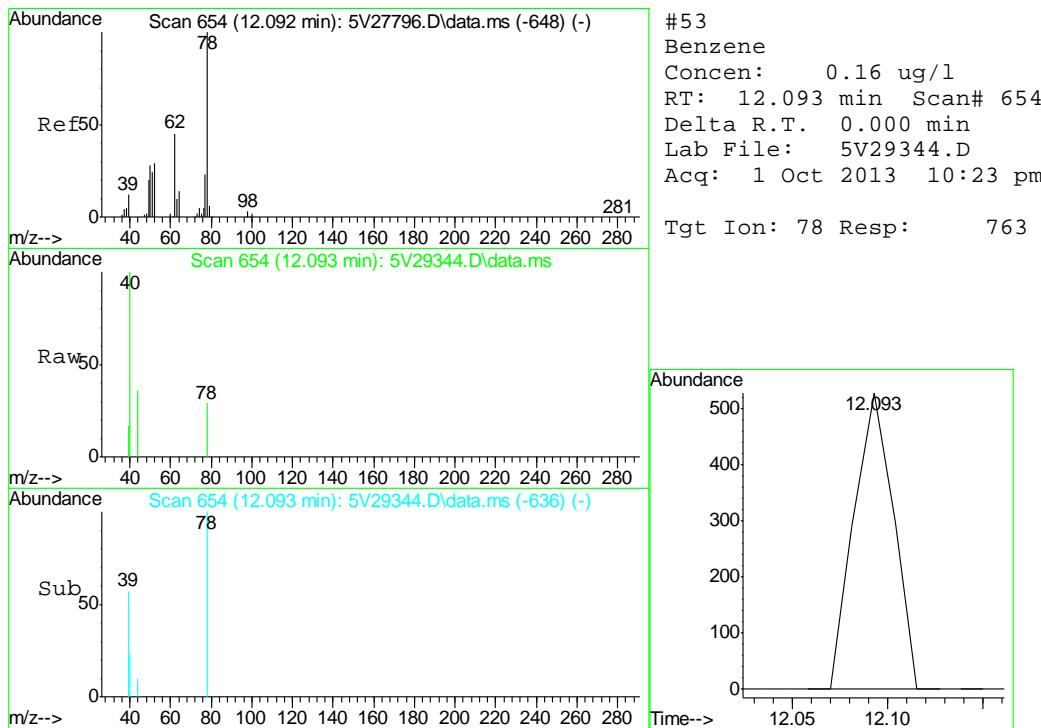
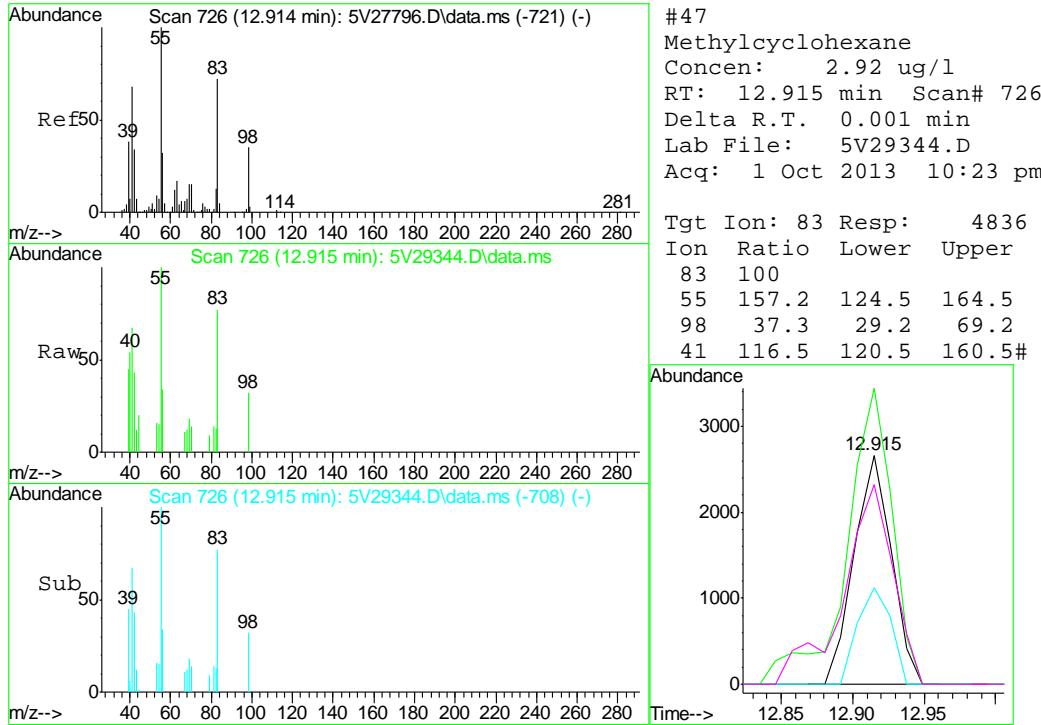
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 Response via : Initial Calibration

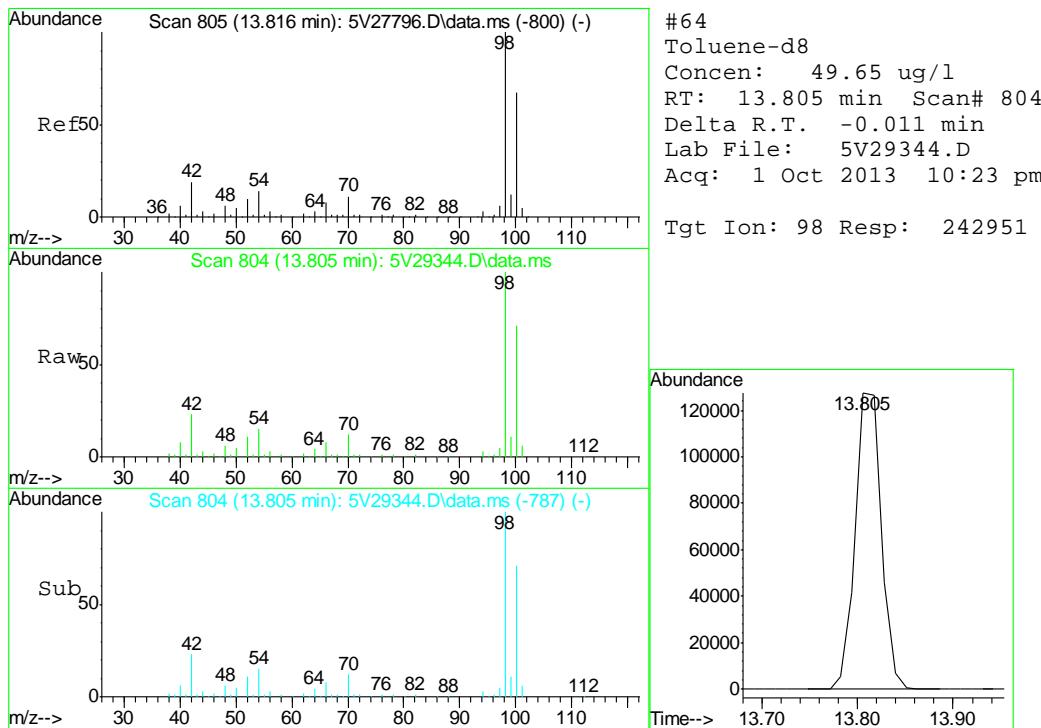
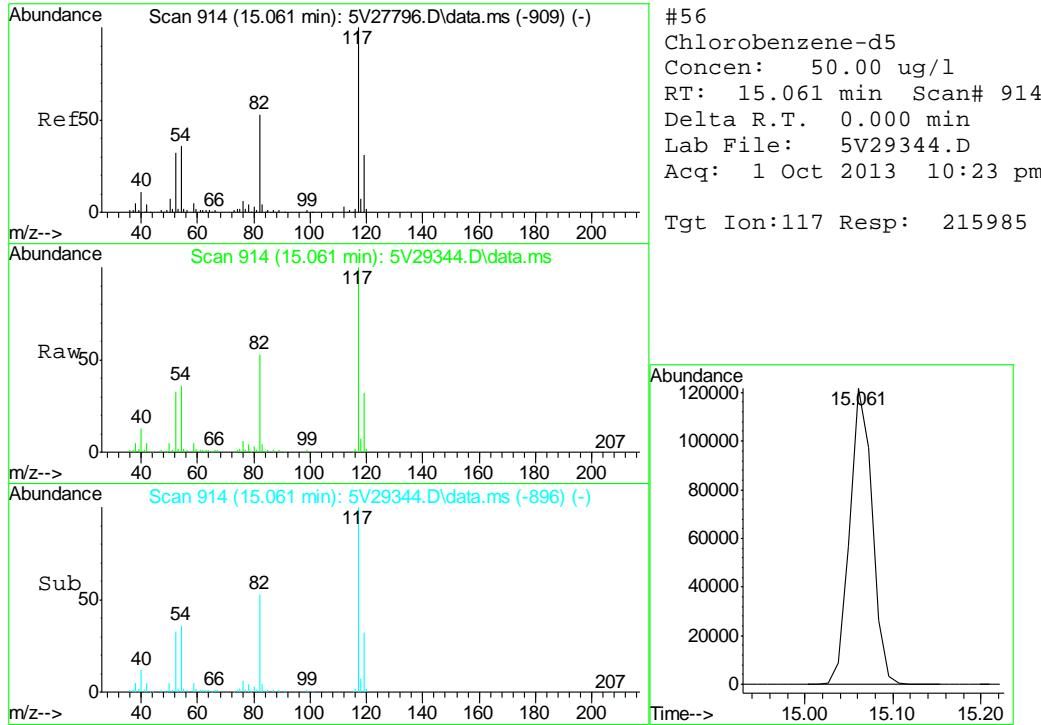


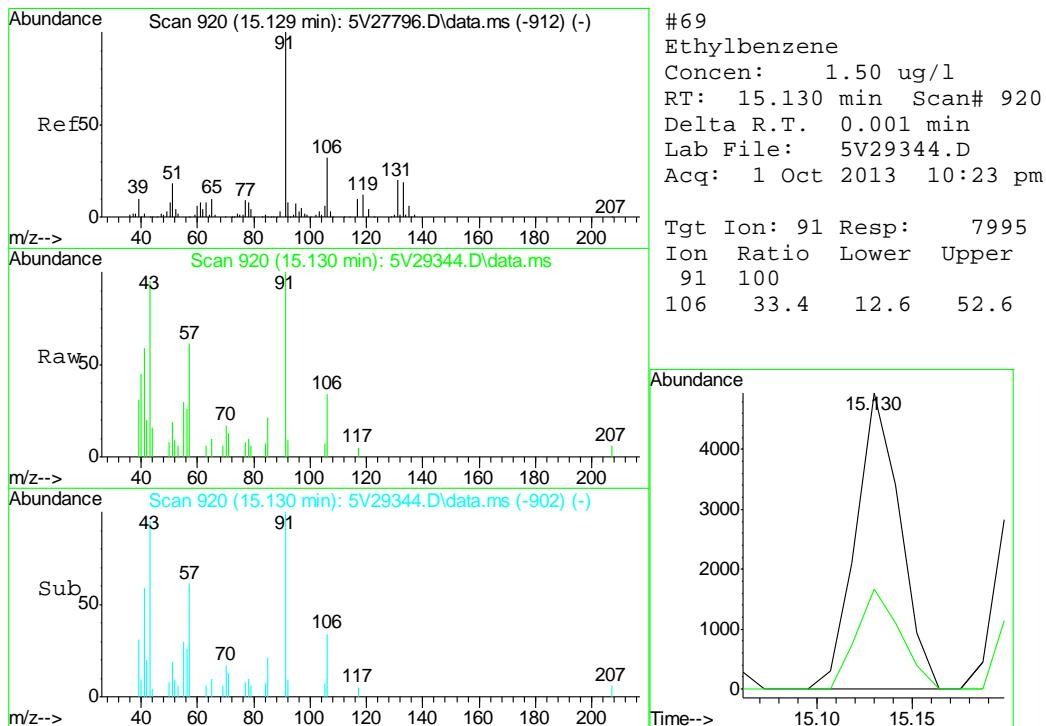
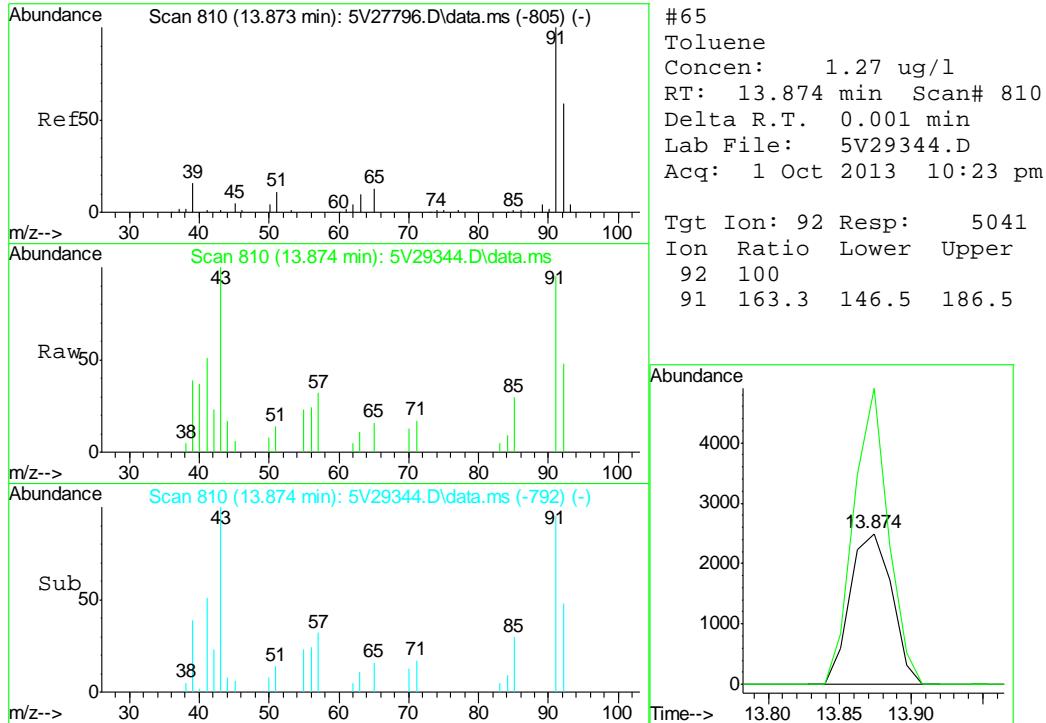


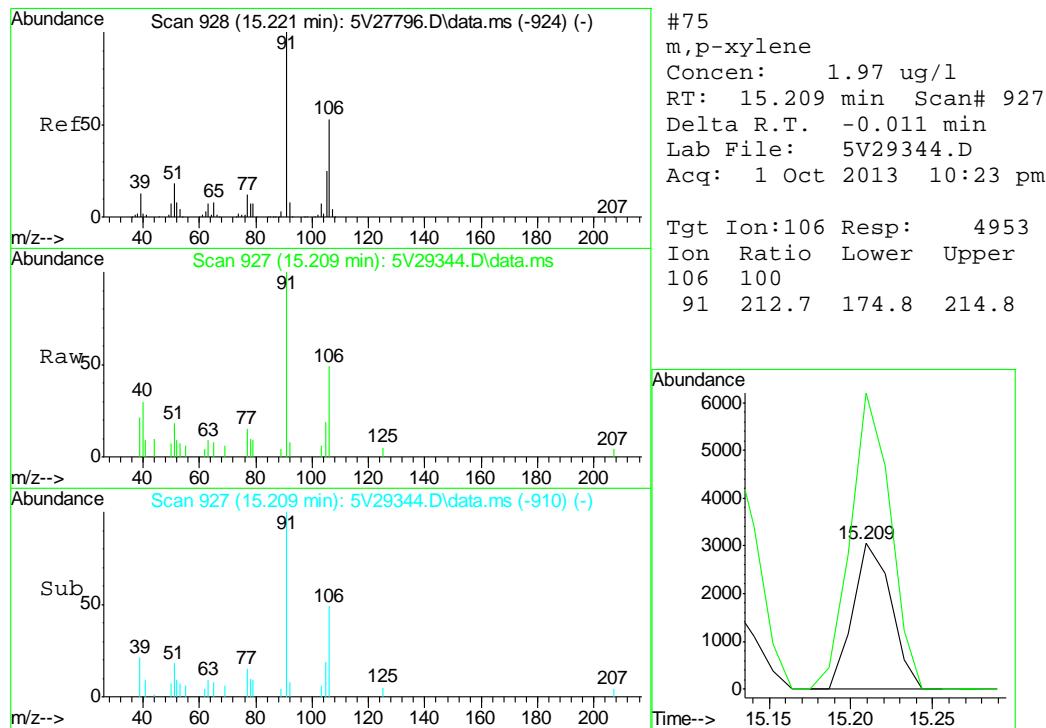
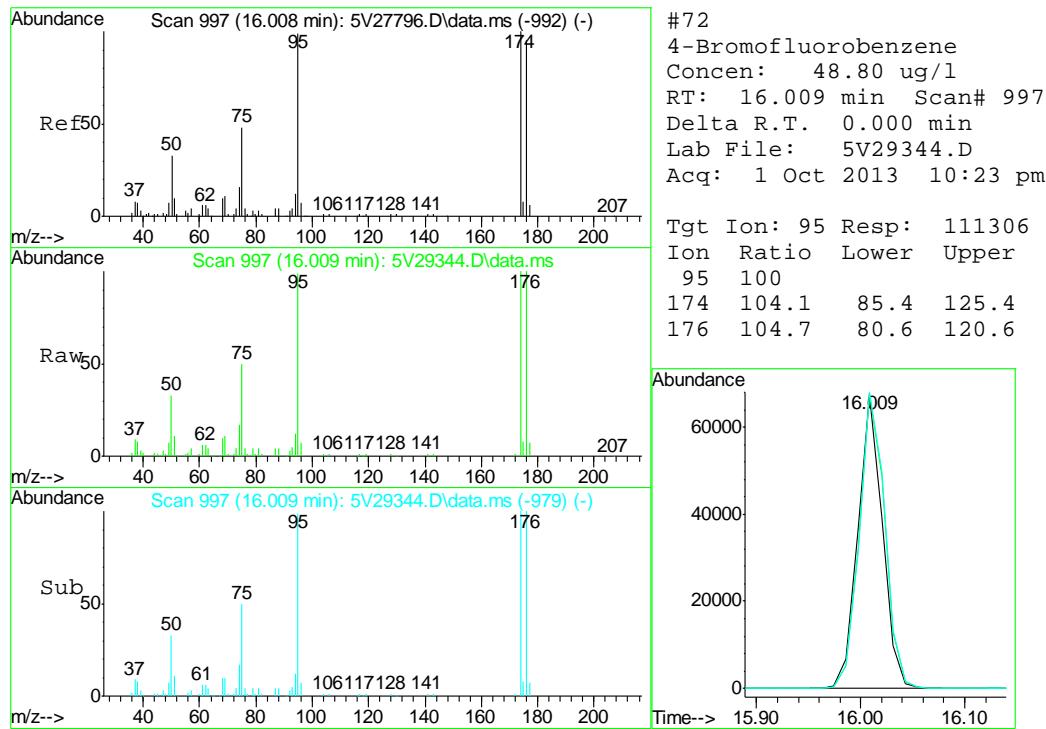


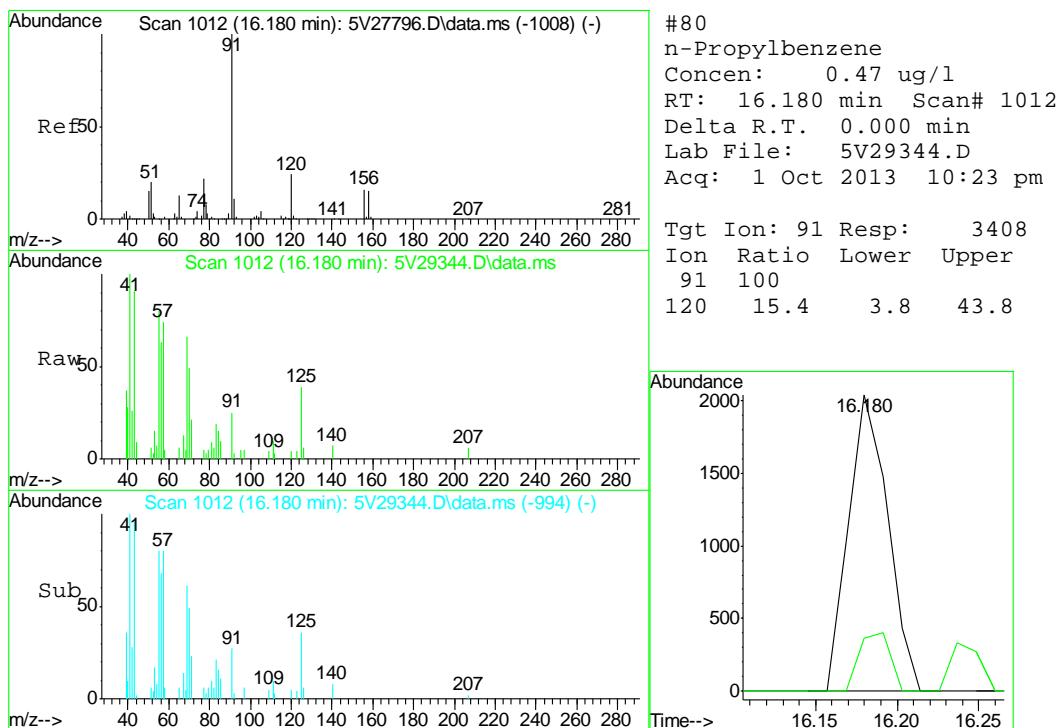
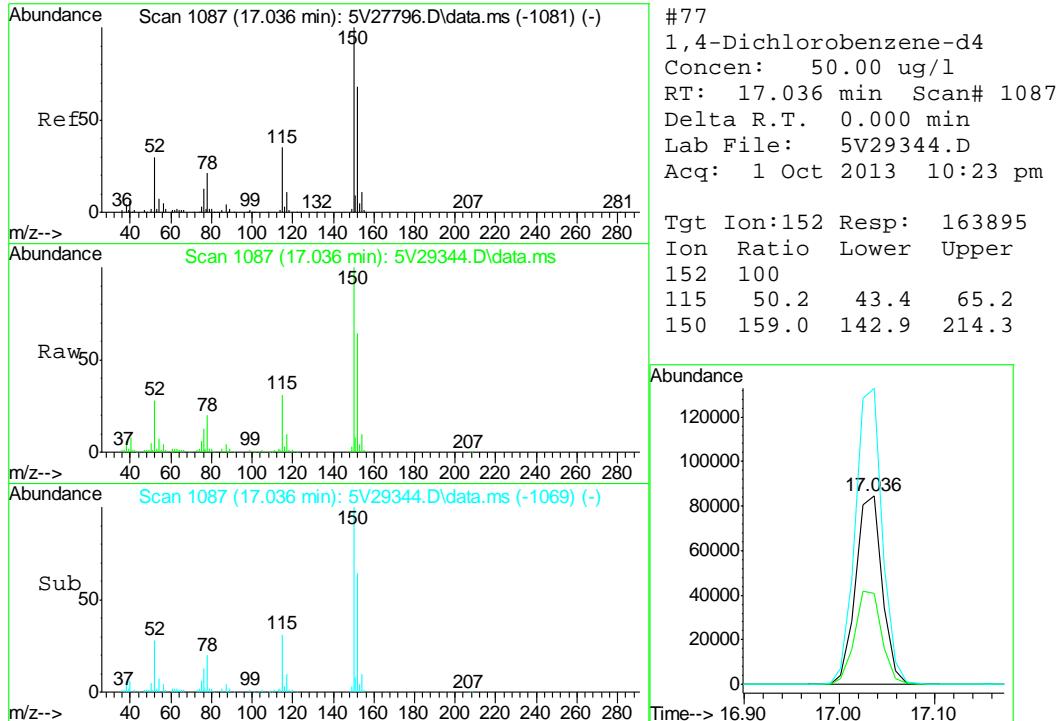


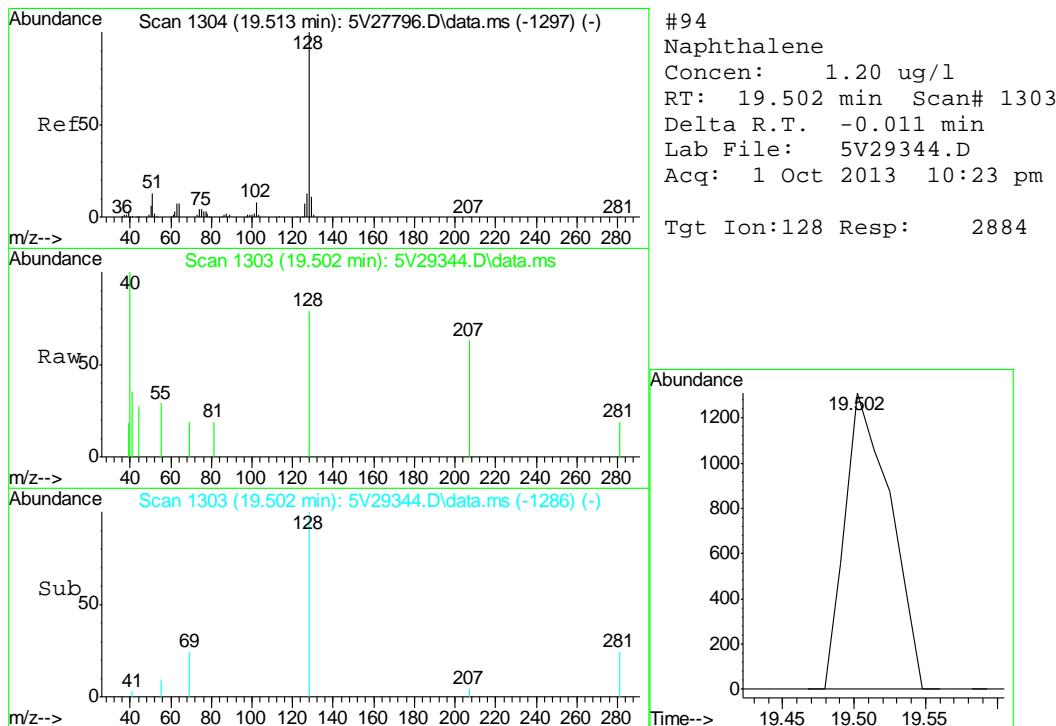
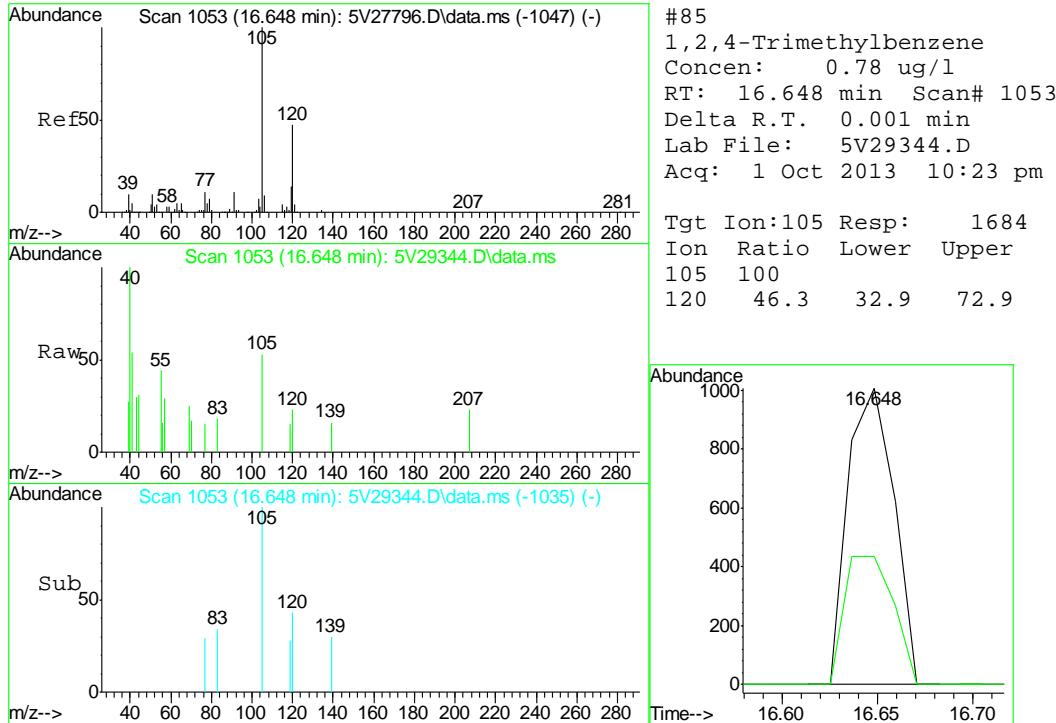


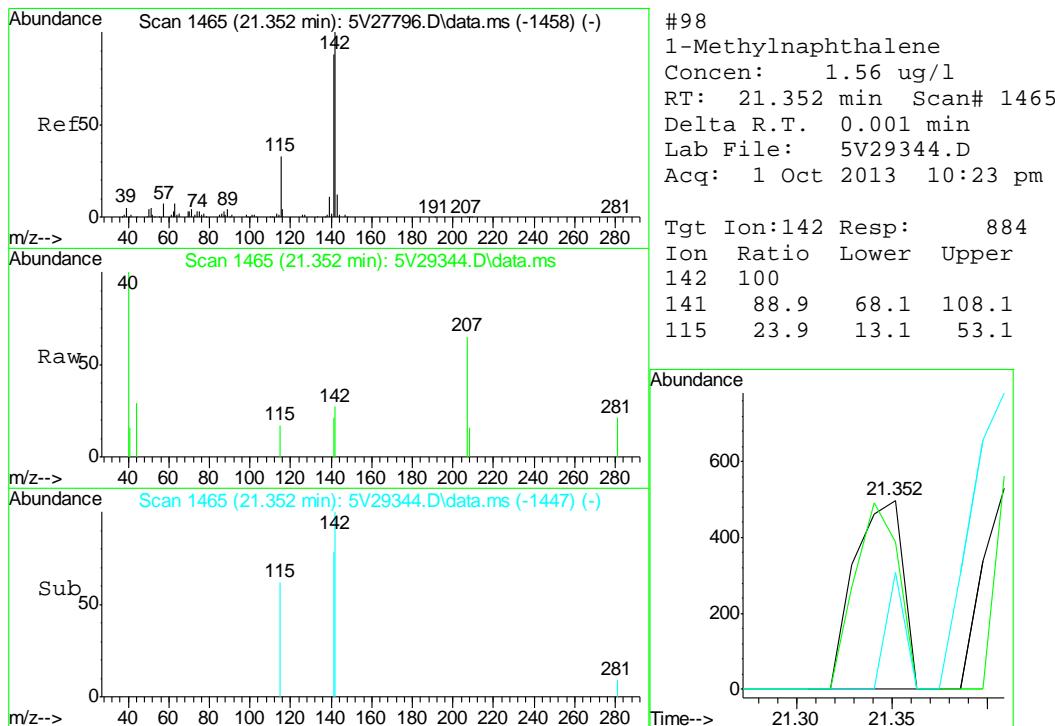
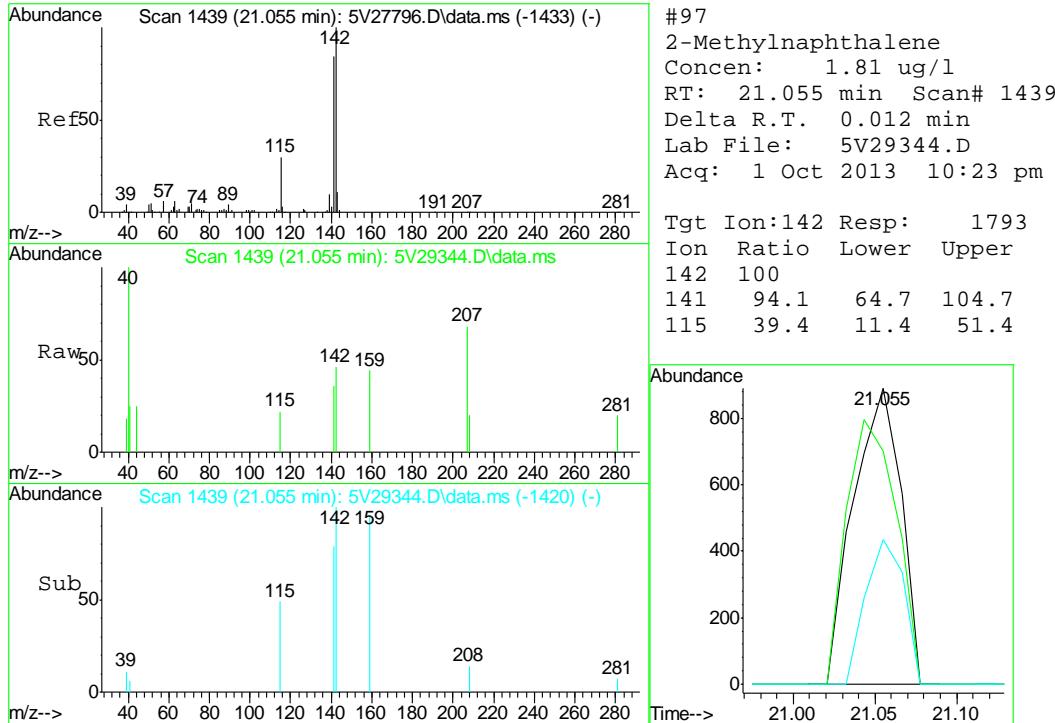












Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5100113.S\
 Data File : 5V29325.D
 Acq On : 1 Oct 2013 12:31 pm
 Operator : BRETD
 Sample : MB
 Misc : MS6474,V5V1763,5.000,,100,5,1
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Oct 02 08:53:24 2013
 Quant Method : C:\msdchem\1\METHODS\V5AP1728TVH1728.M
 Quant Title : 8260
 QLast Update : Tue Aug 20 09:59:22 2013
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.613	168	154496	50.00	ug/l	0.00
37) 1,4-Difluorobenzene	12.412	114	212347	50.00	ug/l	0.00
56) Chlorobenzene-d5	15.061	117	210906	50.00	ug/l	0.00
77) 1,4-Dichlorobenzene-d4	17.024	152	145782	50.00	ug/l	-0.01

System Monitoring Compounds						
35) 1,2-Dichloroethane-d4	12.012	102	16515	52.60	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	105.20%
64) Toluene-d8	13.816	98	233743	48.92	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	97.84%
72) 4-Bromofluorobenzene	16.008	95	96915	43.52	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	87.04%

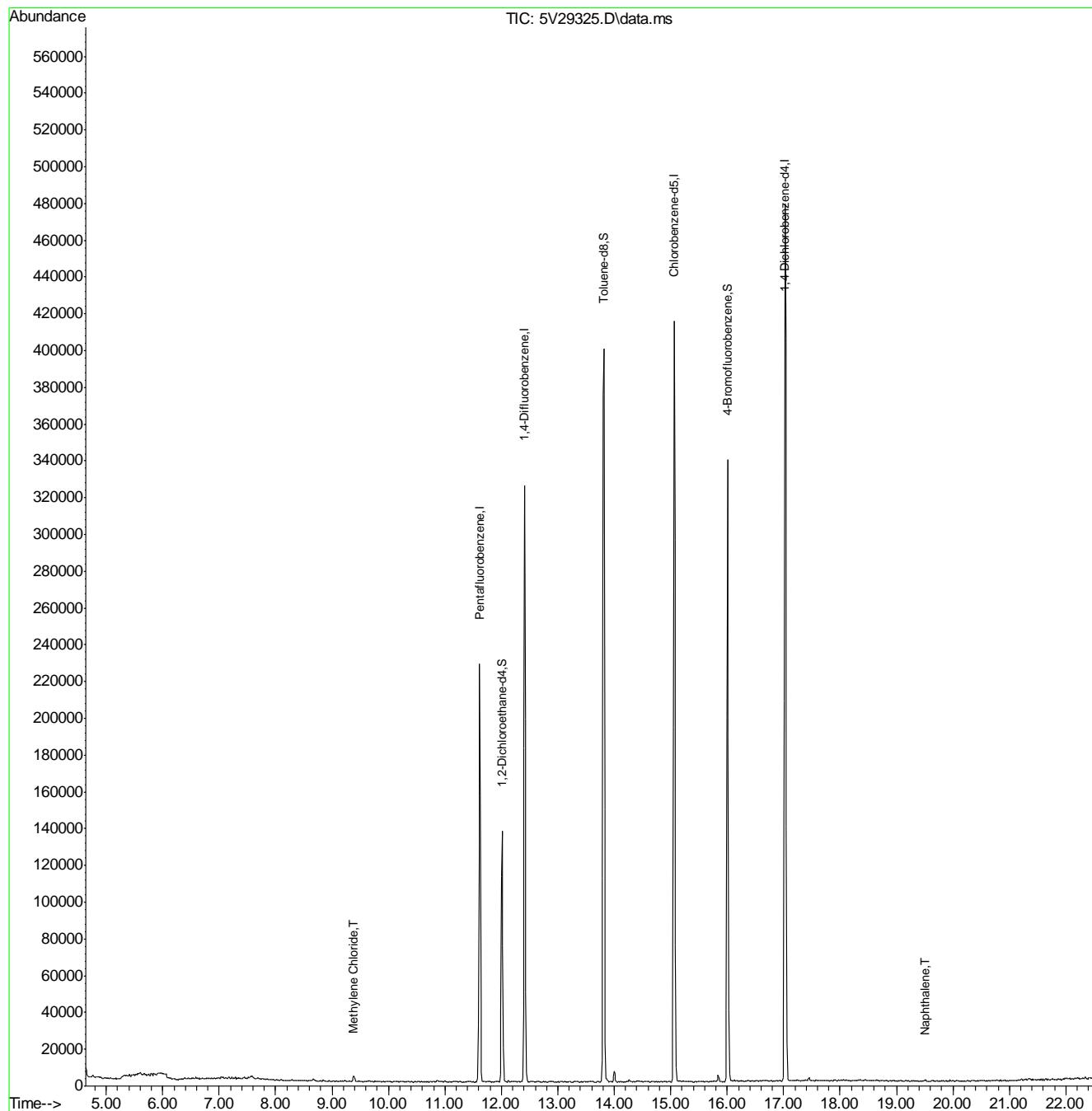
Target Compounds					Qvalue
1) TVH-Gasoline	13.006	TIC	-5250m	57.41	ug/l
18) Methylene Chloride	9.375	84	967	0.75	ug/l # 89
94) Naphthalene	19.502	128	671	0.91	ug/l 100

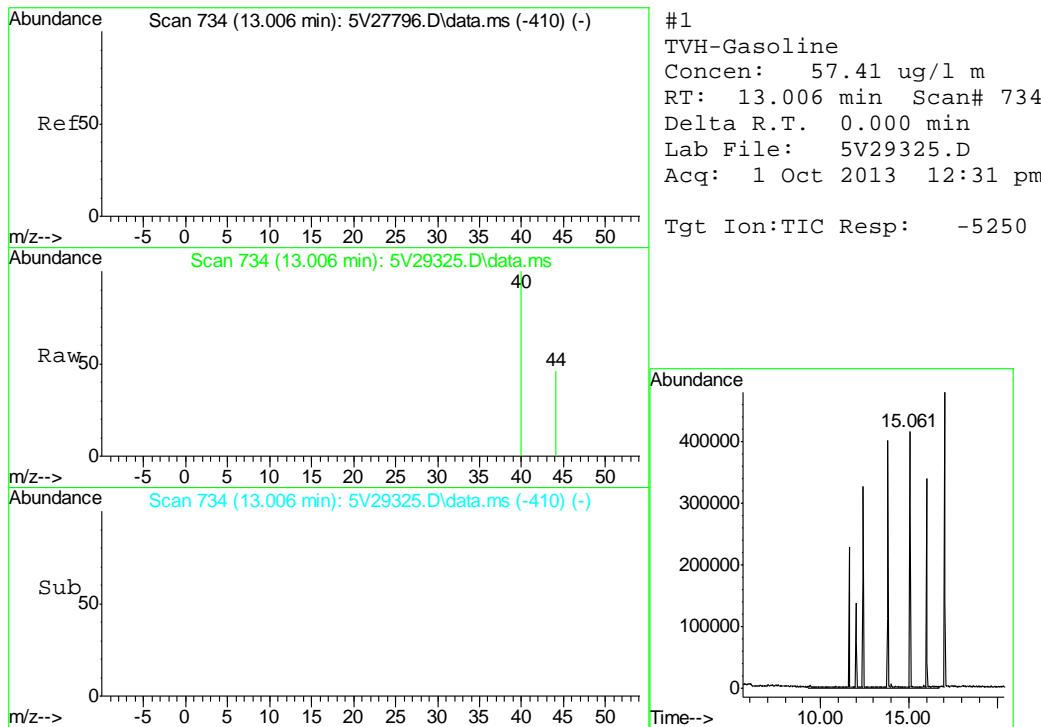
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

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 ALS Vial : 3 Sample Multiplier: 1

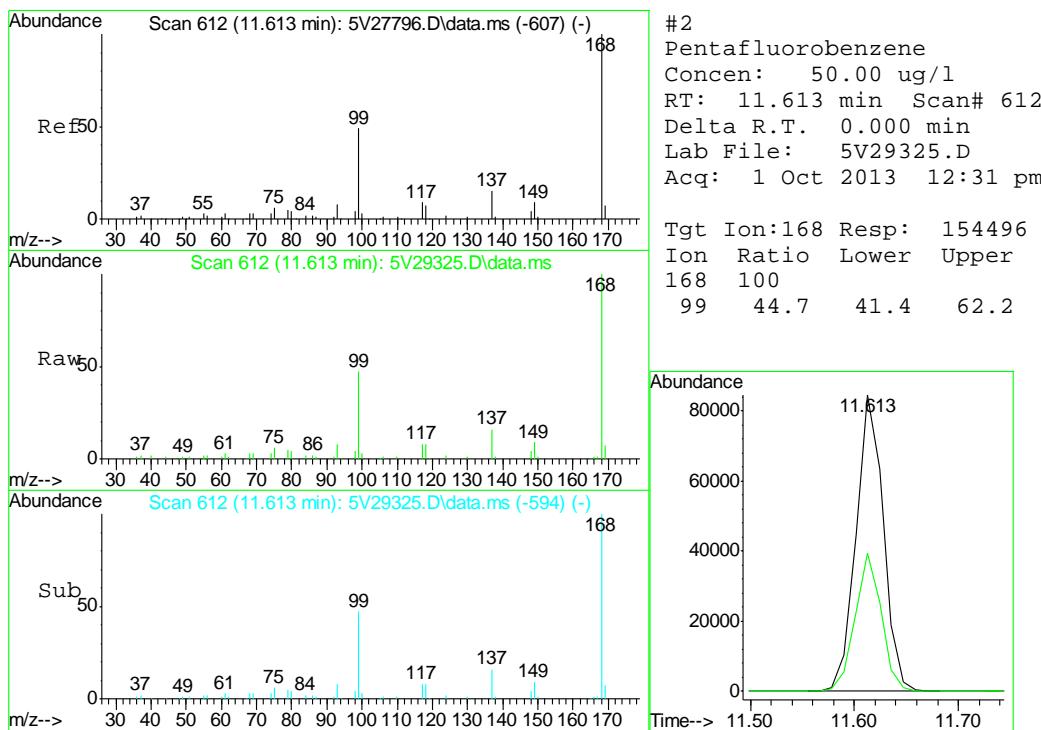
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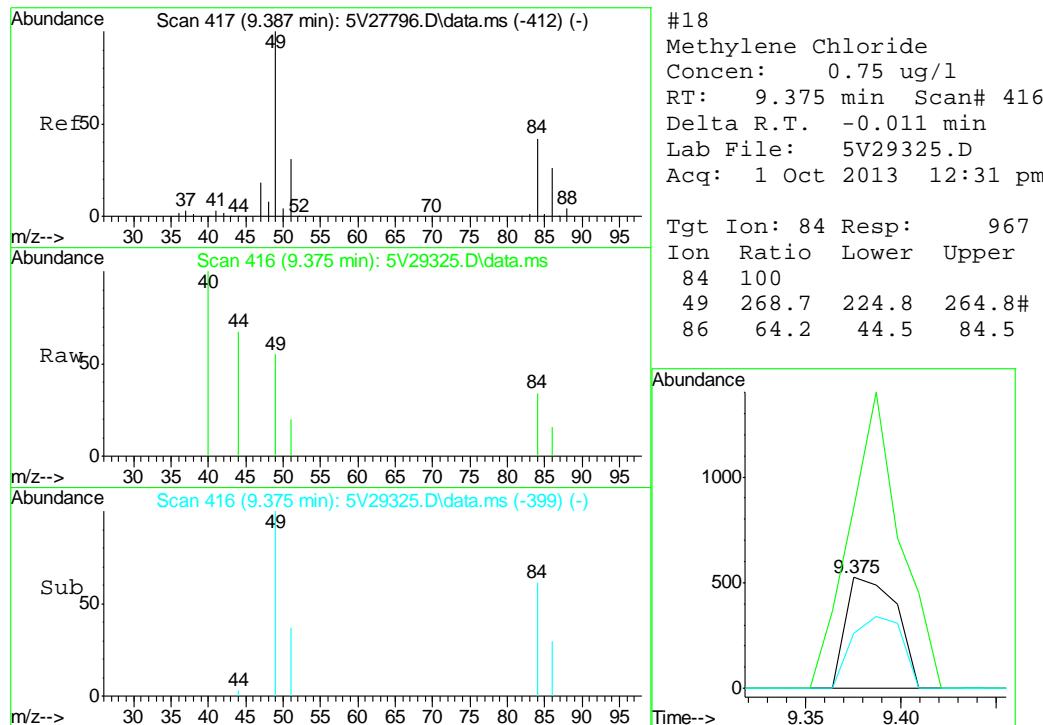




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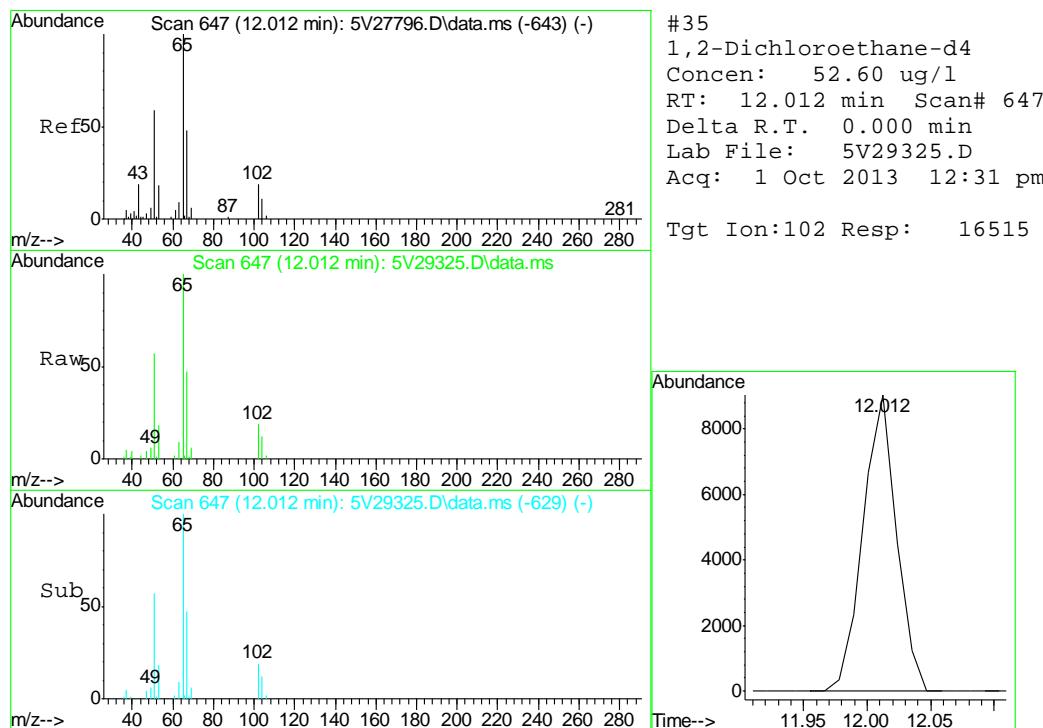
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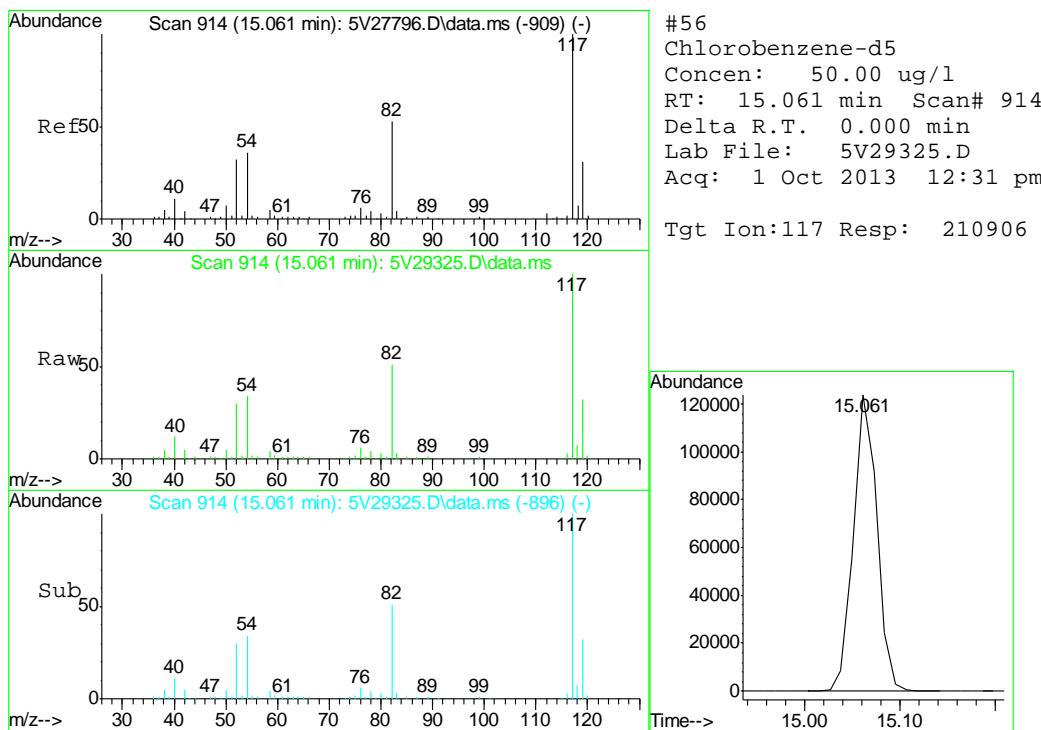
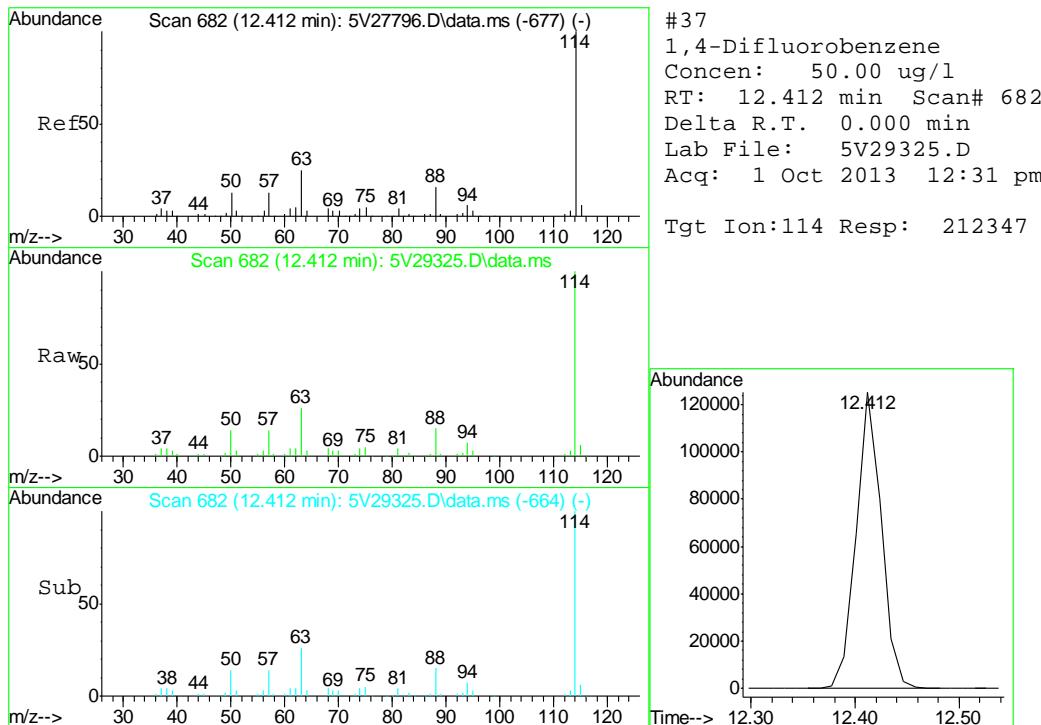


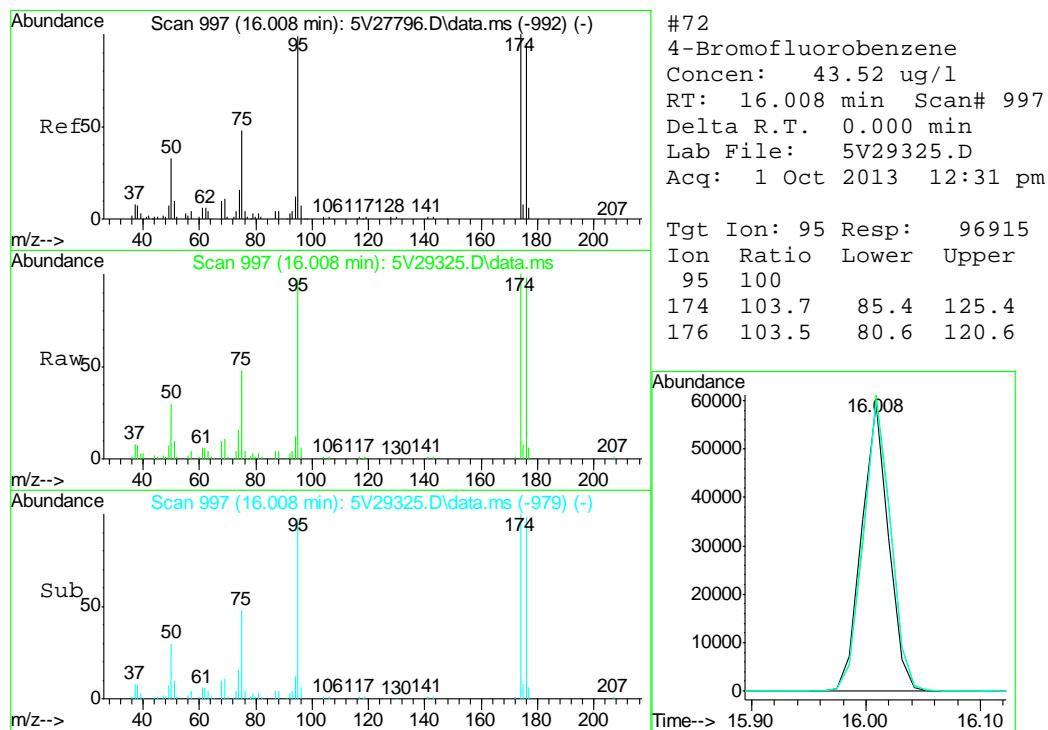
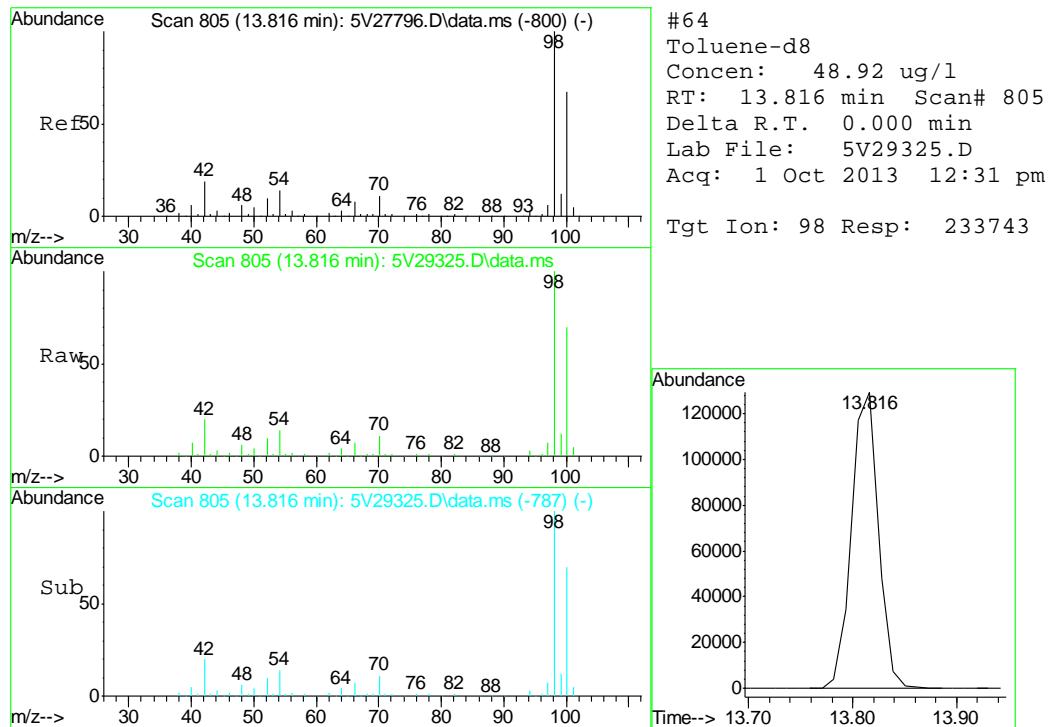


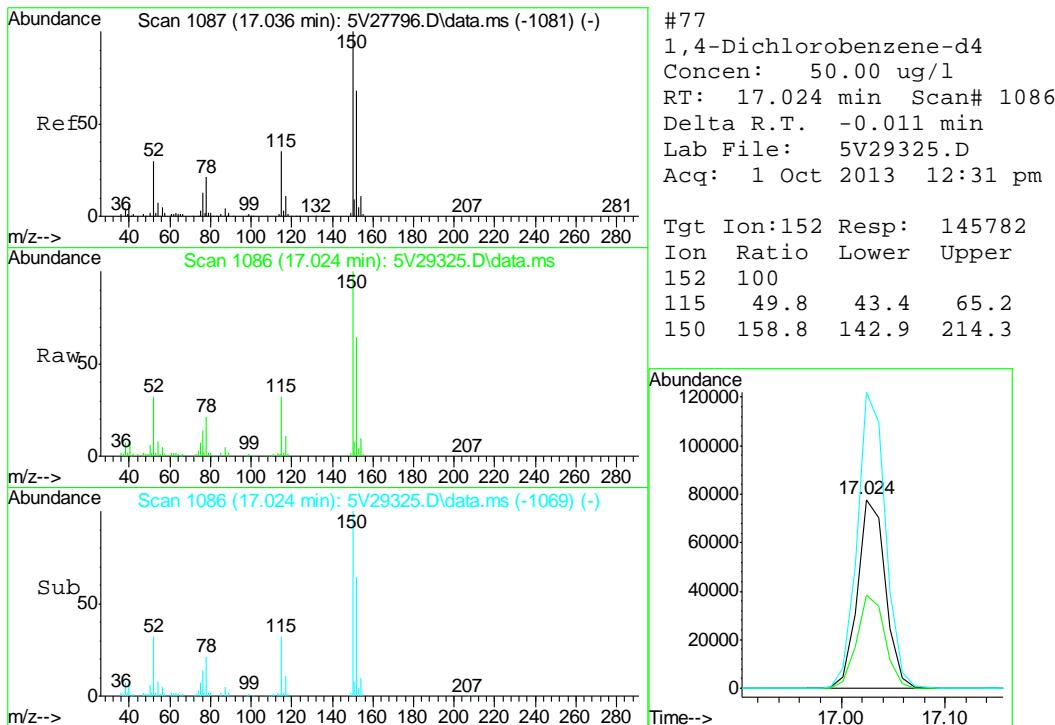
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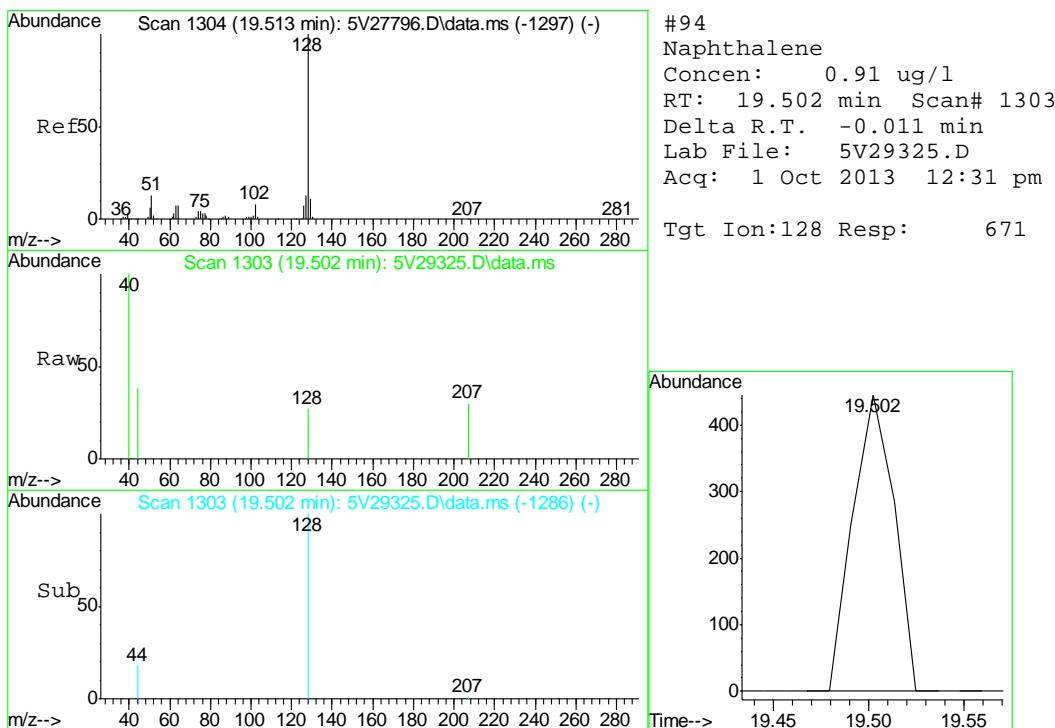






7.2.1

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GC/MS Semi-volatiles

QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Job Number: D51123
Account: XTOKWR XTO Energy
Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8670-MB	3G16517.D	1	10/03/13	DC	10/03/13	OP8670	E3G817

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D51123-1

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	8.3	4.3	ug/kg	
120-12-7	Anthracene	ND	8.3	4.3	ug/kg	
56-55-3	Benzo(a)anthracene	ND	8.3	4.3	ug/kg	
205-99-2	Benzo(b)fluoranthene	ND	8.3	4.3	ug/kg	
207-08-9	Benzo(k)fluoranthene	ND	8.3	4.3	ug/kg	
50-32-8	Benzo(a)pyrene	ND	8.3	4.3	ug/kg	
218-01-9	Chrysene	ND	8.3	4.3	ug/kg	
53-70-3	Dibenzo(a,h)anthracene	ND	8.3	4.3	ug/kg	
206-44-0	Fluoranthene	ND	8.3	4.3	ug/kg	
86-73-7	Fluorene	ND	8.3	5.0	ug/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	8.3	4.3	ug/kg	
91-20-3	Naphthalene	ND	12	10	ug/kg	
129-00-0	Pyrene	ND	8.3	4.3	ug/kg	

CAS No.	Surrogate Recoveries	Limits
4165-60-0	Nitrobenzene-d5	97%
321-60-8	2-Fluorobiphenyl	89%
1718-51-0	Terphenyl-d14	112%

Blank Spike Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8670-BS	3G16518.D	1	10/03/13	DC	10/03/13	OP8670	E3G817

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D51123-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	71.6	86	55-130
120-12-7	Anthracene	83.3	65.3	78	60-130
56-55-3	Benzo(a)anthracene	83.3	70.3	84	62-130
205-99-2	Benzo(b)fluoranthene	83.3	77.2	93	55-130
207-08-9	Benzo(k)fluoranthene	83.3	57.0	68	59-130
50-32-8	Benzo(a)pyrene	83.3	63.2	76	64-130
218-01-9	Chrysene	83.3	68.3	82	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	63.1	76	56-130
206-44-0	Fluoranthene	83.3	62.6	75	59-130
86-73-7	Fluorene	83.3	73.5	88	58-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	62.9	75	60-130
91-20-3	Naphthalene	83.3	69.2	83	56-130
129-00-0	Pyrene	83.3	72.0	86	65-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	100%	10-175%
321-60-8	2-Fluorobiphenyl	91%	25-130%
1718-51-0	Terphenyl-d14	105%	41-133%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8670-MS1	3G16520.D	1	10/03/13	DC	10/03/13	OP8670	E3G817
OP8670-MSD1	3G16521.D	1	10/03/13	DC	10/03/13	OP8670	E3G817
D51039-1	3G16519.D	1	10/03/13	DC	10/03/13	OP8670	E3G817

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D51123-1

CAS No.	Compound	D51039-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND	102	79.3	78	74.9	73	6	29-139/30	
120-12-7	Anthracene	ND	102	76.5	75	80.3	79	5	10-182/30	
56-55-3	Benzo(a)anthracene	ND	102	84.3	82	89.9	88	6	35-149/30	
205-99-2	Benzo(b)fluoranthene	ND	102	68.5	67	72.1	71	5	22-174/30	
207-08-9	Benzo(k)fluoranthene	ND	102	82.8	81	89.3	88	8	10-185/30	
50-32-8	Benzo(a)pyrene	ND	102	72.7	71	76.7	75	5	10-168/30	
218-01-9	Chrysene	ND	102	77.0	75	83.1	81	8	10-168/30	
53-70-3	Dibenzo(a,h)anthracene	ND	102	69.4	68	73.0	72	5	12-160/30	
206-44-0	Fluoranthene	ND	102	75.4	74	79.9	78	6	20-156/30	
86-73-7	Fluorene	ND	102	86.3	84	84.4	83	2	10-164/30	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	102	69.6	68	72.7	71	4	29-136/30	
91-20-3	Naphthalene	ND	102	76.3	75	68.8	67	10	10-258/30	
129-00-0	Pyrene	ND	102	86.6	85	93.6	92	8	10-196/30	

CAS No.	Surrogate Recoveries	MS	MSD	D51039-1	Limits
4165-60-0	Nitrobenzene-d5	86%	80%	64%	10-175%
321-60-8	2-Fluorobiphenyl	77%	75%	67%	25-130%
1718-51-0	Terphenyl-d14	94%	107%	102%	41-133%

* = Outside of Control Limits.

8.3.1
8



GC/MS Semi-volatiles

Raw Data

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\100313\
 Data File : 3g16541.D
 Acq On : 3 Oct 2013 10:30 pm
 Operator : DONC
 Sample : D51123-1
 Misc : OP8670,E3G817,30.11,,,1,1
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Oct 04 15:08:02 2013
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G810.M
 Quant Title : PAHSIM BASE
 QLast Update : Tue Sep 24 08:29:29 2013
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.682	136	179046	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.398	164	103824	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.881	188	150047	4.0000	ug/mL	0.00
19) Chrysene-d12	11.508	240	117598	4.0000	ug/mL	0.00
24) Perylene-d12	12.887	264	89349	4.0000	ug/mL	0.02

System Monitoring Compounds						
2) Nitrobenzene-d5	4.996	82	775741	34.4445	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 68.88%	
7) 2-Fluorobiphenyl	6.736	172	1511248	37.3602	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 74.72%	
21) Terphenyl-d14	10.472	244	1261635	56.7024	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 113.40%	

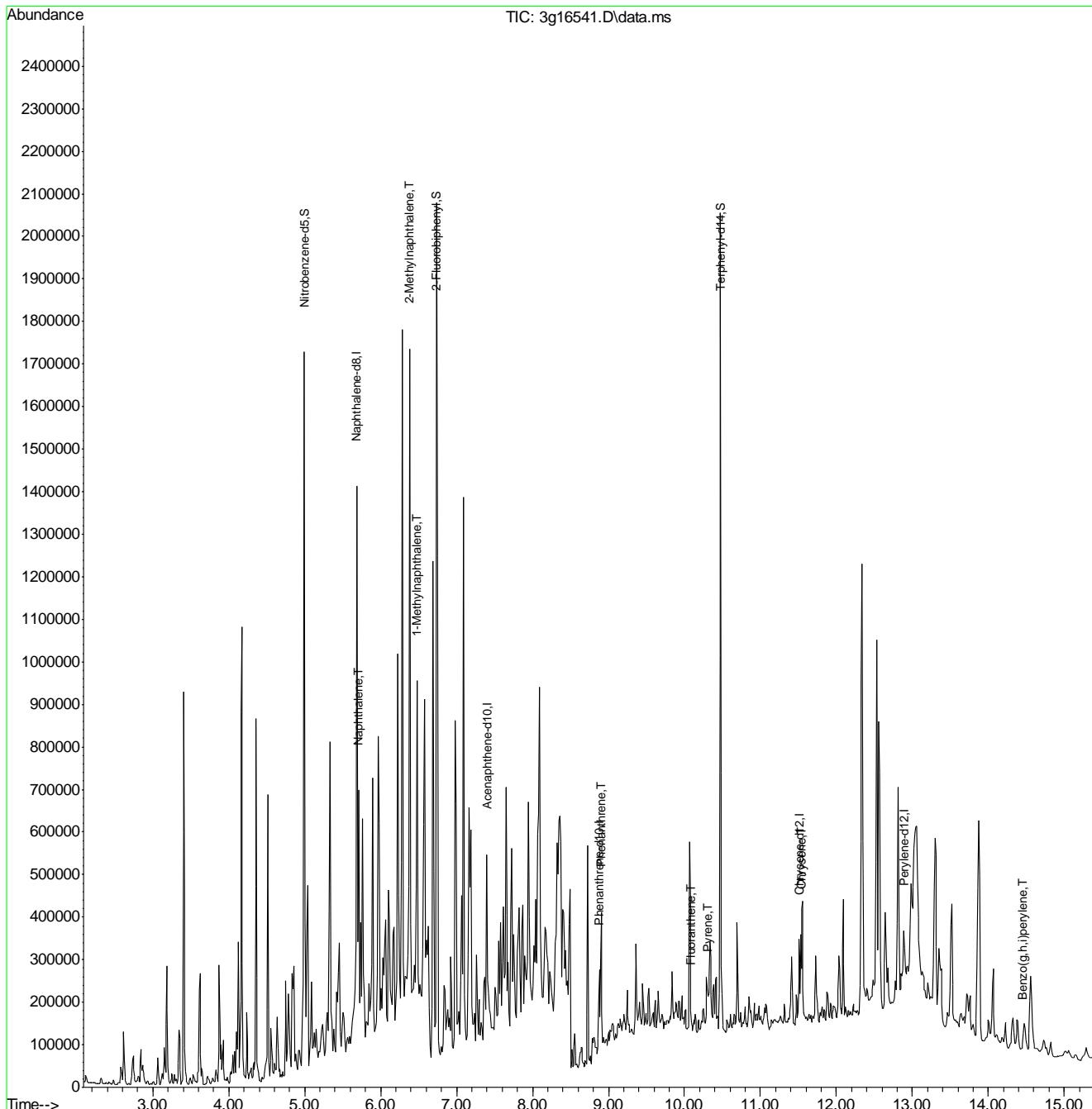
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.414	74	74	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.707	128	403589	6.2309 ug/mL	67
8) 2-Methylnaphthalene	6.380	142	589082	14.1266 ug/mL	93
9) 1-Methylnaphthalene	6.480	142	298917	8.4510 ug/mL	92
10) Acenaphthylene	0.000	152	0	N.D. d	
11) Acenaphthene	0.000	154	0	N.D. d	
12) Dibenzofuran	0.000	168	0	N.D. d	
13) Fluorene	0.000	166	0	N.D. d	
14) Diphenylamine	0.000	169	0	N.D. d	
16) Phenanthrene	8.905	178	212937	3.4878 ug/mL	70
17) Anthracene	0.000	178	0	N.D. d	
18) Fluoranthene	10.084	202	21135	0.3464 ug/mL#	37
20) Pyrene	10.306	202	30681	0.5583 ug/mL#	44
22) Benzo(a)anthracene	0.000	228	0	N.D. d	
23) Chrysene	11.535	228	45336	0.8558 ug/mL	70
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d	
26) Benzo(k)fluoranthene	0.000	252	0	N.D. d	
27) Benzo(a)pyrene	0.000	252	0	N.D. d	
28) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D. d	
29) Dibenz(a,h)anthracene	0.000	278	0	N.D. d	
30) Benzo(g,h,i)perylene	14.453	276	9194	0.2656 ug/mL#	1

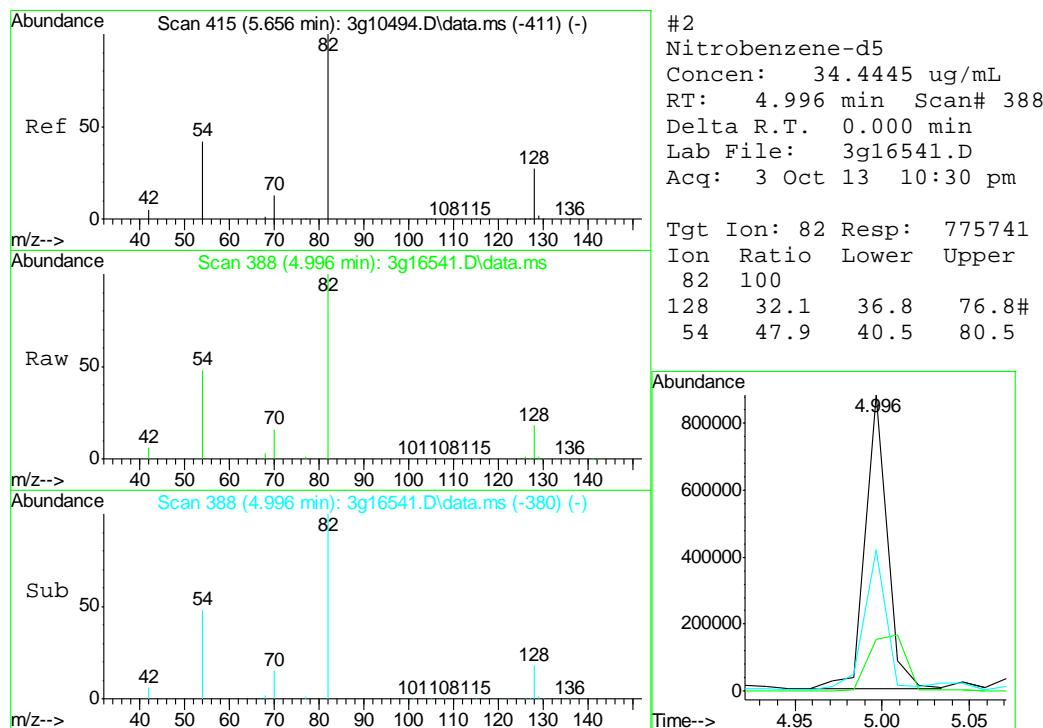
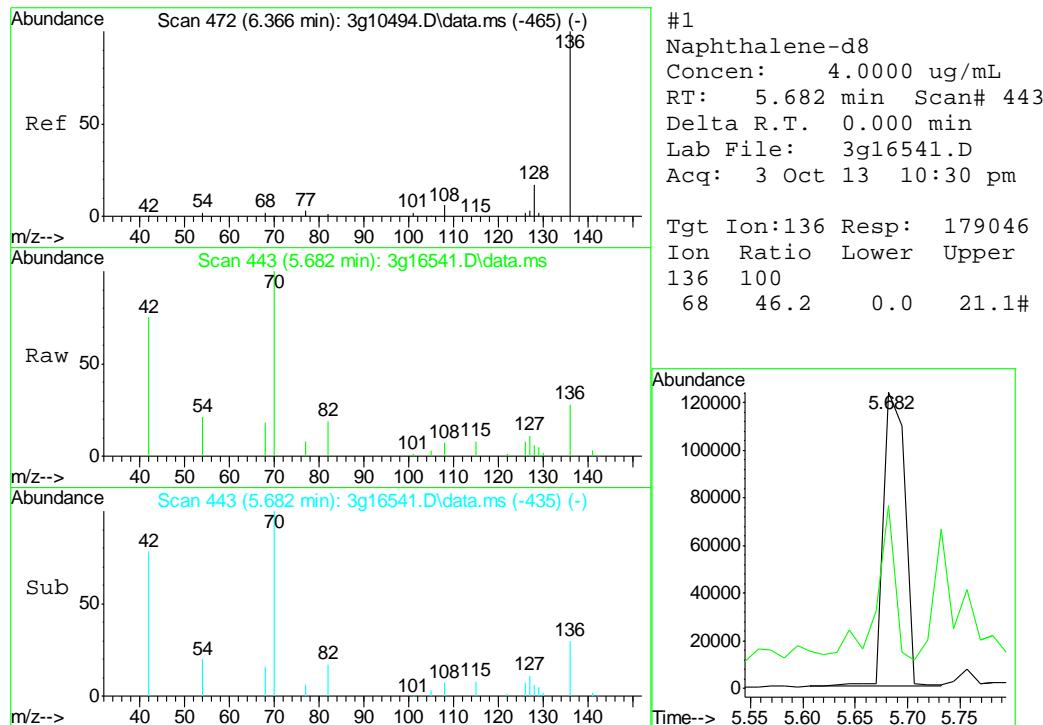
(#) = qualifier out of range (m) = manual integration (+) = signals summed

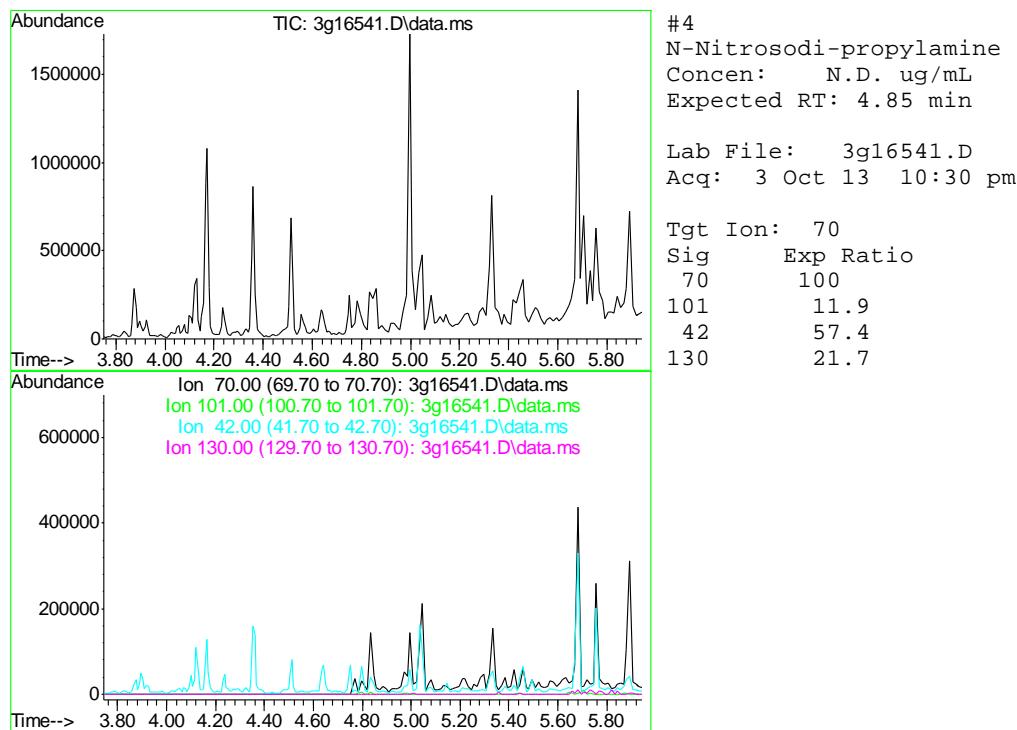
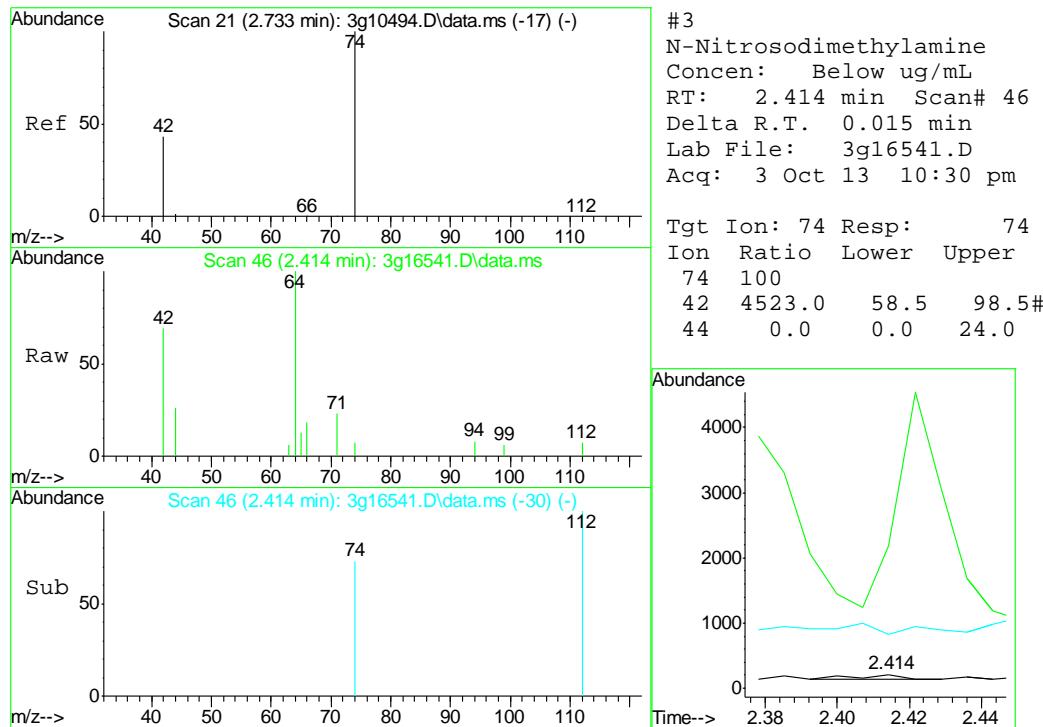
Quantitation Report (QT Reviewed)

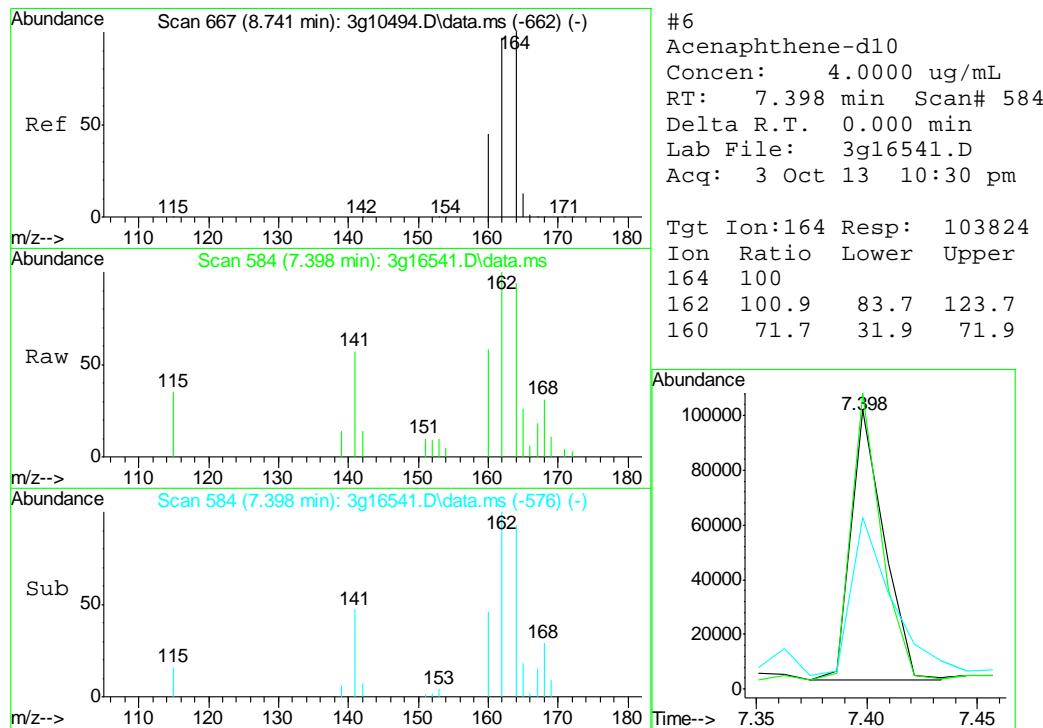
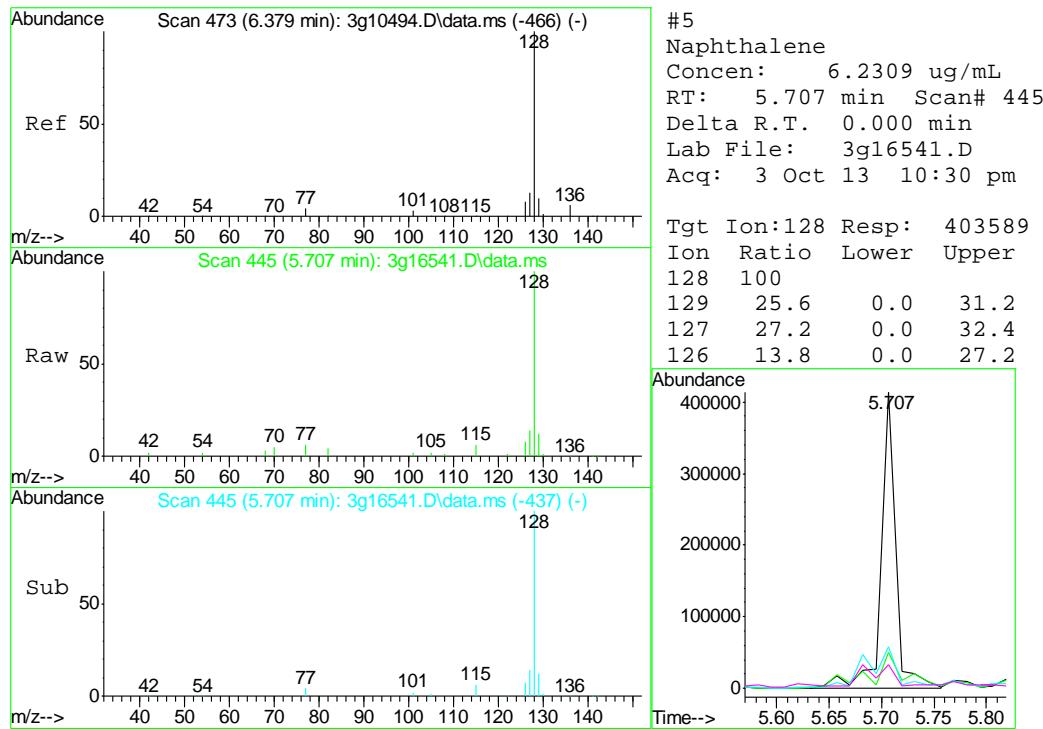
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 Operator : DONC
 Sample : D51123-1
 Misc : OP8670,E3G817,30.11,,,1,1
 ALS Vial : 28 Sample Multiplier: 1

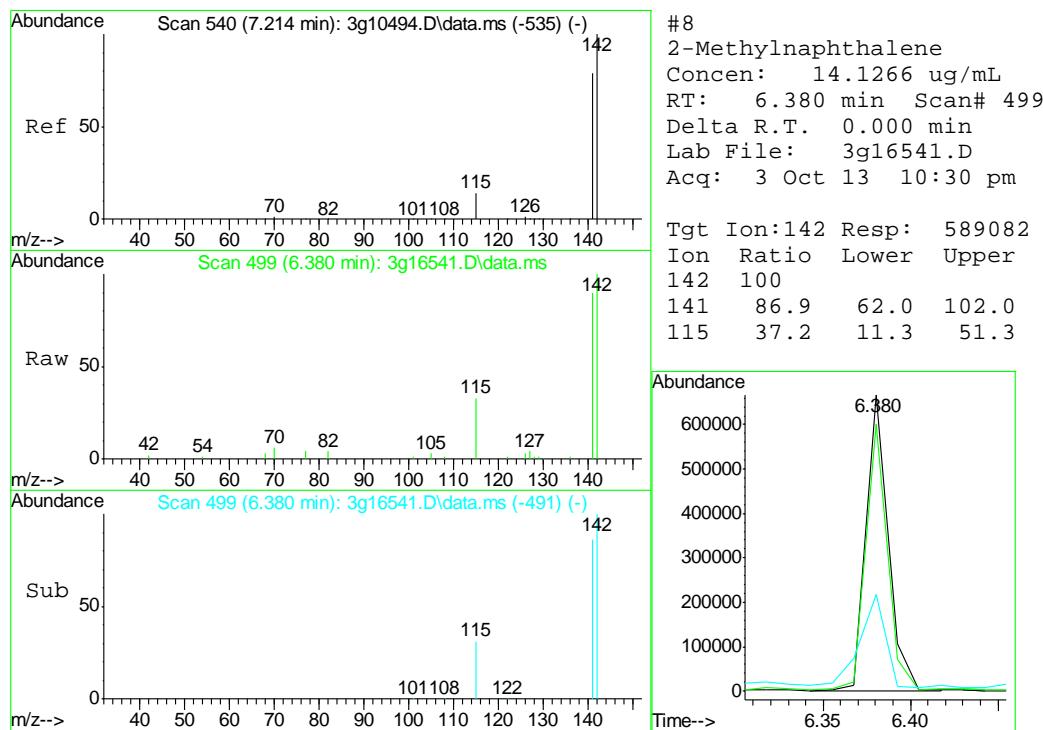
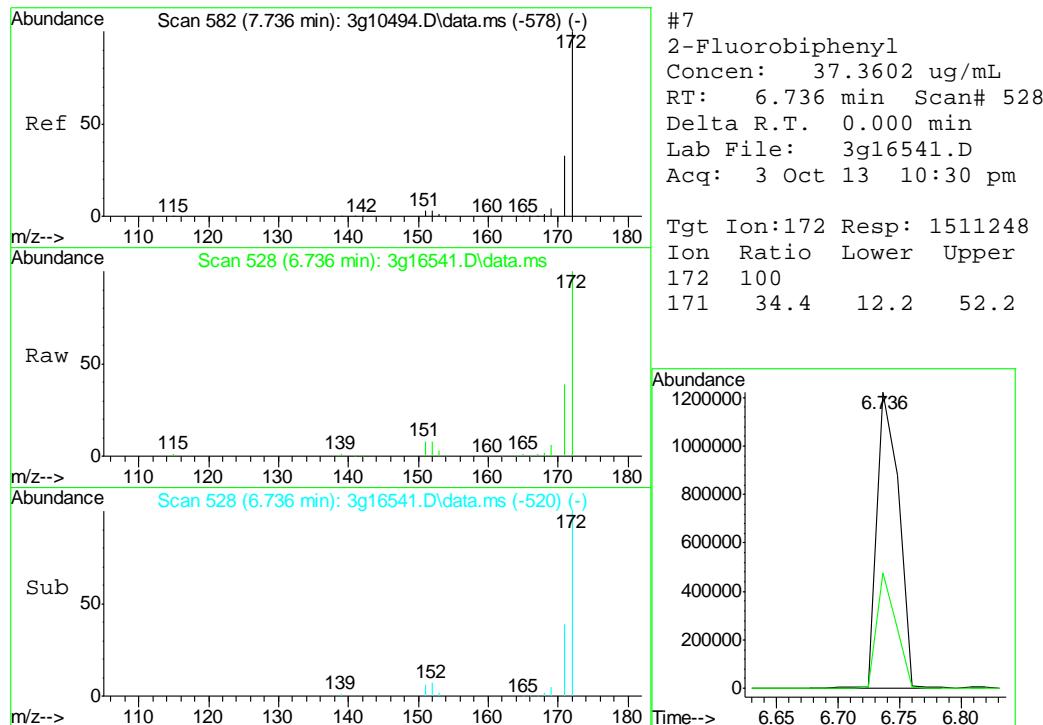
Quant Time: Oct 04 15:08:02 2013
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G810.M
 Quant Title : PAHSIM BASE
 QLast Update : Tue Sep 24 08:29:29 2013
 Response via : Initial Calibration

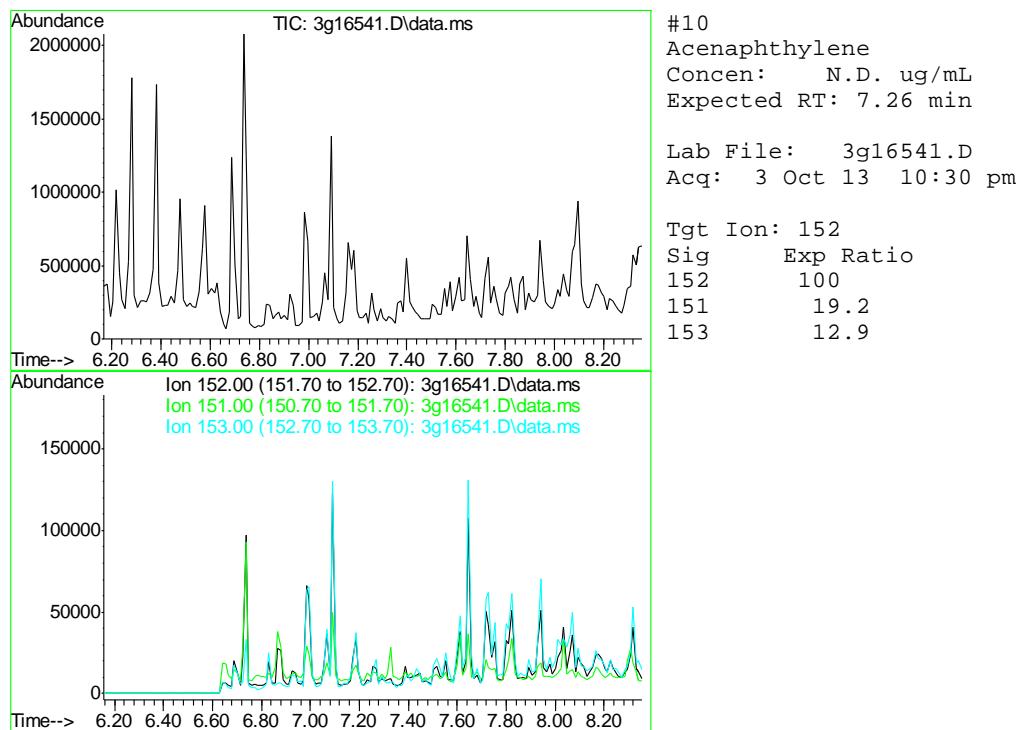
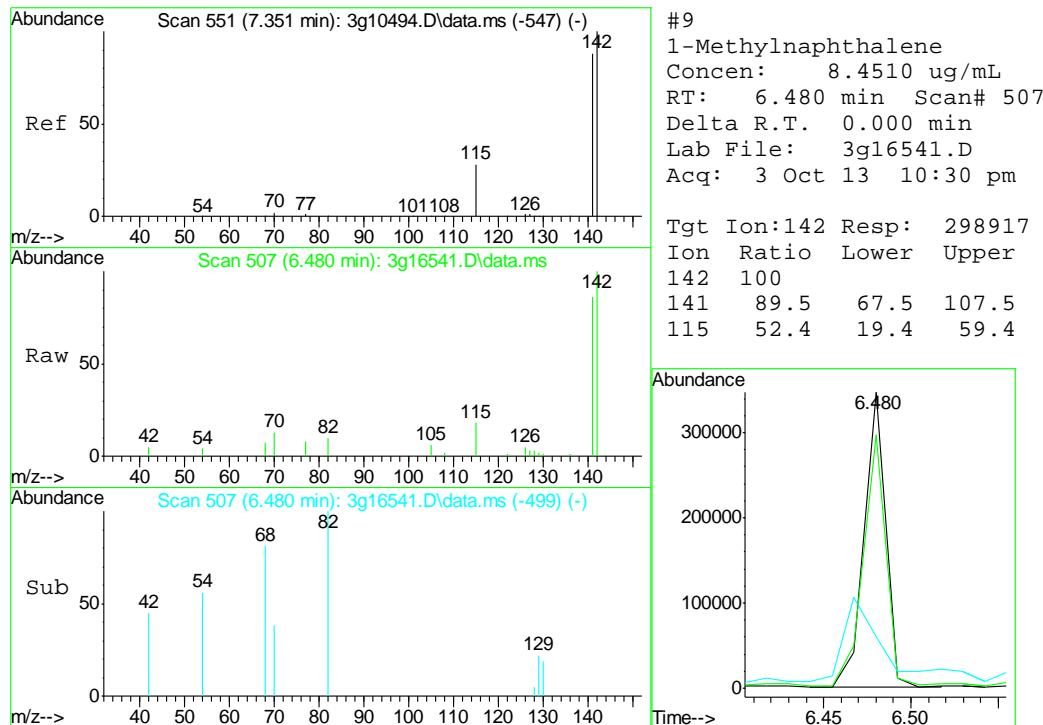


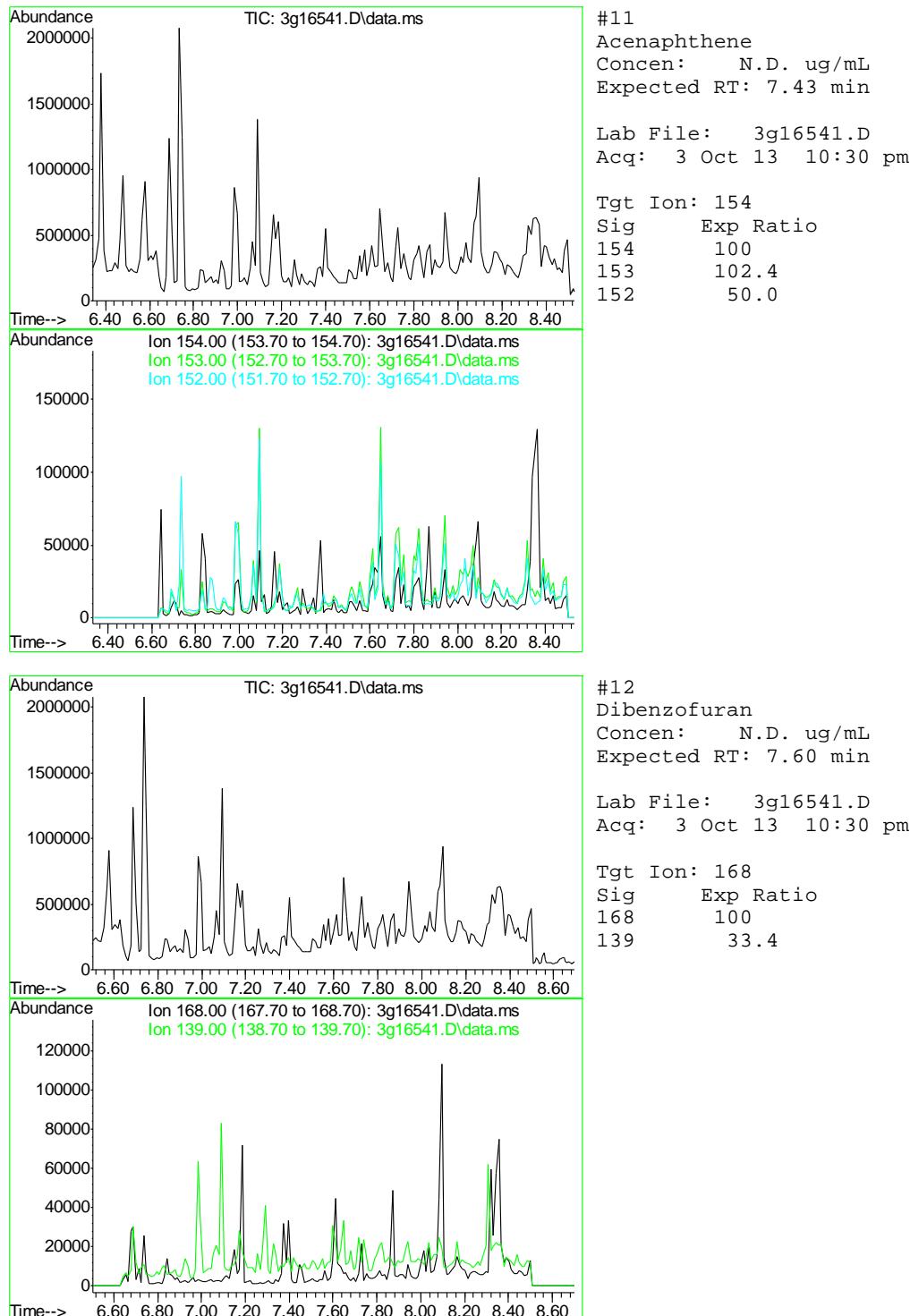


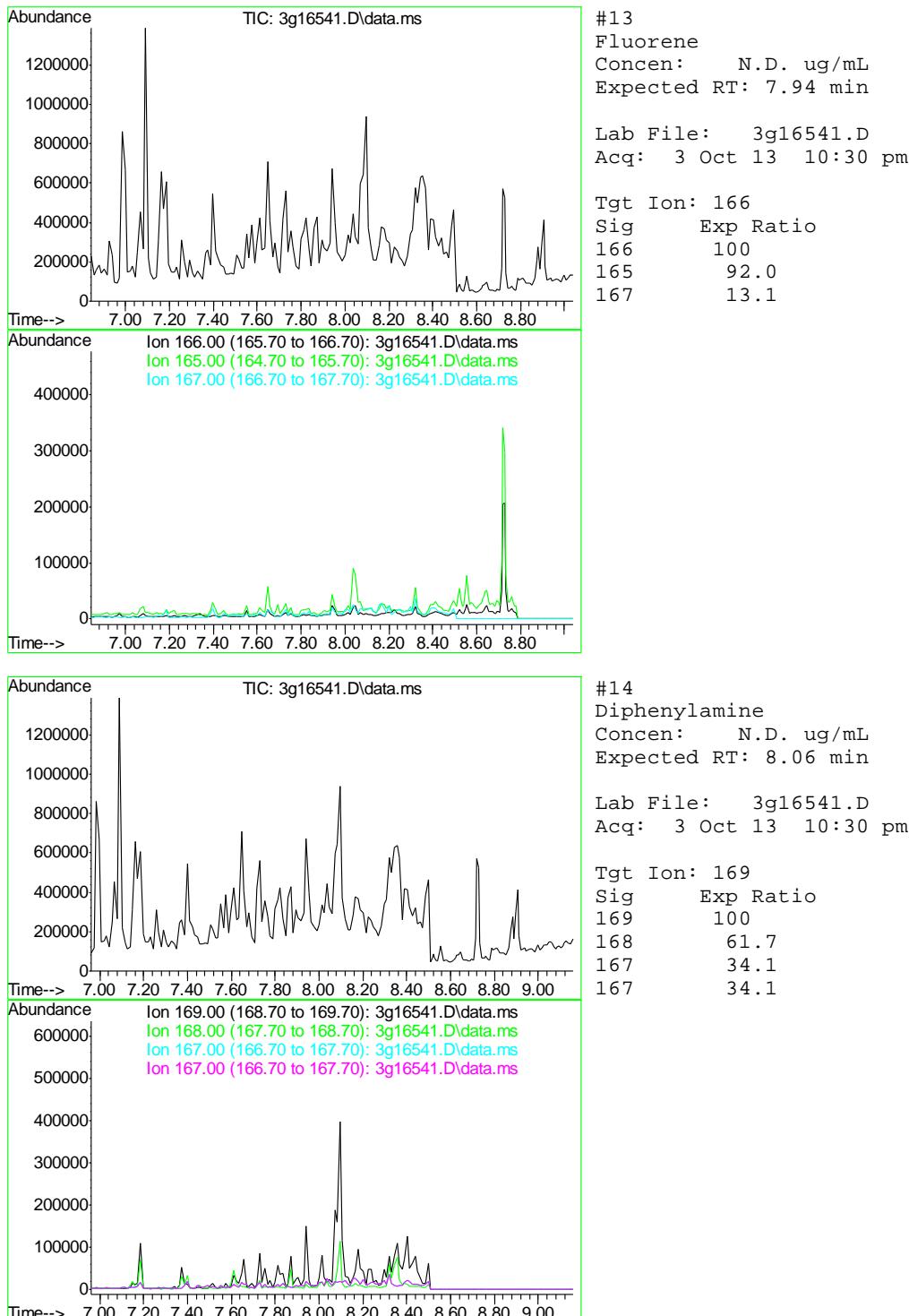


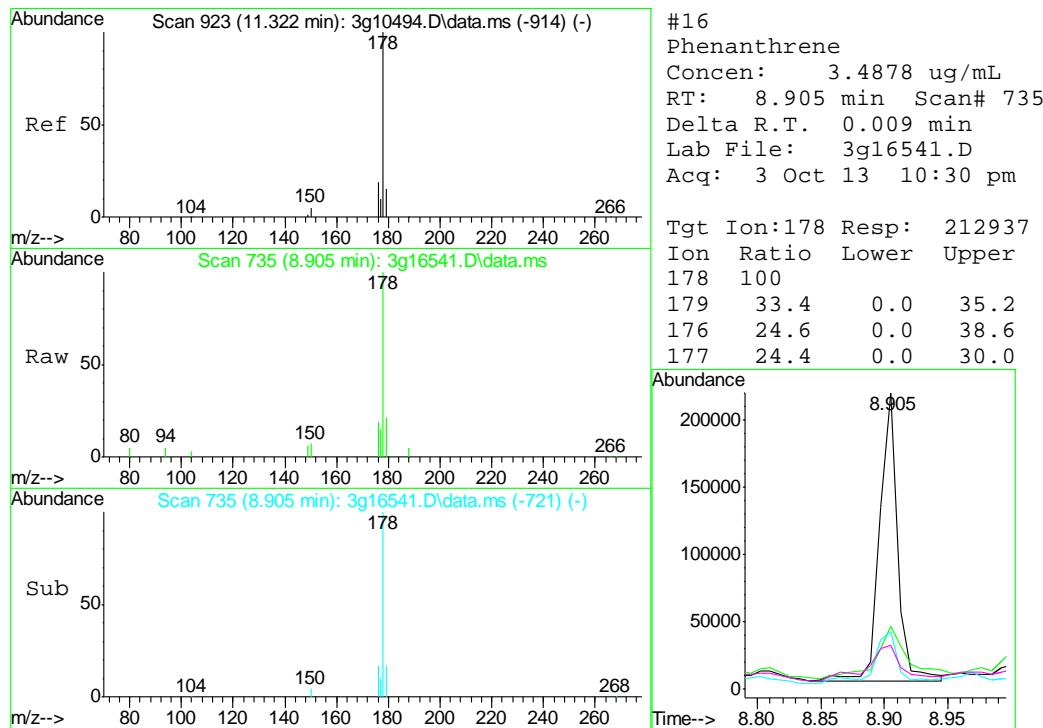
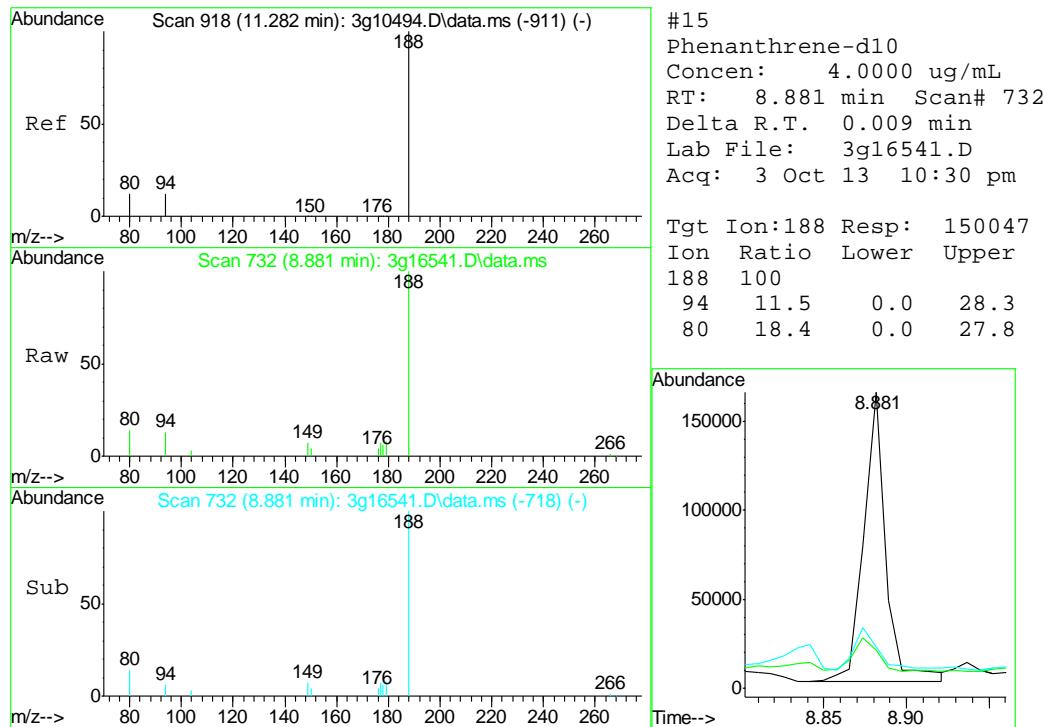


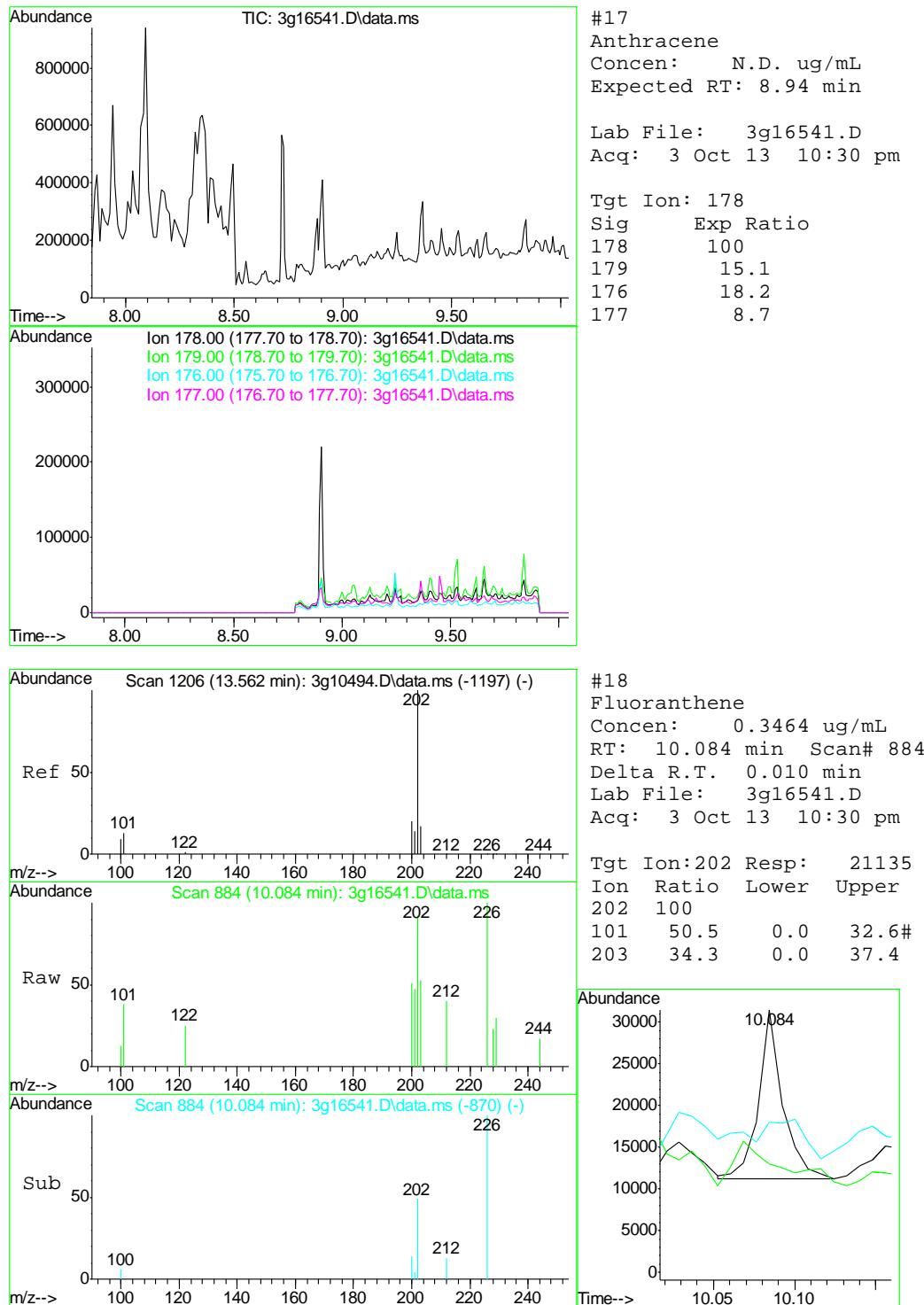


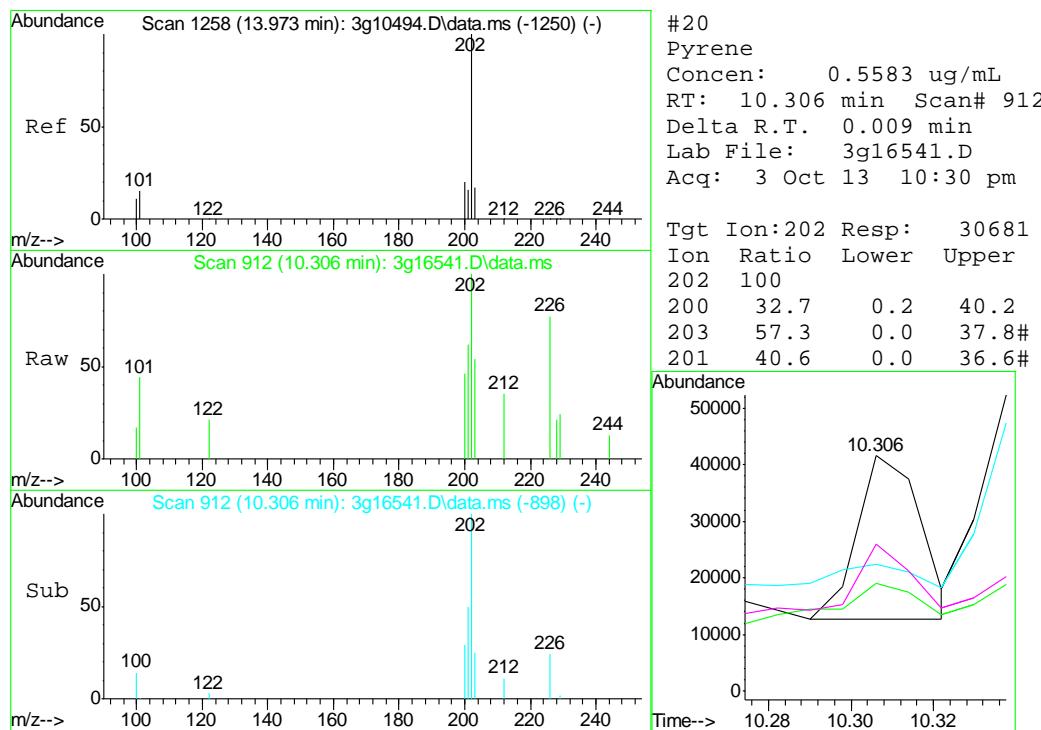
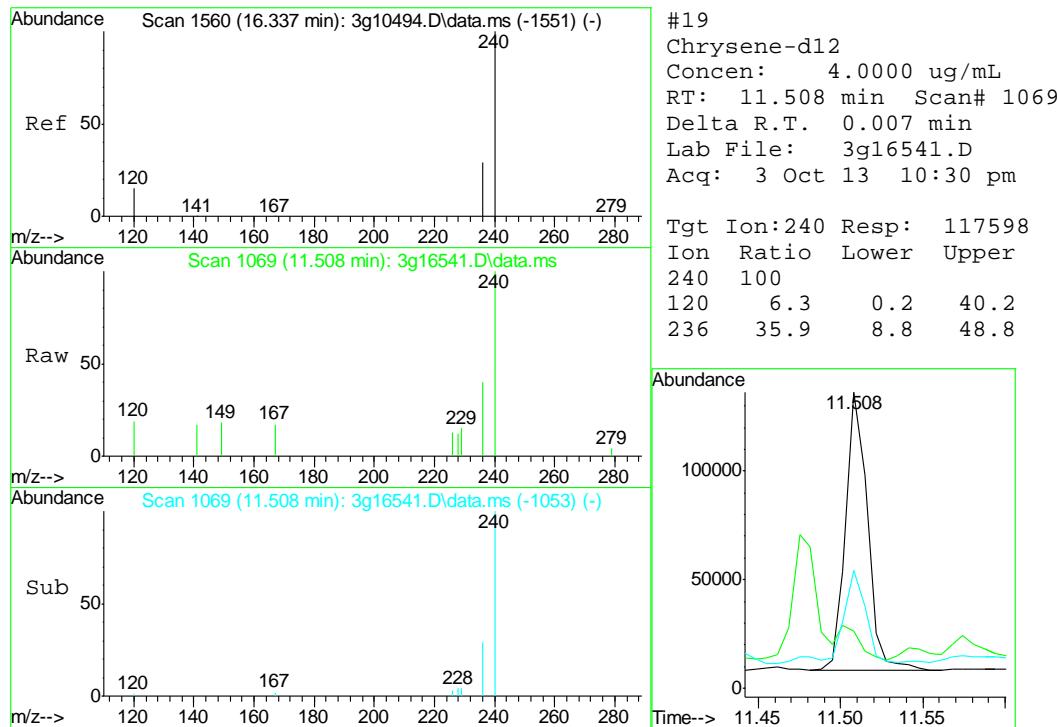


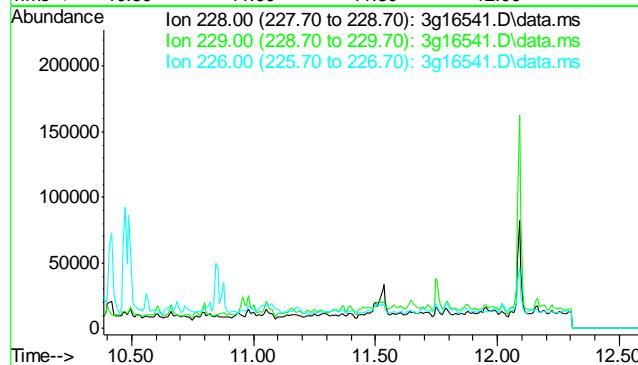
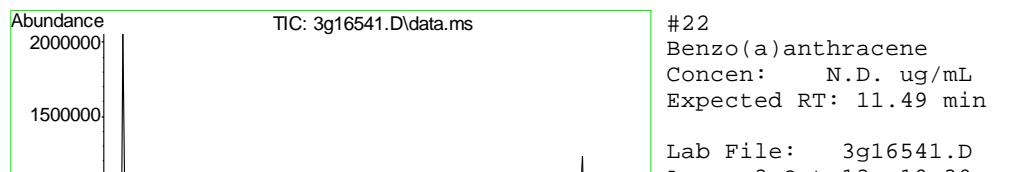
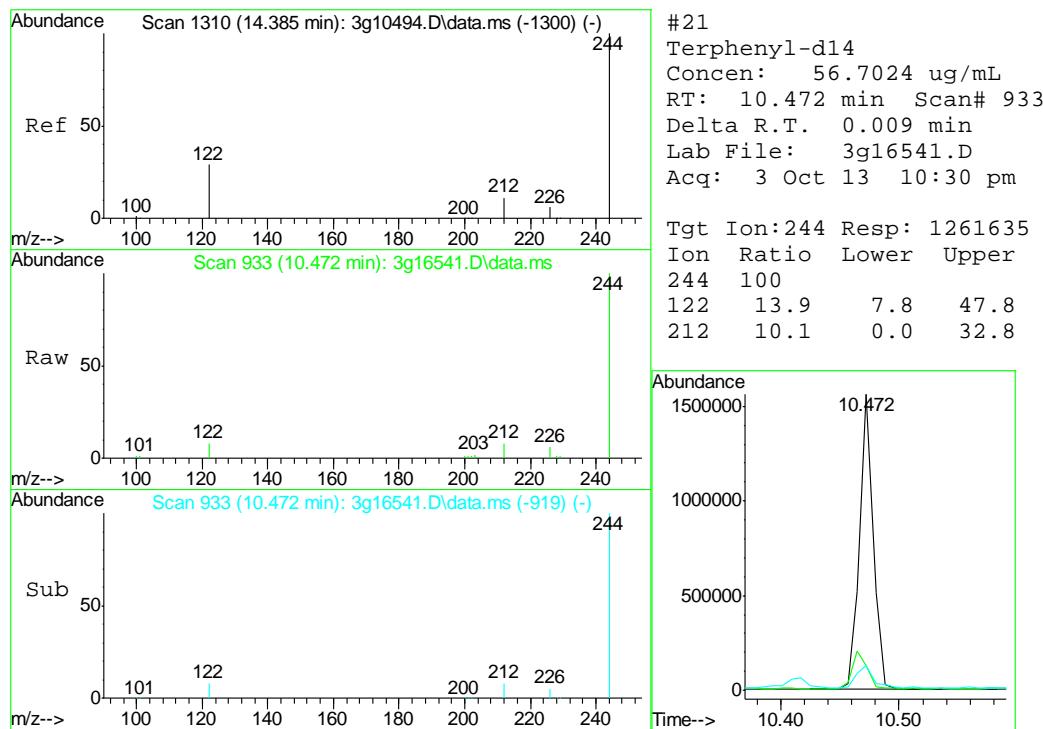


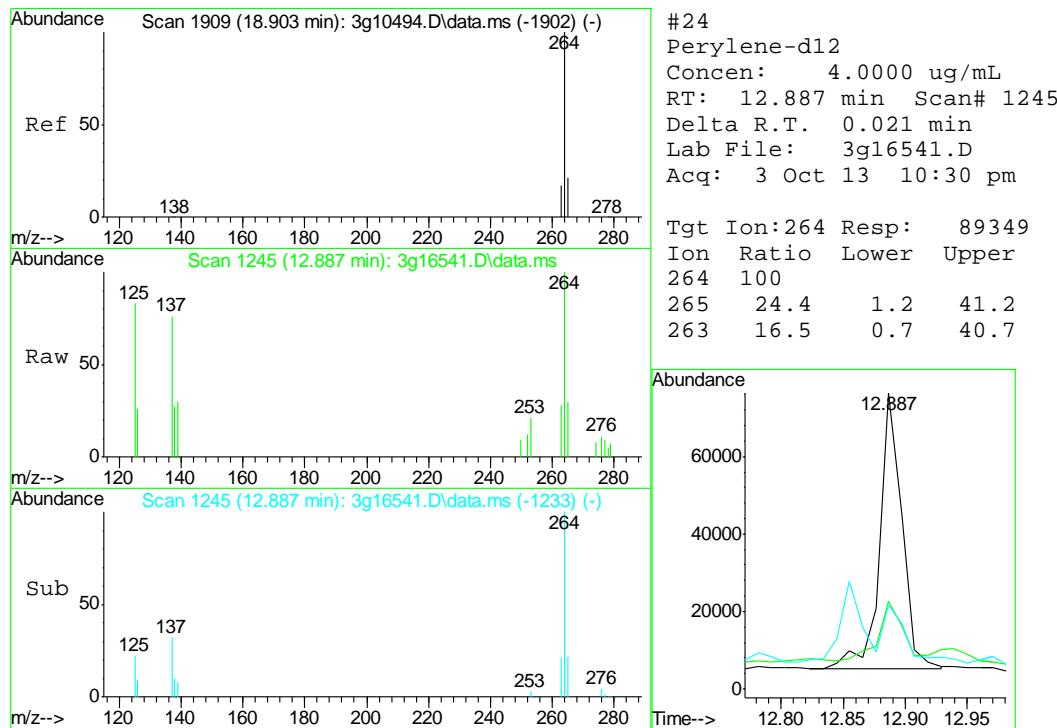
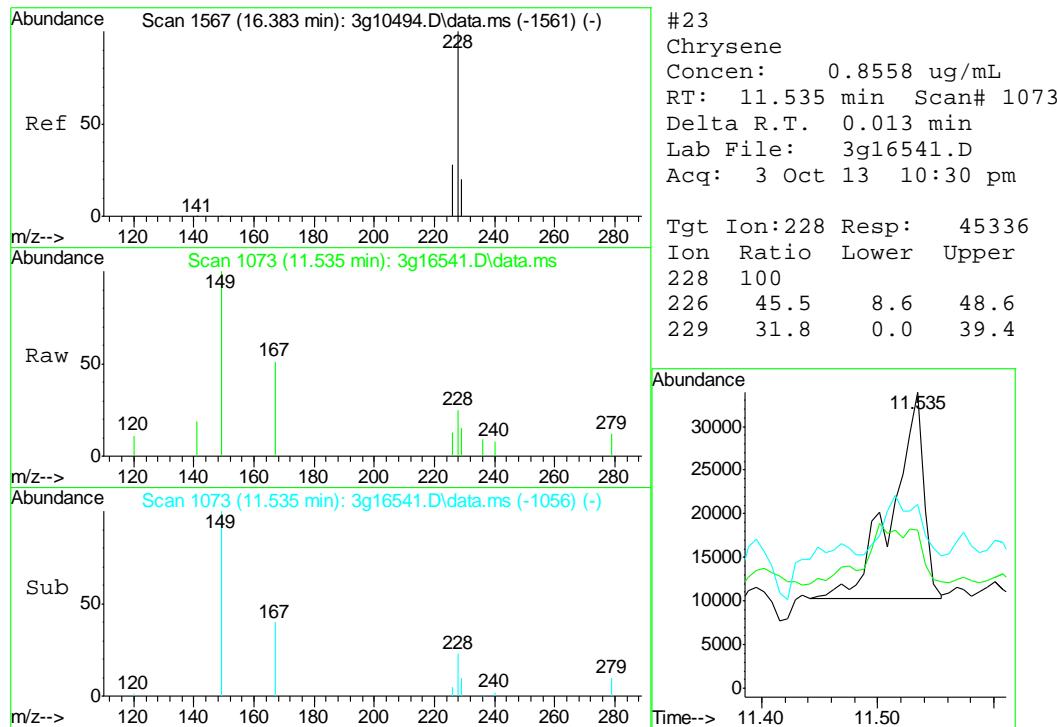


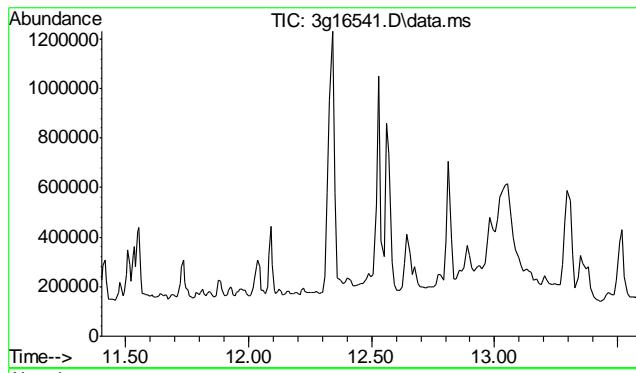








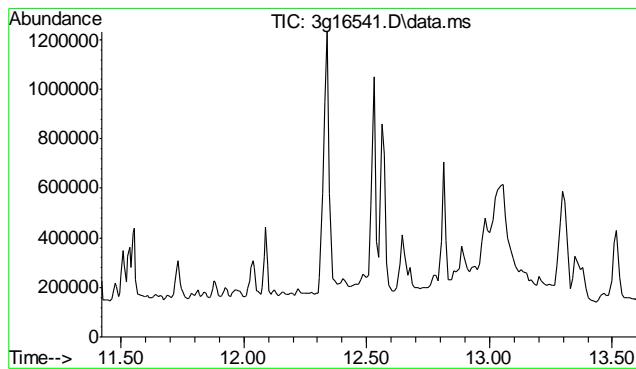
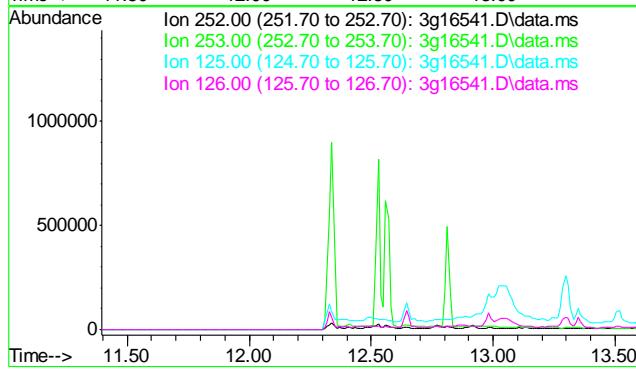




#25
 Benzo(b)fluoranthene
 Concen: N.D. ug/mL
 Expected RT: 12.50 min

Lab File: 3g16541.D
 Acq: 3 Oct 13 10:30 pm

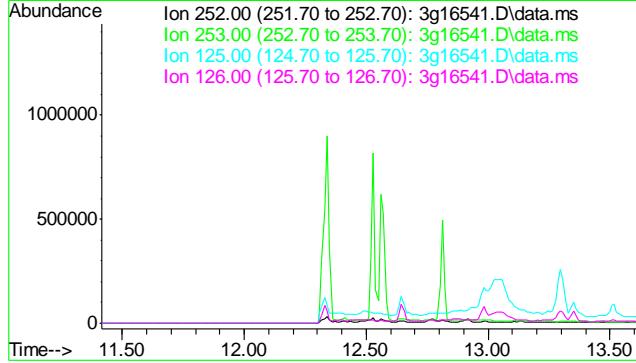
Tgt Ion:	Sig	Exp Ratio
252	100	
253	51.5	
125	13.2	
126	46.9	

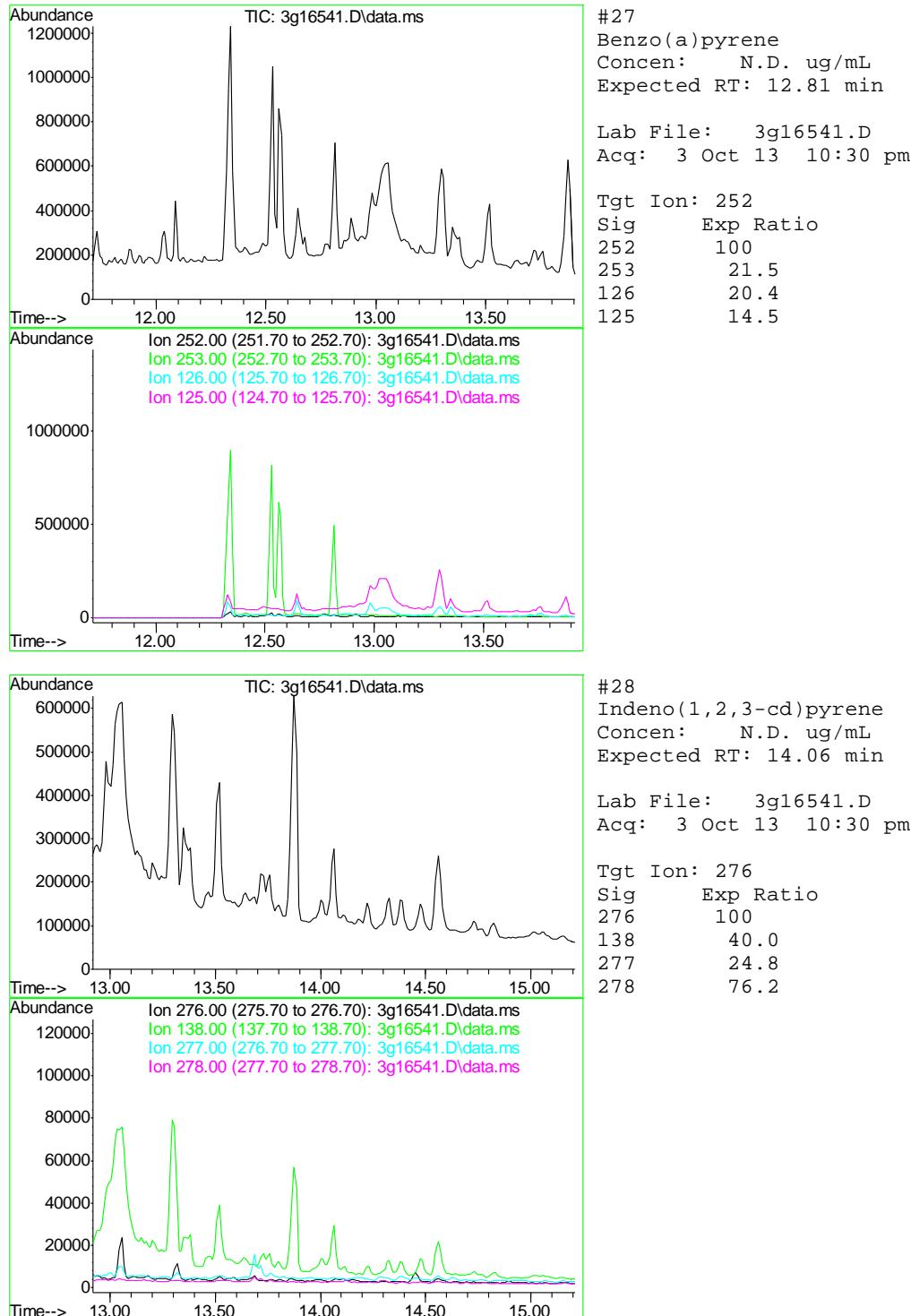


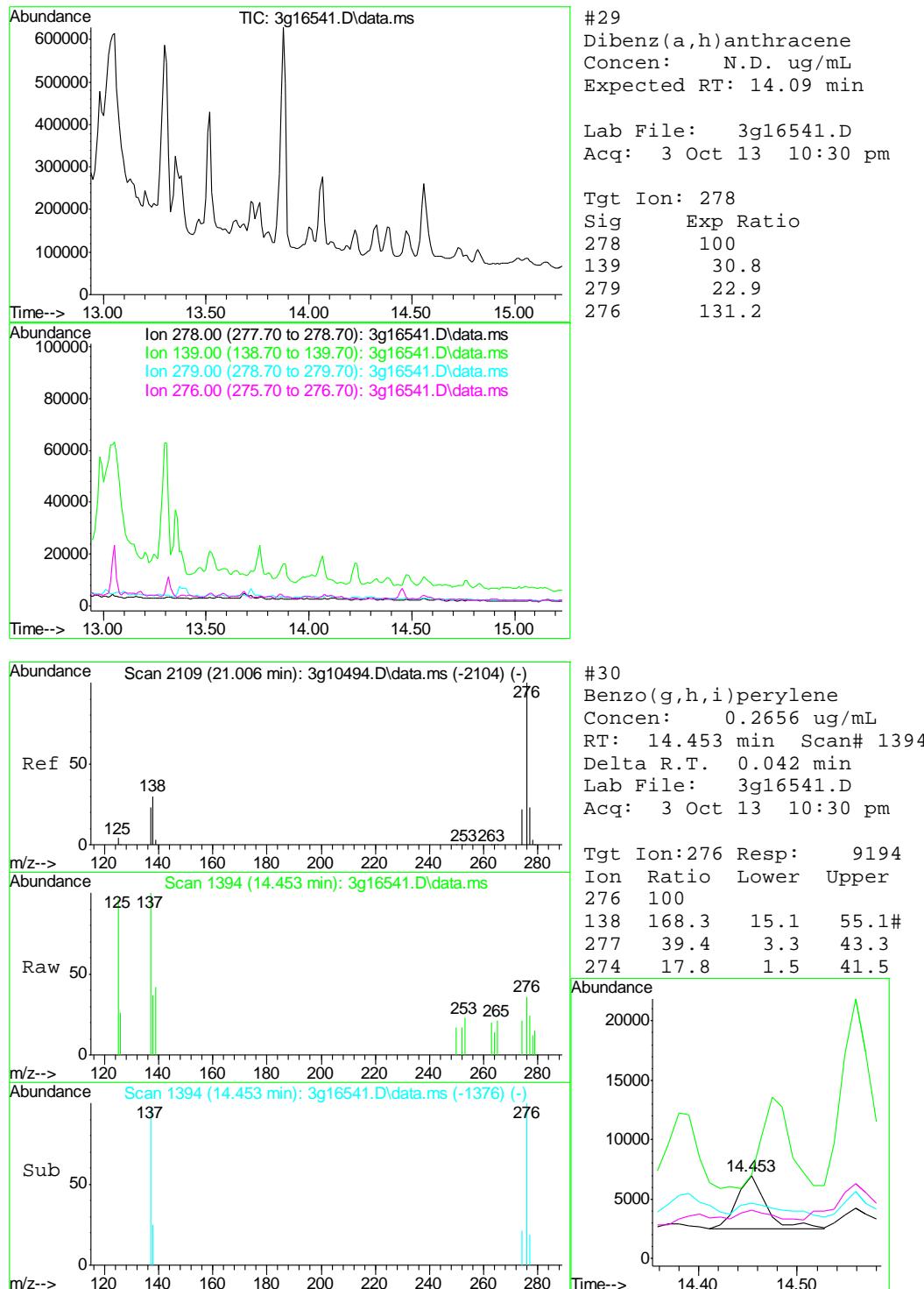
#26
 Benzo(k)fluoranthene
 Concen: N.D. ug/mL
 Expected RT: 12.52 min

Lab File: 3g16541.D
 Acq: 3 Oct 13 10:30 pm

Tgt Ion:	Sig	Exp Ratio
252	100	
253	37.3	
125	9.6	
126	34.1	







Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\100313\
 Data File : 3g16517.D
 Acq On : 3 Oct 2013 12:54 pm
 Operator : DONC
 Sample : OP8670-MB
 Misc : OP8670,E3G817,30.00,,,1,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Oct 04 13:28:03 2013
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G810.M
 Quant Title : PAHSIM BASE
 QLast Update : Tue Sep 24 08:29:29 2013
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.682	136	217864	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.398	164	117785	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.873	188	188996	4.0000	ug/mL	0.00
19) Chrysene-d12	11.501	240	166287	4.0000	ug/mL	0.00
24) Perylene-d12	12.865	264	135532	4.0000	ug/mL	0.00

System Monitoring Compounds						
2) Nitrobenzene-d5	4.996	82	1331104	48.5730	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 97.14%	
7) 2-Fluorobiphenyl	6.736	172	2051275	44.6998	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 89.40%	
21) Terphenyl-d14	10.464	244	1754721	55.7722	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 111.54%	

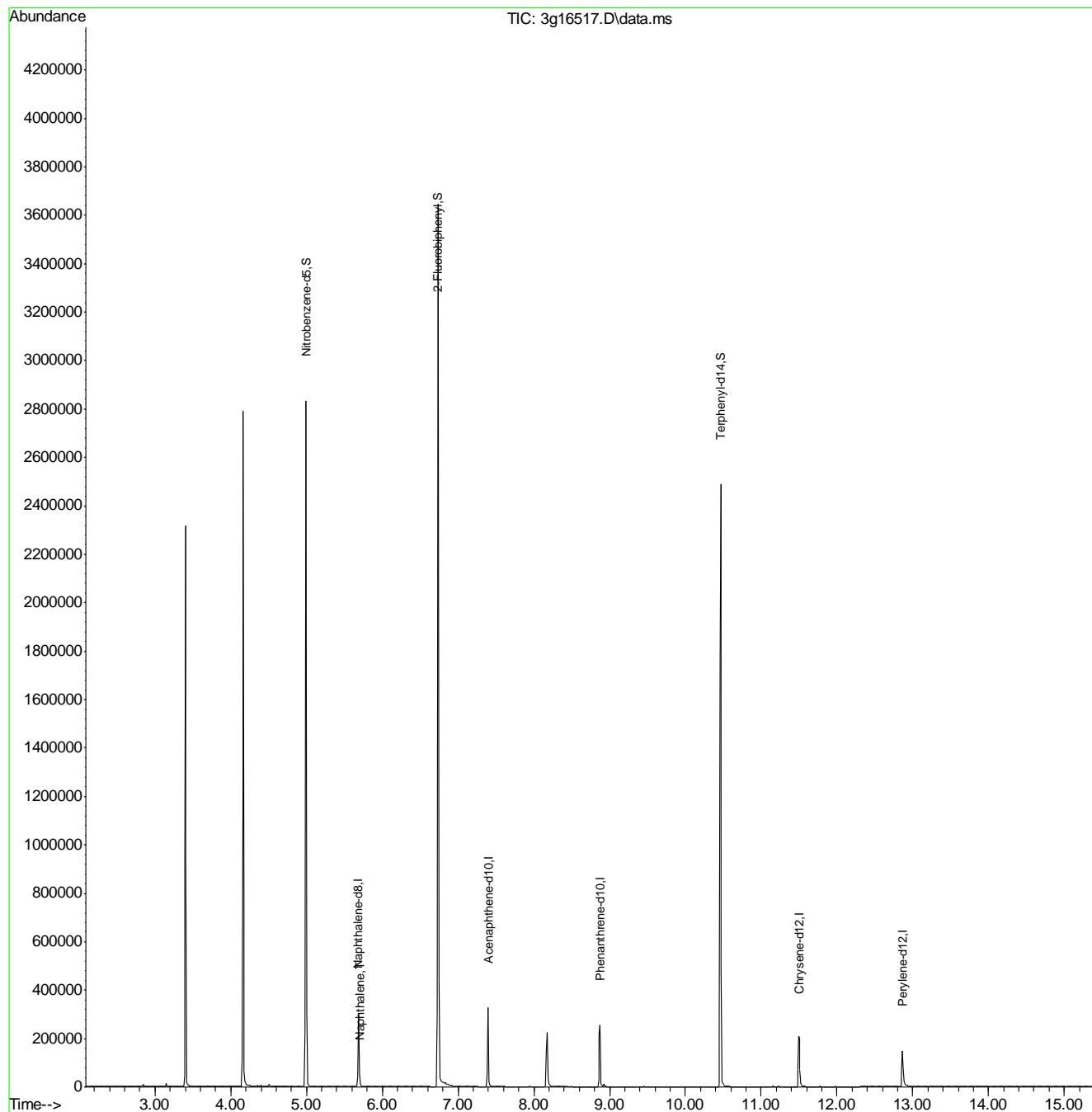
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D.	d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D.	d
5) Naphthalene	5.707	128	6089	0.0773	ug/mL 98
8) 2-Methylnaphthalene	6.380	142	1089	N.D.	
9) 1-Methylnaphthalene	6.480	142	479	N.D.	
10) Acenaphthylene	7.256	152	63	N.D.	
11) Acenaphthene	7.422	154	1060	N.D.	
12) Dibenzofuran	7.599	168	528	N.D.	
13) Fluorene	7.941	166	540	N.D.	
14) Diphenylamine	0.000	169	0	N.D.	d
16) Phenanthrene	8.889	178	943	N.D.	
17) Anthracene	0.000	178	0	N.D.	d
18) Fluoranthene	0.000	202	0	N.D.	d
20) Pyrene	10.298	202	337	N.D.	
22) Benzo(a)anthracene	11.495	228	896	N.D.	
23) Chrysene	11.495	228	896	N.D.	
25) Benzo(b)fluoranthene	0.000	252	0	N.D.	d
26) Benzo(k)fluoranthene	0.000	252	0	N.D.	d
27) Benzo(a)pyrene	0.000	252	0	N.D.	d
28) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D.	d
29) Dibenz(a,h)anthracene	0.000	278	0	N.D.	d
30) Benzo(g,h,i)perylene	0.000	276	0	N.D.	d

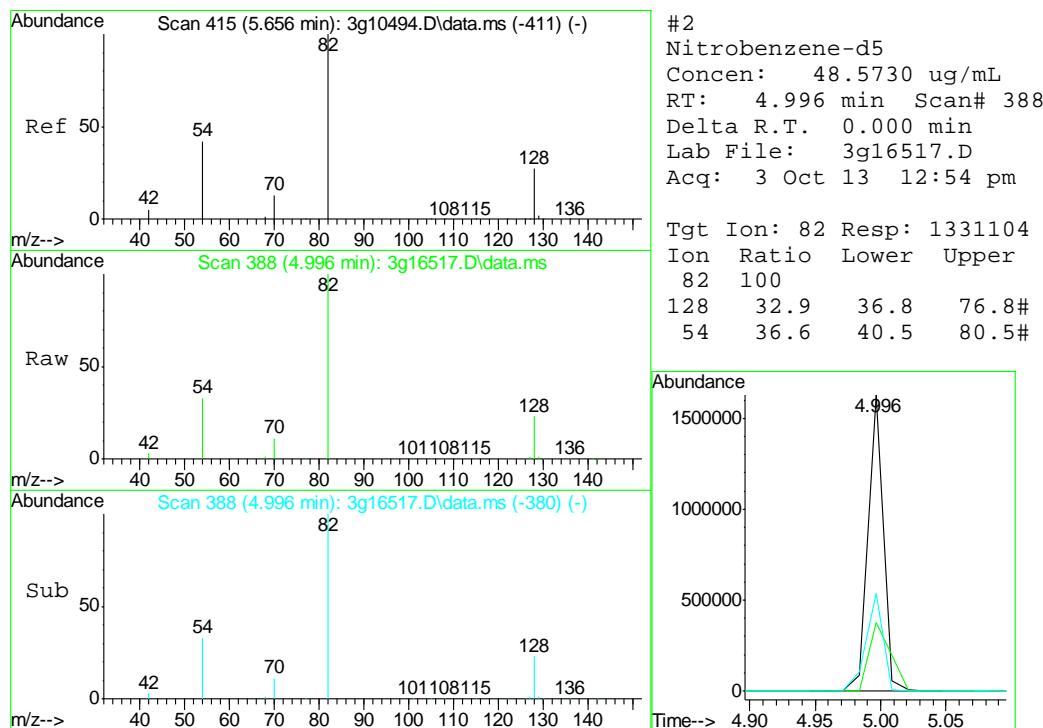
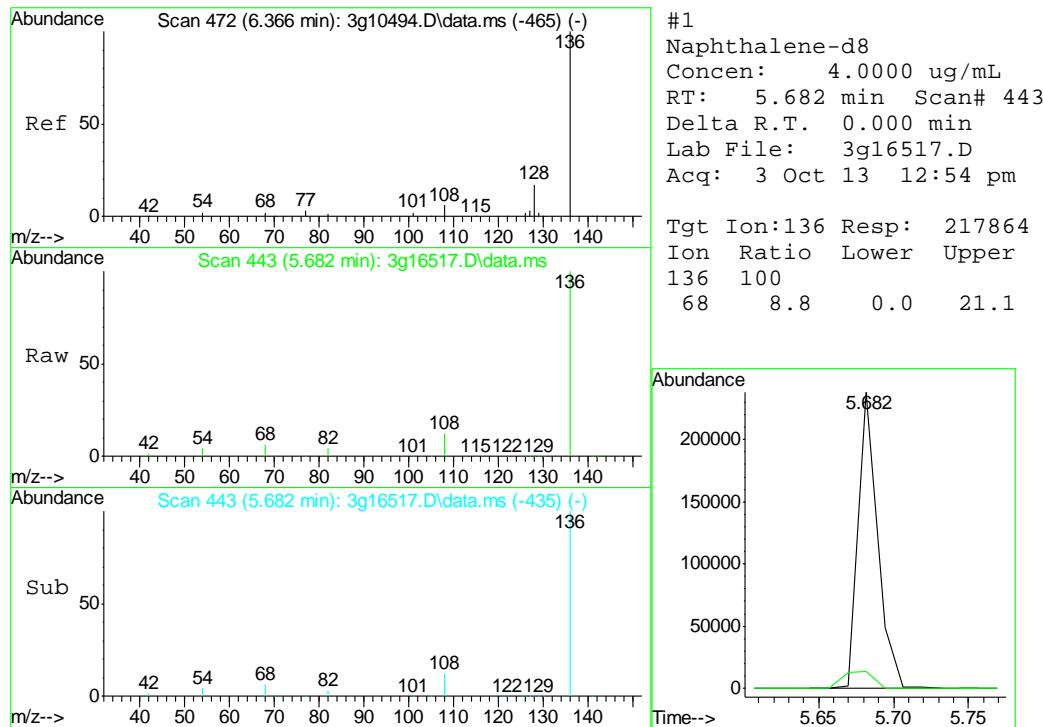
(#) = qualifier out of range (m) = manual integration (+) = signals summed

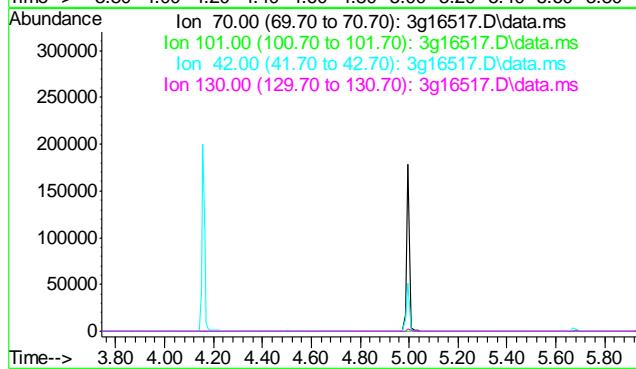
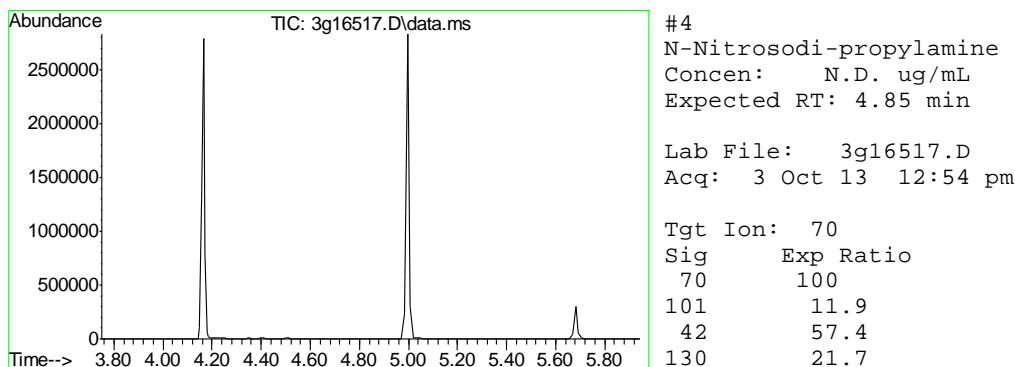
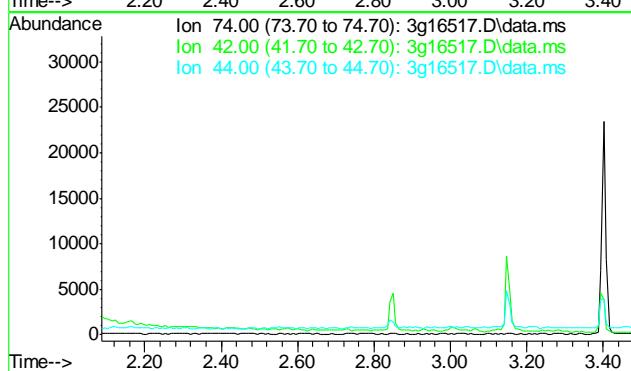
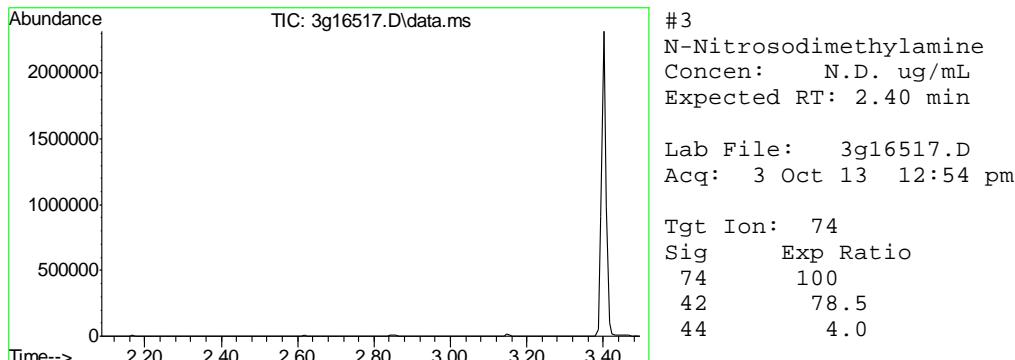
Quantitation Report (QT Reviewed)

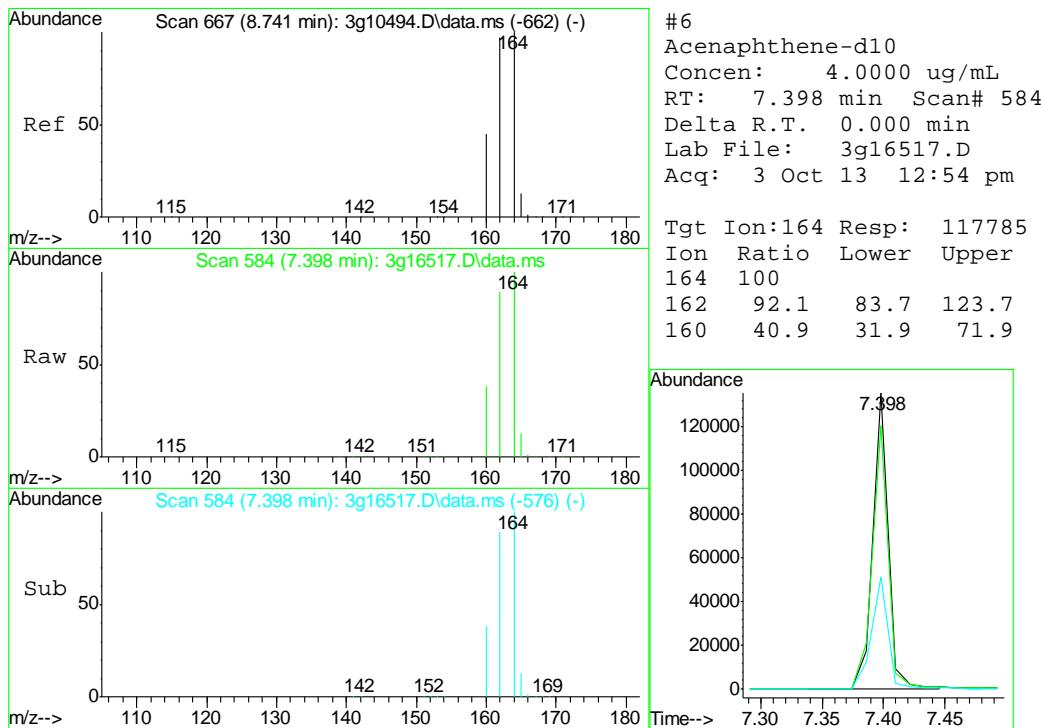
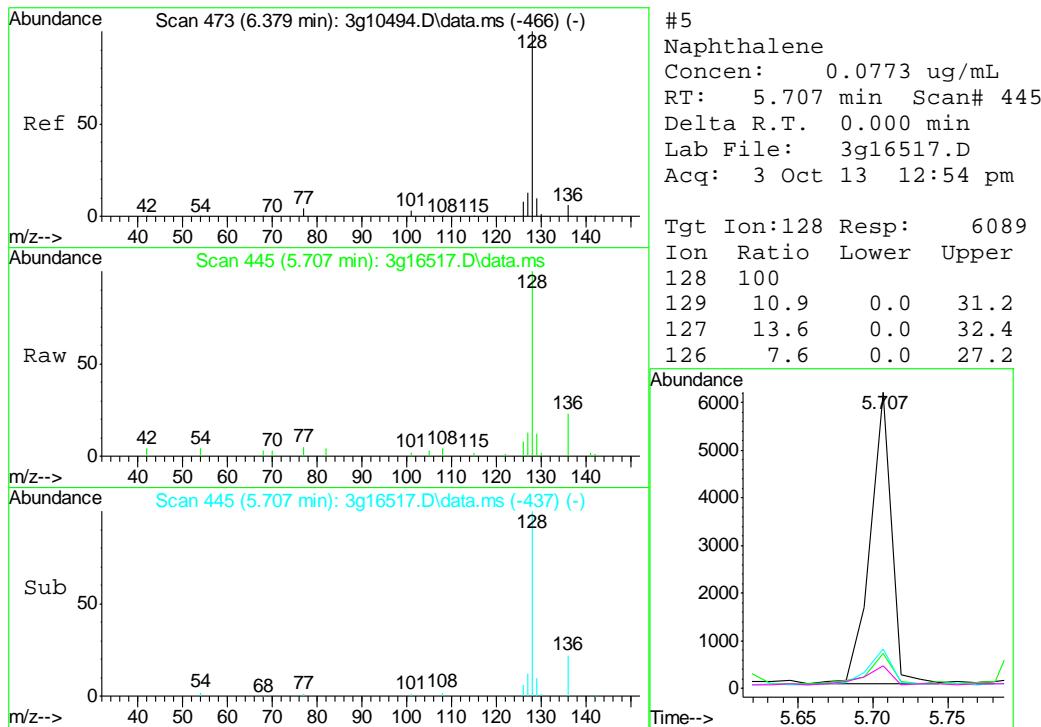
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 ALS Vial : 4 Sample Multiplier: 1

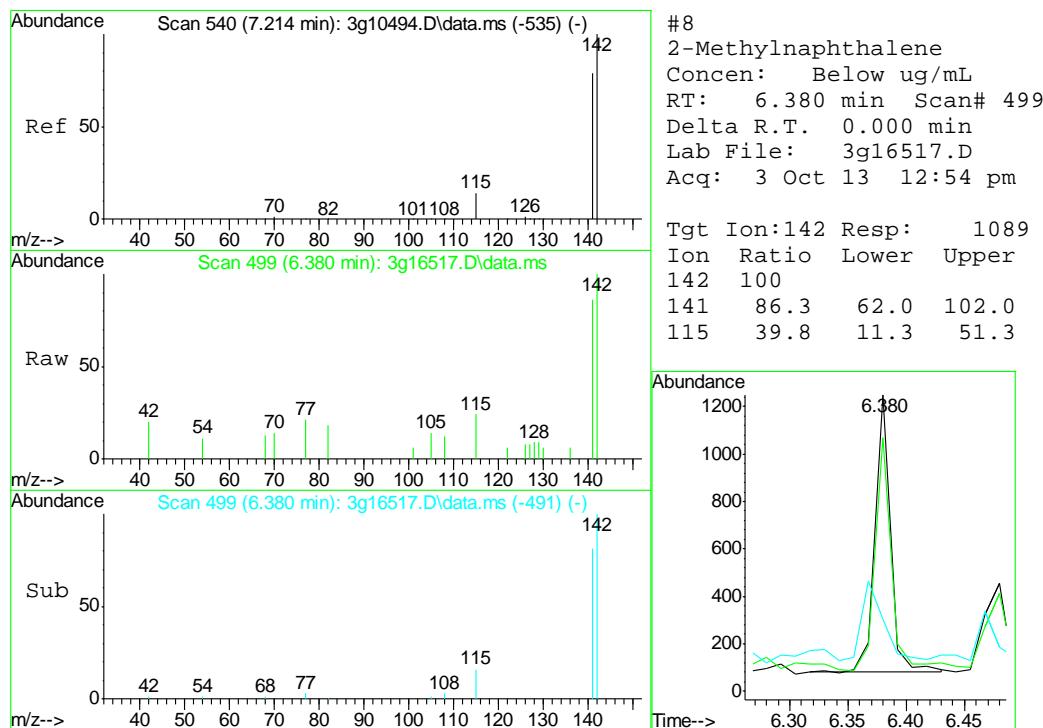
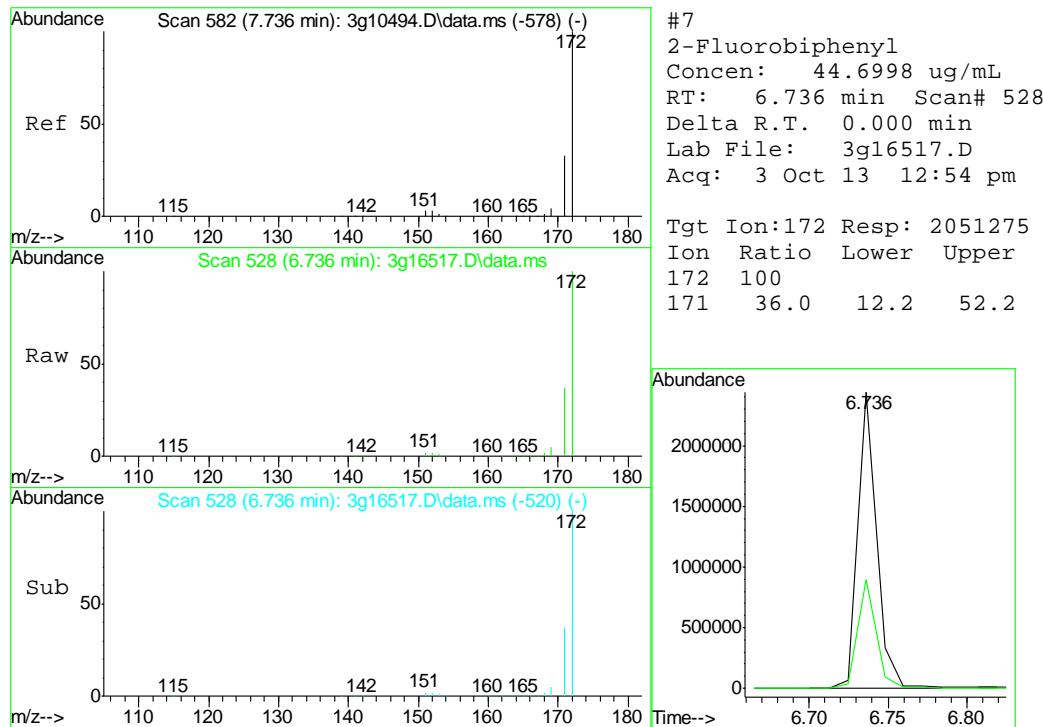
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 Response via : Initial Calibration

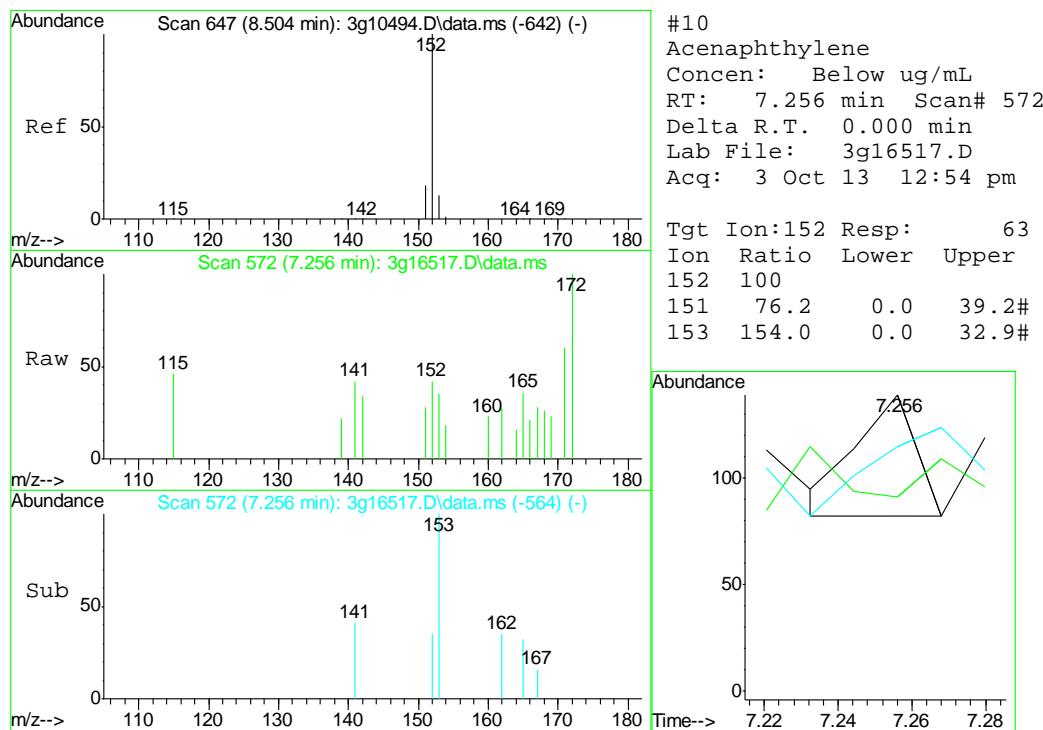
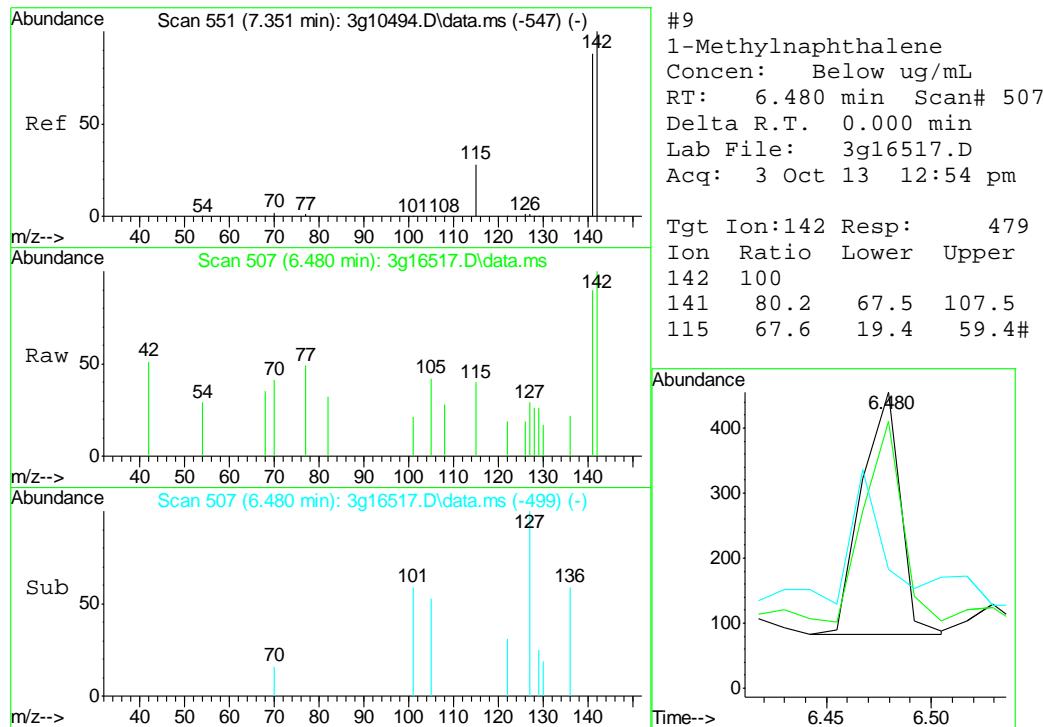


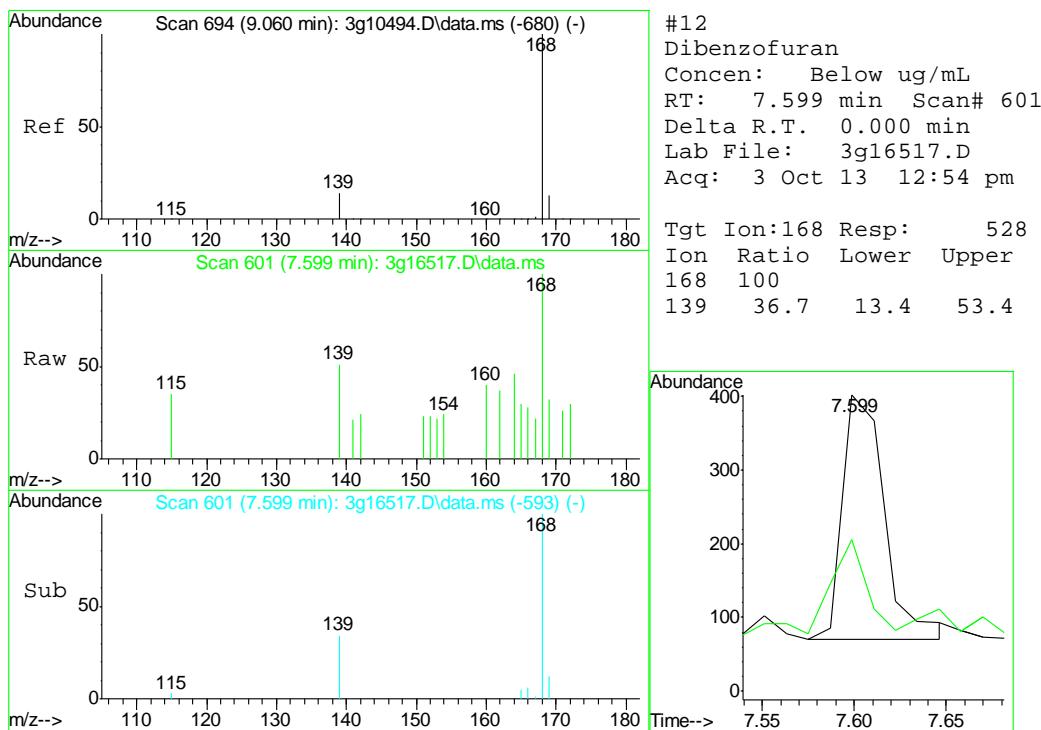
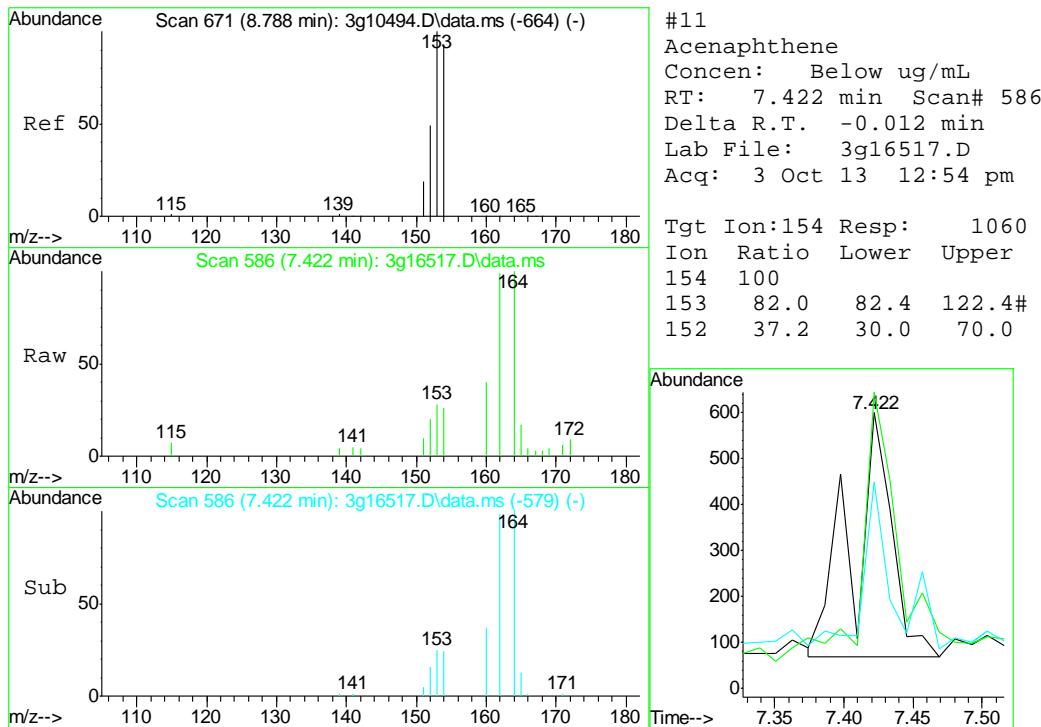


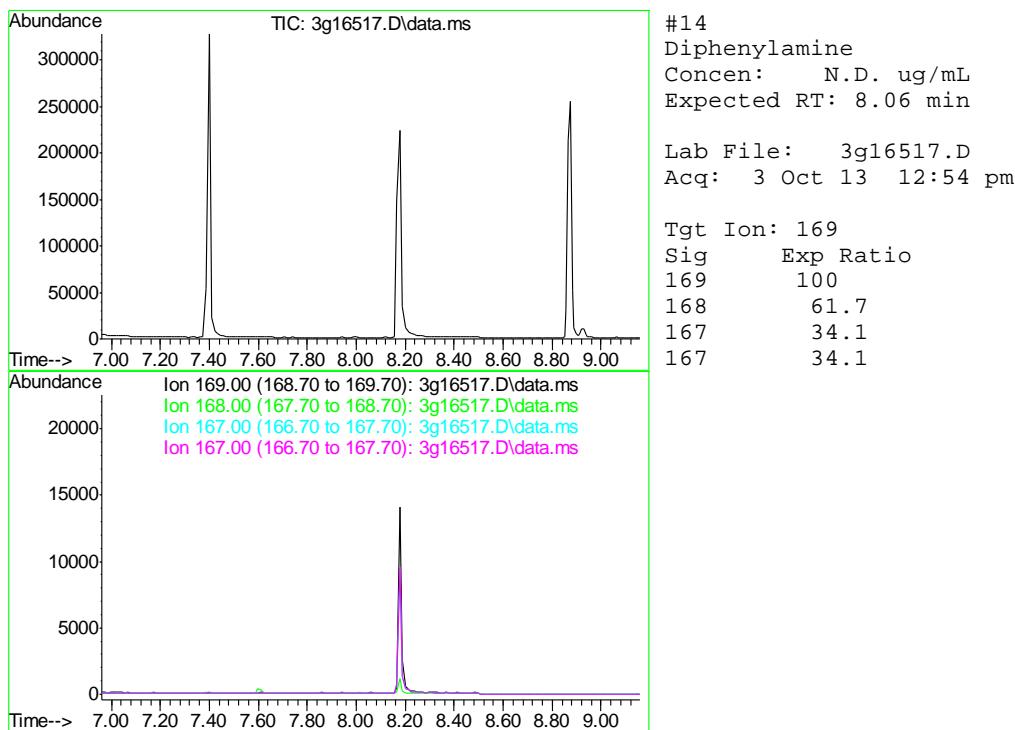
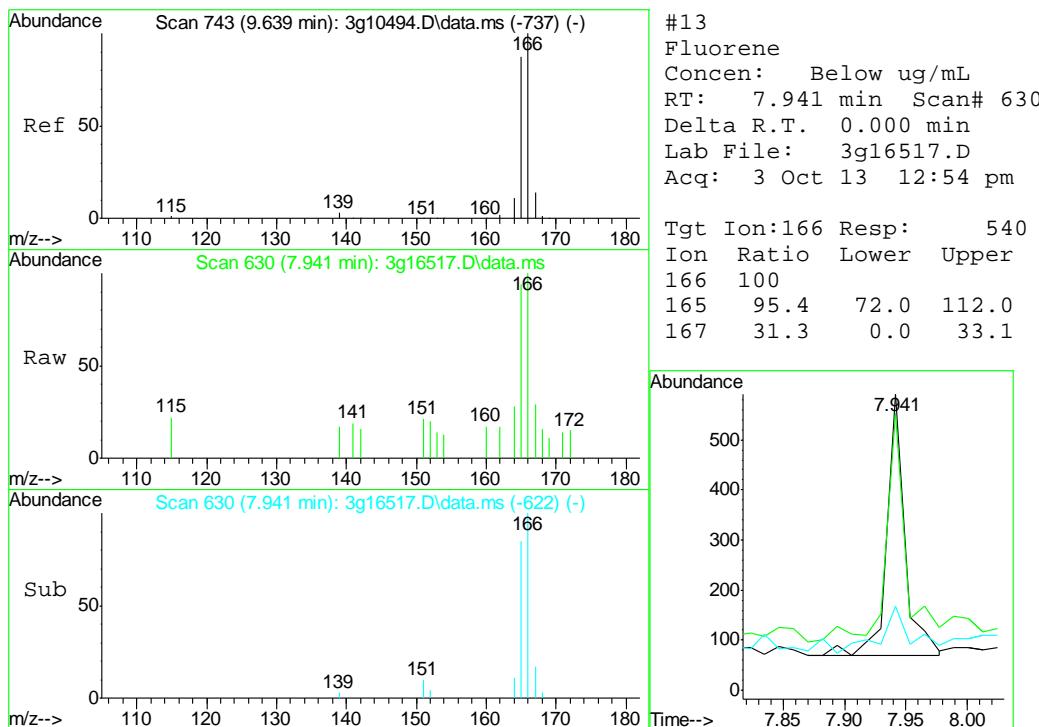


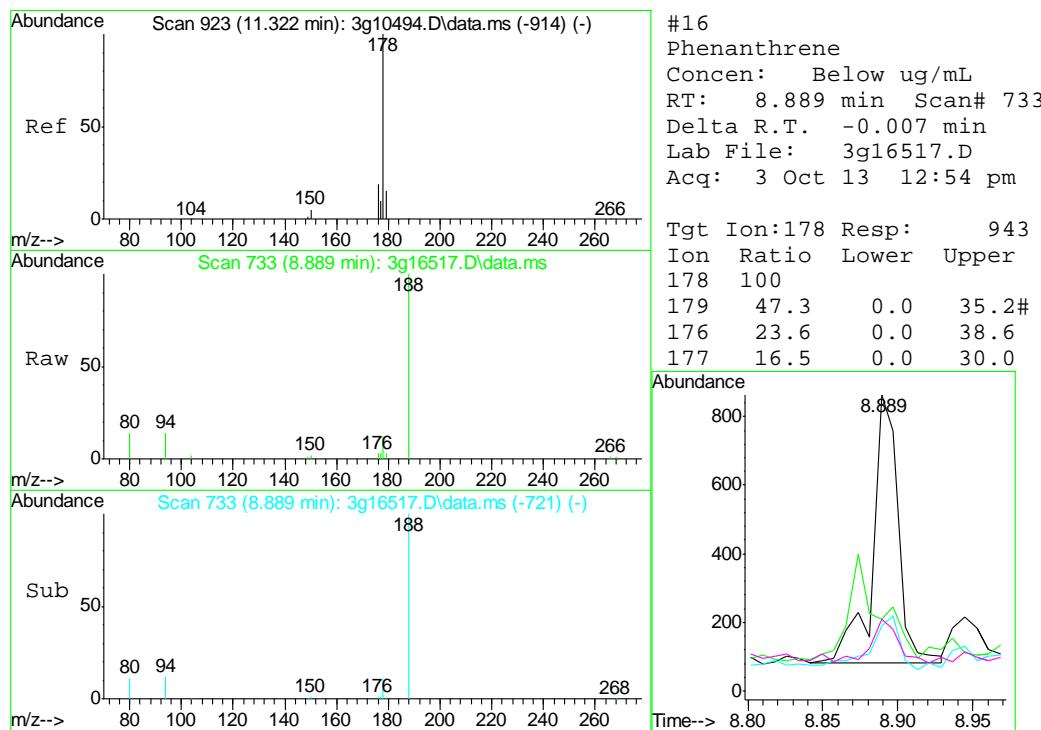
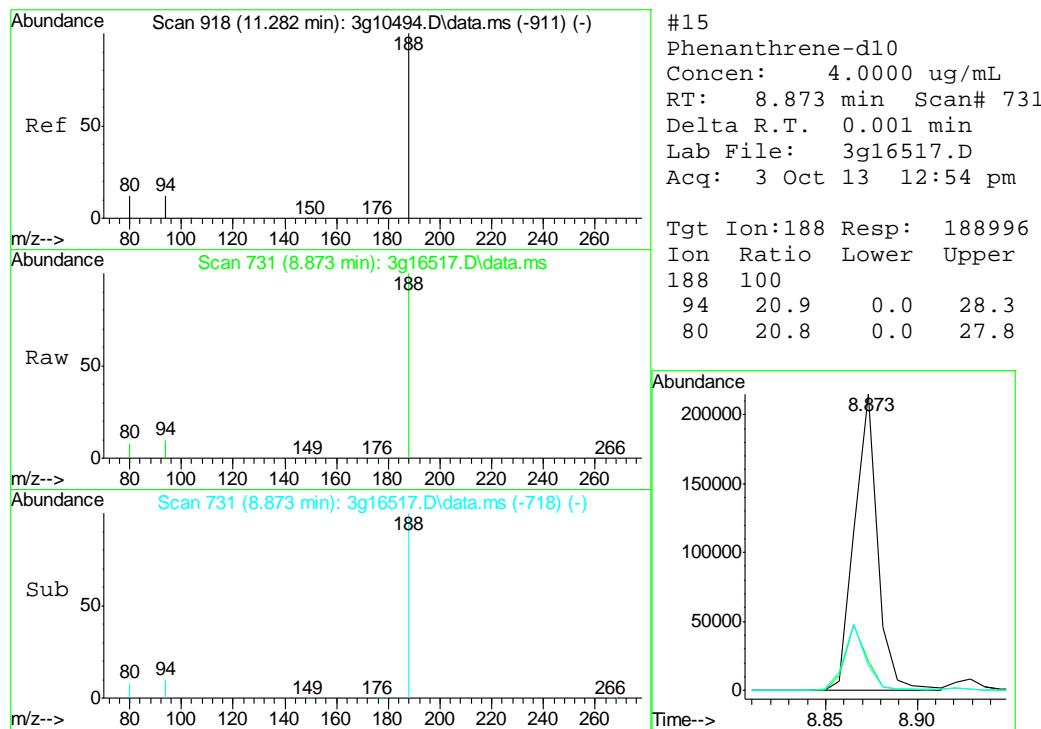


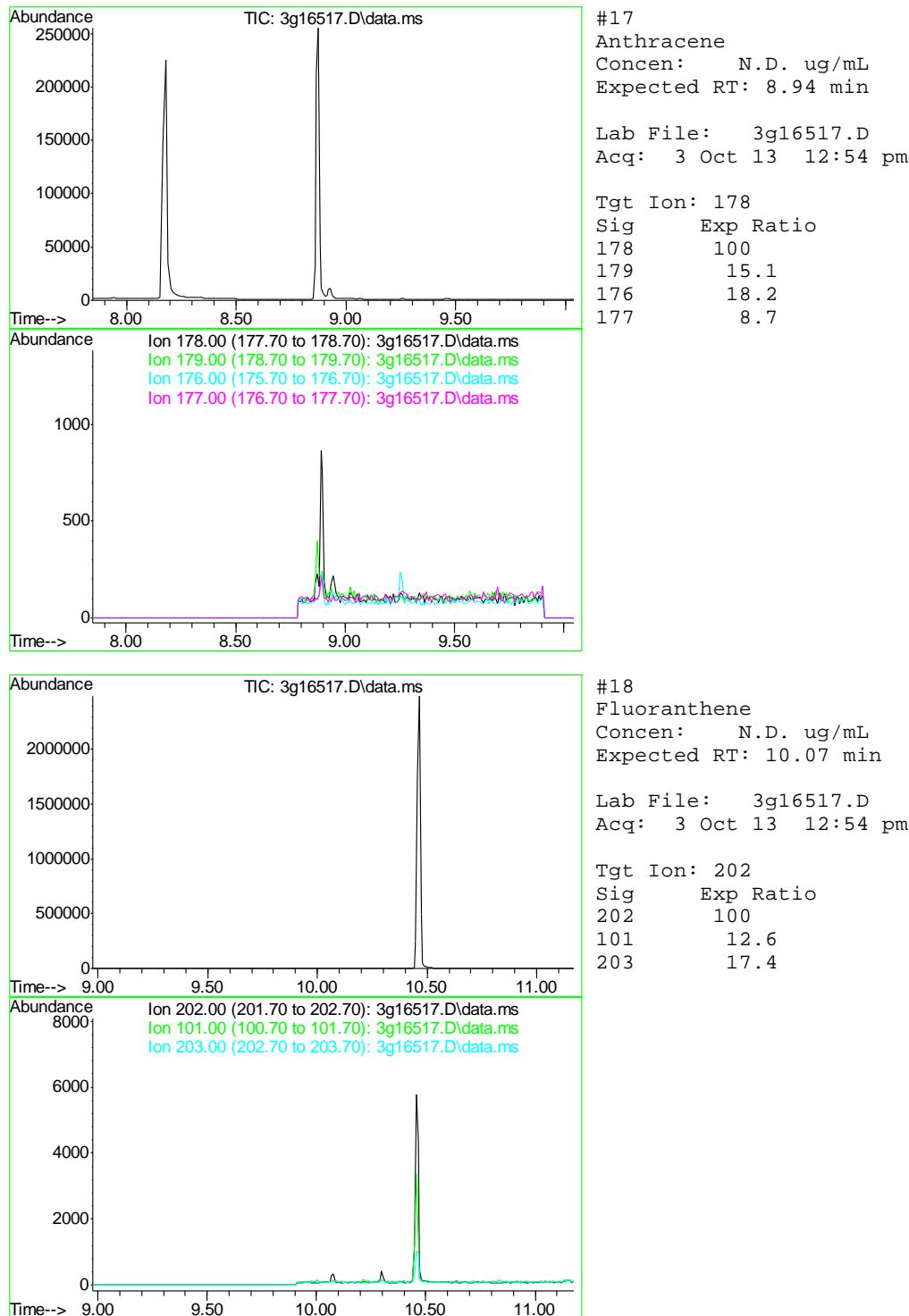


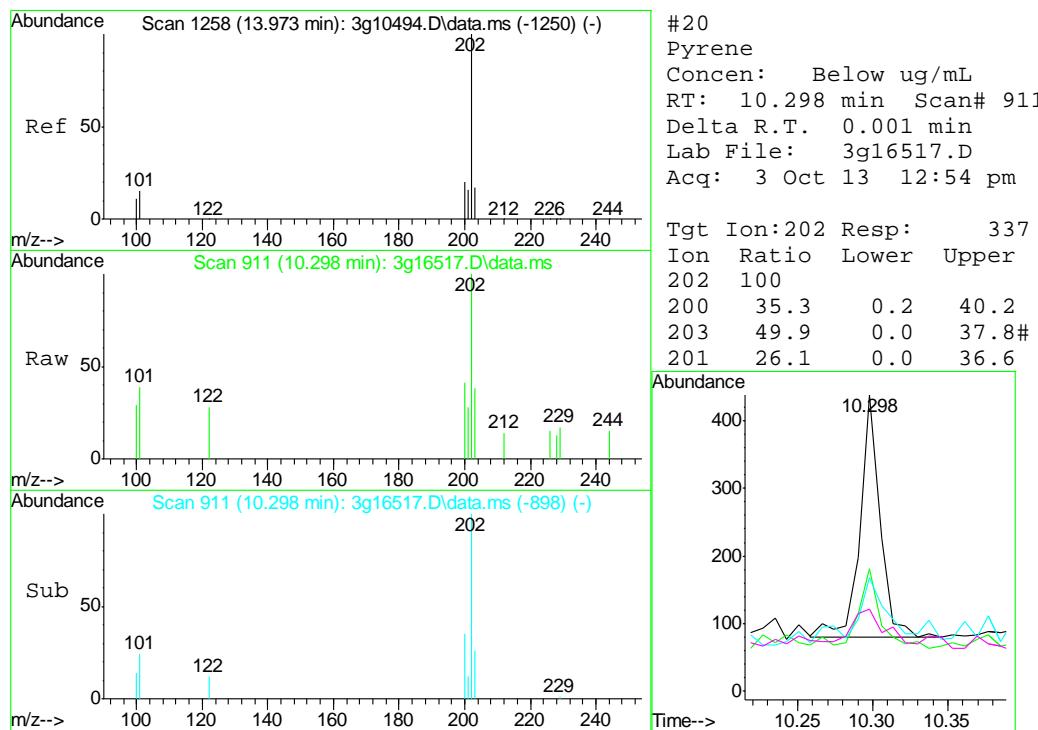
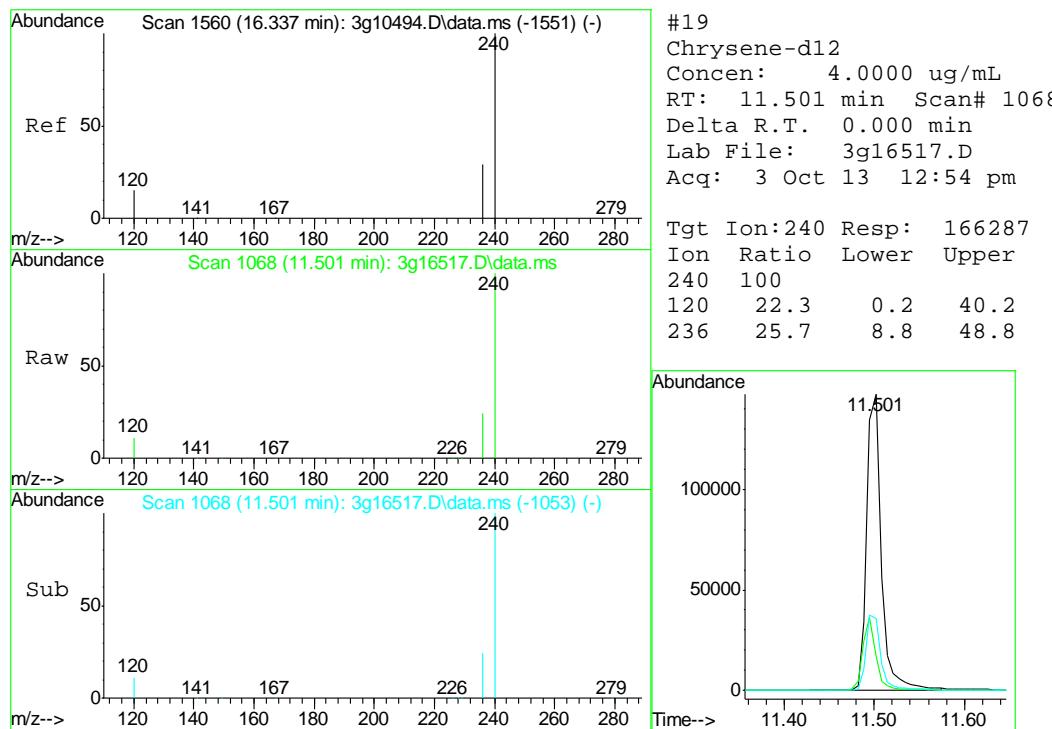


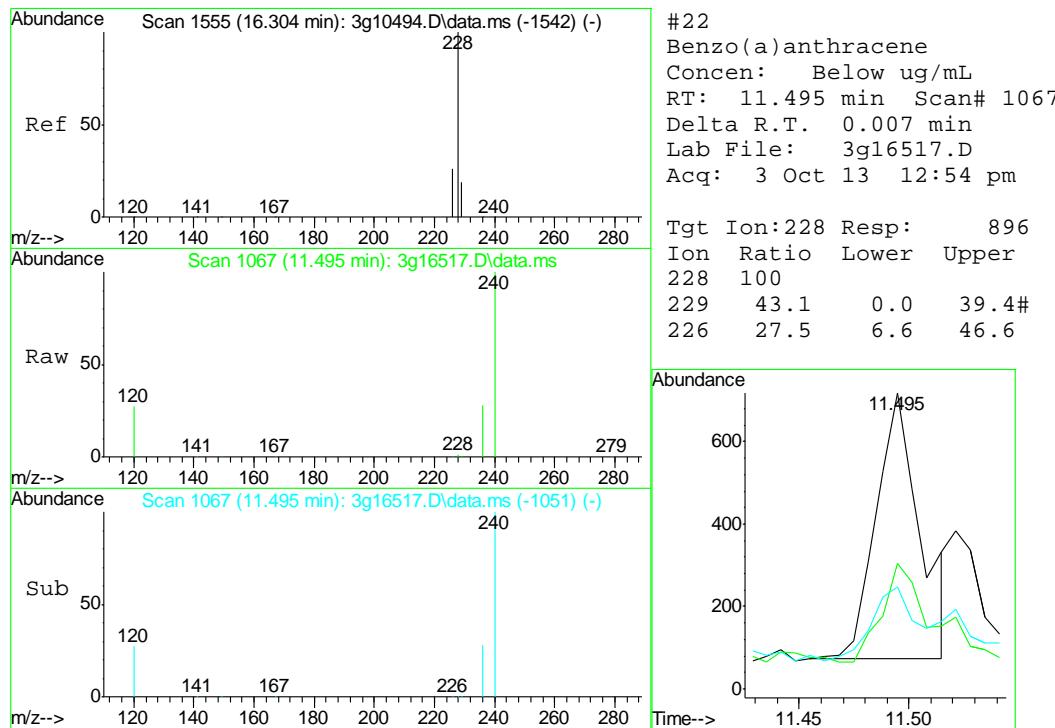
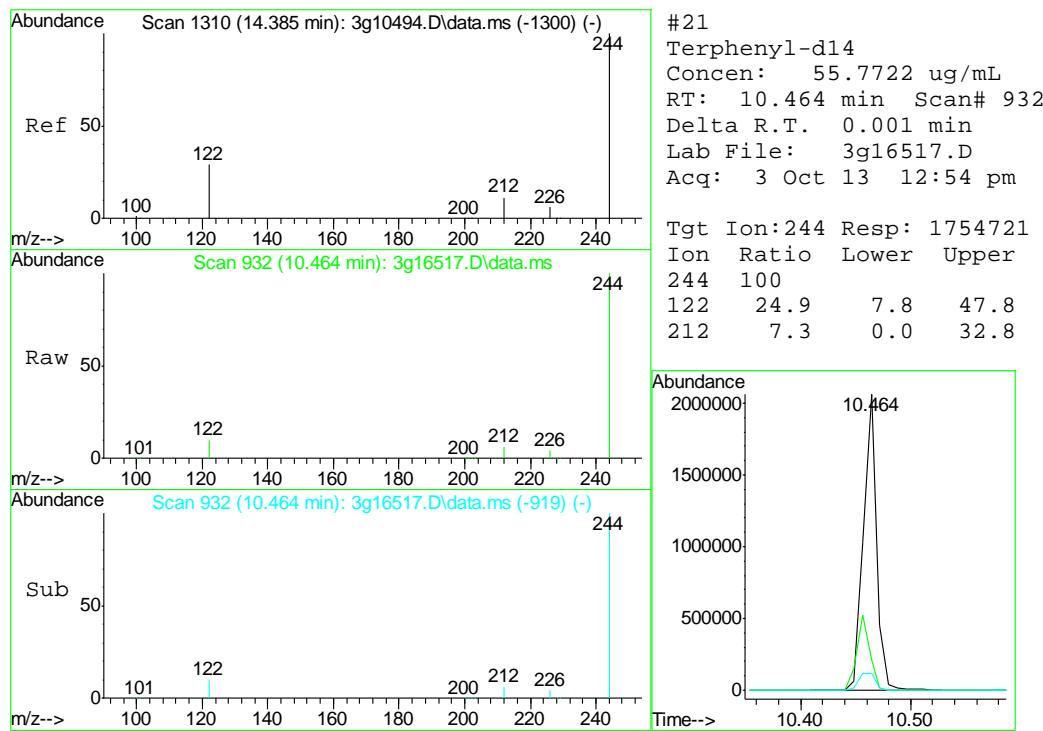


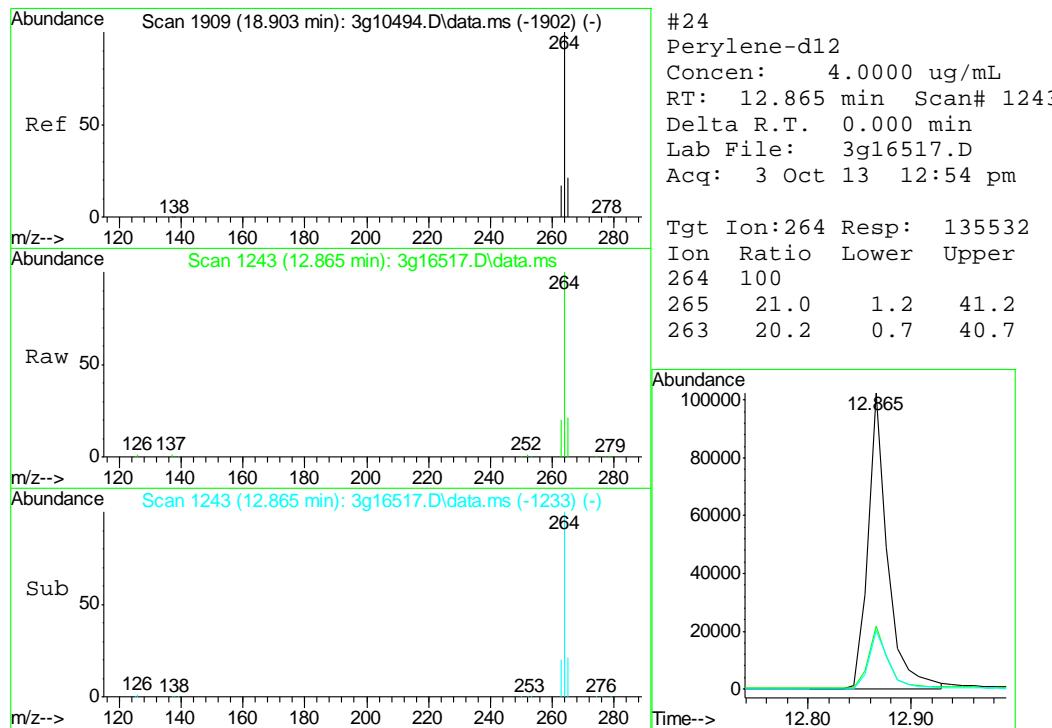
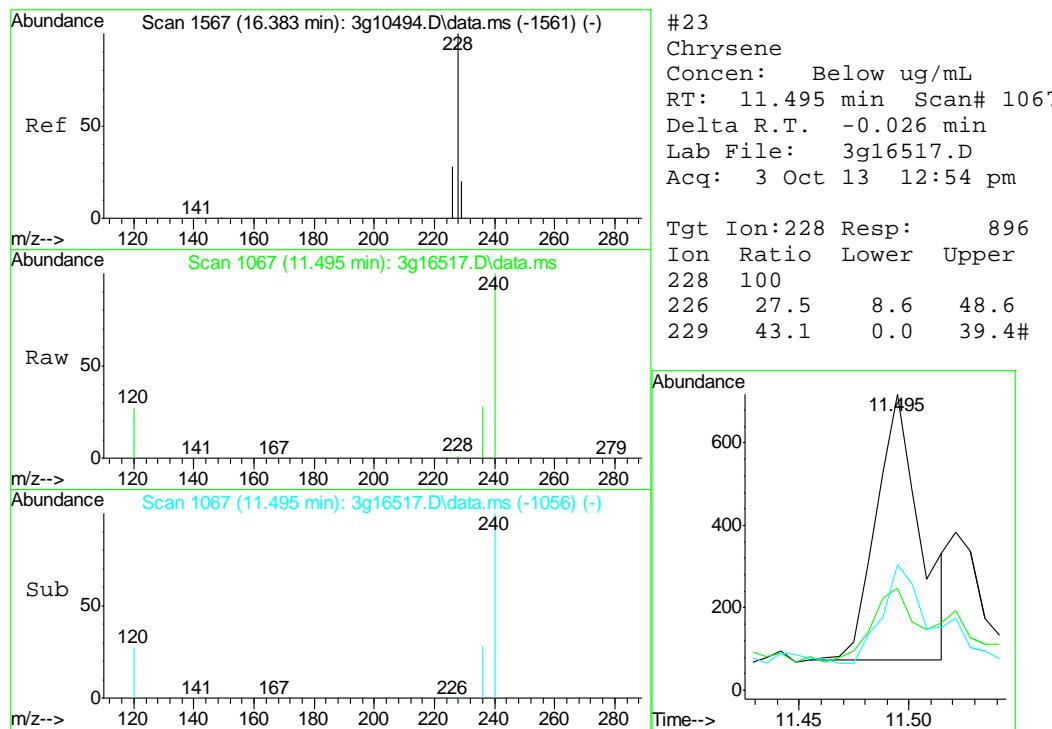


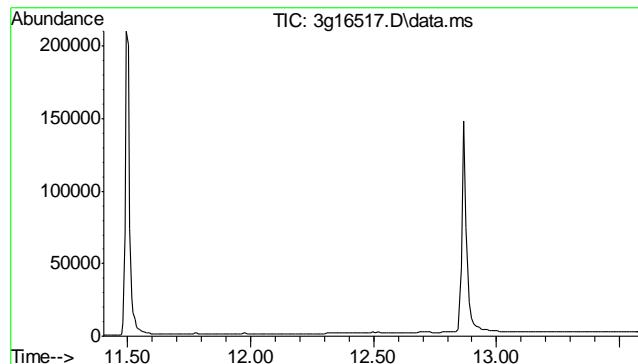








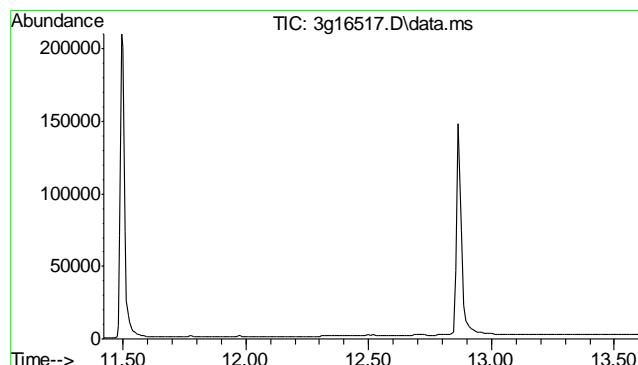
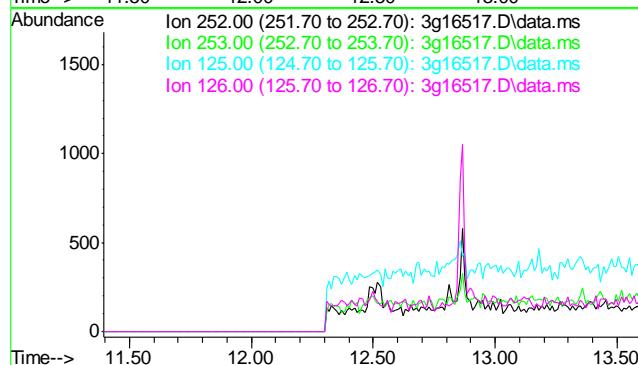




```
#25
Benzo(b)fluoranthene
Concen: N.D. ug/mL
Expected RT: 12.50 min

Lab File: 3g16517.D
Acq: 3 Oct 13 12:54 pm

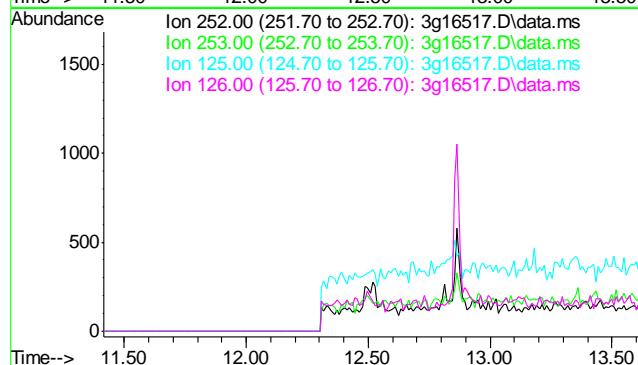
Tgt Ion: 252
Sig      Exp Ratio
252      100
253      51.5
125      13.2
126      46.9
```



```
#26
Benzo(k)fluoranthene
Concen: N.D. ug/mL
Expected RT: 12.52 min

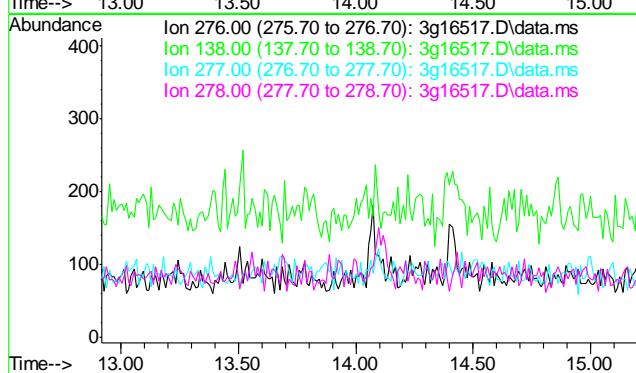
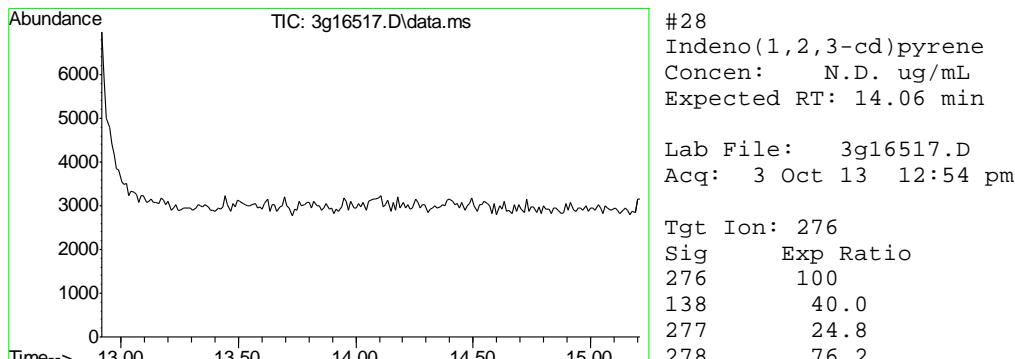
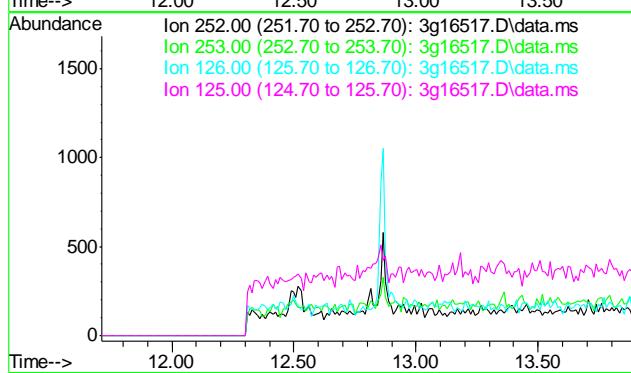
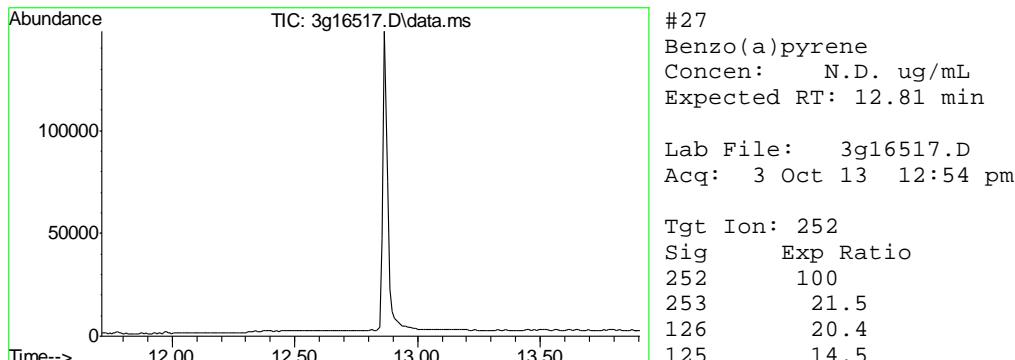
Lab File: 3g16517.D
Acq: 3 Oct 13 12:54 pm

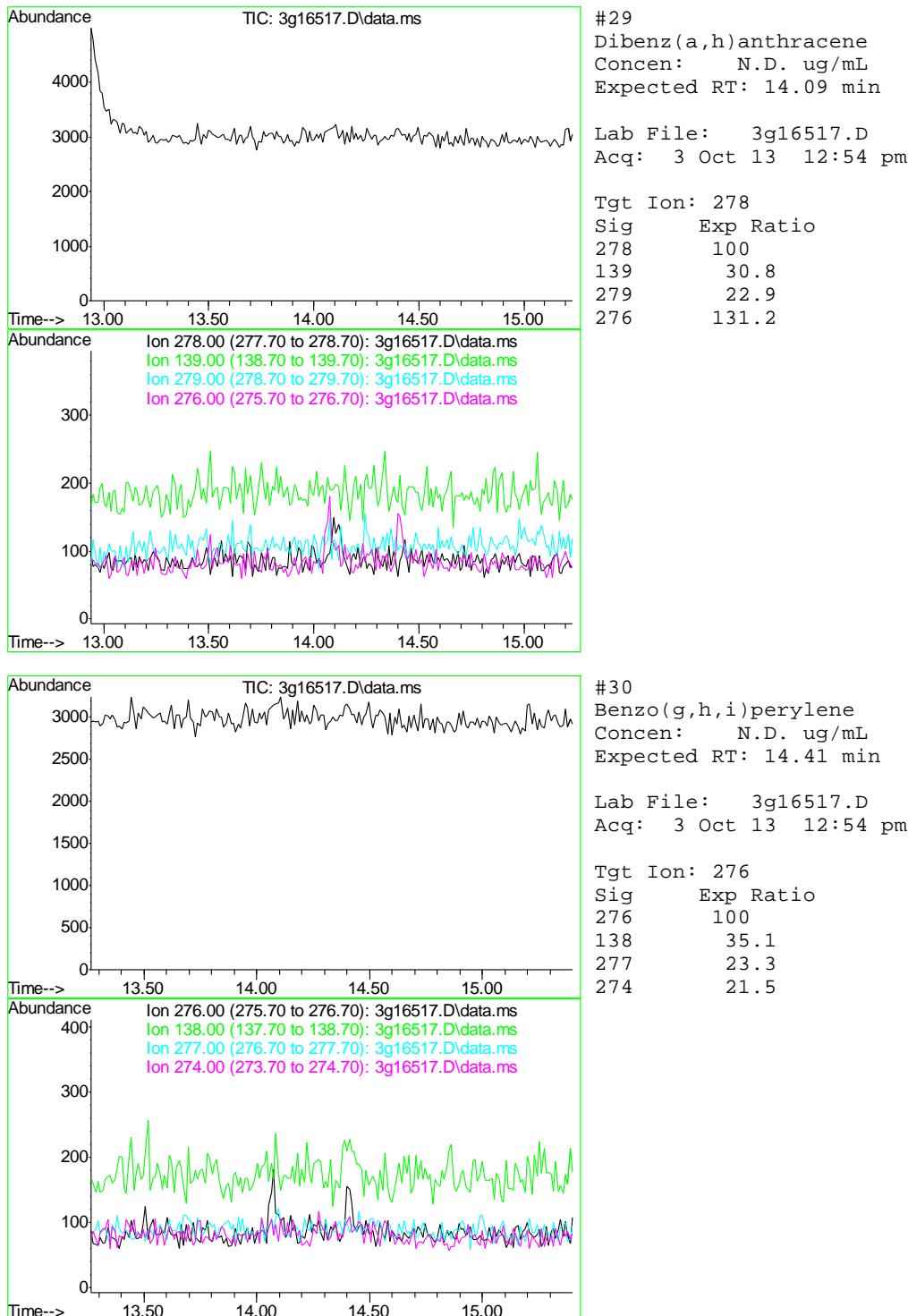
Tgt Ion: 252
Sig      Exp Ratio
252      100
253      37.3
125      9.6
126      34.1
```



9.2.1

9







GC Volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Job Number: D51123
Account: XTOKWR XTO Energy
Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1230-MB	GB22358.D	1	10/01/13	EV	n/a	n/a	GGB1230

The QC reported here applies to the following samples:

Method: SW846 8015B

D51123-1

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	10	5.0	mg/kg	

CAS No.	Surrogate Recoveries	Limits
120-82-1	1,2,4-Trichlorobenzene	85% 60-140%

10.1.1

10

Blank Spike Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1230-BS	GB22359.D	1	10/01/13	EV	n/a	n/a	GGB1230

The QC reported here applies to the following samples:

Method: SW846 8015B

D51123-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	112	102	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	94%	60-140%

10.2.1
10

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D51008-1MS	GB22361.D	1	10/01/13	EV	n/a	n/a	GGB1230
D51008-1MSD	GB22362.D	1	10/01/13	EV	n/a	n/a	GGB1230
D51008-1	GB22360.D	1	10/01/13	EV	n/a	n/a	GGB1230

The QC reported here applies to the following samples:

Method: SW846 8015B

D51123-1

CAS No.	Compound	D51008-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		138	141	102	141	102	0	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D51008-1	Limits
120-82-1	1,2,4-Trichlorobenzene	92%	94%	84%	60-140%

* = Outside of Control Limits.

10.3.1
10



GC Volatiles

Raw Data

Manual Integrations
APPROVED
(compounds with "m" flag)

Jennifer Laidlaw
10/02/13 14:07

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100113\GB22383.D\FID1A.CH Vial: 29
 Signal #2 : Y:\1\DATA\100113\GB22383.D\FID2B.CH
 Acq On : 2 Oct 2013 1:44 am Operator: ELISEV
 Sample : D51123-1 Inst : GC/MS Ins
 Misc : GC3911,GGB1230,5.035,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 02 08:44:19 2013 Quant Results File: TB1125GB1125SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Wed Oct 02 08:35:36 2013
 Response via : Initial Calibration
 DataAcq Meth : TVB4.M

Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.36	2295629	75.986 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	10893218	82.494 %	m

Target Compounds

1) H	TVH-Gasoline	7.31	26349645	0.376 mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	4.14	109459	0.285 ug/L m
6) T	Toluene	7.66	1110639	3.001 ug/L m
7) T	Ethylbenzene	10.28	1041407	3.341 ug/L m
8) T	m,p-Xylene	10.46	2397056	6.350 ug/L m
9) T	o-Xylene	10.96	110918	0.355 ug/L m
11) T	Naphthalene	14.55	72529	0.421 ug/L m

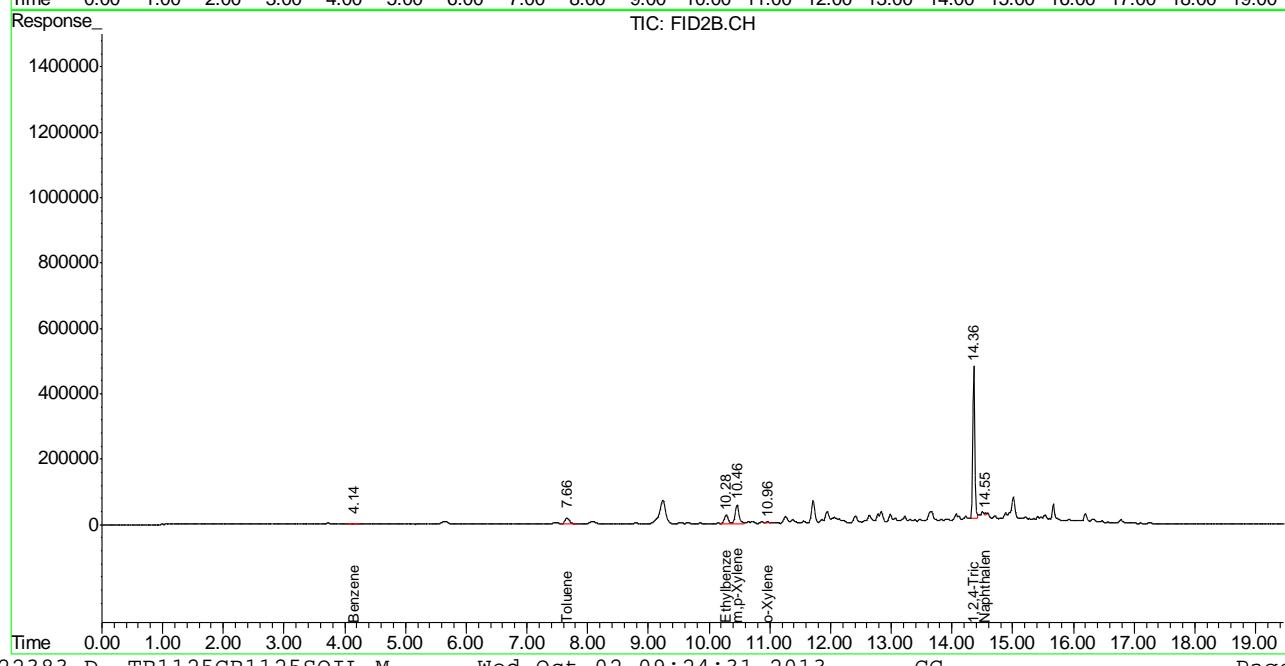
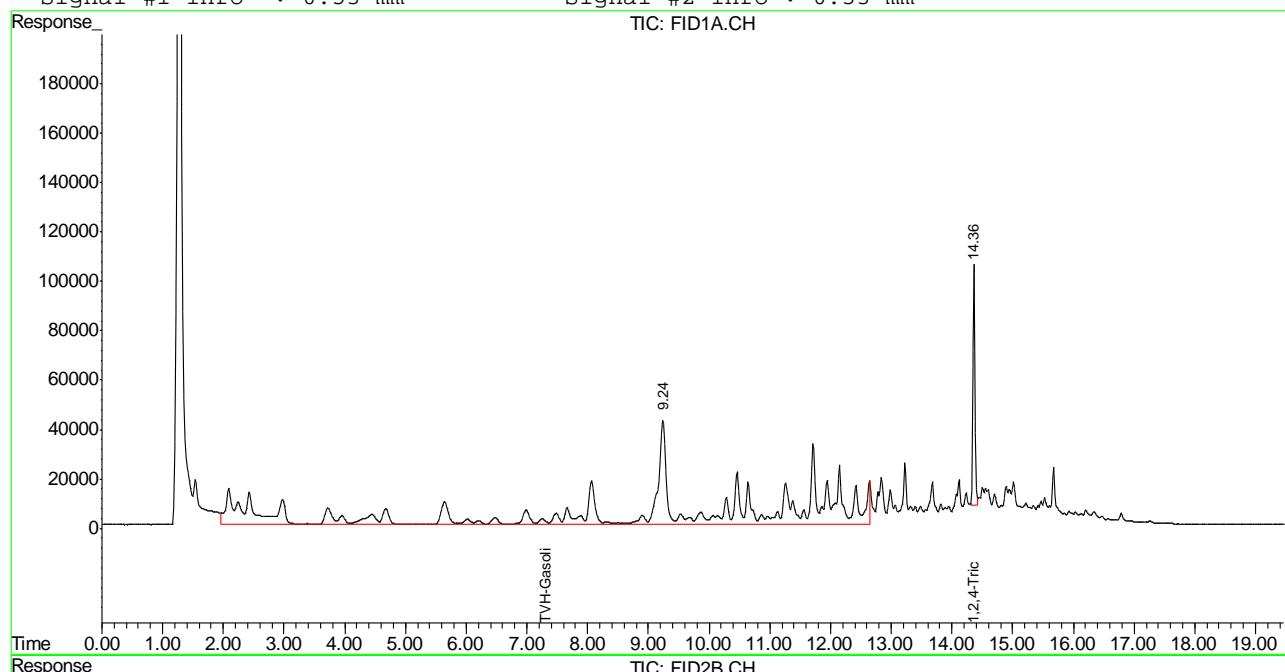
 (f)=RT Delta > 1/2 Window (m)=manual int.
 GB22383.D TB1125GB1125SOIL.M Wed Oct 02 09:24:31 2013 GC

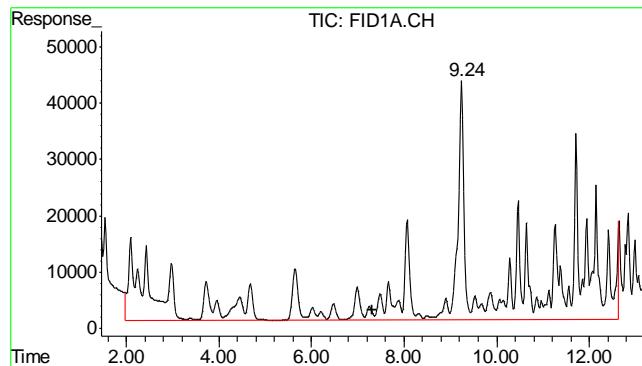
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100113\GB22383.D\FID1A.CH Vial: 29
 Signal #2 : Y:\1\DATA\100113\GB22383.D\FID2B.CH
 Acq On : 2 Oct 2013 1:44 am Operator: ELISEV
 Sample : D51123-1 Inst : GC/MS Ins
 Misc : GC3911,GGB1230,5.035,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 2 9:23 2013 Quant Results File: TB1125GB1125SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Wed Oct 02 08:35:36 2013
 Response via : Multiple Level Calibration
 DataAcq Meth : TVB4.M

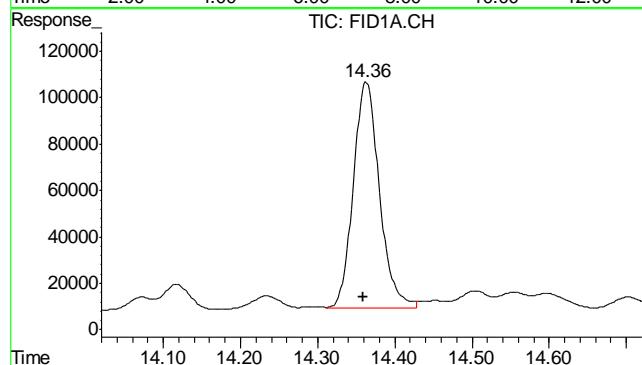
Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





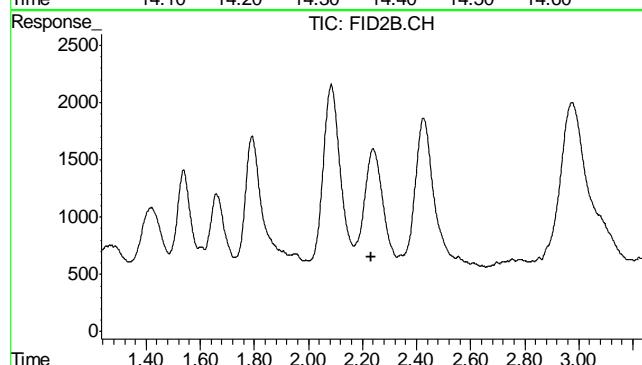
#1 TVH-Gasoline

R.T.: 7.310 min
Delta R.T.: 0.000 min
Response: 26349645
Conc: 0.38 mg/L m



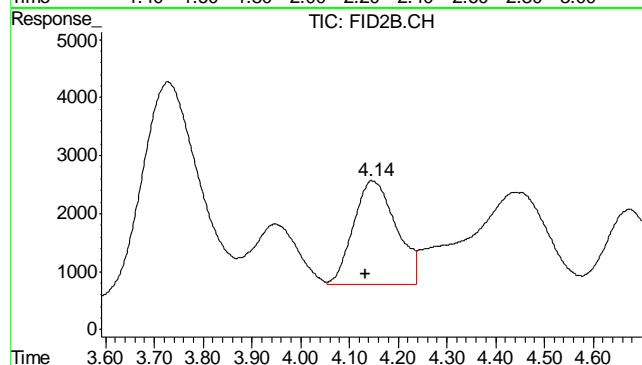
#2 1,2,4-Trichlorobenzene

R.T.: 14.362 min
Delta R.T.: 0.004 min
Response: 2295629
Conc: 75.99 % m



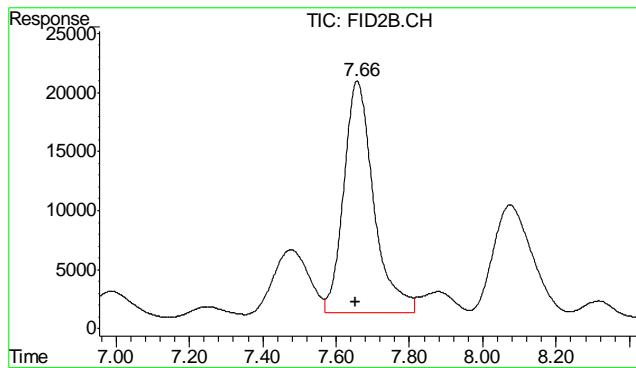
#4 Methyl-t-butyl-ether

R.T.: 0.000 min
Exp R.T.: 2.234 min
Response: 0
Conc: N.D.



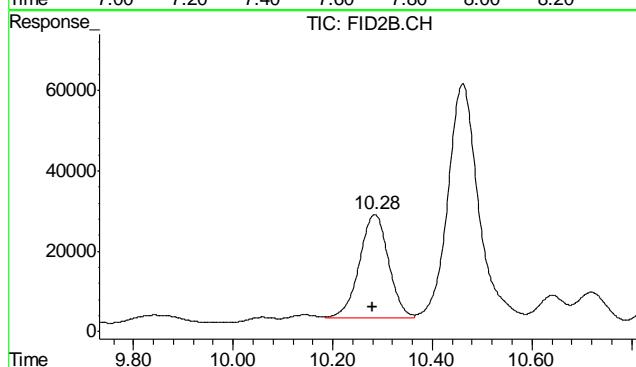
#5 Benzene

R.T.: 4.145 min
Delta R.T.: 0.011 min
Response: 109459
Conc: 0.29 ug/L m



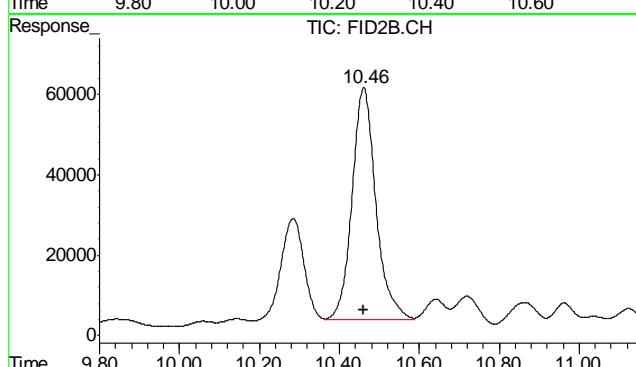
#6 Toluene

R.T.: 7.657 min
Delta R.T.: 0.004 min
Response: 1110639
Conc: 3.00 ug/L m



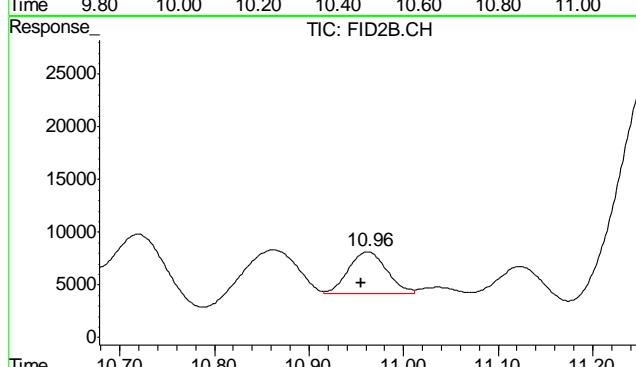
#7 Ethylbenzene

R.T.: 10.284 min
Delta R.T.: 0.005 min
Response: 1041407
Conc: 3.34 ug/L m



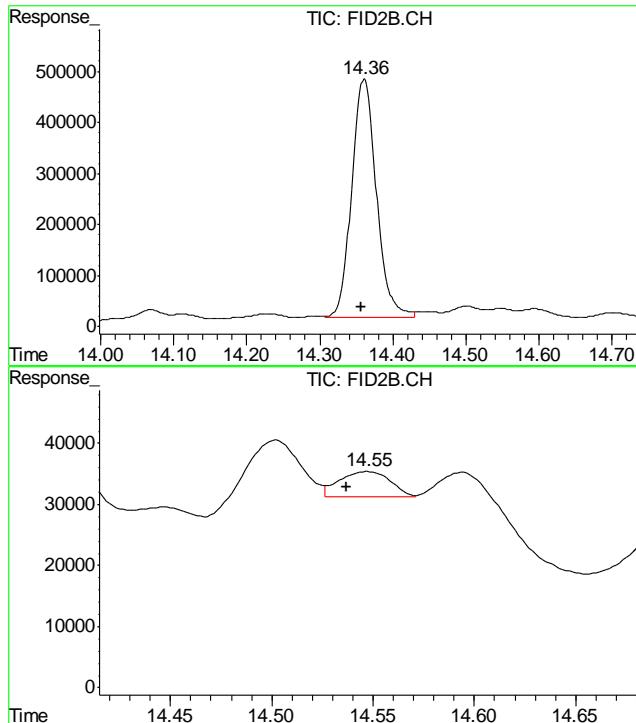
#8 m,p-Xylene

R.T.: 10.461 min
Delta R.T.: 0.002 min
Response: 2397056
Conc: 6.35 ug/L m



#9 o-Xylene

R.T.: 10.962 min
Delta R.T.: 0.006 min
Response: 110918
Conc: 0.36 ug/L m



#10 1,2,4-Trichlorobenzene (P)

R.T.: 14.360 min
 Delta R.T.: 0.004 min
 Response: 10893218
 Conc: 82.49 % m

#11 Naphthalene

R.T.: 14.547 min
 Delta R.T.: 0.010 min
 Response: 72529
 Conc: 0.42 ug/L m

11.1

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100113\GB22358.D\FID1A.CH Vial: 4
 Signal #2 : Y:\1\DATA\100113\GB22358.D\FID2B.CH
 Acq On : 1 Oct 2013 10:44 am Operator: ELISEV
 Sample : MB, S Inst : GC/MS Ins
 Misc : GC3911,GGB1230,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 02 08:42:47 2013 Quant Results File: TB1125GB1125SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Wed Oct 02 08:35:36 2013
 Response via : Initial Calibration
 DataAcq Meth : TVB4.M

Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm

Compound	R.T.	Response	Conc	Units
----------	------	----------	------	-------

System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.35	2552889	84.502 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.35	11821067	89.520 %	m

Target Compounds

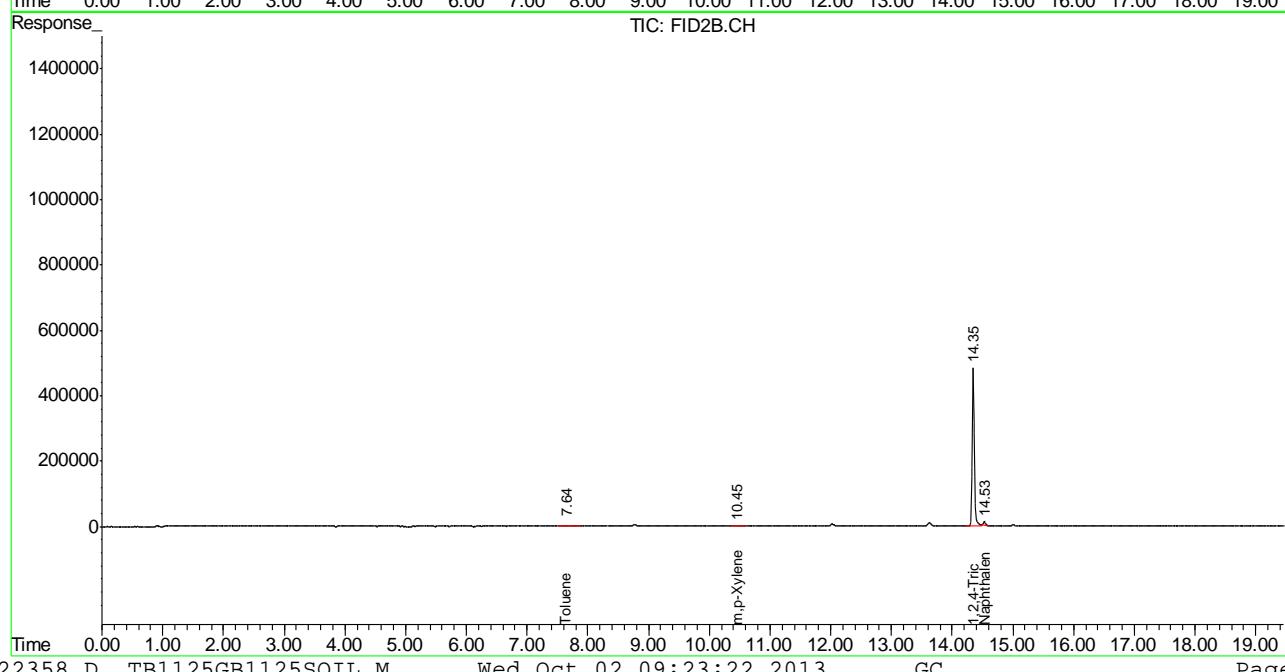
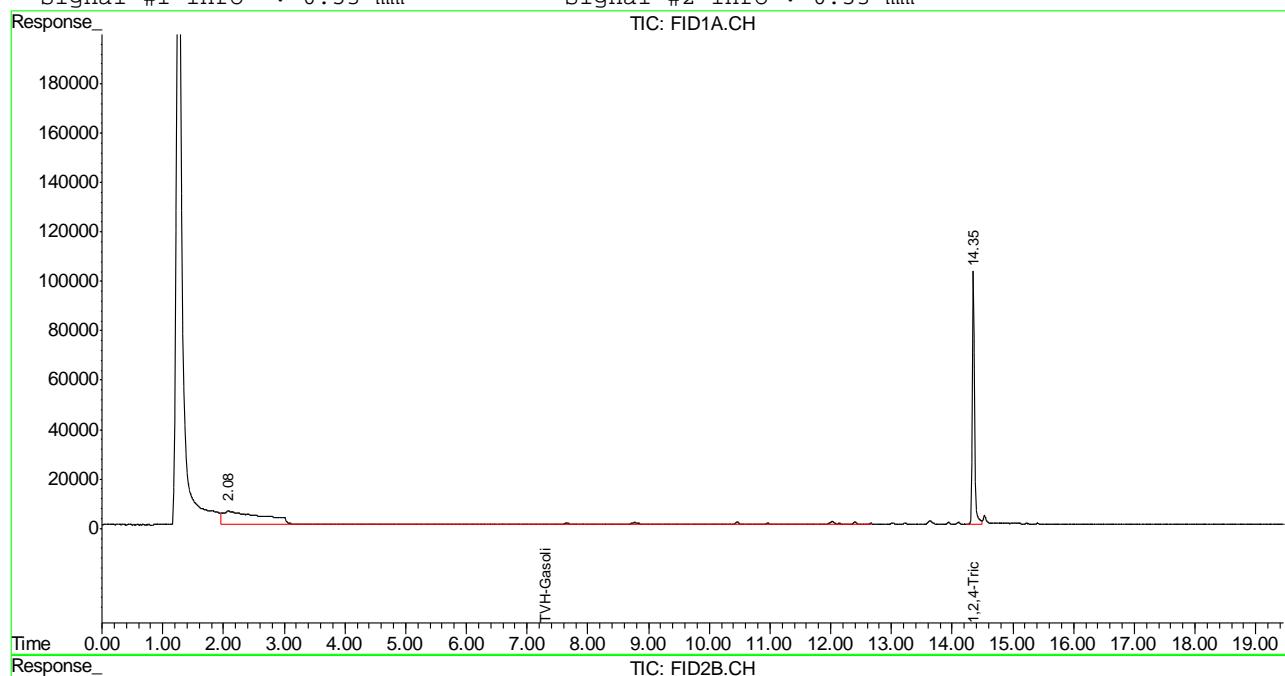
1) H	TVH-Gasoline	7.31	3539480	0.050 mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	0.00	0	N.D. ug/L d
6) T	Toluene	7.64	115021	0.311 ug/L m
7) T	Ethylbenzene	0.00	0	N.D. ug/L d
8) T	m,p-Xylene	10.46	136889	0.363 ug/L
9) T	o-Xylene	0.00	0	N.D. ug/L d
11) T	Naphthalene	14.53	259461	1.506 ug/L m

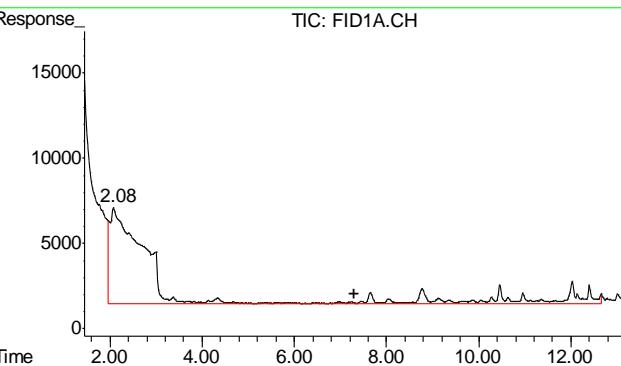
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100113\GB22358.D\FID1A.CH Vial: 4
 Signal #2 : Y:\1\DATA\100113\GB22358.D\FID2B.CH
 Acq On : 1 Oct 2013 10:44 am Operator: ELISEV
 Sample : MB, S Inst : GC/MS Ins
 Misc : GC3911, GGB1230, 5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 2 8:54 2013 Quant Results File: TB1125GB1125SOIL.RES

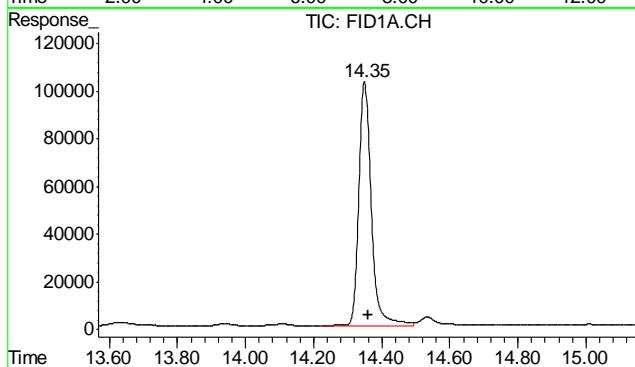
Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Wed Oct 02 08:35:36 2013
 Response via : Multiple Level Calibration
 DataAcq Meth : TVB4.M

Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm

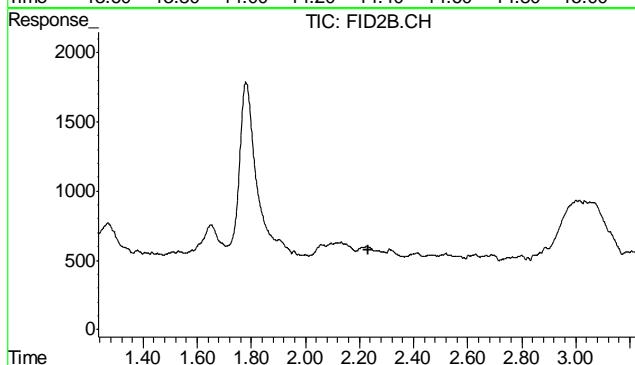




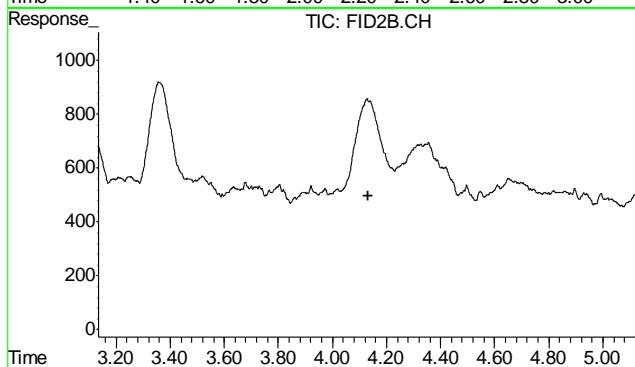
#1 TVH-Gasoline
R.T.: 7.310 min
Delta R.T.: 0.000 min
Response: 3539480
Conc: 0.05 mg/L m



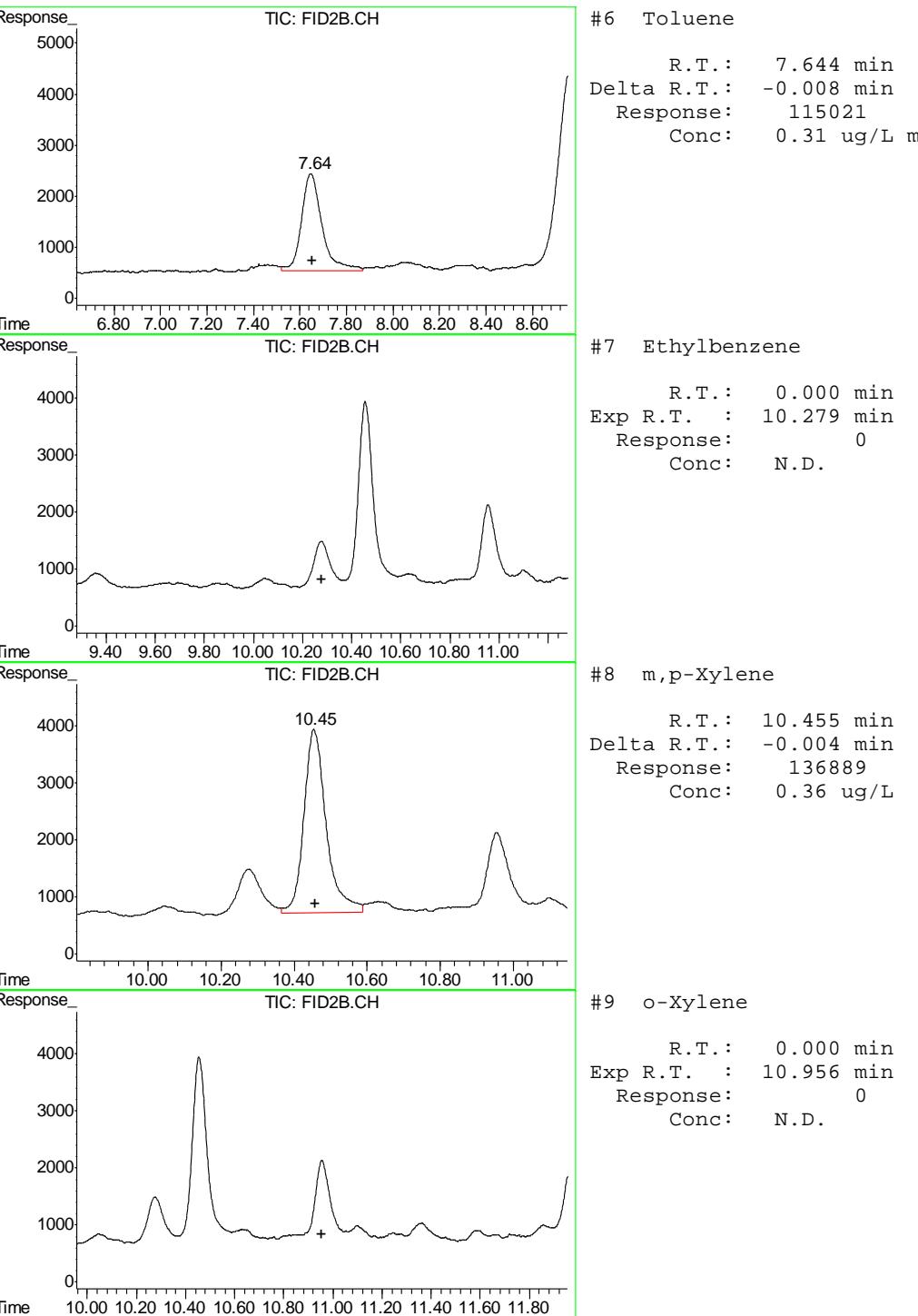
#2 1,2,4-Trichlorobenzene
R.T.: 14.349 min
Delta R.T.: -0.009 min
Response: 2552889
Conc: 84.50 % m



#4 Methyl-t-butyl-ether
R.T.: 0.000 min
Exp R.T. : 2.234 min
Response: 0
Conc: N.D.

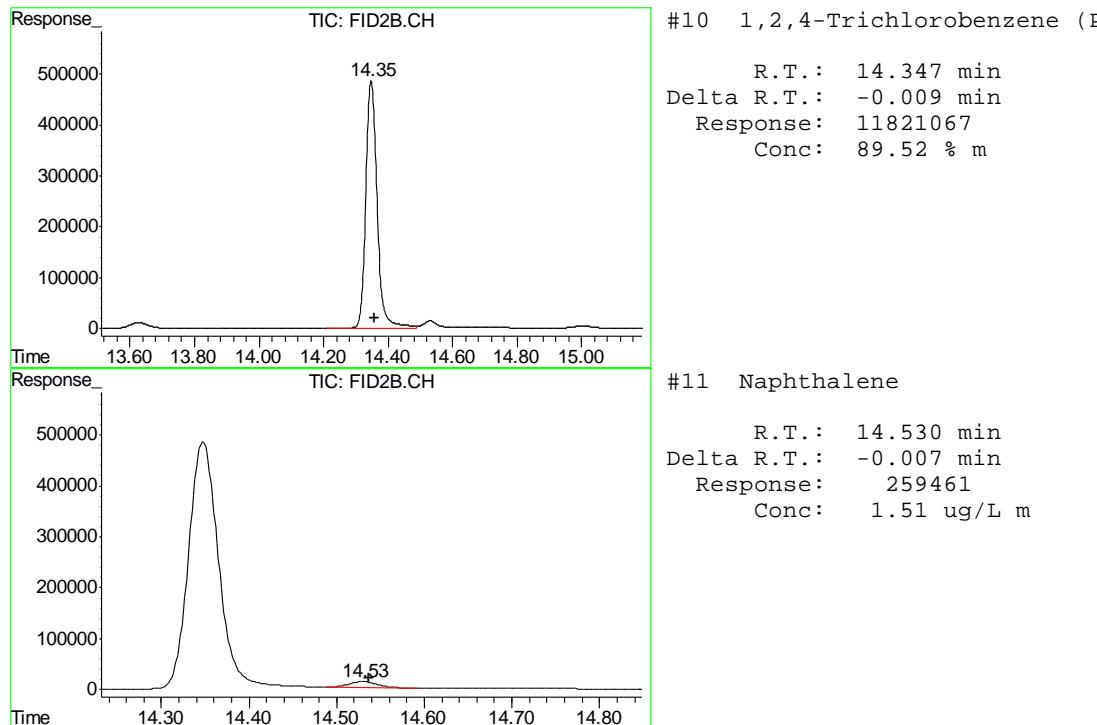


#5 Benzene
R.T.: 0.000 min
Exp R.T. : 4.134 min
Response: 0
Conc: N.D.



11.2.1

11



11.2.1

11



GC Semi-volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8666-MB	FH013648.D	1	10/03/13	TU	10/02/13	OP8666	GFH720

The QC reported here applies to the following samples:

Method: SW846-8015B

D51123-1

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	ND	6.7	5.0	mg/kg	

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	71% 20-130%

12.1.1

12

Blank Spike Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8666-BS	FH013650.D	1	10/03/13	TU	10/02/13	OP8666	GFH720

The QC reported here applies to the following samples:

Method: SW846-8015B

D51123-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	422	63	42-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	64%	20-130%

* = Outside of Control Limits.

12.2.1

12

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51123

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8666-MS	FH013672.D	10	10/03/13	TU	10/02/13	OP8666	GFH722
OP8666-MSD	FH013673.D	10	10/03/13	TU	10/02/13	OP8666	GFH722
D51127-1	FH013674.D	10	10/03/13	TU	10/02/13	OP8666	GFH722

The QC reported here applies to the following samples:

Method: SW846-8015B

D51123-1

CAS No.	Compound	D51127-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	10500		727	11100	83	11800	179* a	6	20-150/30

CAS No.	Surrogate Recoveries	MS	MSD	D51127-1	Limits
84-15-1	o-Terphenyl	82%	92%	74%	20-130%

(a) Outside control limits due to high level in sample relative to spike amount.

* = Outside of Control Limits.



GC Semi-volatiles

Raw Data

Quantitation Report (QT Reviewed)

Data File : C:\FID6_DATA\FI100313.SEC\FI09508.D Vial: 64
 Acq On : 3 Oct 2013 9:02 pm Operator: TIMU
 Sample : D51123-1 Inst : Fid6
 Misc : OP8666,GFI637,30.10,,,1,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Oct 04 08:32:59 2013 Quant Results File: ORO-RR-GFI585.RES

Quant Method : C:\MSDCHEM\1...\ORO-RR-GFI585.M (Chemstation Integrator)
 Title : 8015B TEH Front detector
 Last Update : Thu Oct 03 11:49:33 2013
 Response via : Initial Calibration
 DataAcq Meth : DUAL_B2.M

Volume Inj. : 1ul
 Signal Phase : RTX-5
 Signal Info : 530um

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	13.56	120979241	1669.086 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (C10-C28)	11.71	1296321928	20227.288 mg/L
3) H TPH-ORO (>C28-C40)	23.52	753218054	18053.122 mg/L

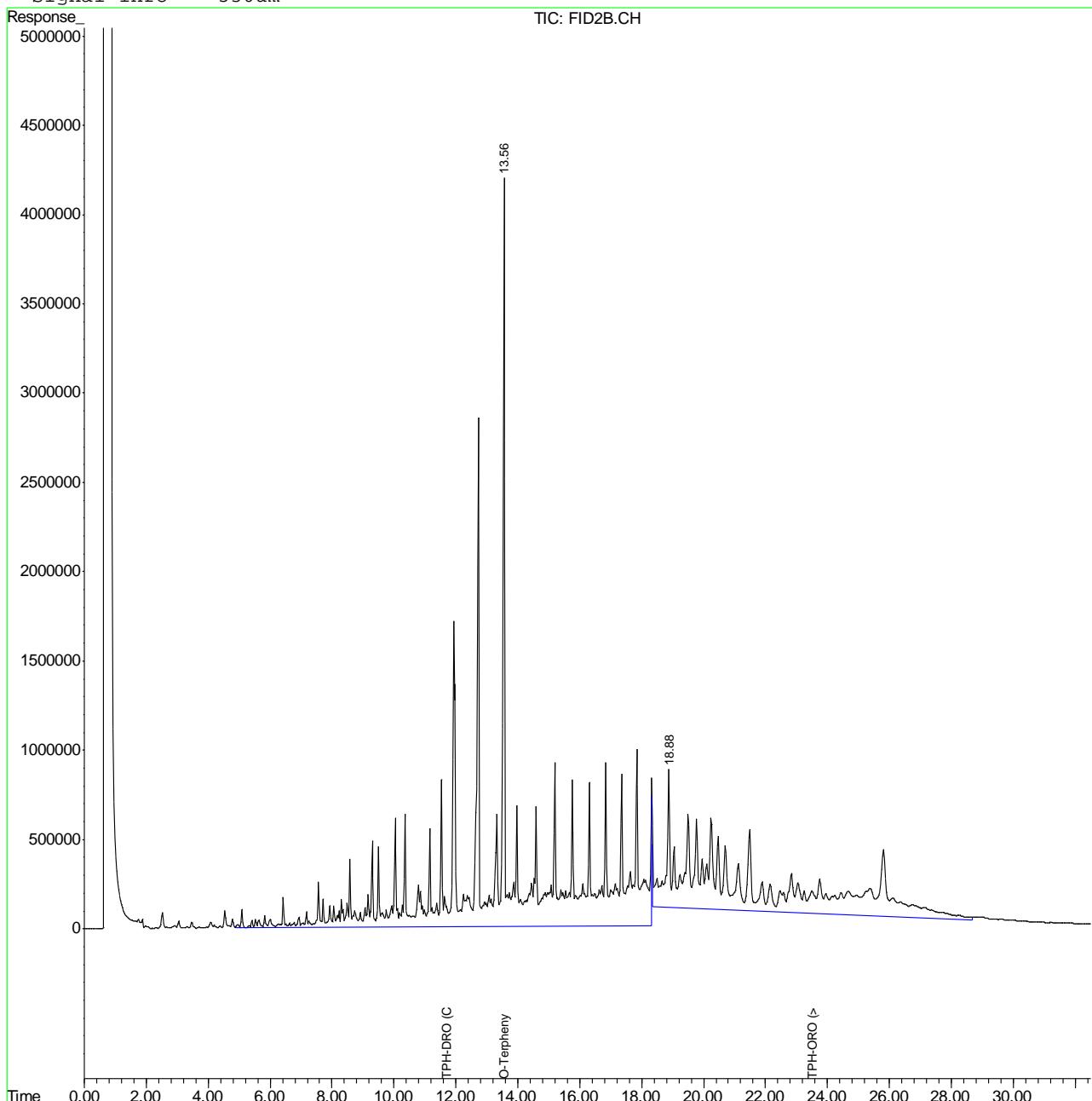
(f)=RT Delta > 1/2 Window (m)=manual int.
 FI09508.D ORO-RR-GFI585.M Fri Oct 04 08:40:15 2013 TEH

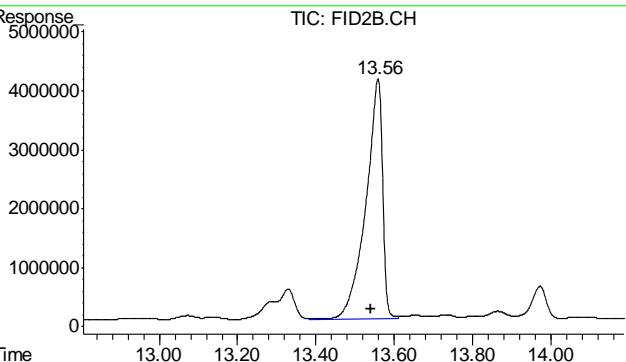
Quantitation Report (QT Reviewed)

Data File : C:\FID6_DATA\FI100313.SEC\FI09508.D Vial: 64
 Acq On : 3 Oct 2013 9:02 pm Operator: TIMU
 Sample : D51123-1 Inst : Fid6
 Misc : OP8666,GFI637,30.10,,,1,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Oct 4 8:37 2013 Quant Results File: ORO-RR-GFI585.RES

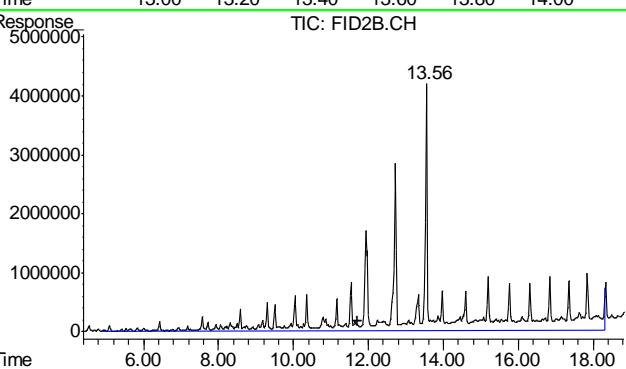
Quant Method : C:\MSDCHEM\1...\ORO-RR-GFI585.M (Chemstation Integrator)
 Title : 8015B TEH Front detector
 Last Update : Thu Oct 03 11:49:33 2013
 Response via : Multiple Level Calibration
 DataAcq Meth : DUAL_B2.M

Volume Inj. : 1ul
 Signal Phase : RTX-5
 Signal Info : 530um

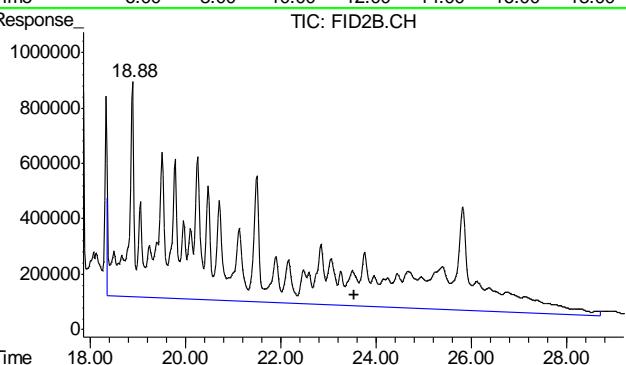




#1 O-Terphenyl
R.T.: 13.558 min
Delta R.T.: 0.018 min
Response: 120979241
Conc: 1669.09 mg/L



#2 TPH-DRO (C10-C28)
R.T.: 11.710 min
Delta R.T.: 0.000 min
Response: 1296321928
Conc: 20227.29 mg/L m



#3 TPH-ORO (>C28-C40)
R.T.: 23.523 min
Delta R.T.: 0.000 min
Response: 753218054
Conc: 18053.12 mg/L m

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH100213.SEC\
 Data File : FH013648.D
 Signal(s) : FID2B.ch
 Acq On : 3 Oct 2013 4:58 am
 Operator : TIMU
 Sample : OP8666-MB
 Misc : OP8666,GFH720,30.00,,,1,1
 ALS Vial : 78 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Oct 03 08:21:13 2013
 Quant Method : C:\msdchem\1\METHODS\DRD-GFH689R.M
 Quant Title : DRO-ORO REAR
 QLast Update : Wed Sep 11 09:58:51 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
1) S o-Terphenyl	12.169	2478226606	1428.275	ug/ml
<hr/>				
Target Compounds				
2) H TPH-DRO (C10-C28)	9.818	150985896	107.343	ug/ml
<hr/>				

(f)=RT Delta > 1/2 Window

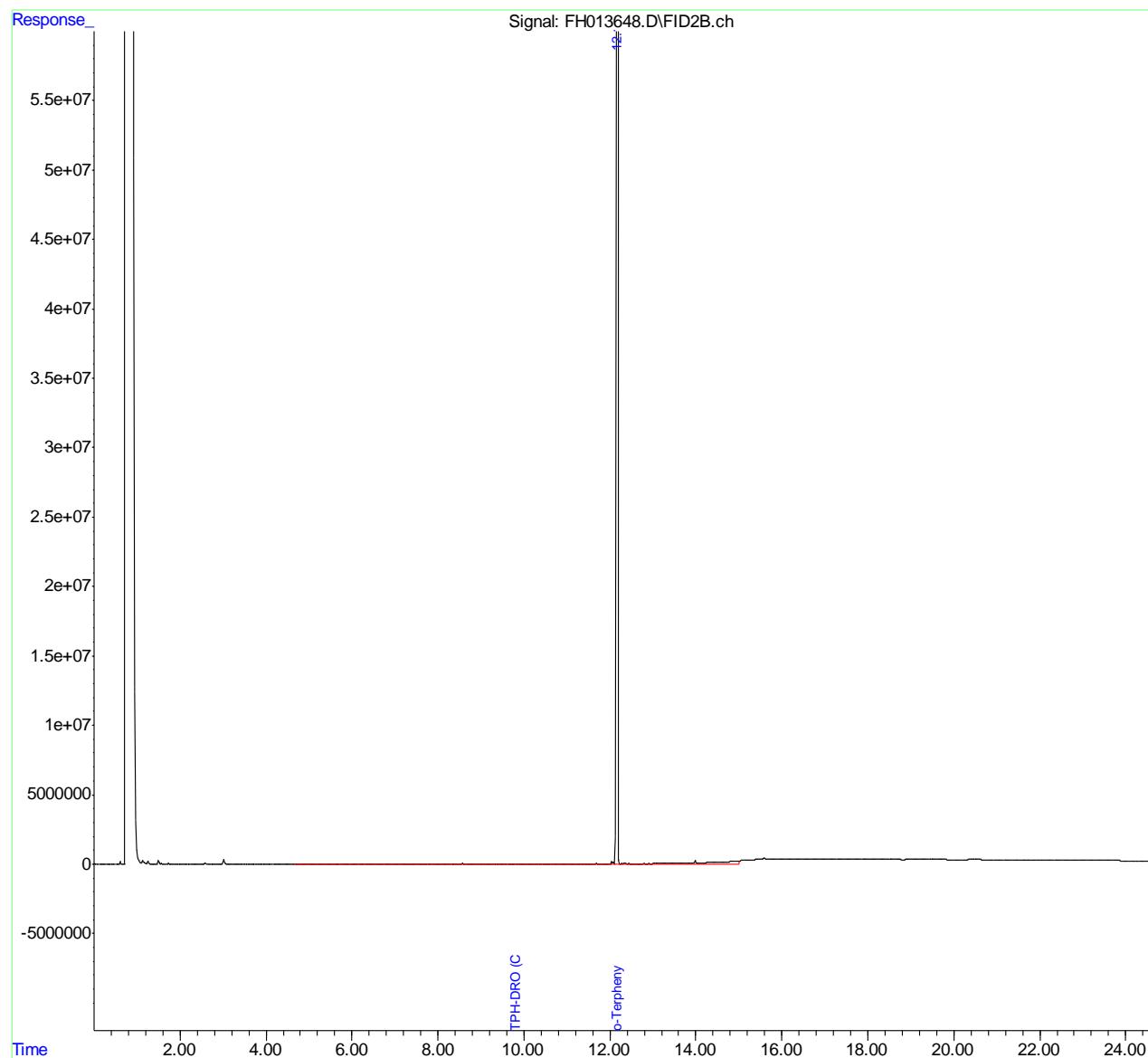
(m)=manual int.

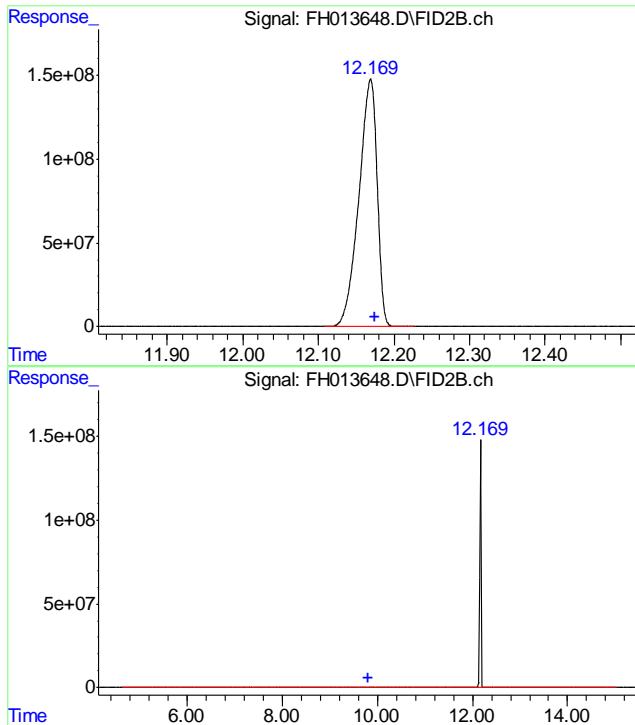
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH100213.SEC\
 Data File : FH013648.D
 Signal(s) : FID2B.ch
 Acq On : 3 Oct 2013 4:58 am
 Operator : TIMU
 Sample : OP8666-MB
 Misc : OP8666,GFH720,30.00,,,1,1
 ALS Vial : 78 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Oct 03 08:21:13 2013
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH689R.M
 Quant Title : DRO-ORO REAR
 QLast Update : Wed Sep 11 09:58:51 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :





#1 o-Terphenyl

R.T.: 12.169 min
Delta R.T.: -0.006 min
Response: 2478226606
Conc: 1428.27 ug/ml

#2 TPH-DRO (C10-C28)

R.T.: 9.818 min
Delta R.T.: 0.000 min
Response: 150985896
Conc: 107.34 ug/ml



Metals Analysis

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Matrix Spike and Duplicate Summaries
- Blank Spike and Lab Control Sample Summaries
- Serial Dilution Summaries

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11267
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

10/02/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	1.1	1.8		
Antimony	3.0	.21	.5		
Arsenic	2.5	.38	.63		
Barium	1.0	.02	.36	0.040	<1.0
Beryllium	1.0	.09	.06		
Boron	5.0	.08	.16		
Cadmium	1.0	.02	.28	0.030	<1.0
Calcium	40	.24	6.8		
Chromium	1.0	.03	.03	0.040	<1.0
Cobalt	0.50	.05	.039		
Copper	1.0	.08	.13	-0.020	<1.0
Iron	7.0	.15	1.8		
Lead	5.0	.21	.25	0.080	<5.0
Lithium	0.50	.04	.13		
Magnesium	20	.68	1.8		
Manganese	0.50	.05	.038		
Molybdenum	1.0	.04	.13		
Nickel	3.0	.05	.07	-0.020	<3.0
Phosphorus	10	1.5	1.2		
Potassium	200	9.9	12		
Selenium	5.0	.71	1.1	-0.47	<5.0
Silicon	5.0	.47	1.1		
Silver	3.0	.03	.05	-0.030	<3.0
Sodium	40	.73	3.7		
Strontium	5.0	.001	.022		
Thallium	1.0	.18	.46		
Tin	5.0	1.2	2.3		
Titanium	1.0	.01	.46		
Uranium	5.0	.29	.31		
Vanadium	1.0	.04	.043		
Zinc	3.0	.04	.16	0.34	<3.0

Associated samples MP11267: D51123-1

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11267
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11267
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 10/02/13

Metal	D51122-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	2780	2690	248	-36.3(a) 75-125
Beryllium				
Boron				
Cadmium	1.0	51.4	62	81.2 75-125
Calcium				
Chromium	23.5	71.7	62	77.7 75-125
Cobalt				
Copper	30.0	87.0	62	91.9 75-125
Iron				
Lead	17.7	118	124	80.8 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	14.5	64.3	62	80.3 75-125
Phosphorus				
Potassium				
Selenium	2.8	115	124	90.4 75-125
Silicon				
Silver	0.036	23.0	24.8	92.5 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	50.9	101	62	80.8 75-125

Associated samples MP11267: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11267
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11267
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

10/02/13

Metal	D51122-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	anr					
Barium	2780	3100	251	127.7(a)	14.2	20
Beryllium						
Boron						
Cadmium	1.0	51.5	62.7	80.6	0.2	20
Calcium						
Chromium	23.5	75.6	62.7	83.1	5.3	20
Cobalt						
Copper	30.0	81.0	62.7	81.4	7.1	20
Iron						
Lead	17.7	117	125	79.2	0.9	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	14.5	63.0	62.7	77.4	2.0	20
Phosphorus						
Potassium						
Selenium	2.8	113	125	87.9	1.8	20
Silicon						
Silver	0.036	23.1	25.1	92.0	0.4	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	50.9	97.7	62.7	74.7N(b)	3.3	20

Associated samples MP11267: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11267
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- (b) Spike recovery indicates possible matrix interference.

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11267
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

10/02/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	192	200	96.0	80-120
Beryllium				
Boron				
Cadmium	45.5	50	91.0	80-120
Calcium				
Chromium	48.8	50	97.6	80-120
Cobalt				
Copper	47.2	50	94.4	80-120
Iron				
Lead	94.1	100	94.1	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	46.7	50	93.4	80-120
Phosphorus				
Potassium				
Selenium	96.1	100	96.1	80-120
Silicon				
Silver	20.2	20	101.0	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	44.9	50	89.8	80-120

Associated samples MP11267: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11267
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51123
 Account: XTOKWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11267
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: ug/l

Prep Date:

10/02/13

Metal	D51122-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	22800	24900	9.0	0-10
Beryllium				
Boron				
Cadmium	8.40	7.00	16.7 (a)	0-10
Calcium				
Chromium	193	207	7.3	0-10
Cobalt				
Copper	247	246	0.5	0-10
Iron				
Lead	146	160	9.5	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	119	134	12.4*(b)	0-10
Phosphorus				
Potassium				
Selenium	23.4	0.00	100.0(a)	0-10
Silicon				
Silver	0.300	3.50	1066.7(a)	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	419	483	15.3*(b)	0-10

Associated samples MP11267: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11267
Matrix Type: SOLID

Methods: SW846 6010C
Units: ug/l

Prep Date:

Metal

- (anr) Analyte not requested
(a) Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
(b) Serial dilution indicates possible matrix interference.

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11268
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date:

10/02/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.55	.75		
Antimony	0.20	.0011	.029		
Arsenic	0.10	.0085	.024	0.0070	<0.10
Barium	1.0	.008	.16		
Beryllium	0.10	.008	.049		
Boron	20	.25	.07		
Cadmium	0.050	.018	.038		
Calcium	200	2.8	13		
Chromium	1.0	.027	.11		
Cobalt	0.10	.0025	.0085		
Copper	1.0	.03	.1		
Iron	5.0	1.8	1.8		
Lead	0.25	.004	.0075		
Magnesium	50	.65	.65		
Manganese	0.50	.06	.07		
Molybdenum	0.50	.025	.046		
Nickel	1.0	.0044	.17		
Phosphorus	30	1.3	4.9		
Potassium	100	1.5	2.5		
Selenium	0.20	.03	.13		
Silver	0.050	.00095	.01		
Sodium	250	2.5	5.5		
Strontium	10	.005	.027		
Thallium	0.10	.0012	.0075		
Tin	5.0	.032	2.3		
Titanium	1.0	.03	.085		
Uranium	0.25	.00085	.0015		
Vanadium	2.0	.019	.11		
Zinc	5.0	.11	1.4		

Associated samples MP11268: D51123-1

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

14.2.1
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
 Account: XTOKWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11268
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

10/02/13

Metal	D51122-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	10.7	132	124	97.8 75-125
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP11268: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (N) Matrix Spike Rec. outside of QC limits
 (anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11268
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

10/02/13

Metal	D51122-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	10.7	132	125	96.8	0.0	20
Barium						
Beryllium						
Boron						
Cadmium						
Calcium						
Chromium						
Cobalt						
Copper						
Iron						
Lead						
Magnesium						
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium						
Silver						
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP11268: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (N) Matrix Spike Rec. outside of QC limits
 (anr) Analyte not requested

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11268
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 10/02/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	94.6	100	94.6	80-120
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP11268: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (anr) Analyte not requested

14.2.3
14

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51123
 Account: XTOKWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11268
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: ug/l

Prep Date:

10/02/13

Metal	D51122-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	88.0	82.1	6.6	0-10
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP11268: D51123-1

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (anr) Analyte not requested

14.2.4
14

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11269
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 10/04/13

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.008	-0.0011	<0.10

Associated samples MP11269: D51123-1

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11269
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 10/04/13

Metal	D51122-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.034	0.50	0.428	108.9 75-125

Associated samples MP11269: D51123-1

Results < IDL are shown as zero for calculation purposes

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11269
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

10/04/13

Metal	D51122-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.034	0.47	0.407	107.2	6.2	20

Associated samples MP11269: D51123-1

Results < IDL are shown as zero for calculation purposes

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

(anr) Analyte not requested

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11269
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

10/04/13

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.42	0.4	105.0	80-120

Associated samples MP11269: D51123-1

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

14.3.3
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BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11305
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
Units: ug/l

Prep Date:

10/04/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	55	210		
Antimony	150	11	95		
Arsenic	130	19	28		
Barium	50	1	7		
Beryllium	50	4.5	6		
Boron	250	4	33		
Cadmium	50	1	1.8		
Calcium	2000	12	210	-29	<2000
Chromium	50	1.5	2		
Cobalt	25	2.5	2.9		
Copper	50	4	9.5		
Iron	350	7.5	48		
Lead	250	11	110		
Lithium	25	2	14		
Magnesium	1000	34	95	16.0	<1000
Manganese	25	2.5	2.3		
Molybdenum	50	2	4.2		
Nickel	150	2.5	4.4		
Phosphorus	500	75	100		
Potassium	5000	500	1400		
Selenium	250	36	55		
Silicon	250	24	26		
Silver	150	1.5	3		
Sodium	2000	37	850	-140	<2000
Strontium	25	.05	.6		
Thallium	50	9	20		
Tin	250	60	80		
Titanium	50	.5	11		
Uranium	250	15	28		
Vanadium	50	2	2		
Zinc	150	2	16		

Associated samples MP11305: D51123-1A

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11305
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
 Account: XTOKWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11305
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
 Units: ug/l

Prep Date:

10/04/13

Metal	D51224-6A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	19600	148000	125000	102.7
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	5170	127000	125000	97.5
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	16500	137000	125000	96.4
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP11305: D51123-1A

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11305
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
Units: ug/l

Prep Date:

Metal

(N) Matrix Spike Rec. outside of QC limits
(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11305
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
 Units: ug/l

Prep Date: 10/04/13

Metal	D51224-6A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	19600	149000	125000	103.5	0.7
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	5170	128000	125000	98.3	0.8
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	16500	137000	125000	96.4	0.0
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP11305: D51123-1A

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11305
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
Units: ug/l

Prep Date:

Metal

(N) Matrix Spike Rec. outside of QC limits
(anr) Analyte not requested

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51123
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11305
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
 Units: ug/l

Prep Date: 10/04/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	130000	125000	104.0	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	122000	125000	97.6	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	121000	125000	96.8	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP11305: D51123-1A

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11305
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51123
 Account: XTOKWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11305
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
 Units: ug/l

Prep Date:

10/04/13

Metal	D51224-6A	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	3920	3920		0.2	0-10
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	1030	1100		6.3	0-10
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	3300	3260		0.9	0-10
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP11305: D51123-1A

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51123
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11305
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested



General Chemistry

QC Data Summaries

Includes the following where applicable:

- Method Blank and Blank Spike Summaries
- Duplicate Summaries
- Matrix Spike Summaries

METHOD BLANK AND SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP11063/GN22129	1.0	0.0	mg/kg	106	101	94.7	80-120%
Specific Conductivity	GP11068/GN22136			umhos/cm	9979	9840	98.6	90-110%
pH	GN22154			su	8.00	8.03	100.4	99.3-100.7%

Associated Samples:

Batch GN22154: D51123-1

Batch GP11063: D51123-1

Batch GP11068: D51123-1

(*) Outside of QC limits

DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP11063/GN22129 GN22168	D51041-1 D51122-1	mg/kg mv	0.12 133	0.0 130	47.2(a) 2.3	0-20% 0-20%

Associated Samples:

Batch GN22168: D51123-1

Batch GP11063: D51123-1

(*) Outside of QC limits

(a) RPD acceptable due to low duplicate and sample concentrations.

MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP11063/GN22129	D51041-1	mg/kg	0.12	40.0	36.3	90.9	75-125%

Associated Samples:

Batch GP11063: D51123-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51123
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP11063/GN22129	D51041-1	mg/kg	0.12	40.0	37.6	3.5	20%

Associated Samples:

Batch GP11063: D51123-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits