



10/09/13

Technical Report for

XTO Energy

FRU 197-31A

1111-02A Cut 1 Subliner Comp

Accutest Job Number: D51203

Sampling Date: 09/30/13



Report to:

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Total number of pages in report: 141



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)

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Test results relate only to samples analyzed.

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Sample Summary

XTO Energy

Job No: D51203

FRU 197-31A

Project No: 1111-02A Cut 1 Subliner Comp

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D51203-1	09/30/13	11:45 DS	10/02/13	SO	Soil	CUT 1 SUBLINER COMP
D51203-1A	09/30/13	11:45 DS	10/02/13	SO	Soil	CUT 1 SUBLINER COMP

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



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: XTO Energy

Job No D51203

Site: FRU 197-31A

Report Date 10/9/2013 1:45:52 PM

On 10/02/2013, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 3.6 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D51203 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: V5V1764

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51084-1MS, D51084-1MSD 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 and 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.

Volatiles by GC By Method SW846 8015B

Matrix: SO

Batch ID: GGB1231

- All samples were analyzed within the recommended method holding time.
- Sample(s) D51200-1MS, D51200-1MSD 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: OP8682

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D51224-3MS, D51224-3MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- The matrix spike (MS) and matrix spike duplicate (MSD) recovery(s) of TPH-DRO (C10-C28) are outside control limits. Probable cause due to matrix interference.

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: MP11290

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51202-1MS, D51202-1MSD, D51202-1SDL were used as the QC samples for the metals analysis.
- The serial dilution RPD(s) for Cadmium, Copper, Lead, Silver are outside control limits for sample MP11290-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP11290-SD1 for Barium, Chromium, Nickel, Zinc : Serial dilution indicates possible matrix interference.

Metals By Method SW846 6020A

Matrix: SO

Batch ID: MP11291

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51202-1MS, D51202-1MSD, D51202-1SDL were used as the QC samples for the metals analysis.

Metals By Method SW846 7471B

Matrix: SO

Batch ID: MP11292

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51148-1MS, D51148-1MSD 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: GN22130

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

Wet Chemistry By Method SW846 3060A/7196A

Matrix: SO

Batch ID: GP11117

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D51202-1MS, D51202-1MSD, D51202-1DUP were used as the QC samples for the Chromium, Hexavalent analysis.
- The duplicate RPD(s) for Chromium, Hexavalent are outside control limits for sample GP11117-D1. RPD acceptable due to low duplicate and sample concentrations.

Wet Chemistry By Method SW846 3060A/7196A M

Matrix: SO

Batch ID: R18969

- The data for SW846 3060A/7196A M meets quality control requirements.
- D51203-1 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: D51203-1

Wet Chemistry By Method USDA HANDBOOK 60**Matrix:** SO**Batch ID:** MP11305

- D51203-1A 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: D51203
Account: XTO Energy
Project: FRU 197-31A
Collected: 09/30/13

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Lab Sample ID	Client Sample ID	Result/ Analyte	Qual	RL	MDL	Units	Method
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D51203-1 CUT 1 SUBLINER COMP

TPH-DRO (C10-C28)	26.7	7.5	5.6	mg/kg	SW846-8015B
Arsenic	5.0	0.11		mg/kg	SW846 6020A
Barium	361	1.1		mg/kg	SW846 6010C
Chromium	41.3	1.1		mg/kg	SW846 6010C
Copper	8.5	1.1		mg/kg	SW846 6010C
Lead	9.2	5.4		mg/kg	SW846 6010C
Nickel	14.8	3.3		mg/kg	SW846 6010C
Zinc	36.8	3.3		mg/kg	SW846 6010C
Specific Conductivity	1610	1.0		umhos/cm	SM 2510B-2011 MOD
Chromium, Trivalent ^a	41.0	2.1		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	225			mv	ASTM D1498-76M
pH	10.18			su	SW846 9045D

D51203-1A CUT 1 SUBLINER COMP

Calcium	4.75	2.0	mg/l	SW846 6010C
Magnesium	3.56	1.0	mg/l	SW846 6010C
Sodium	370	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	31.3		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

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Report of Analysis

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Client Sample ID: CUT 1 SUBLINER COMP**Lab Sample ID:** D51203-1**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** FRU 197-31A**Date Sampled:** 09/30/13**Date Received:** 10/02/13**Percent Solids:** 88.5

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	5V29362.D	1	10/02/13	BD	n/a	n/a	V5V1764
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.07 g	5.0 ml	100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	0.062	0.031	mg/kg	
108-88-3	Toluene	ND	0.12	0.062	mg/kg	
100-41-4	Ethylbenzene	ND	0.12	0.024	mg/kg	
1330-20-7	Xylene (total)	ND	0.25	0.12	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	98%		64-130%
460-00-4	4-Bromofluorobenzene	101%		62-131%
17060-07-0	1,2-Dichloroethane-D4	96%		70-130%

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

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Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP			Date Sampled:	09/30/13		
Lab Sample ID:	D51203-1			Date Received:	10/02/13		
Matrix:	SO - Soil			Percent Solids:	88.5		
Method:	SW846 8270C BY SIM SW846 3546						
Project:	FRU 197-31A						
	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G16537.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	0.0094	0.0049	mg/kg	
120-12-7	Anthracene	ND	0.0094	0.0049	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0094	0.0049	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0094	0.0049	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0094	0.0049	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0094	0.0049	mg/kg	
218-01-9	Chrysene	ND	0.0094	0.0049	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.0094	0.0049	mg/kg	
206-44-0	Fluoranthene	ND	0.0094	0.0049	mg/kg	
86-73-7	Fluorene	ND	0.0094	0.0056	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0094	0.0049	mg/kg	
91-20-3	Naphthalene	ND	0.013	0.012	mg/kg	
129-00-0	Pyrene	ND	0.0094	0.0049	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	72%		10-175%		
321-60-8	2-Fluorobiphenyl	71%		25-130%		
1718-51-0	Terphenyl-d14	86%		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

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	09/30/13
Lab Sample ID:	D51203-1	Date Received:	10/02/13
Matrix:	SO - Soil	Percent Solids:	88.5
Method:	SW846 8015B		
Project:	FRU 197-31A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB22395.D	1	10/02/13	EV	n/a	n/a	GGB1231
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.1 g	5.0 ml	100 ul
Run #2			

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	12	6.2	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2		Limits	
120-82-1	1,2,4-Trichlorobenzene	80%			60-140%	

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

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Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	09/30/13
Lab Sample ID:	D51203-1	Date Received:	10/02/13
Matrix:	SO - Soil	Percent Solids:	88.5
Method:	SW846-8015B SW846 3546		
Project:	FRU 197-31A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FH013760.D	1	10/05/13	TU	10/04/13	OP8682	GFH723
Run #2							

	Initial Weight	Final Volume
Run #1	30.0 g	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	26.7	7.5	5.6	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	65%		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

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	09/30/13
Lab Sample ID:	D51203-1	Date Received:	10/02/13
Matrix:	SO - Soil	Percent Solids:	88.5
Project:	FRU 197-31A		

Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	5.0	0.11	mg/kg	5	10/03/13	10/04/13	JB	SW846 6020A ²
Barium	361	1.1	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Cadmium	< 1.1	1.1	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Chromium	41.3	1.1	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Copper	8.5	1.1	mg/kg	1	10/03/13	10/07/13	JM	SW846 6010C ⁴
Lead	9.2	5.4	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Mercury	< 0.097	0.097	mg/kg	1	10/04/13	10/04/13	JB	SW846 7471B ¹
Nickel	14.8	3.3	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Selenium	< 5.4	5.4	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Silver	< 3.3	3.3	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³
Zinc	36.8	3.3	mg/kg	1	10/03/13	10/04/13	JM	SW846 6010C ³

- (1) Instrument QC Batch: MA4035
- (2) Instrument QC Batch: MA4036
- (3) Instrument QC Batch: MA4038
- (4) Instrument QC Batch: MA4043
- (5) Prep QC Batch: MP11290
- (6) Prep QC Batch: MP11291
- (7) Prep QC Batch: MP11292

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	09/30/13
Lab Sample ID:	D51203-1	Date Received:	10/02/13
Matrix:	SO - Soil	Percent Solids:	88.5
Project:	FRU 197-31A		

General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	1610	1.0	umhos/cm	1	10/08/13	JD	SM 2510B-2011 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	10/08/13	RW	SW846 3060A/7196A
Chromium, Trivalent ^a	41.0	2.1	mg/kg	1	10/08/13	RW	SW846 3060A/7196A M
Redox Potential Vs H2	225		mv	1	10/04/13	AK	ASTM D1498-76M
Solids, Percent	88.5		%	1	10/03/13	SWT	SM2540B-2011 M
pH	10.18		su	1	10/03/13 12:30	AK	SW846 9045D

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

RL = Reporting Limit

Report of Analysis

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4.2
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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	09/30/13
Lab Sample ID:	D51203-1A	Date Received:	10/02/13
Matrix:	SO - Soil	Percent Solids:	88.5
Project:	FRU 197-31A		

SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	4.75	2.0	mg/l	1	10/04/13	10/04/13 JM	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	3.56	1.0	mg/l	1	10/04/13	10/04/13 JM	SW846 6010C ¹	SW846 3010A/M ²
Sodium	370	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

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	09/30/13
Lab Sample ID:	D51203-1A	Date Received:	10/02/13
Matrix:	SO - Soil	Percent Solids:	88.5
Project:	FRU 197-31A		

General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	31.3		ratio	1	10/04/13 21:27	JM	USDA HANDBOOK 60

(a) Calculated as: $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})/2]}$

RL = Reporting Limit



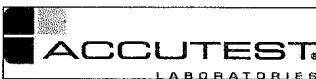
Misc. Forms

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 #
Accutest Quote #	Accutest Job #
D51203	
Requested Analysis (see TEST CODE sheet)	
Matrix Codes	
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	
LAB USE ONLY	

Client / Reporting Information		Project Information		X Table 910														
Company Name KRW Consulting Street Address 8000 West 14th Street; Suite 200 City Lakewood, CO 80214 Project Contact Dwayne Knudson Phone # 970-488-1098 Sampler(s) Name(s) David Sanders 970-488-1098		Project Name: XTO FRV 197-31A Street Billing Information (if different from Report to) Company Name XTO Energy Project # 1111-02A Street Address 21459 CR 5 Client Purchase Order # City Rifle, CO 81650 Project Manager Joe Hess Attention: Jessica Dooling																
Accutest Sample # Field ID / Point of Collection Cut 1 Submer Camp.	MEOH/DI Vial #	Collection				Number of Preserved Bottles												
		Date	Time			Sampled by	Matrix	# of bottles	HCl	NaOH	HNO3	H2SO4	NONE	D/Water	MEOH	ENCORE	Biulatte	
		9/30/13	1145			DS 50	S	X										
Turnaround Time (Business days)		Approved By (Accutest PM): / Date:		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"/>				<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																		
Sample Custody must be documented below each time samples change possession, including courier delivery.																		
Relinquished by Sampler: <i>1 David Sanders</i>	Date Time: <i>10/1/13 1331</i>	Received By: <i>Service Center</i>	Relinquished By: <i>2</i>	Date Time: <i></i>	Received By: <i>2 JACOB Parra</i>	Date Time: <i>13:00 10/2/13</i>												
Relinquished by Sampler: <i>3</i>	Date Time: <i></i>	Received By: <i>3</i>	Relinquished By: <i>4</i>	Date Time: <i></i>	Received By: <i>4</i>	Date Time: <i></i>												
Relinquished by: <i>5</i>	Date Time: <i></i>	Received By: <i>5</i>	Custody Seal # <i>NDLco</i>	Intact <input checked="" type="checkbox"/> Not Intact <input type="checkbox"/>	Preserved where applicable <i>BN</i>	On Ice <input type="checkbox"/>	Cooler Temp. <i>3.6</i>											

D51203: Chain of Custody

Page 1 of 2



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D51203

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 10/2/2013 1:00:00 PM

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

D51203: 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

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1764-MB	5V29351.D	1	10/02/13	BD	n/a	n/a	V5V1764

The QC reported here applies to the following samples:

Method: SW846 8260B

D51203-1

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	101%	64-130%
460-00-4	4-Bromofluorobenzene	84%	62-131%
17060-07-0	1,2-Dichloroethane-D4	101%	70-130%

Blank Spike Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1764-BS	5V29352.D	1	10/02/13	BD	n/a	n/a	V5V1764

The QC reported here applies to the following samples:

Method: SW846 8260B

D51203-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	2500	2560	102	70-130
100-41-4	Ethylbenzene	2500	2700	108	70-130
108-88-3	Toluene	2500	2660	106	70-130
1330-20-7	Xylene (total)	7500	8520	114	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	102%	64-130%
460-00-4	4-Bromofluorobenzene	100%	62-131%
17060-07-0	1,2-Dichloroethane-D4	96%	70-130%

* = Outside of Control Limits.

Blank Spike Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1764-BS	5V29354.D	1	10/02/13	BD	n/a	n/a	V5V1764

The QC reported here applies to the following samples:

Method: SW846 8260B

D51203-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
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CAS No.	Surrogate Recoveries	BSP	Limits
---------	----------------------	-----	--------

2037-26-5	Toluene-D8	102%	64-130%
460-00-4	4-Bromofluorobenzene	91%	62-131%
17060-07-0	1,2-Dichloroethane-D4	90%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D51084-1MS	5V29357.D	1	10/02/13	BD	n/a	n/a	V5V1764
D51084-1MSD	5V29358.D	1	10/02/13	BD	n/a	n/a	V5V1764
D51084-1	5V29356.D	1	10/02/13	BD	n/a	n/a	V5V1764

The QC reported here applies to the following samples:

Method: SW846 8260B

D51203-1

CAS No.	Compound	D51084-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	ND		4170	4230	102	4120	99	3	64-139/30
100-41-4	Ethylbenzene	ND		4170	4180	100	4140	99	1	68-136/30
108-88-3	Toluene	ND		4170	3740	90	3810	91	2	60-130/30
1330-20-7	Xylene (total)	ND		12500	13200	106	13000	104	2	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D51084-1	Limits
2037-26-5	Toluene-D8	98%	101%	102%	64-130%
460-00-4	4-Bromofluorobenzene	119%	118%	125%	62-131%
17060-07-0	1,2-Dichloroethane-D4	92%	96%	91%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D51084-1MS	5V29359.D	1	10/02/13	BD	n/a	n/a	V5V1764
D51084-1MSD	5V29360.D	1	10/02/13	BD	n/a	n/a	V5V1764
D51084-1	5V29356.D	1	10/02/13	BD	n/a	n/a	V5V1764

The QC reported here applies to the following samples:

Method: SW846 8260B

D51203-1

CAS No.	Compound	D51084-1		Spike	MS	MS	MSD	MSD	Limits
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD

CAS No.	Surrogate Recoveries	MS	MSD	D51084-1	Limits
2037-26-5	Toluene-D8	106%	102%	102%	64-130%
460-00-4	4-Bromofluorobenzene	127%	126%	125%	62-131%
17060-07-0	1,2-Dichloroethane-D4	87%	87%	91%	70-130%

* = Outside of Control Limits.



GC/MS Volatiles

Raw Data

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5100213.S\
 Data File : 5V29362.D
 Acq On : 2 Oct 2013 6:51 pm
 Operator : BRETD
 Sample : D51203-1
 Misc : MS6481,V5V1764,5.074,,100,5,1
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Oct 03 08:48:11 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	166253	50.00	ug/l	0.00
37) 1,4-Difluorobenzene	12.412	114	216616	50.00	ug/l	0.00
56) Chlorobenzene-d5	15.061	117	215179	50.00	ug/l	0.00
77) 1,4-Dichlorobenzene-d4	17.036	152	158937	50.00	ug/l	0.00

System Monitoring Compounds						
35) 1,2-Dichloroethane-d4	12.012	102	16252	48.11	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	96.22%
64) Toluene-d8	13.805	98	238910	49.01	ug/l	-0.01
Spiked Amount	50.000	Range	70 - 130	Recovery	=	98.02%
72) 4-Bromofluorobenzene	16.008	95	114398	50.35	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	100.70%

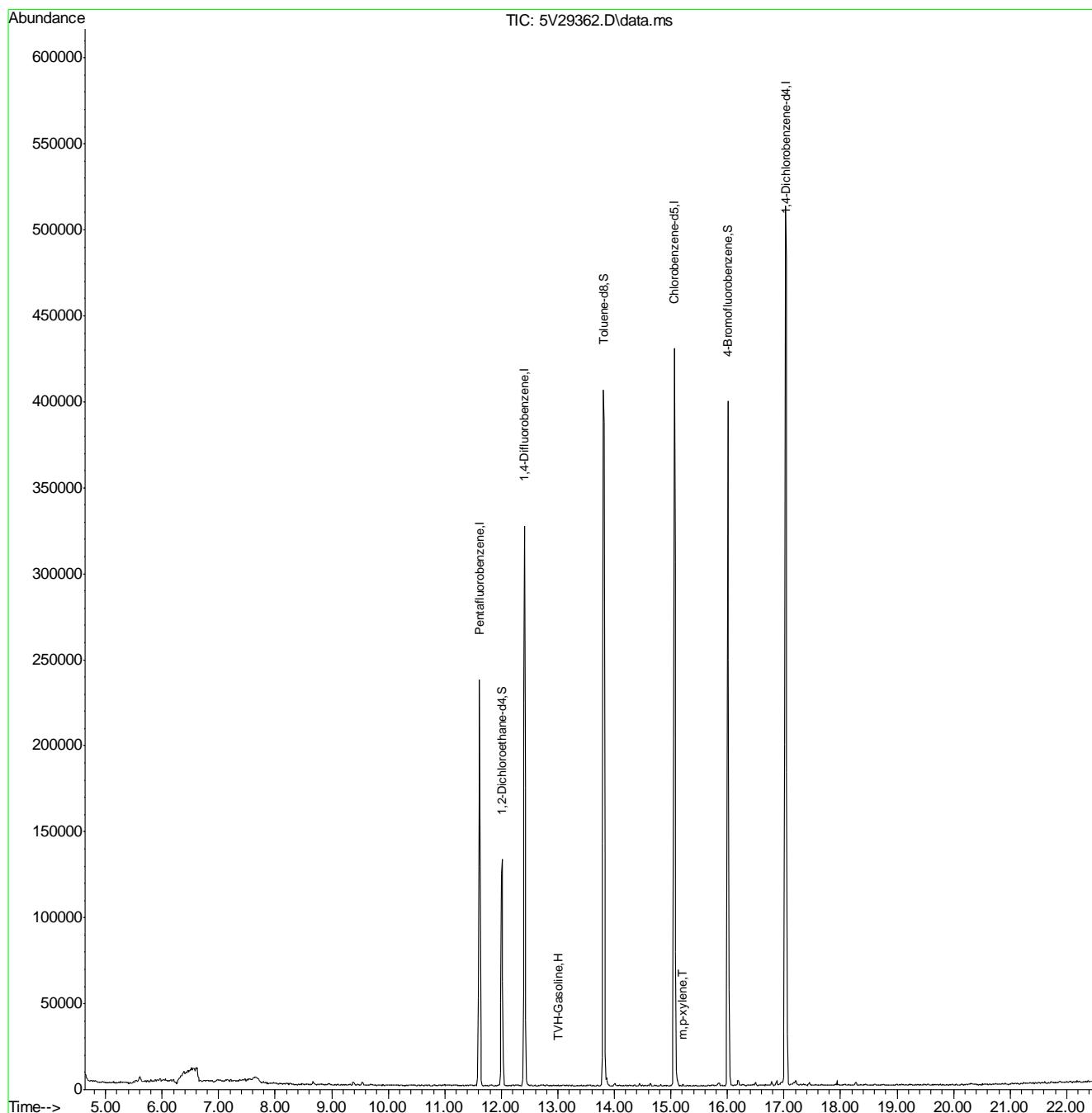
Target Compounds					Qvalue
1) TVH-Gasoline	13.006	TIC	12372m	58.69	ug/l
75) m,p-xylene	15.209	106	373	0.15	ug/l

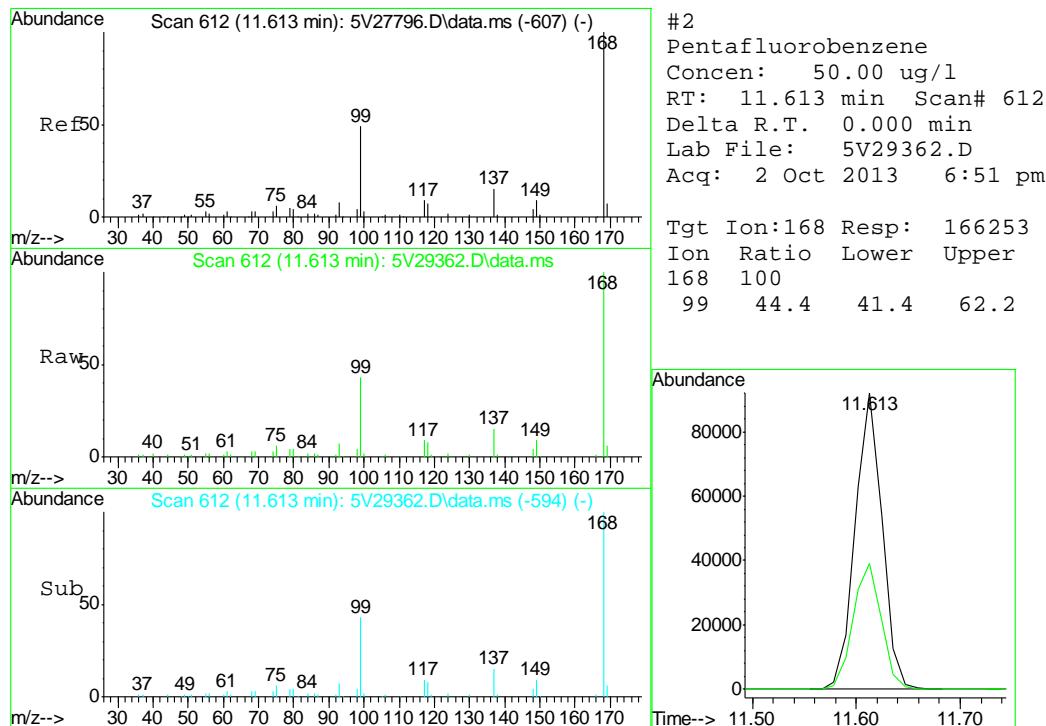
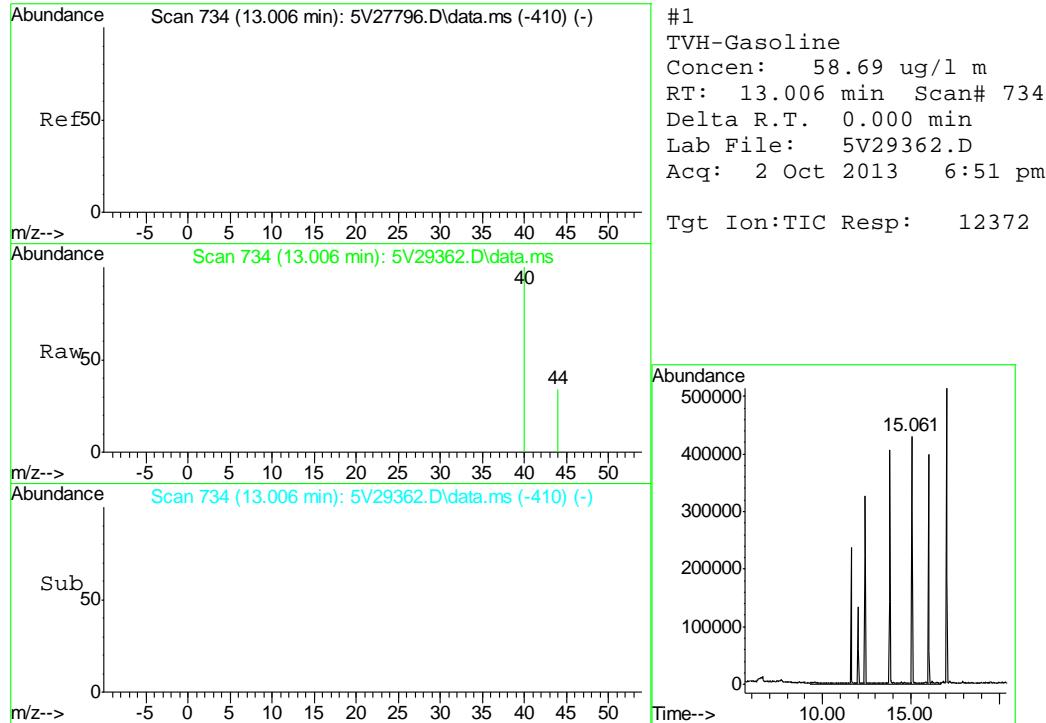
(#) = qualifier out of range (m) = manual integration (+) = signals summed

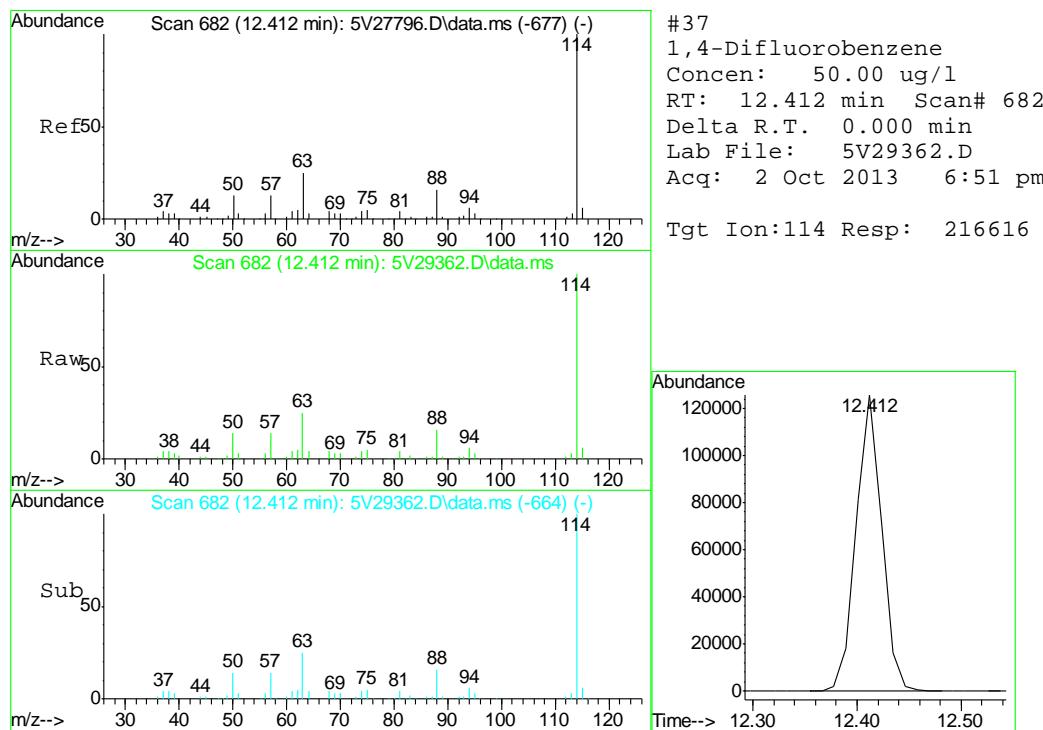
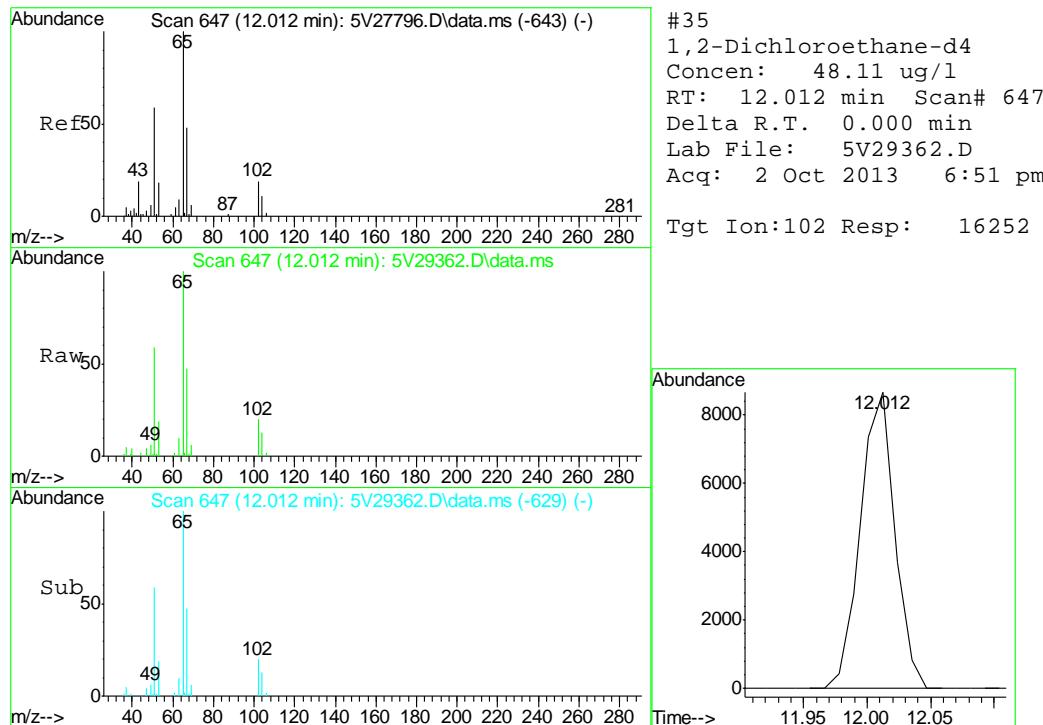
Quantitation Report (QT Reviewed)

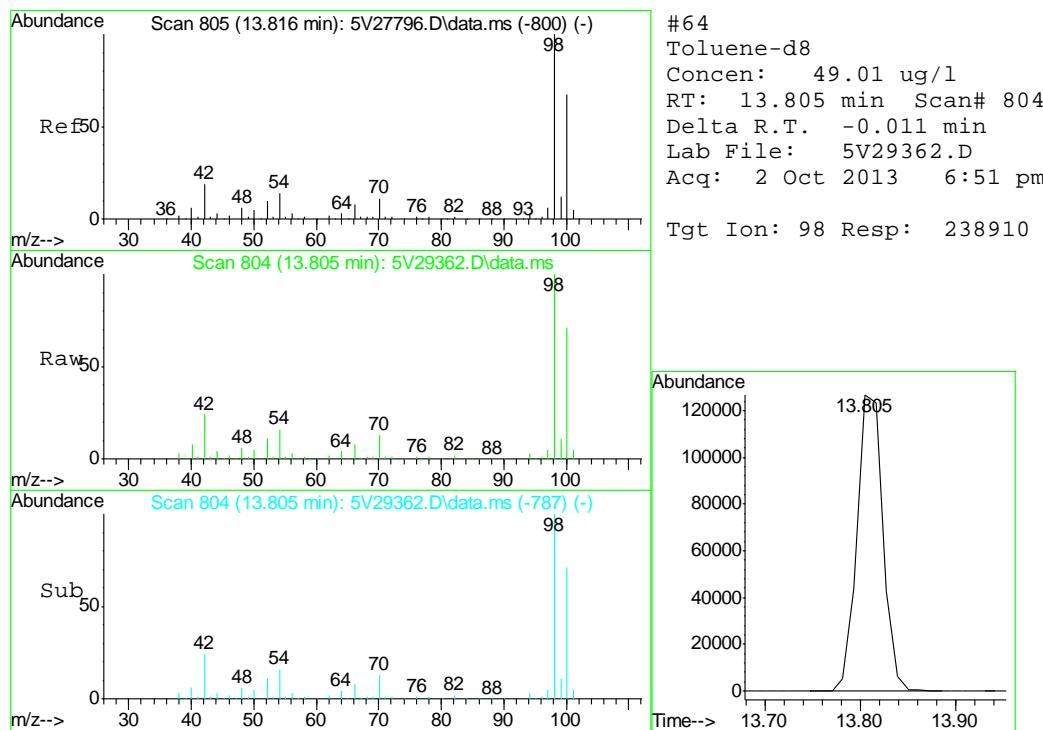
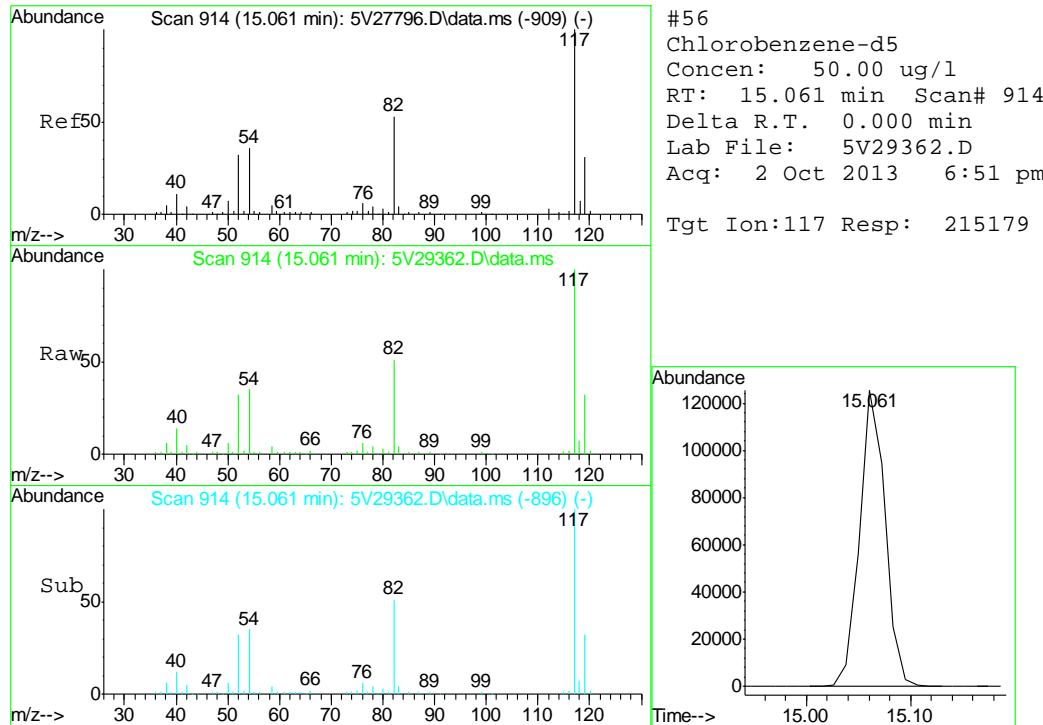
Data Path : C:\msdchem\1\DATA\V5100213.S\
 Data File : 5V29362.D
 Acq On : 2 Oct 2013 6:51 pm
 Operator : BRETD
 Sample : D51203-1
 Misc : MS6481,V5V1764,5.074,,100,5,1
 ALS Vial : 16 Sample Multiplier: 1

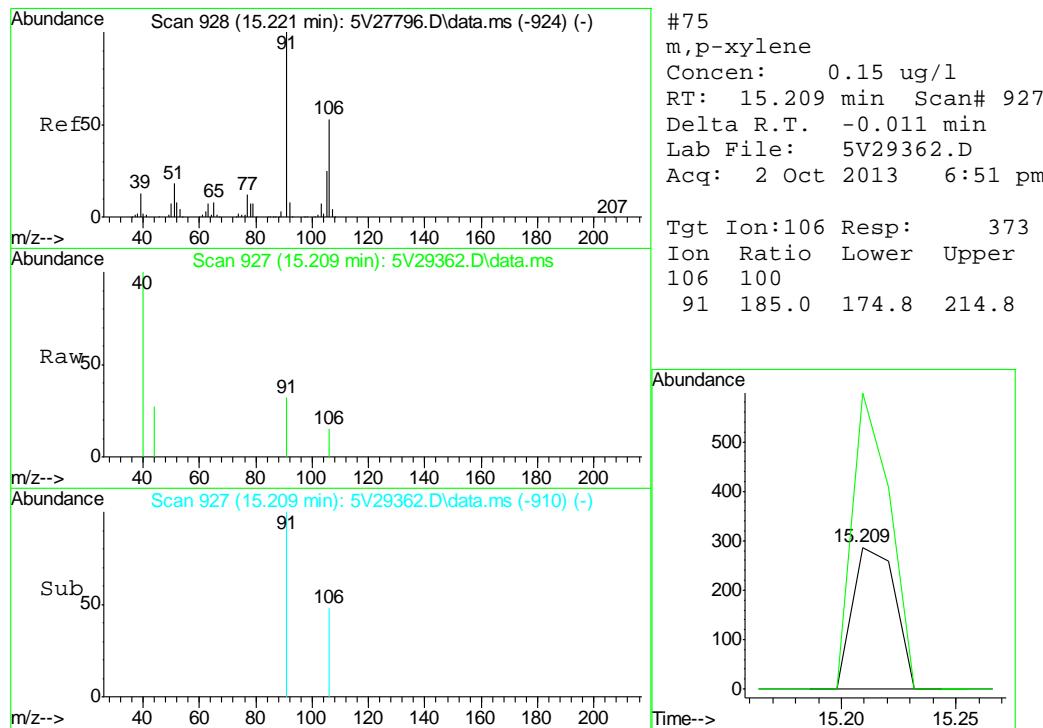
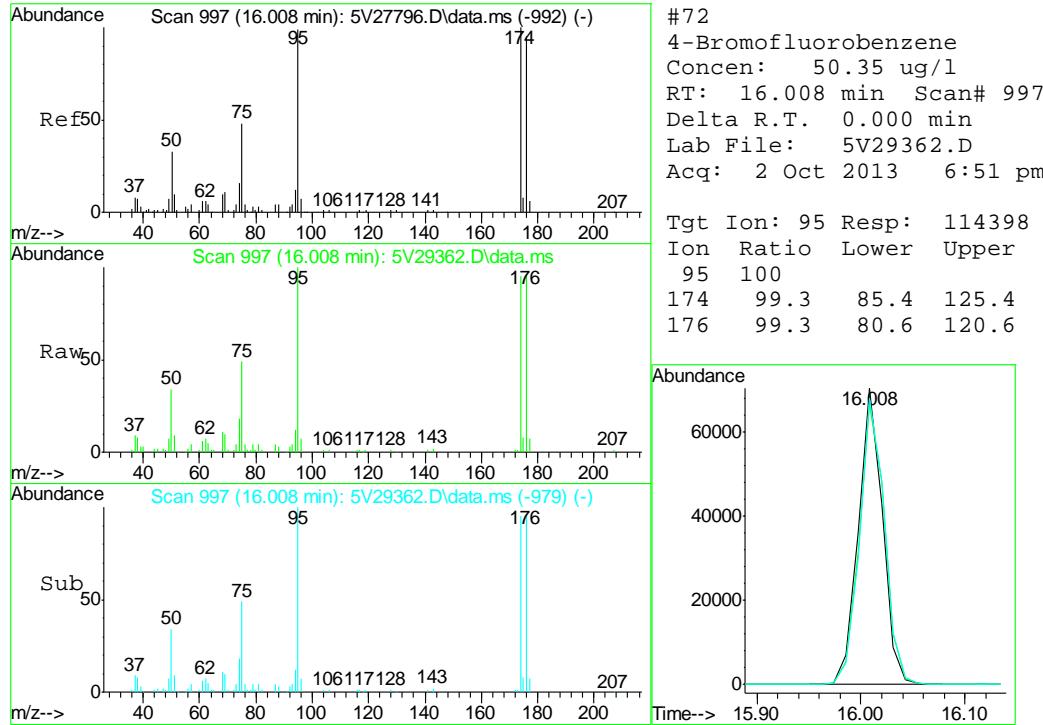
Quant Time: Oct 03 08:48:11 2013
 Quant Method : C:\msdchem\1\METHODS\V5AP1728TVH1728.M
 Quant Title : 8260
 QLast Update : Tue Aug 20 09:59:22 2013
 Response via : Initial Calibration

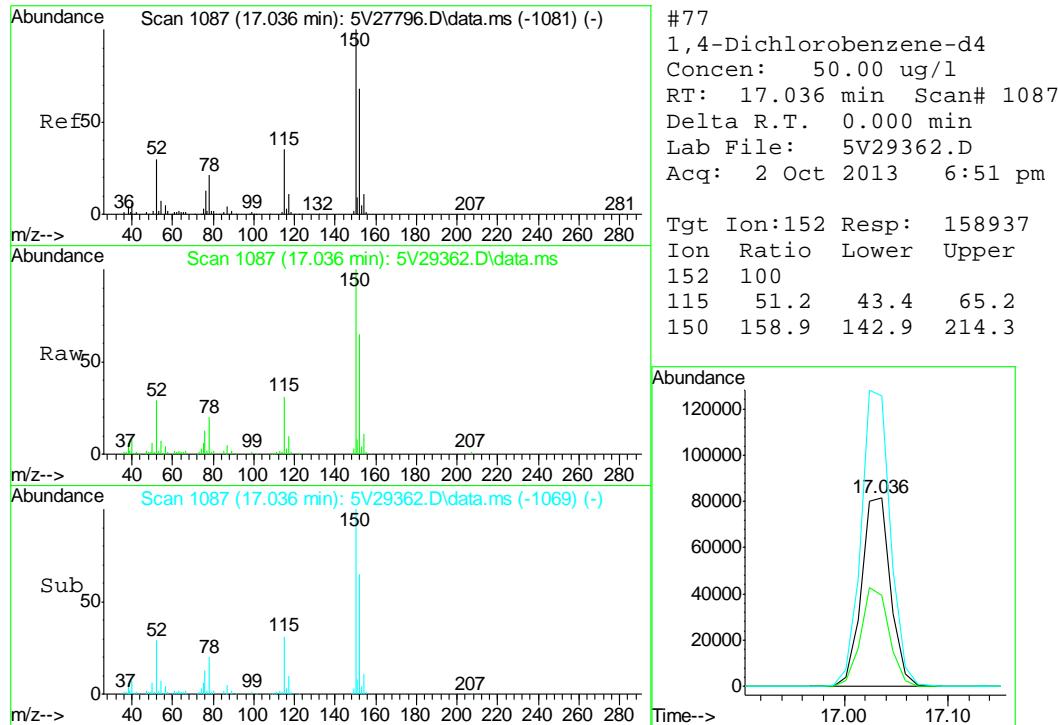












Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5100213.S\
 Data File : 5V29351.D
 Acq On : 2 Oct 2013 1:07 pm
 Operator : BRETD
 Sample : MB
 Misc : MS6481,V5V1764,5.000,,100,5,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Oct 03 09:15:35 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	155231	50.00	ug/l	0.00
37) 1,4-Difluorobenzene	12.412	114	212288	50.00	ug/l	0.00
56) Chlorobenzene-d5	15.061	117	202438	50.00	ug/l	0.00
77) 1,4-Dichlorobenzene-d4	17.025	152	132604	50.00	ug/l	-0.01

System Monitoring Compounds						
35) 1,2-Dichloroethane-d4	12.013	102	15972	50.63	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	101.26%
64) Toluene-d8	13.816	98	231512	50.48	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	100.96%
72) 4-Bromofluorobenzene	16.008	95	89552	41.89	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	83.78%

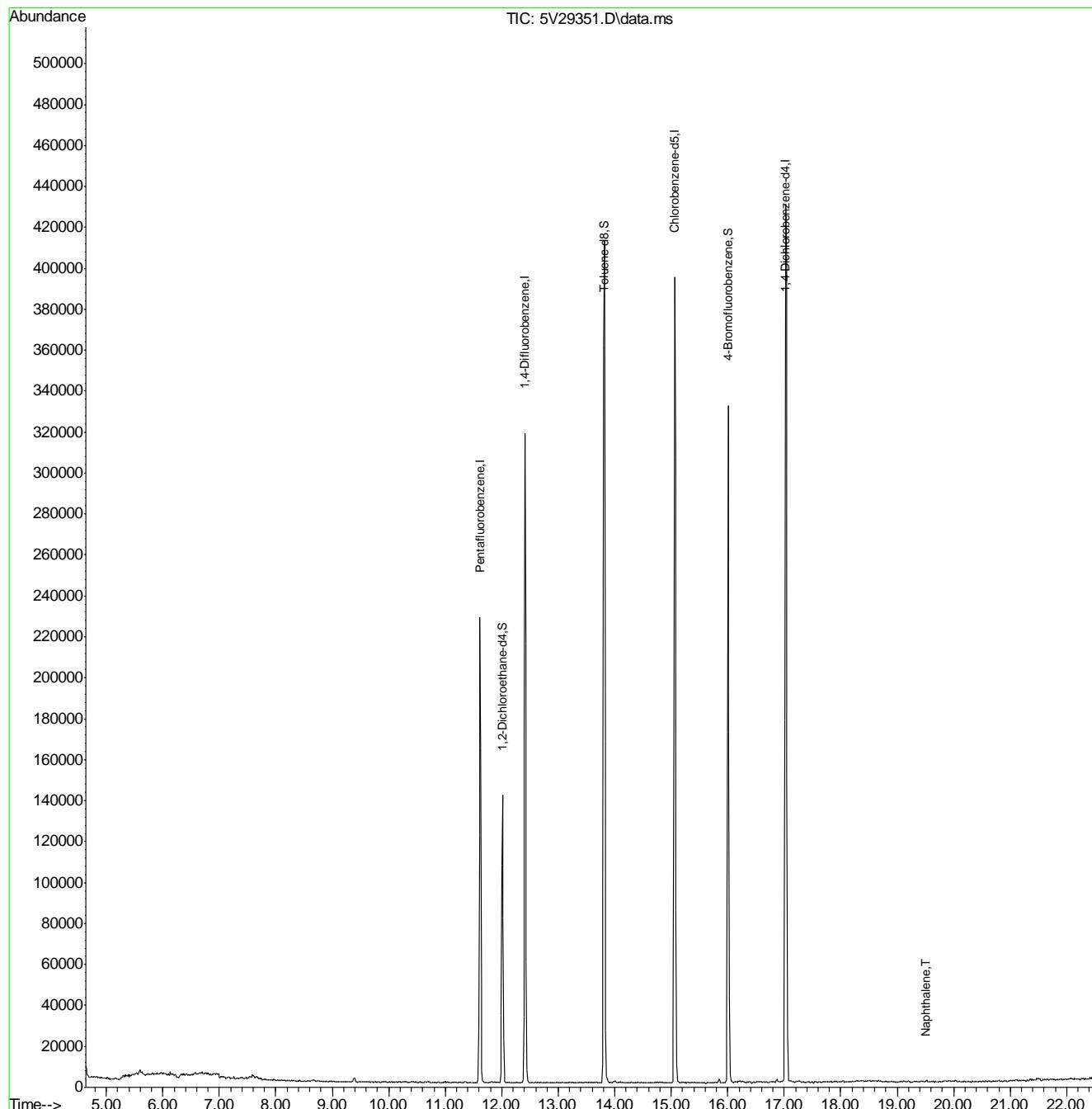
Target Compounds					Qvalue
1) TVH-Gasoline	13.006	TIC	-9922m	57.07	ug/l
94) Naphthalene	19.513	128	995	0.98	ug/l

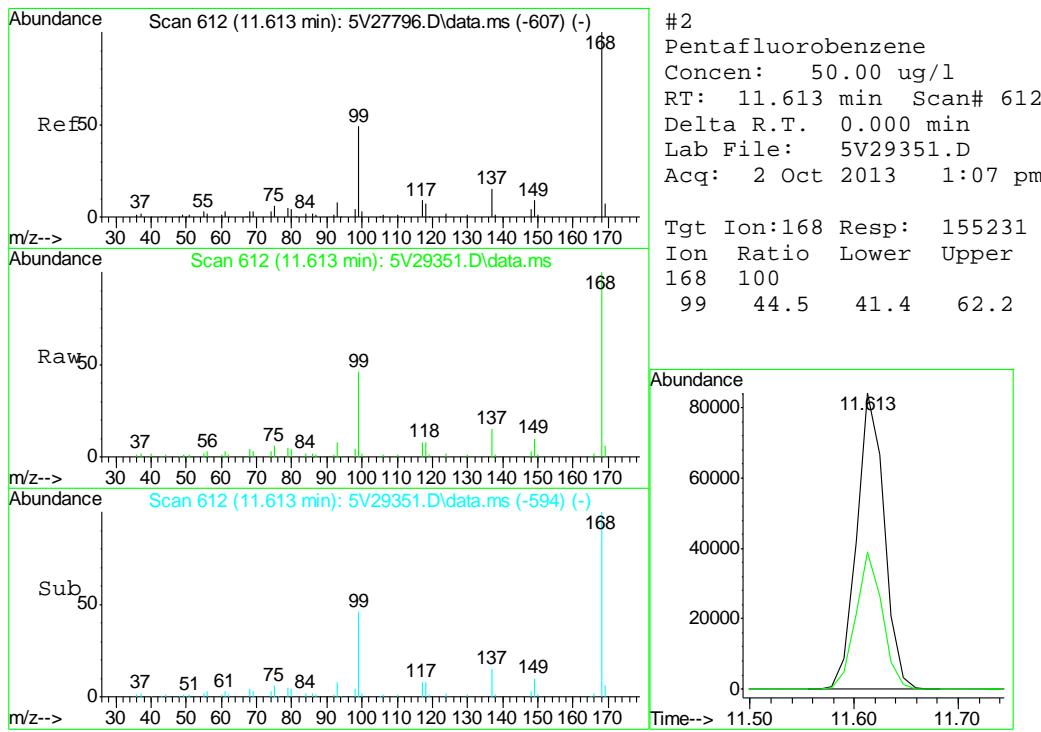
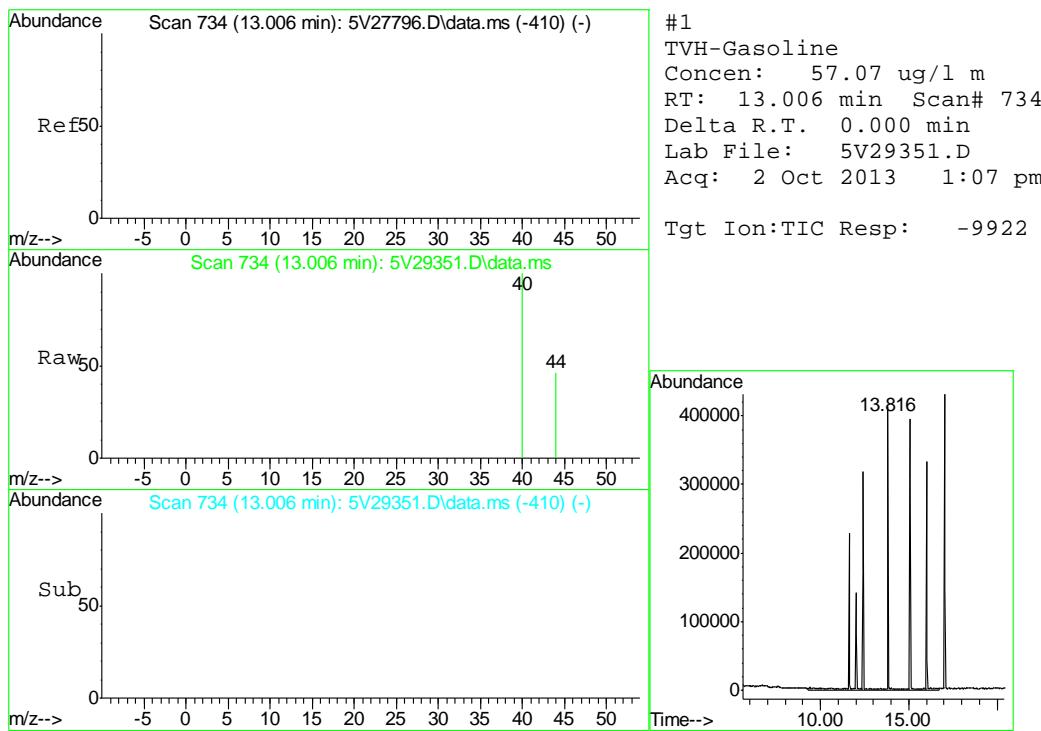
(#) = qualifier out of range (m) = manual integration (+) = signals summed

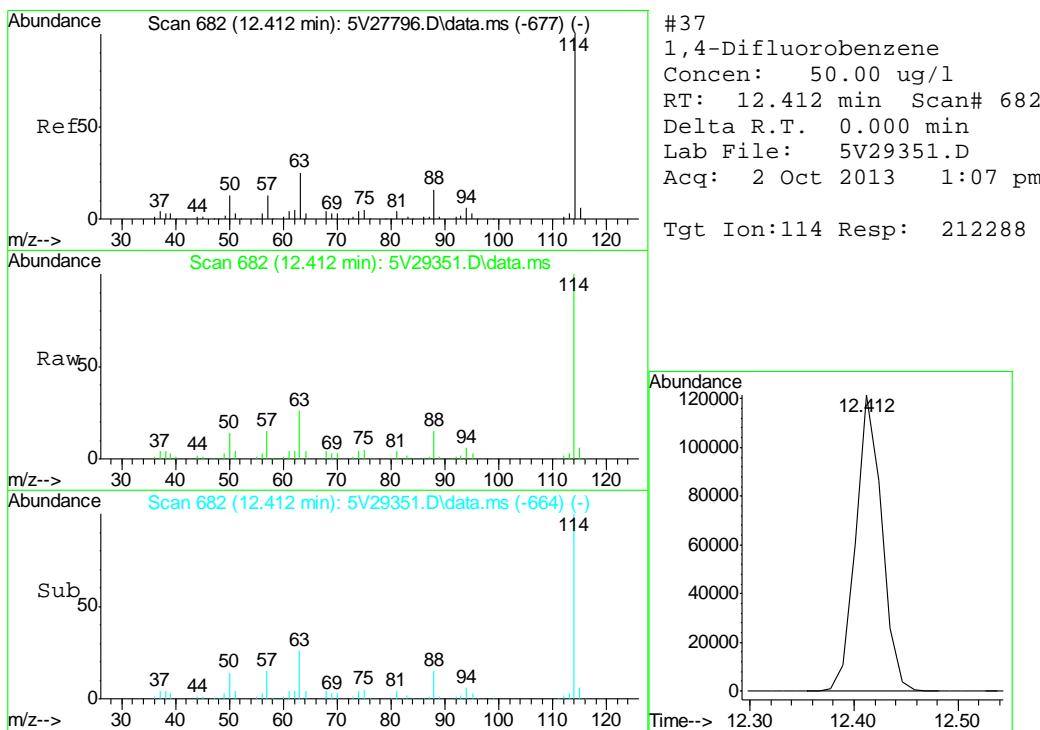
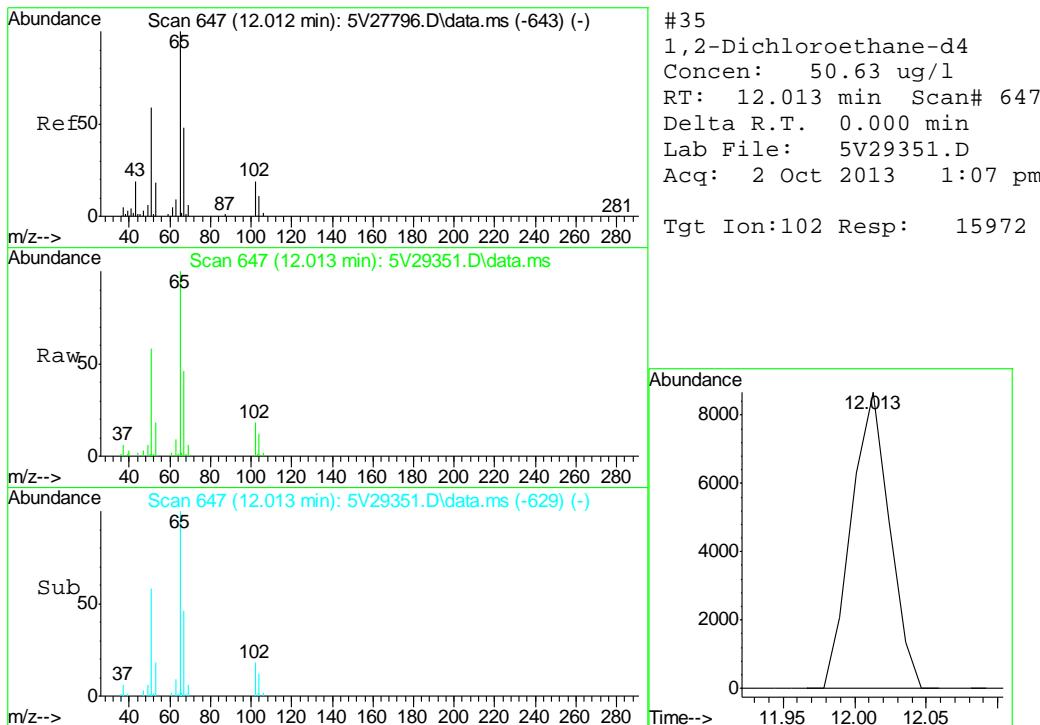
Quantitation Report (QT Reviewed)

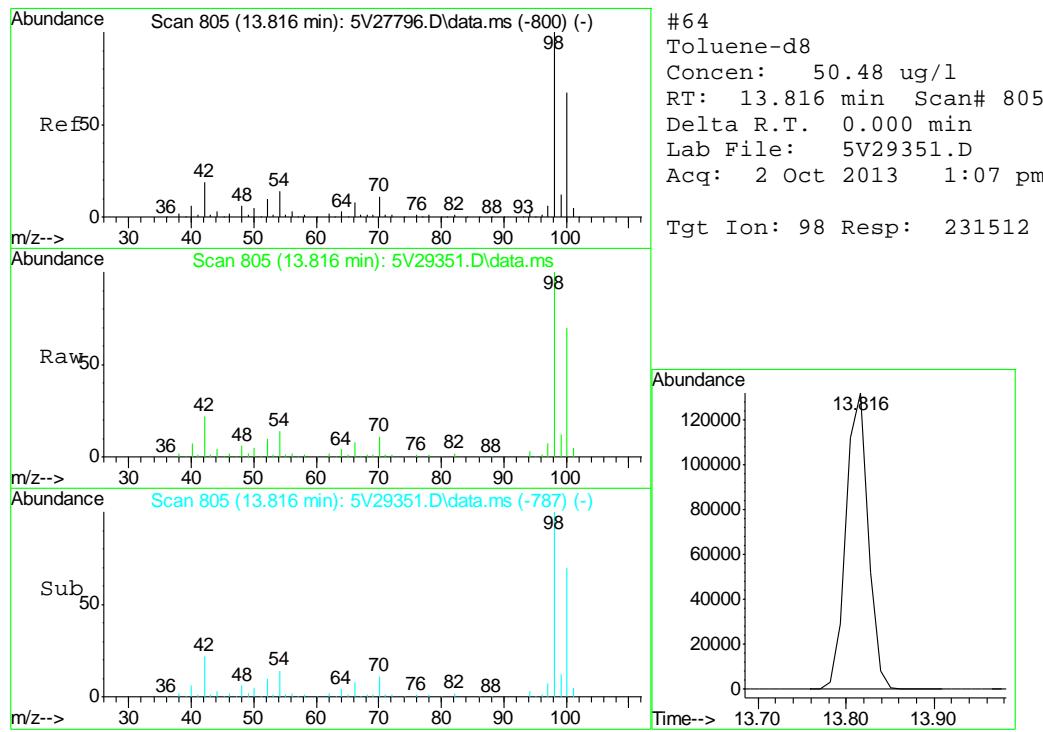
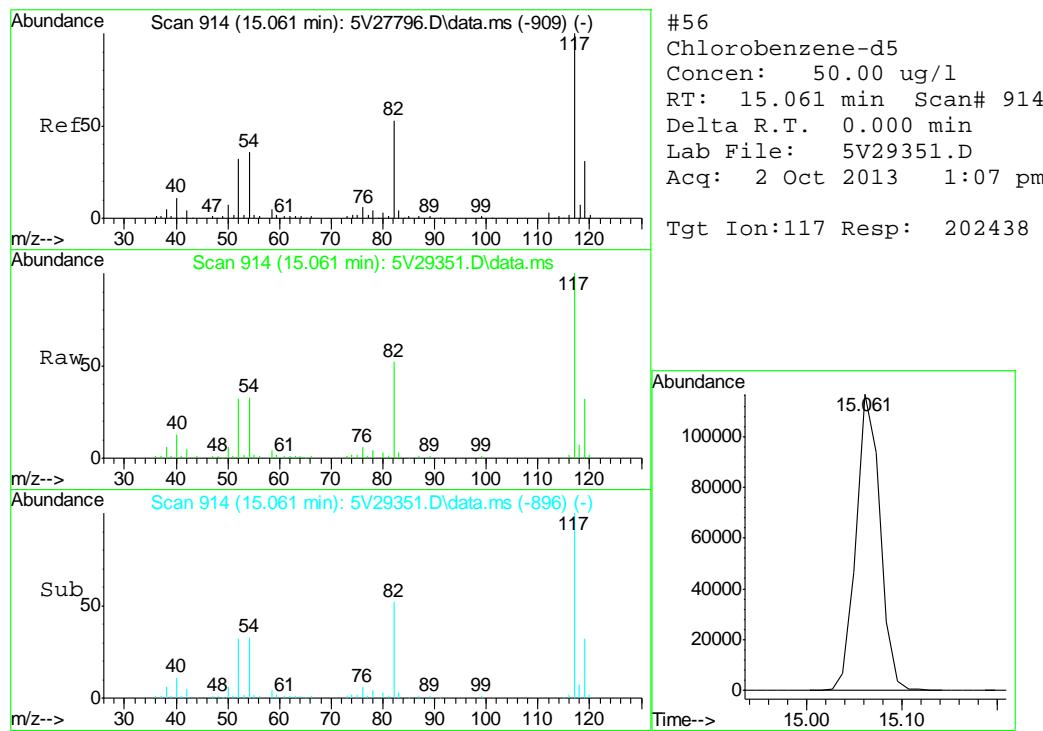
Data Path : C:\msdchem\1\DATA\V5100213.S\
 Data File : 5V29351.D
 Acq On : 2 Oct 2013 1:07 pm
 Operator : BRETD
 Sample : MB
 Misc : MS6481,V5V1764,5.000,,100,5,1
 ALS Vial : 5 Sample Multiplier: 1

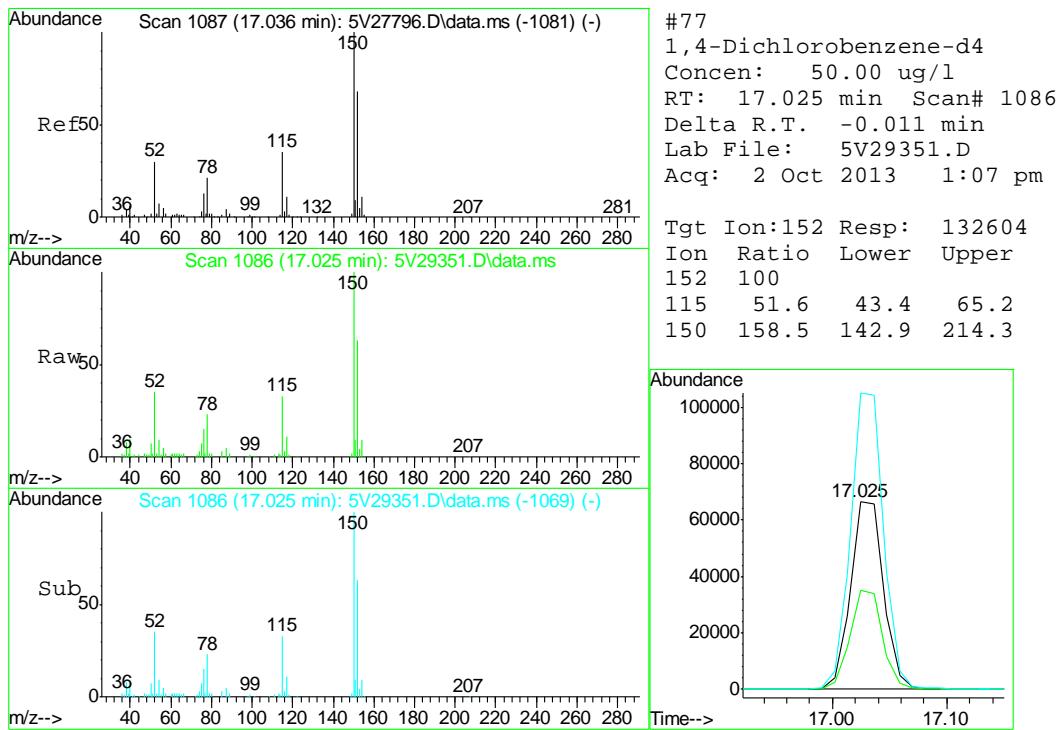
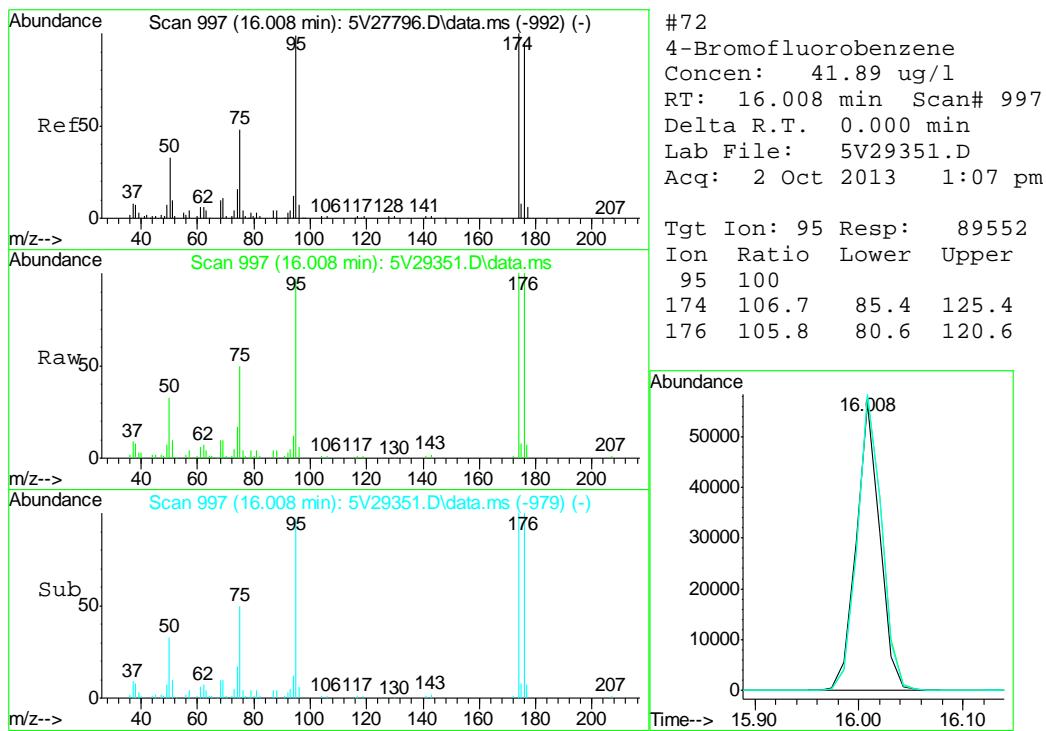
Quant Time: Oct 03 09:15:35 2013
 Quant Method : C:\msdchem\1\METHODS\V5AP1728TVH1728.M
 Quant Title : 8260
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 Response via : Initial Calibration

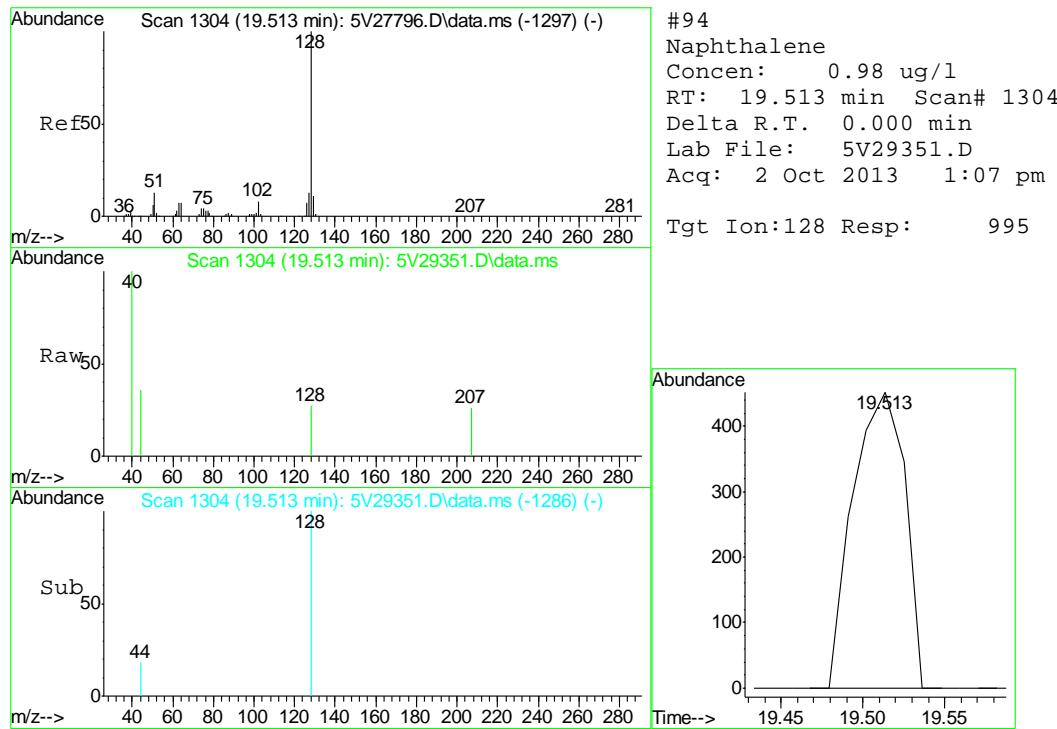












7.2.1

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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: D51203

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

D51203-1

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: D51203
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

D51203-1

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: D51203

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

D51203-1

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 : 3g16537.D
 Acq On : 3 Oct 2013 8:56 pm
 Operator : DONC
 Sample : D51203-1
 Misc : OP8670,E3G817,30.13,,,1,1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Oct 04 15:02:40 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	221041	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.398	164	151002	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.872	188	230669	4.0000	ug/mL	0.00
19) Chrysene-d12	11.501	240	184762	4.0000	ug/mL	0.00
24) Perylene-d12	12.865	264	125572	4.0000	ug/mL	0.00

System Monitoring Compounds						
2) Nitrobenzene-d5	4.996	82	997698	35.8835	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	71.76%
7) 2-Fluorobiphenyl	6.736	172	2088675	35.5026	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	71.00%
21) Terphenyl-d14	10.463	244	1503953	43.0219	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	86.04%

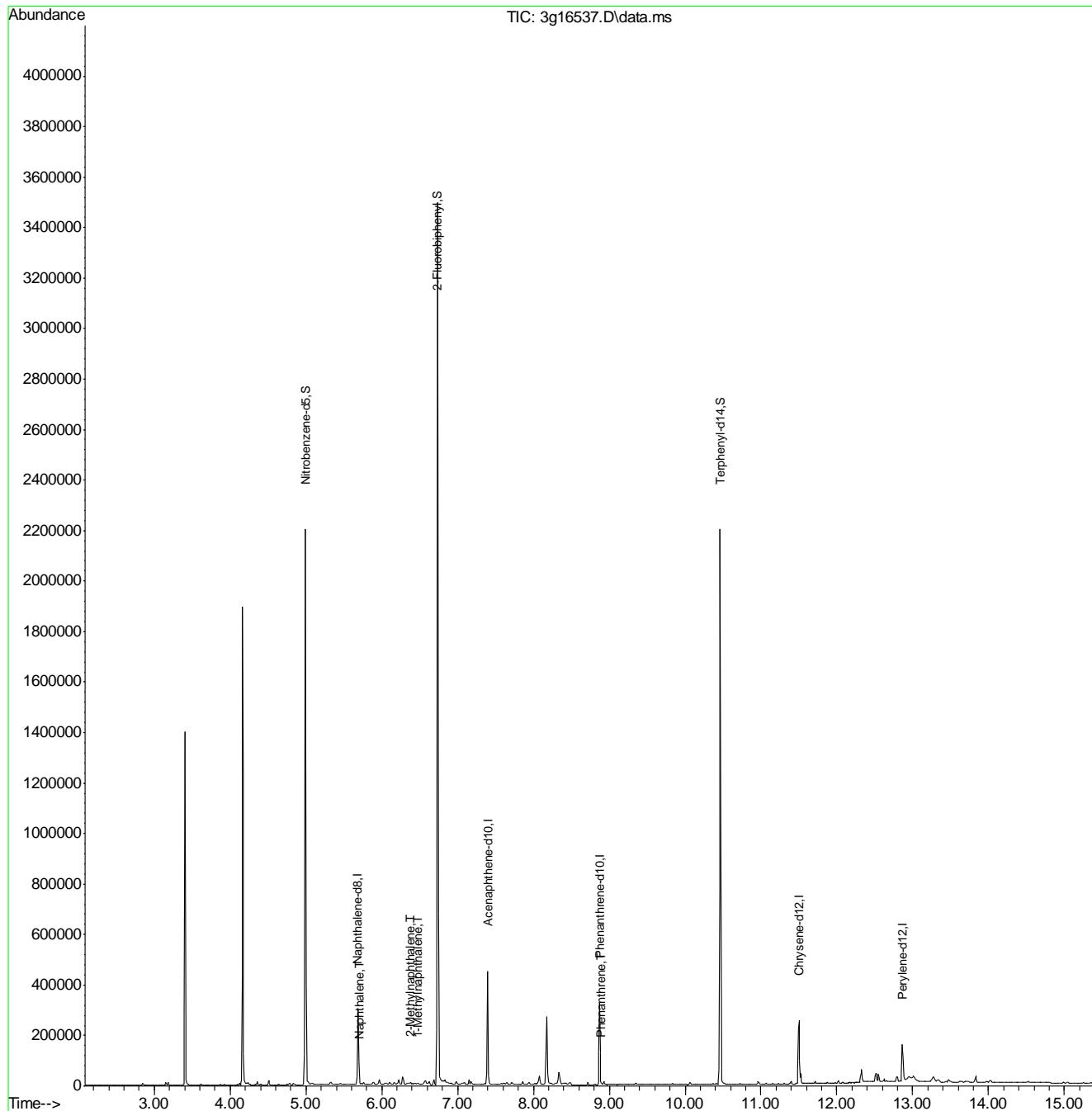
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.407	74	50	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.707	128	2910	0.0364 ug/mL#	79
8) 2-Methylnaphthalene	6.380	142	3503	0.0578 ug/mL	87
9) 1-Methylnaphthalene	6.480	142	2205	0.0429 ug/mL	92
10) Acenaphthylene	7.256	152	645	N.D.	
11) Acenaphthene	7.398	154	863	N.D.	
12) Dibenzofuran	7.610	168	938	N.D.	
13) Fluorene	7.941	166	970	N.D.	
14) Diphenylamine	0.000	169	0	N.D. d	
16) Phenanthrene	8.896	178	3143	0.0335 ug/mL#	72
17) Anthracene	0.000	178	0	N.D. d	
18) Fluoranthene	0.000	202	0	N.D. d	
20) Pyrene	10.305	202	552	N.D.	
22) Benzo(a)anthracene	11.495	228	1278	N.D.	
23) Chrysene	11.495	228	1278	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	

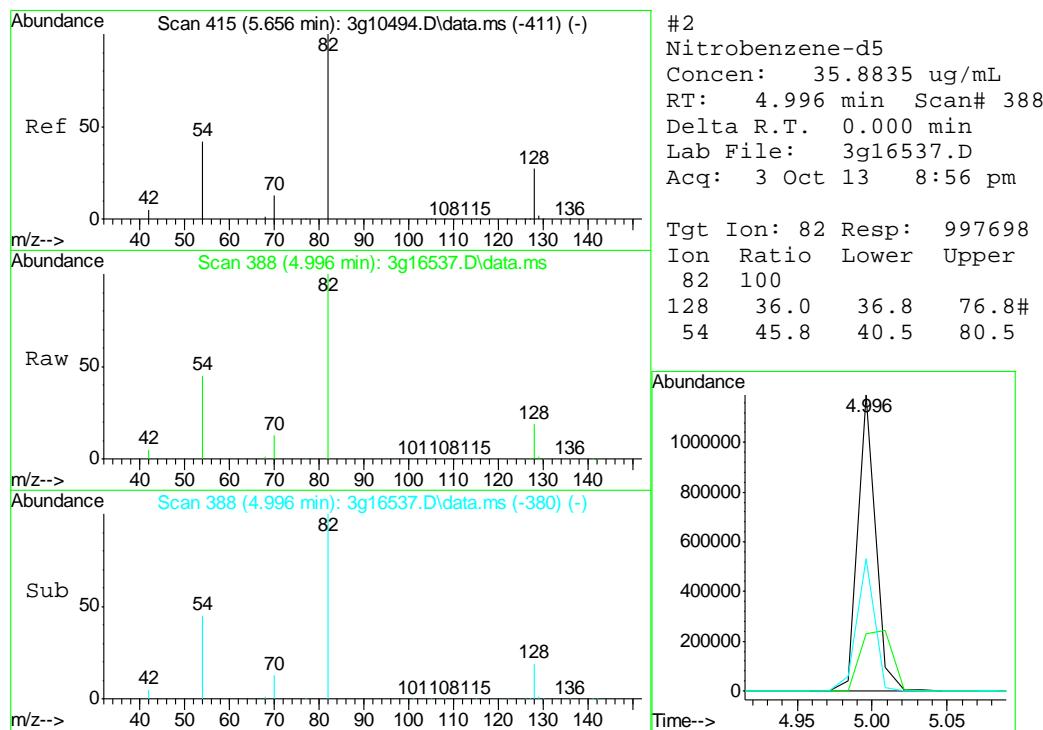
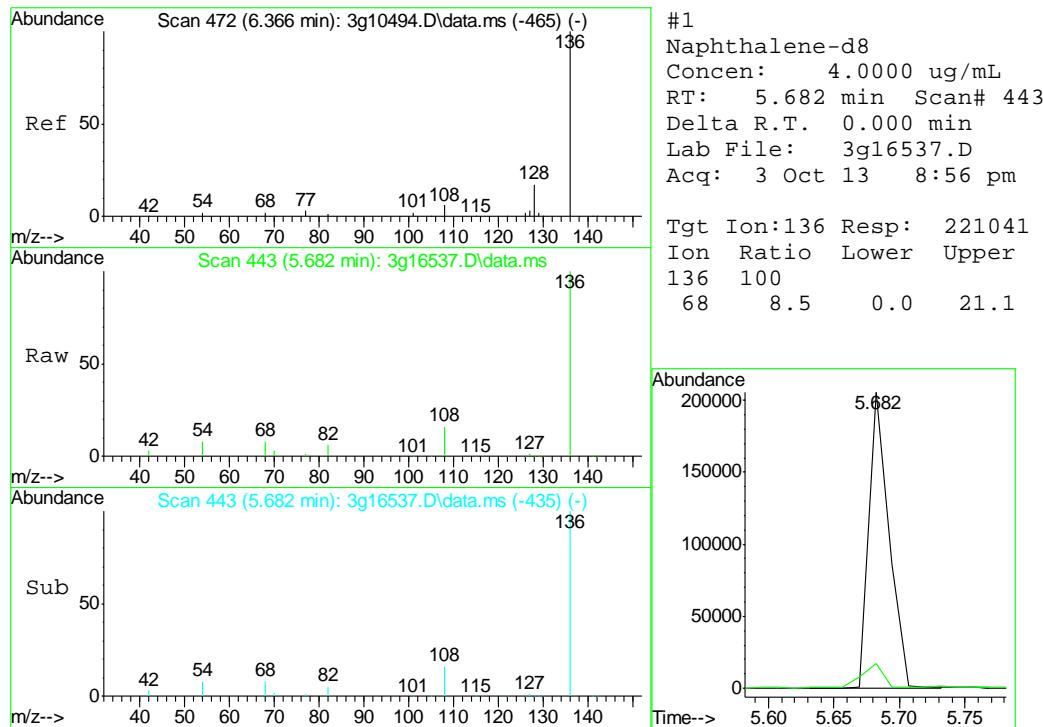
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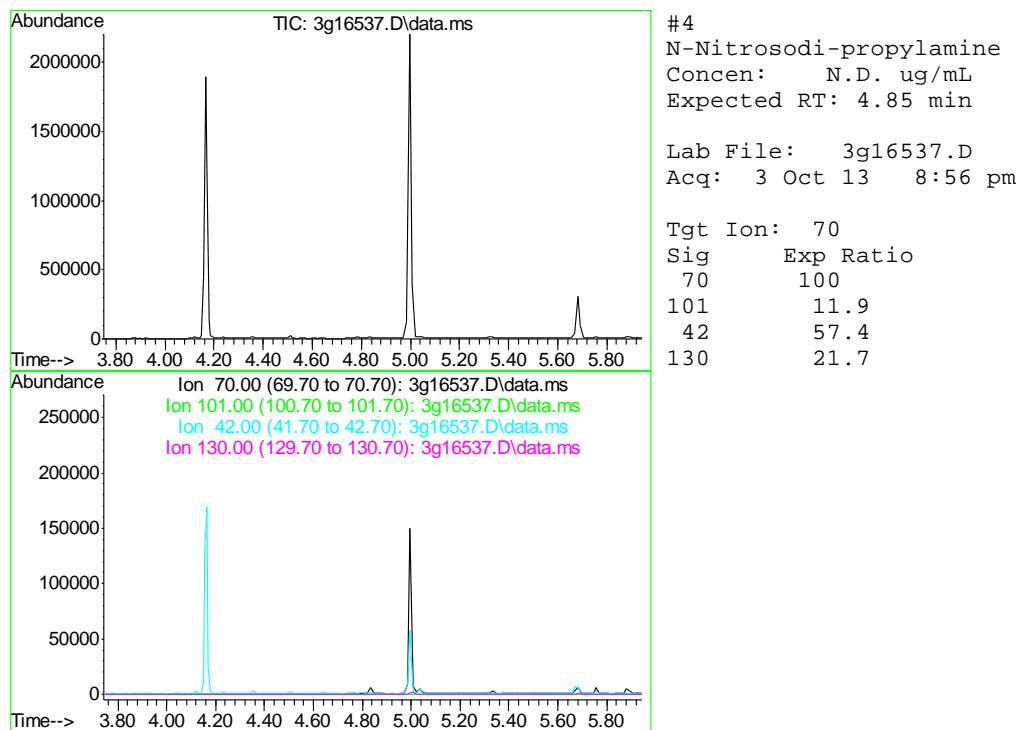
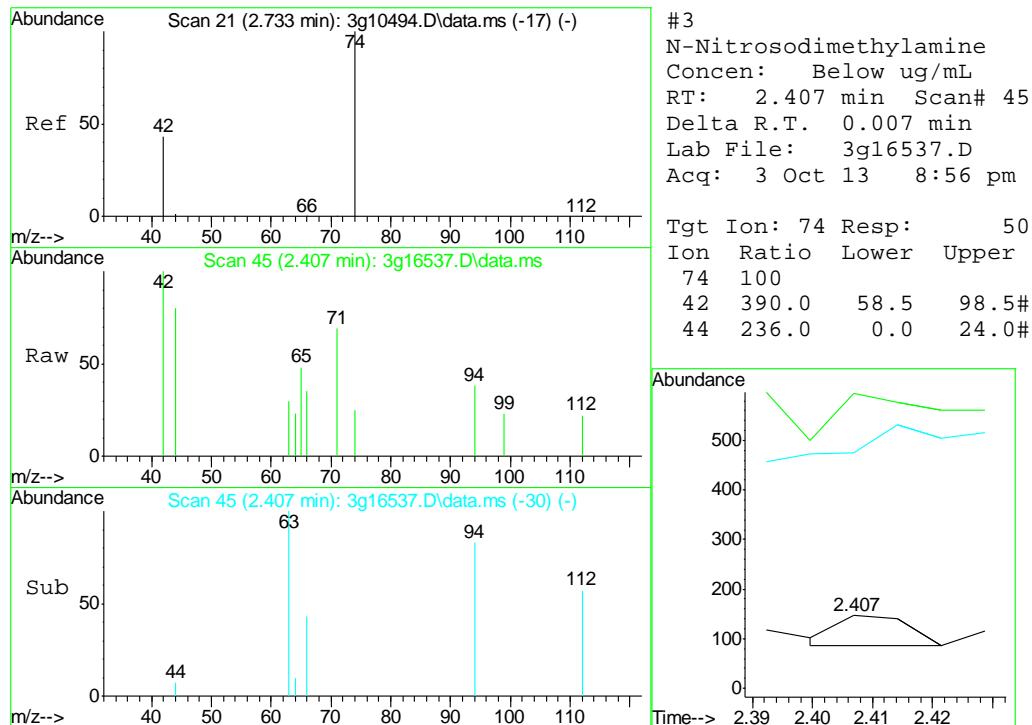
Quantitation Report (QT Reviewed)

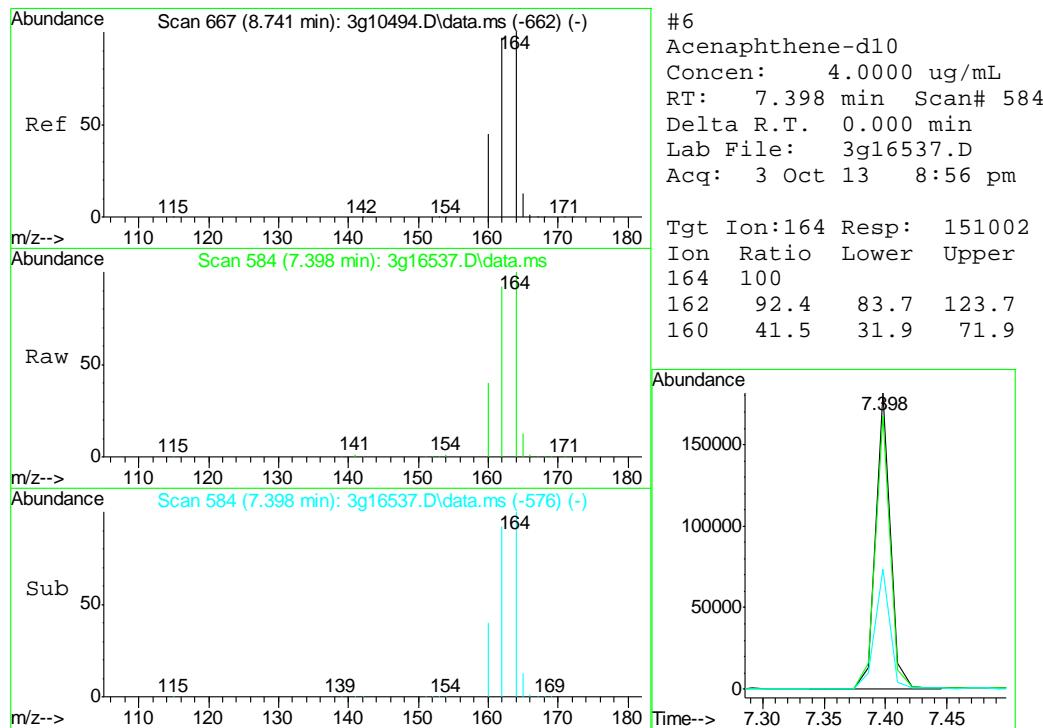
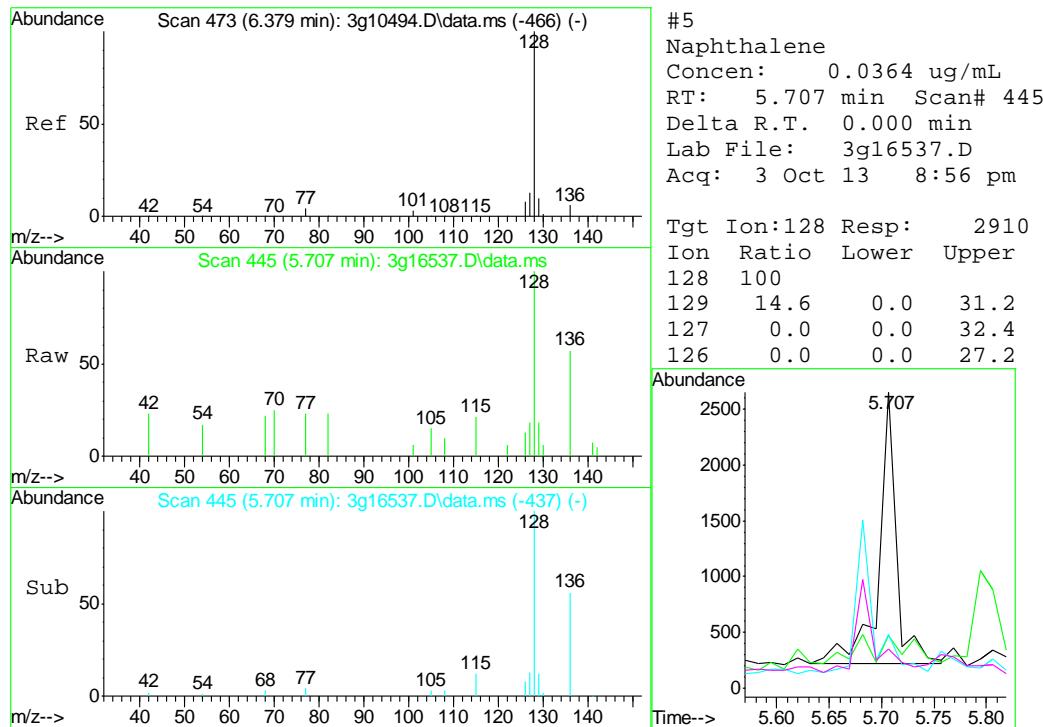
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 Data File : 3g16537.D
 Acq On : 3 Oct 2013 8:56 pm
 Operator : DONC
 Sample : D51203-1
 Misc : OP8670,E3G817,30.13,,,1,1
 ALS Vial : 24 Sample Multiplier: 1

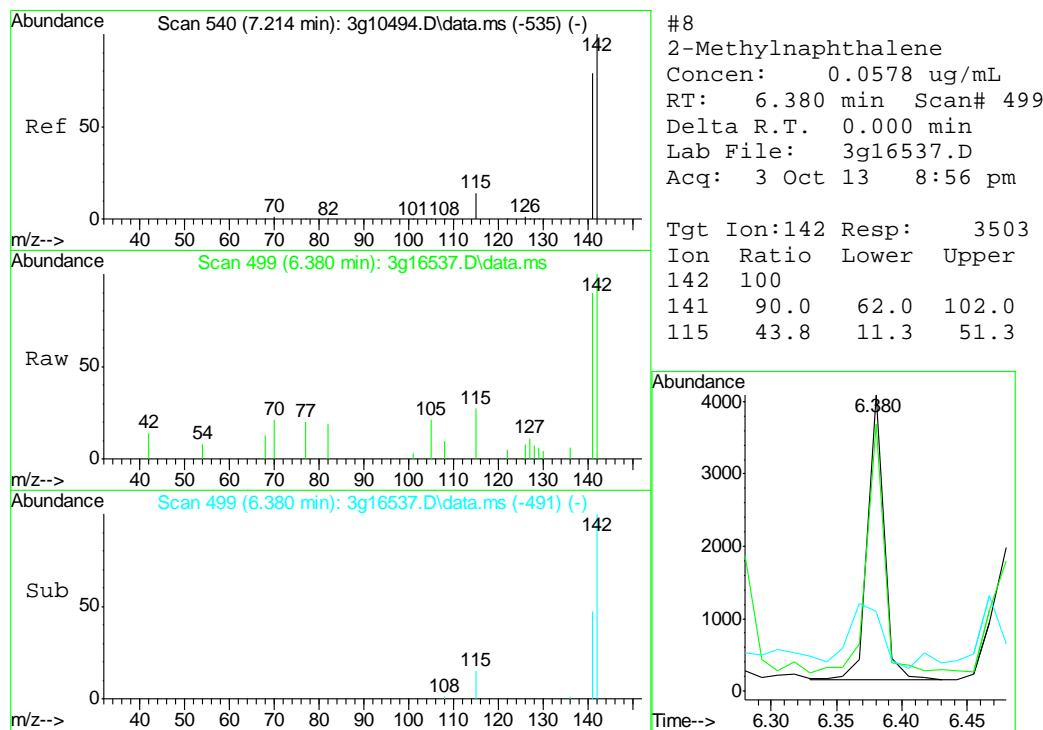
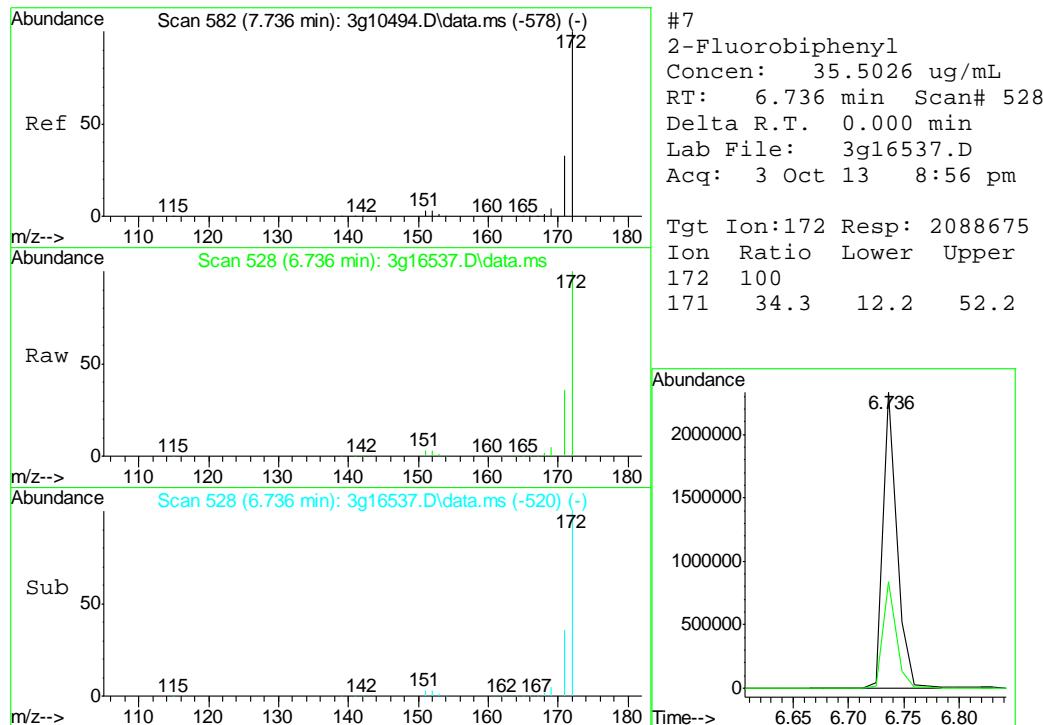
Quant Time: Oct 04 15:02:40 2013
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G810.M
 Quant Title : PAHSIM BASE
 QLast Update : Tue Sep 24 08:29:29 2013
 Response via : Initial Calibration

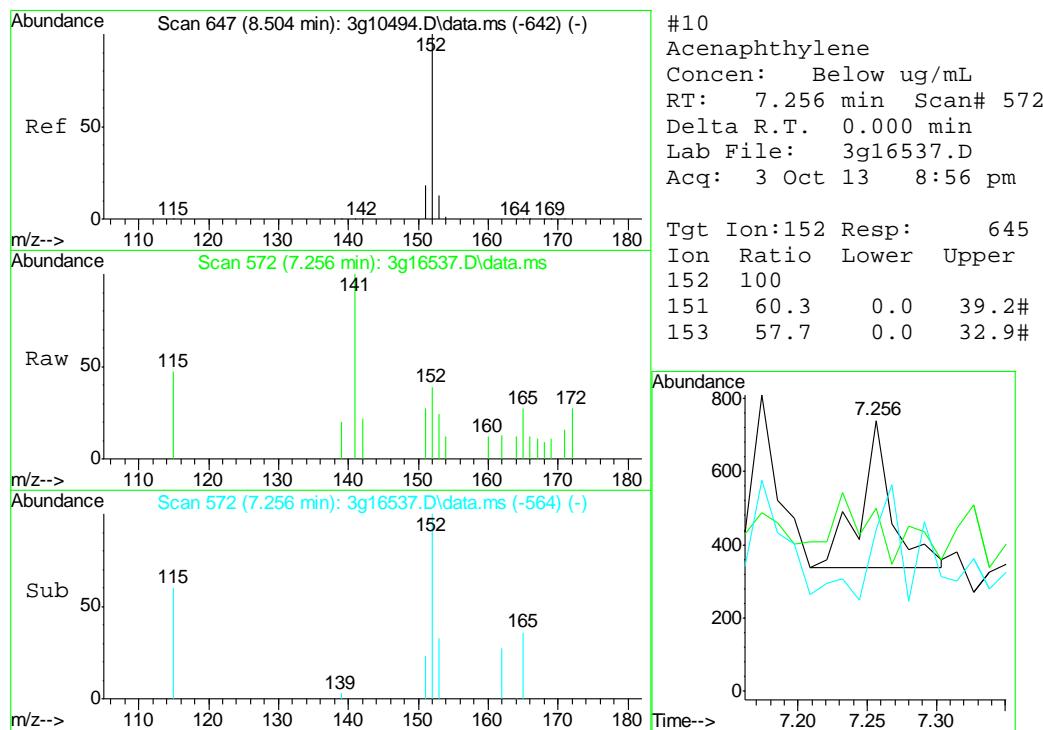
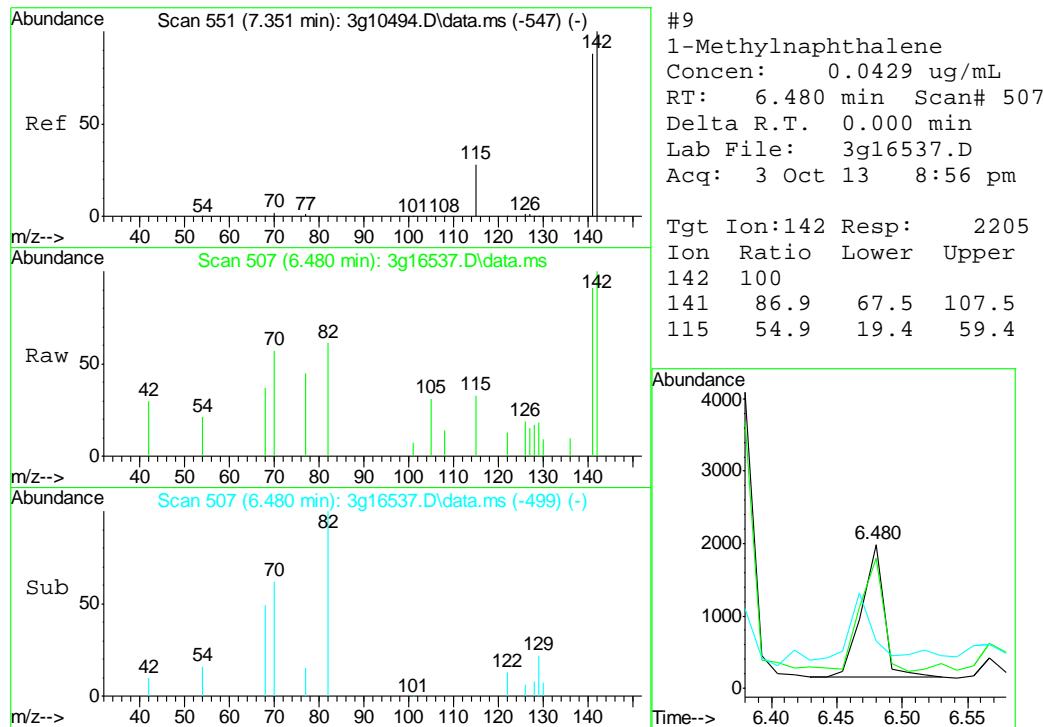


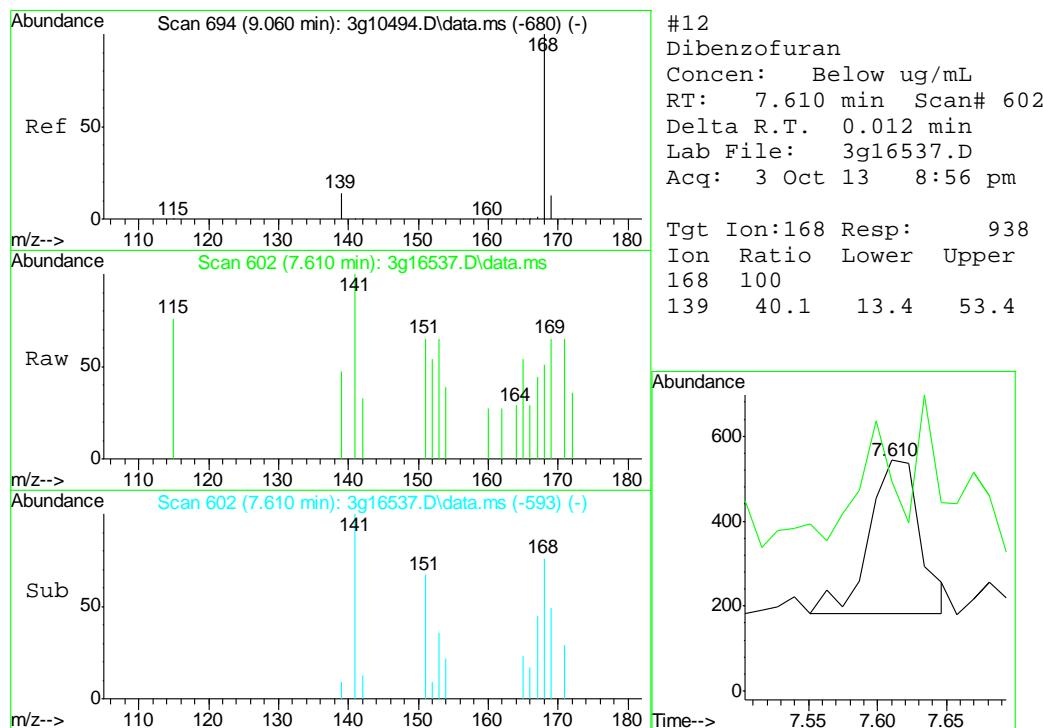
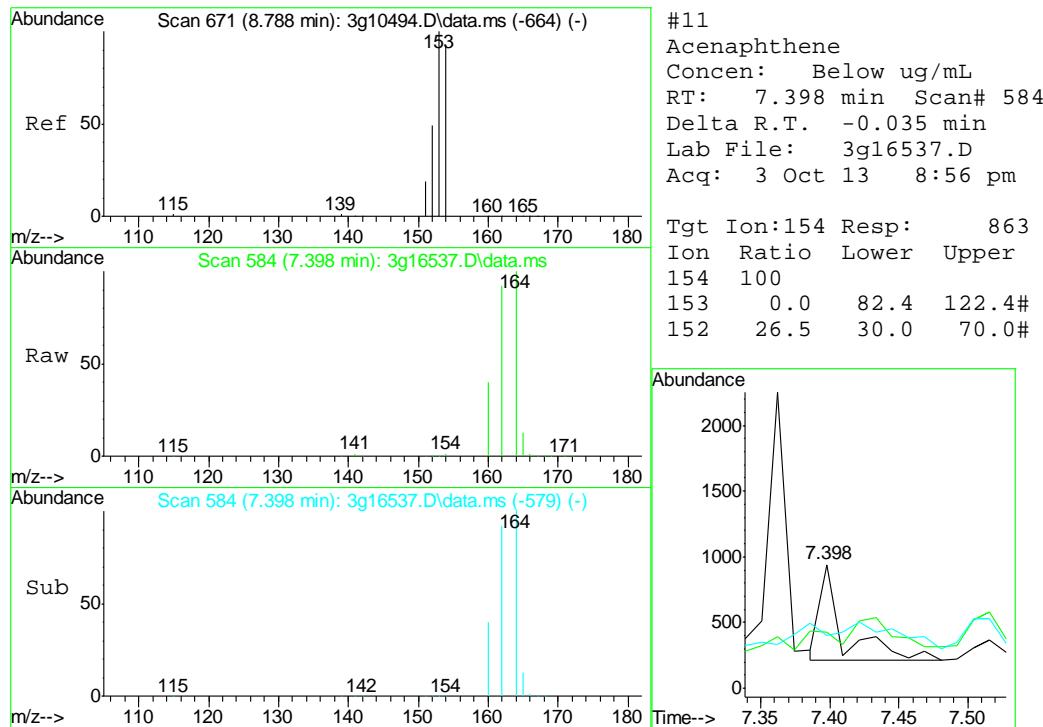


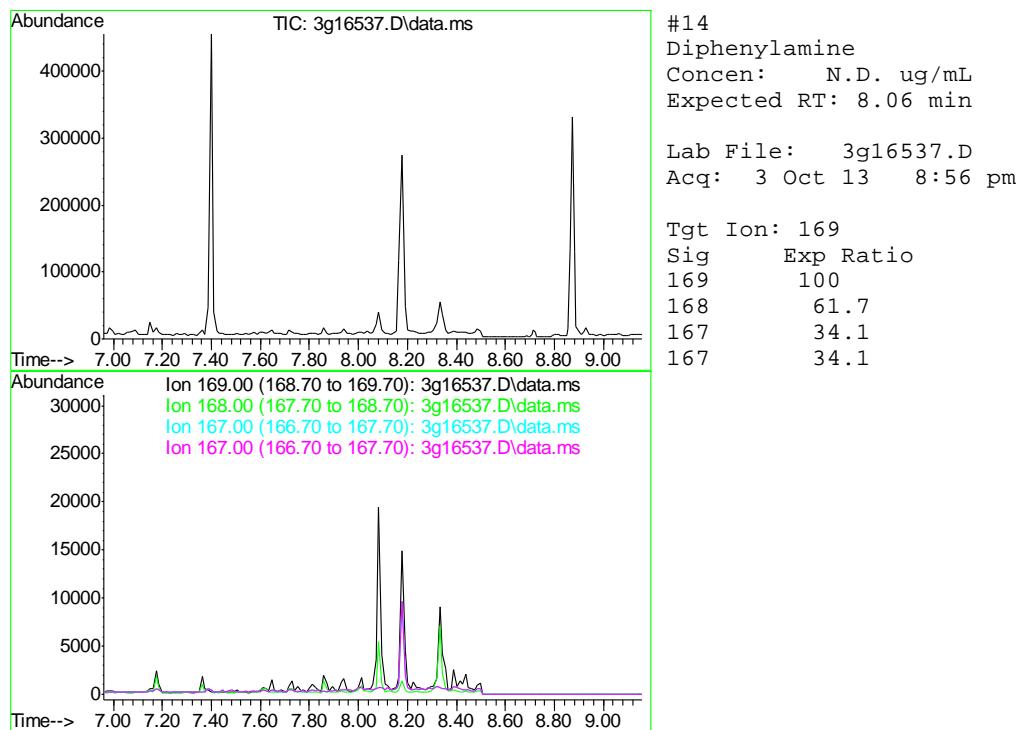
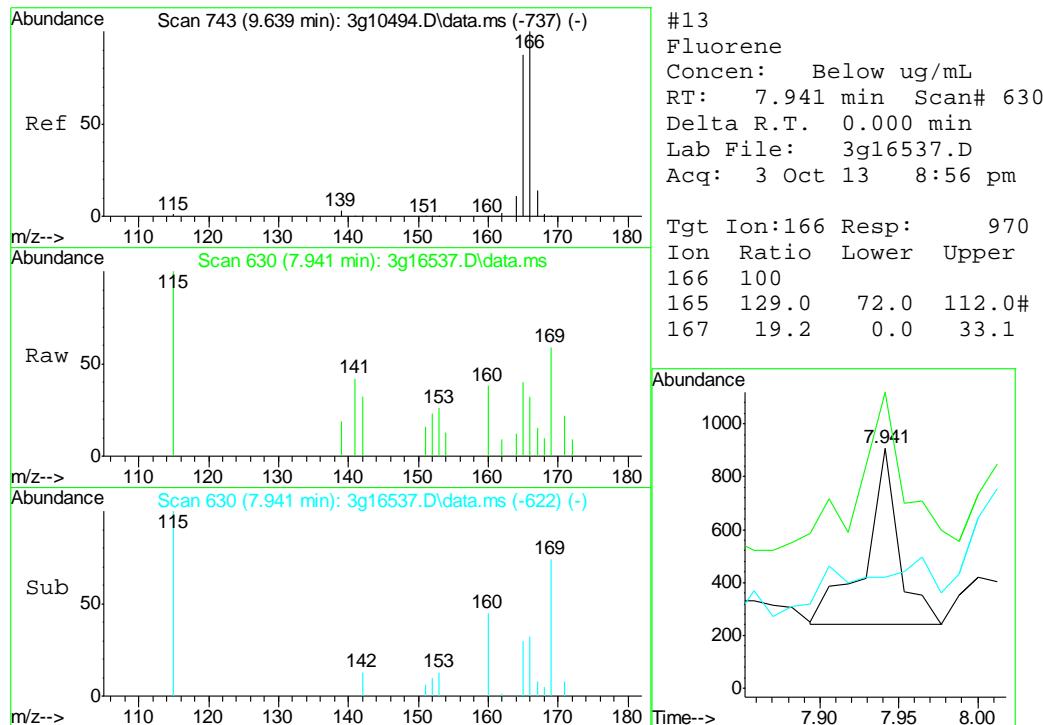


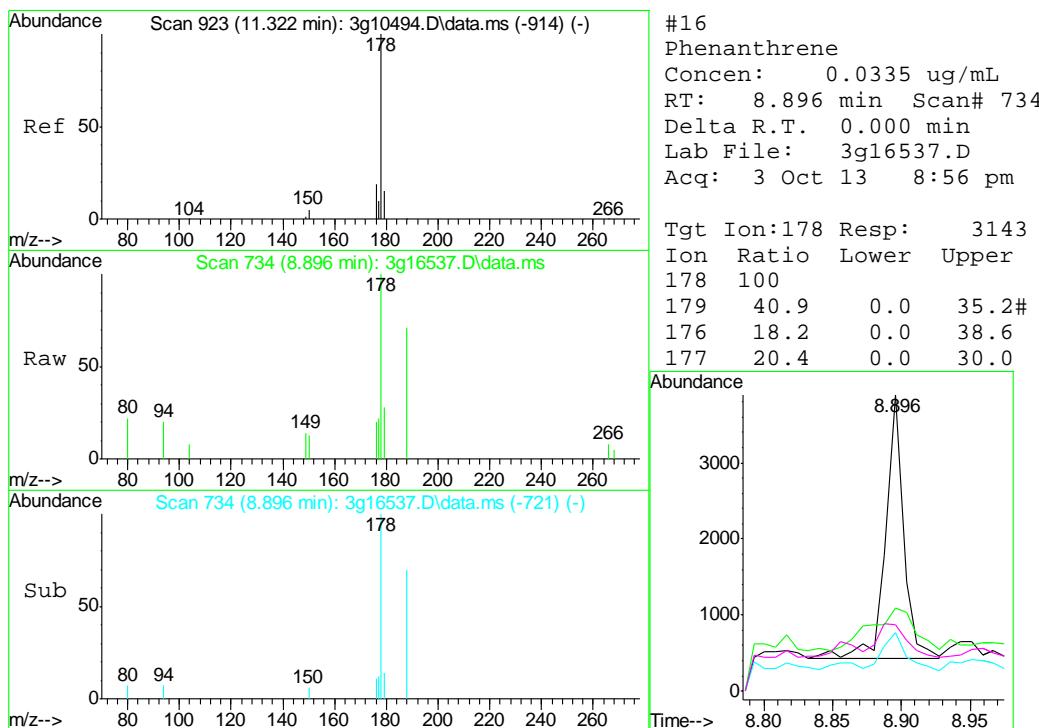
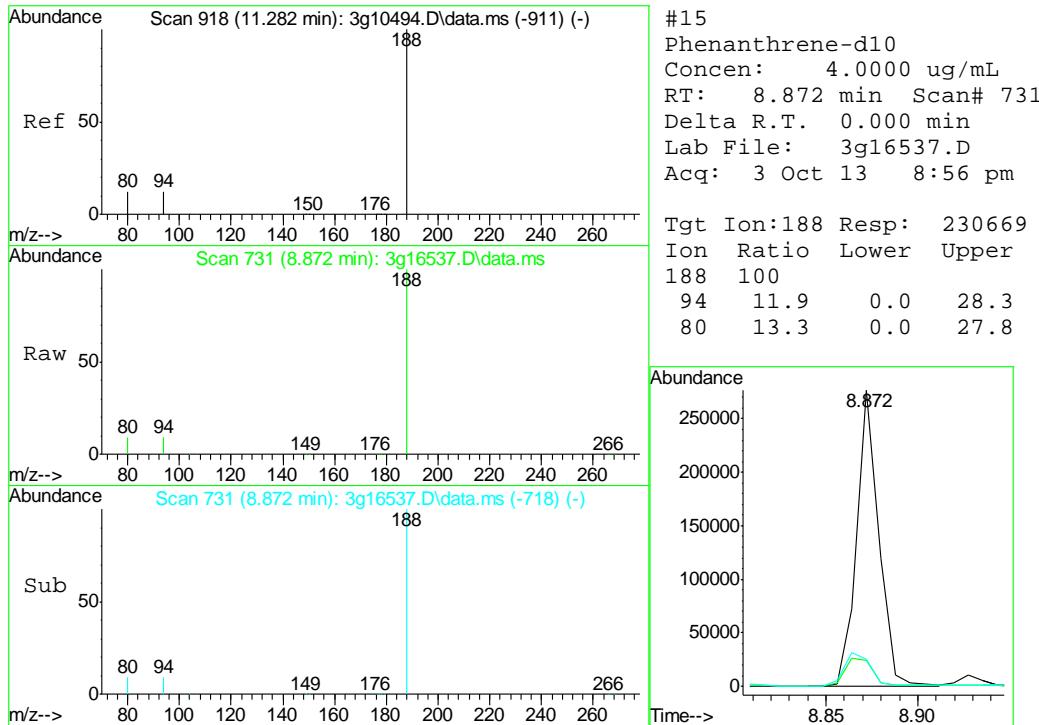


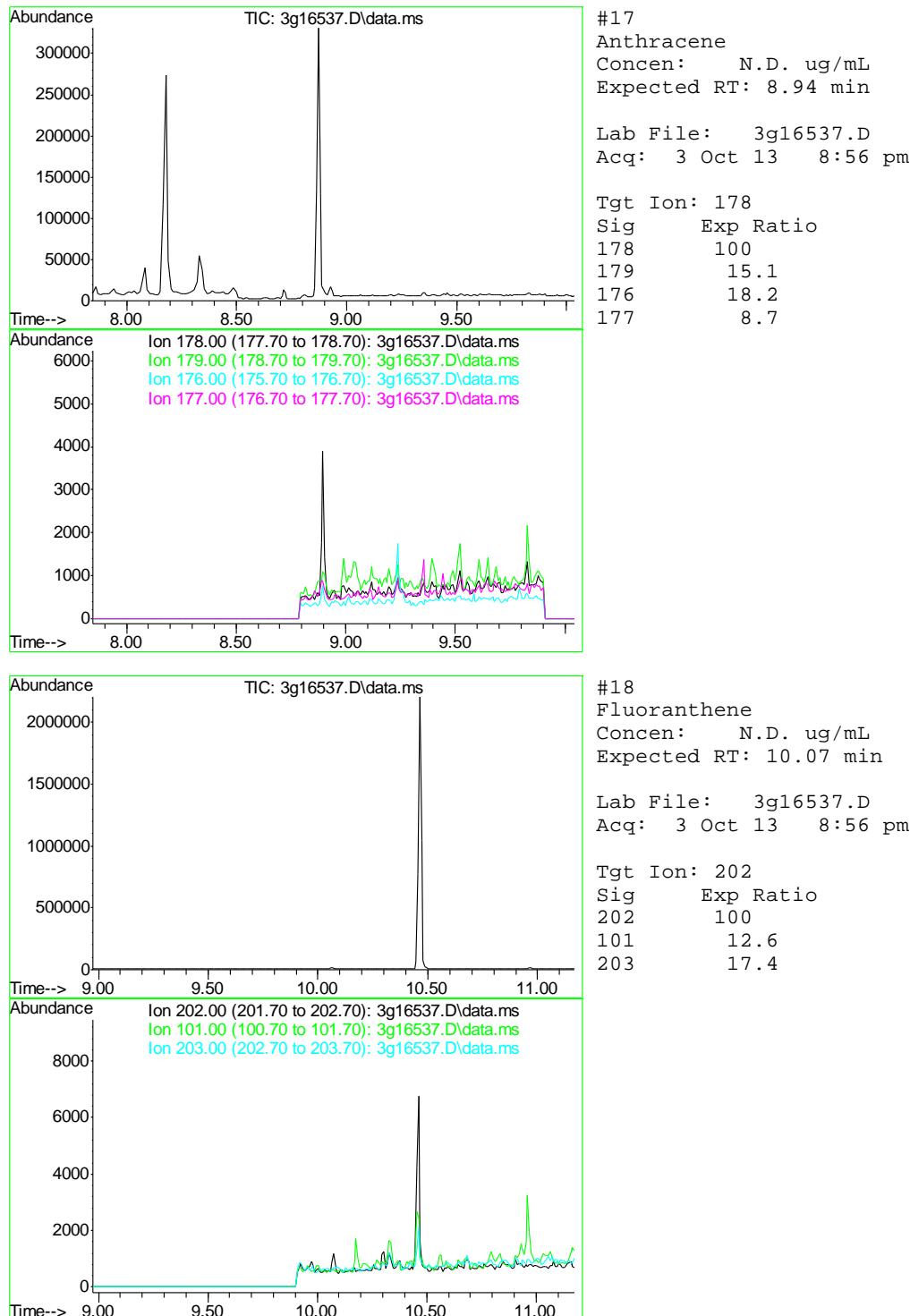


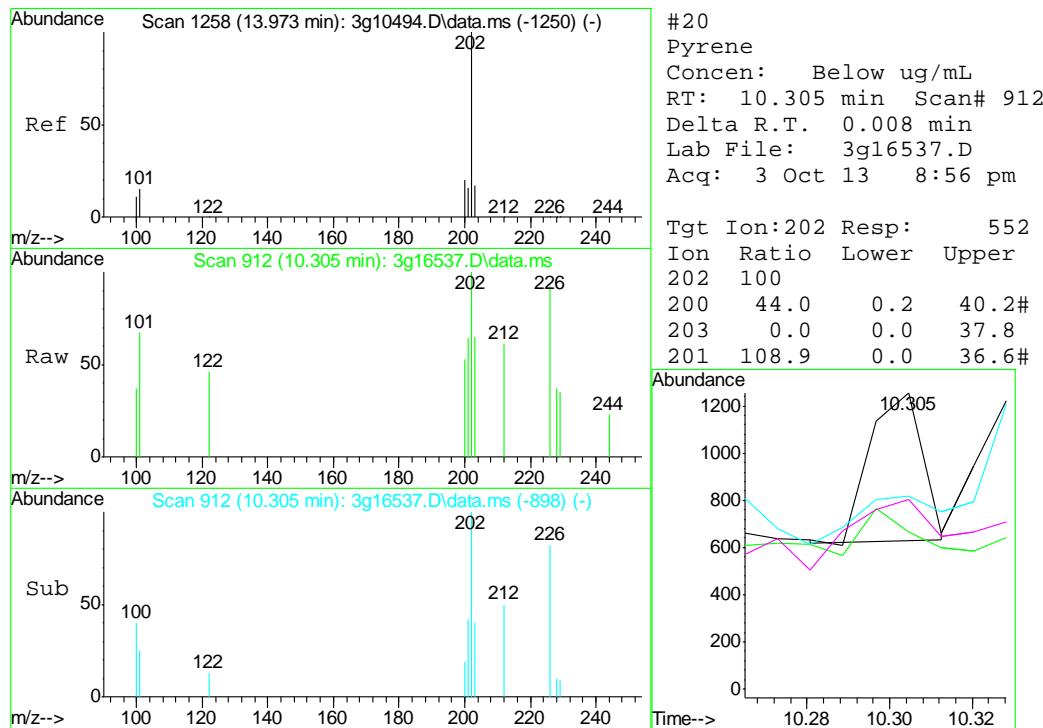
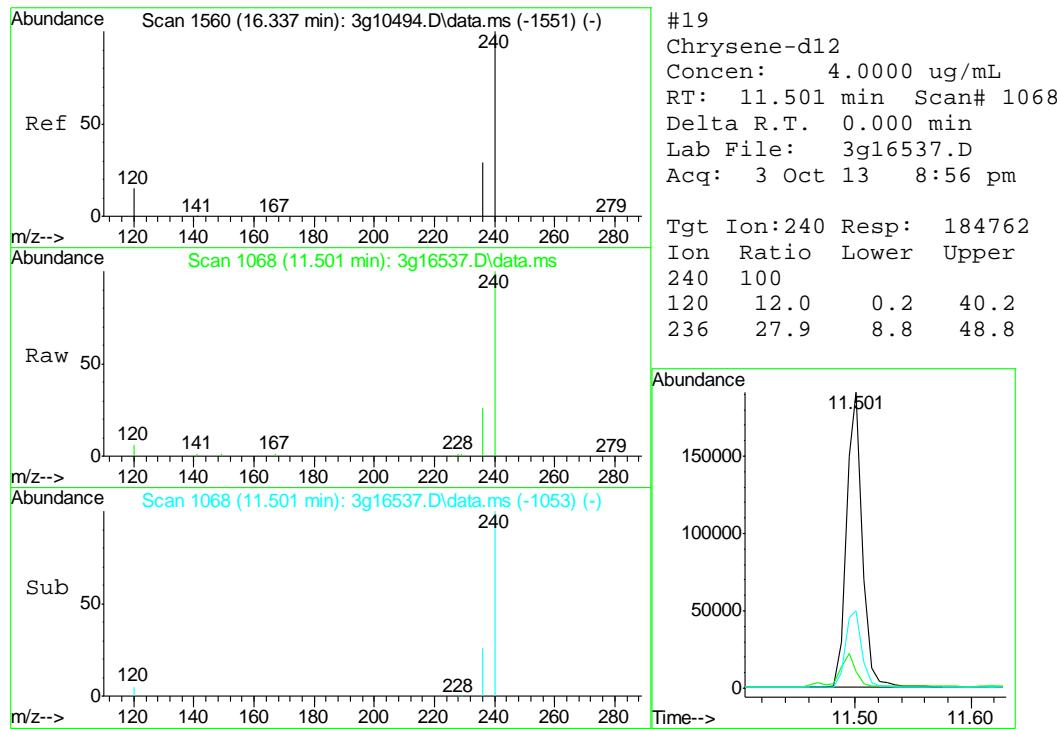


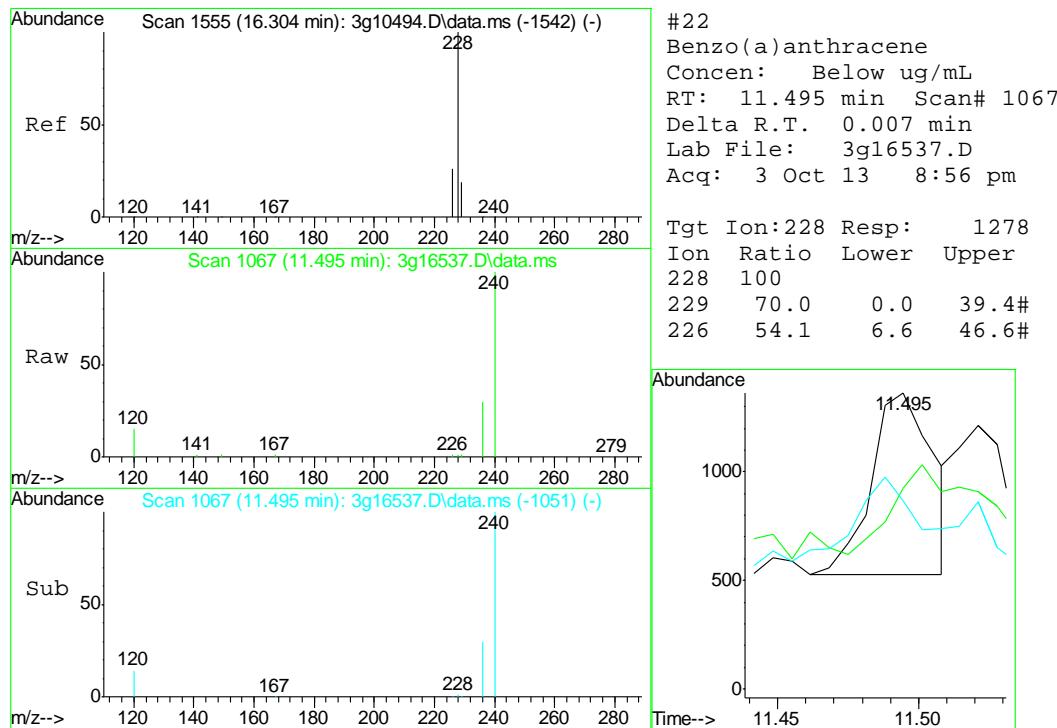
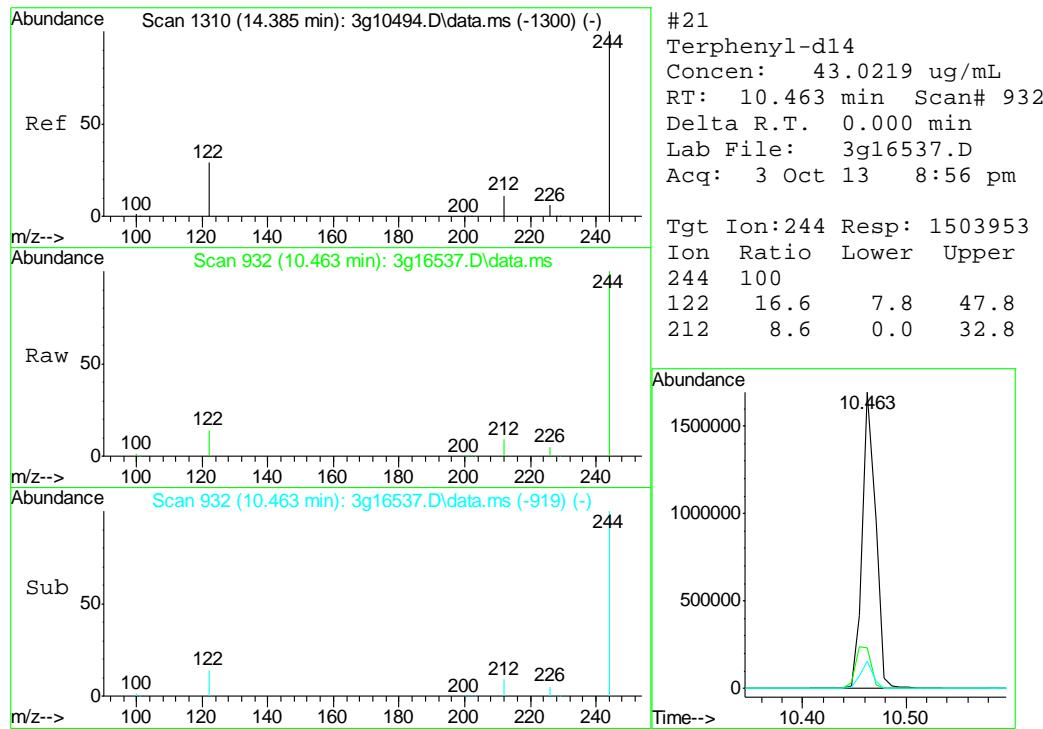


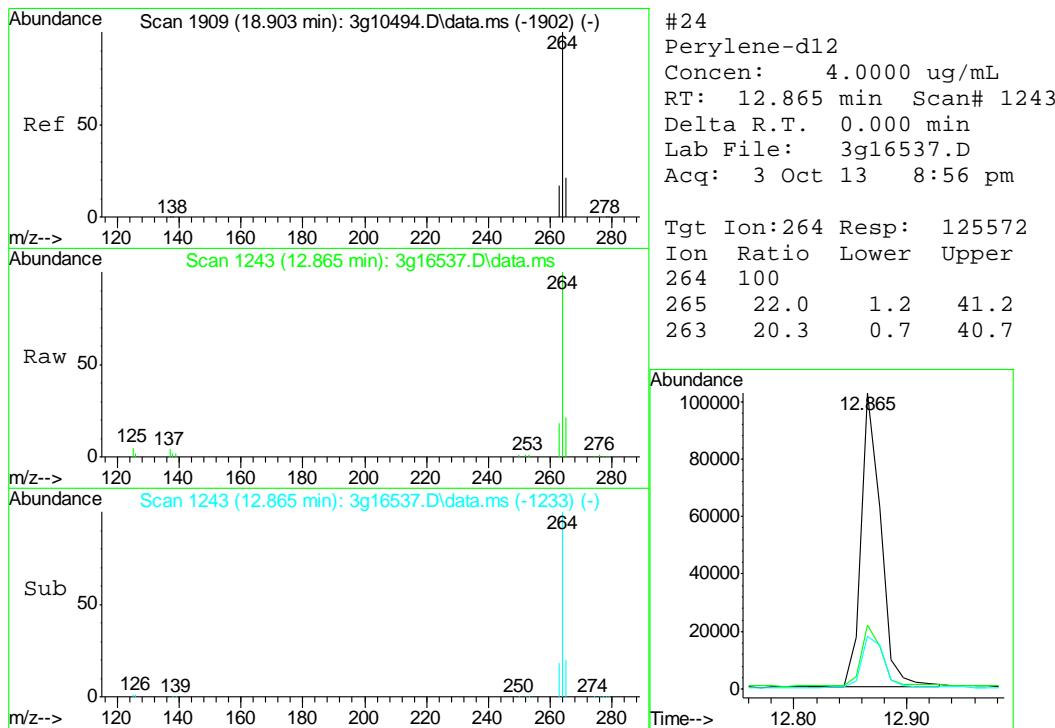
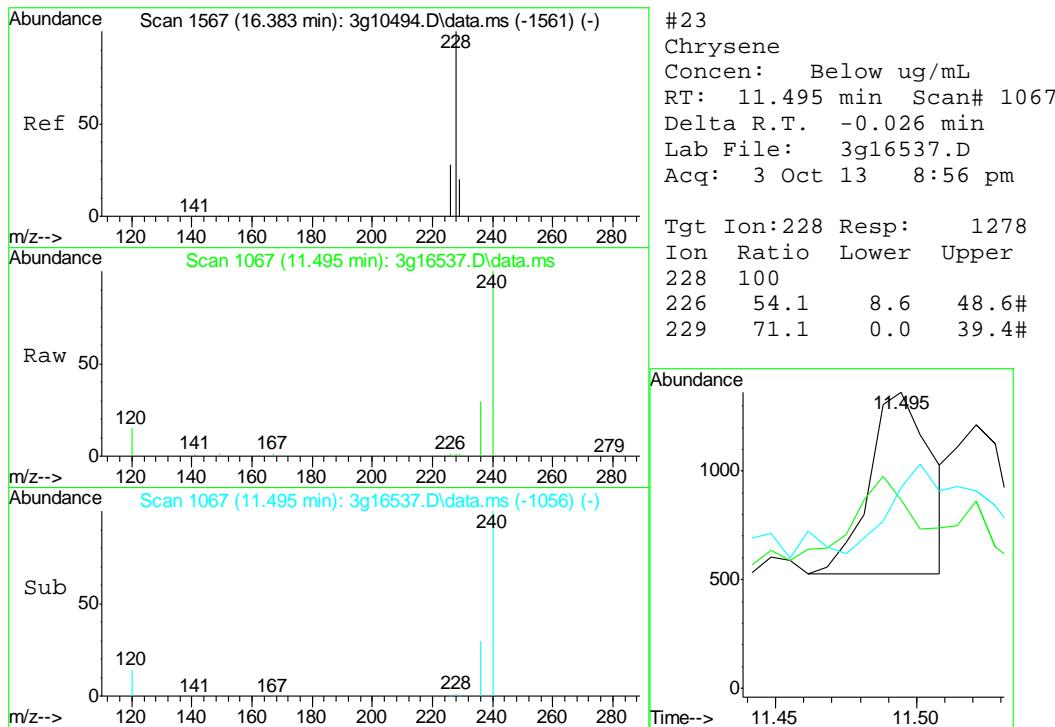


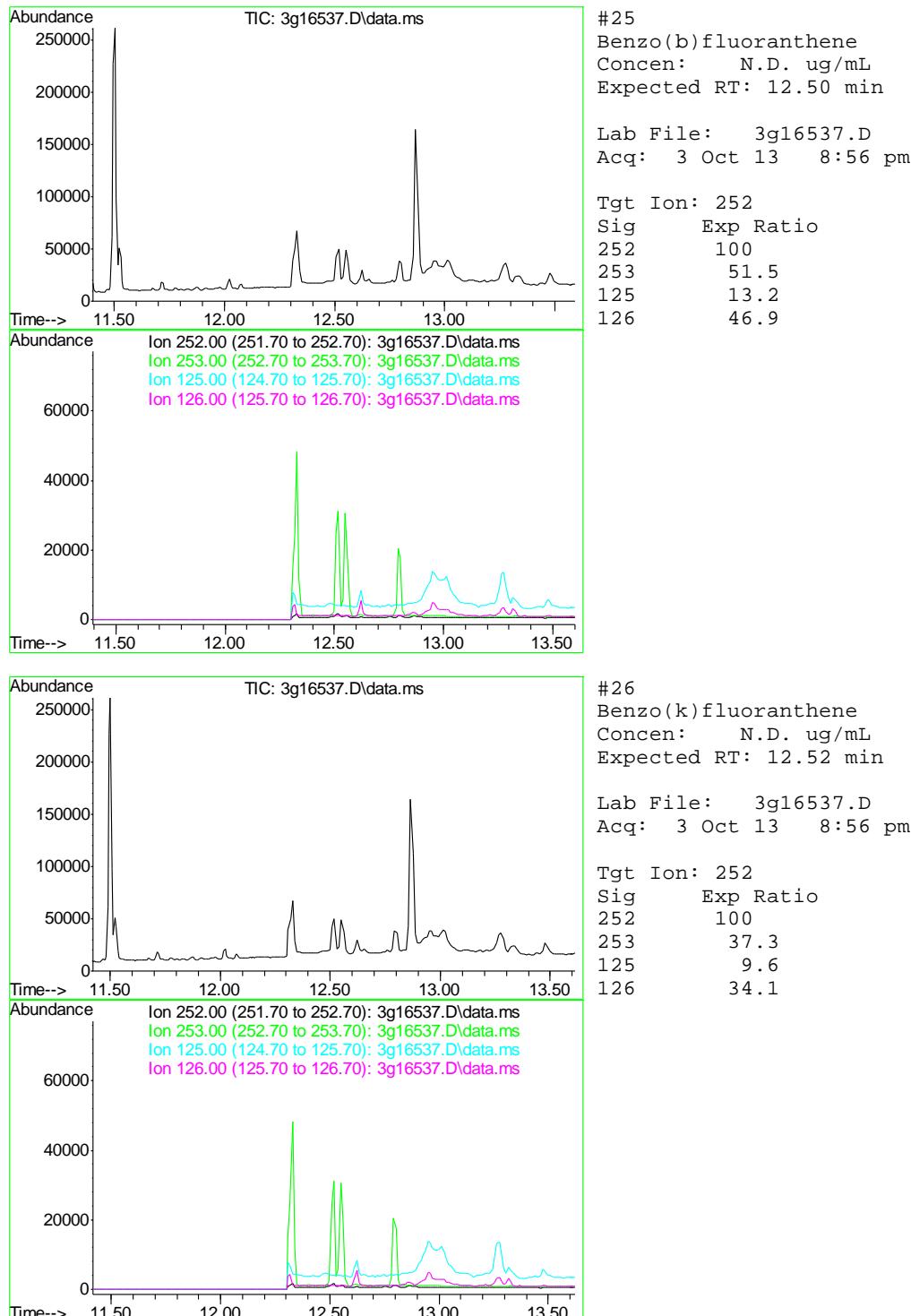


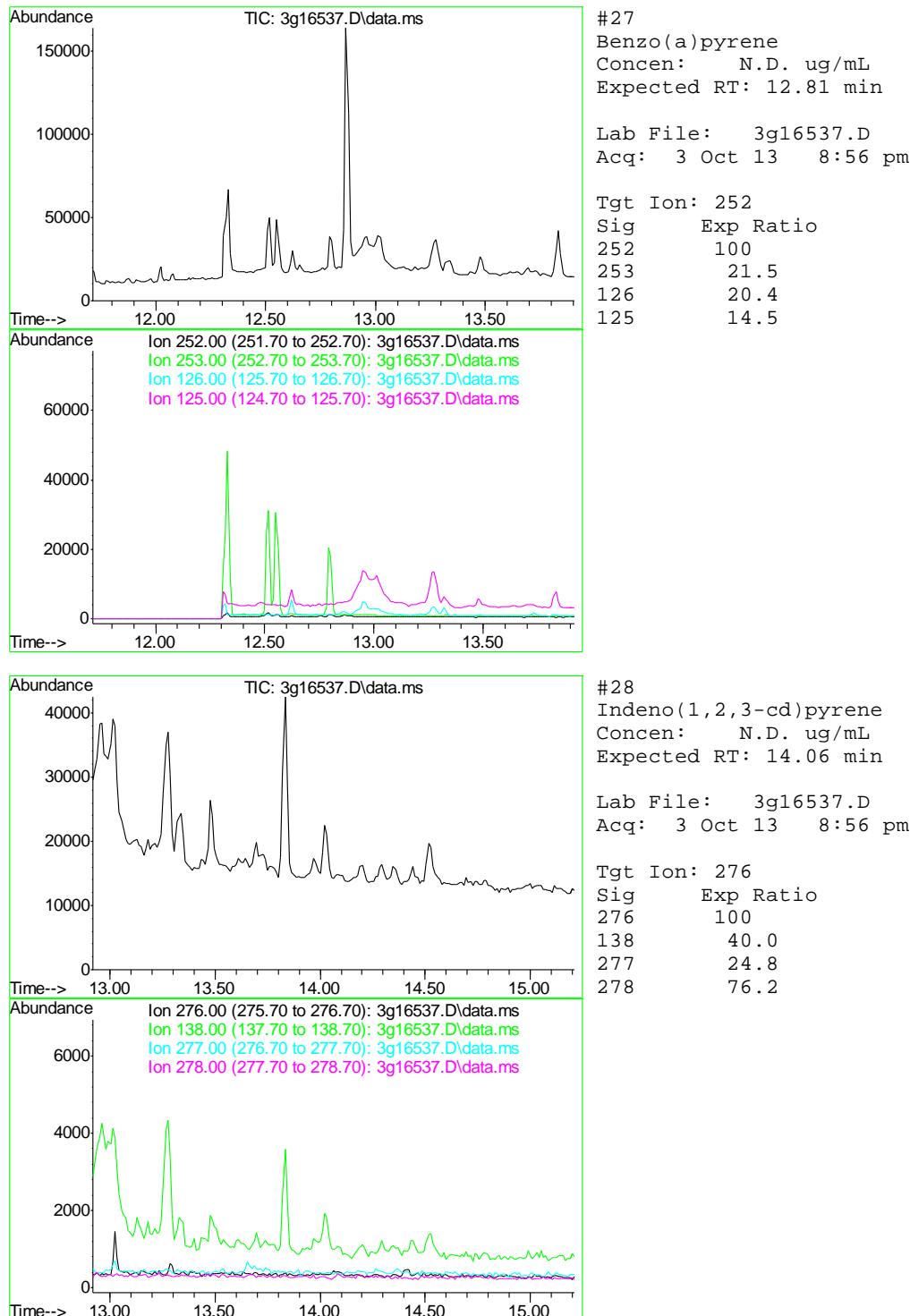


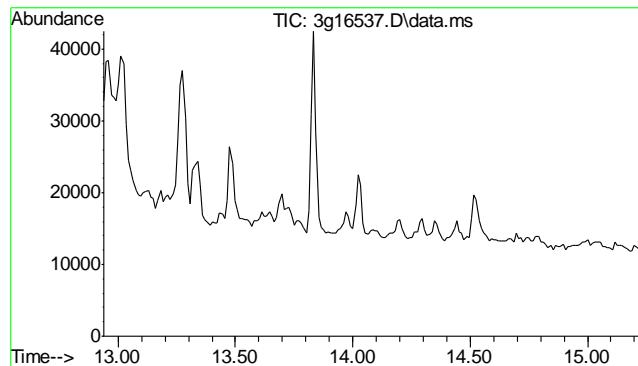








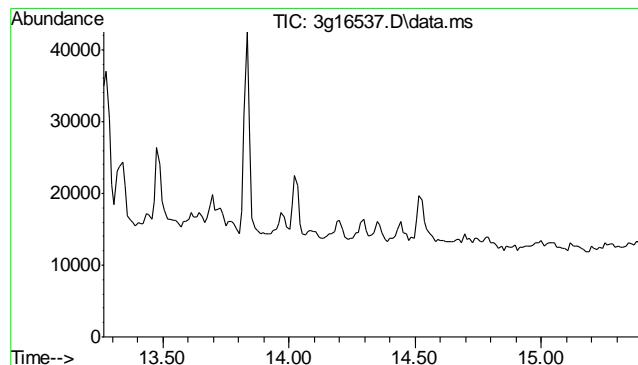
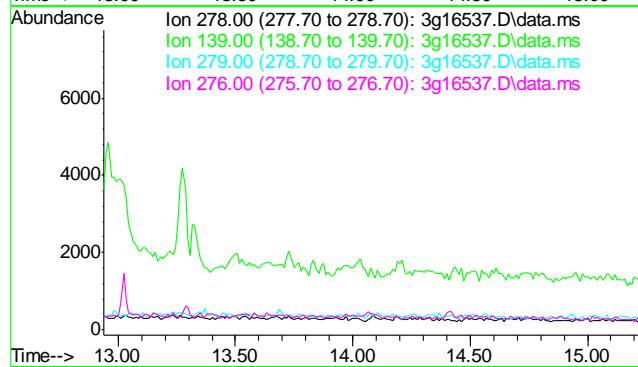




#29
 Dibenz(a,h)anthracene
 Concen: N.D. ug/mL
 Expected RT: 14.09 min

Lab File: 3g16537.D
 Acq: 3 Oct 13 8:56 pm

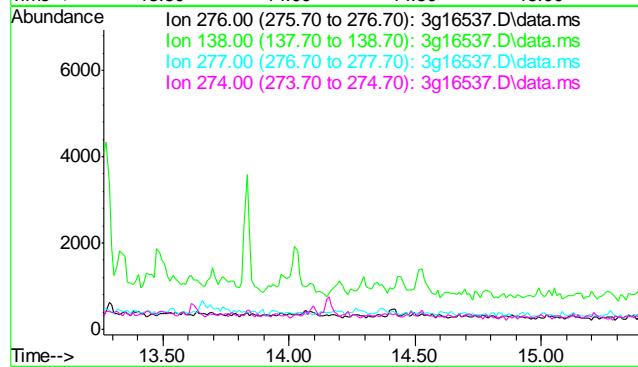
Tgt Ion:	Sig	Exp Ratio
278	100	
139	30.8	
279	22.9	
276	131.2	



#30
 Benzo(g,h,i)perylene
 Concen: N.D. ug/mL
 Expected RT: 14.41 min

Lab File: 3g16537.D
 Acq: 3 Oct 13 8:56 pm

Tgt Ion:	Sig	Exp Ratio
276	100	
138	35.1	
277	23.3	
274	21.5	



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%	

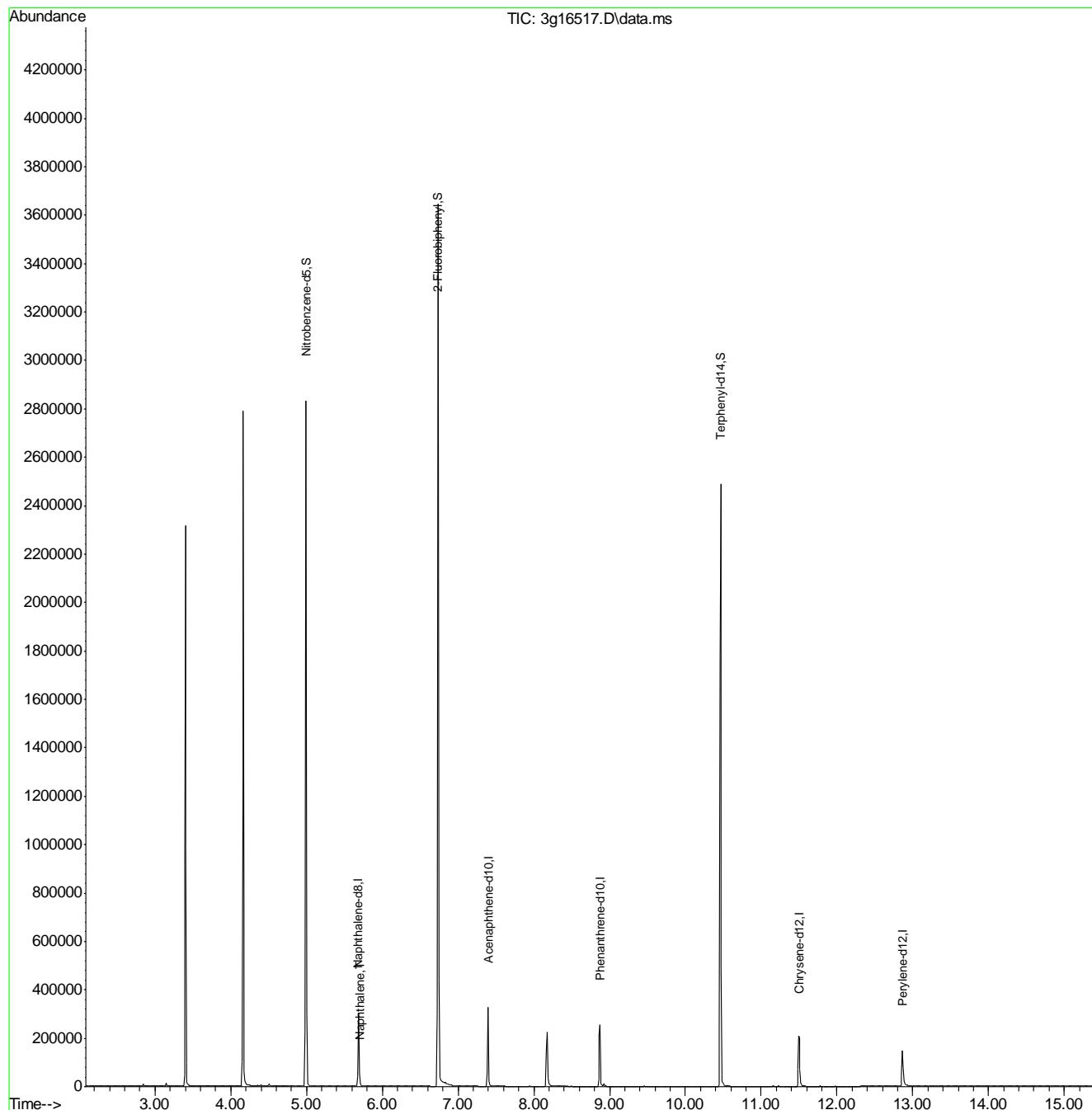
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D.	d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D.	d
5) Naphthalene	5.707	128	6089	0.0773	ug/mL 98
8) 2-Methylnaphthalene	6.380	142	1089	N.D.	
9) 1-Methylnaphthalene	6.480	142	479	N.D.	
10) Acenaphthylene	7.256	152	63	N.D.	
11) Acenaphthene	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

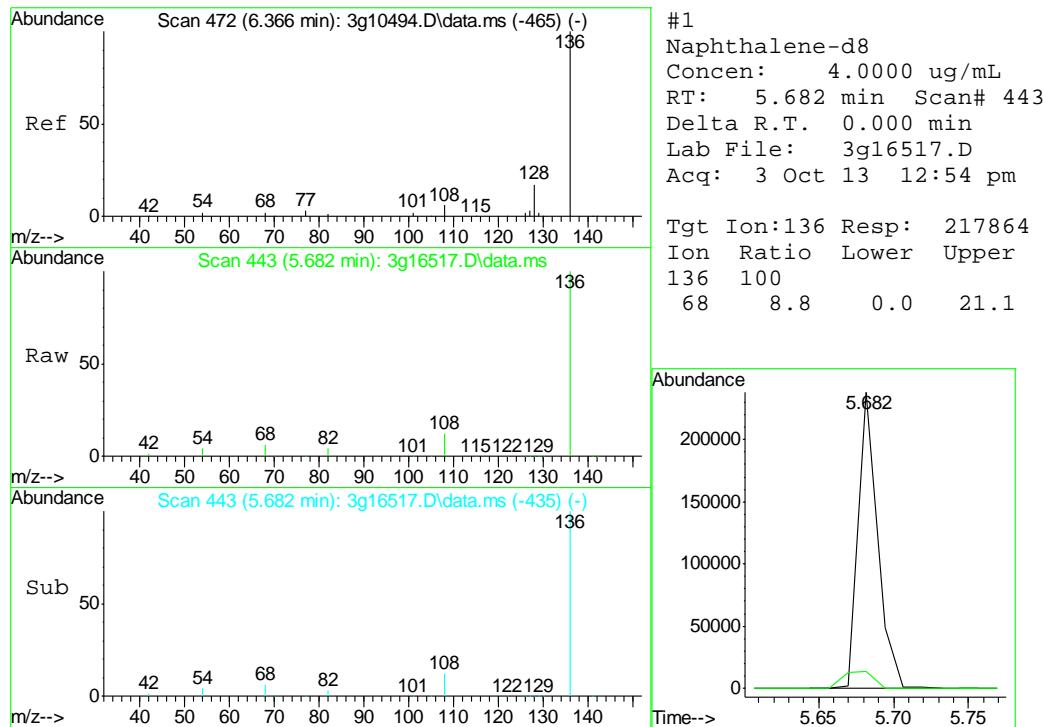
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

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

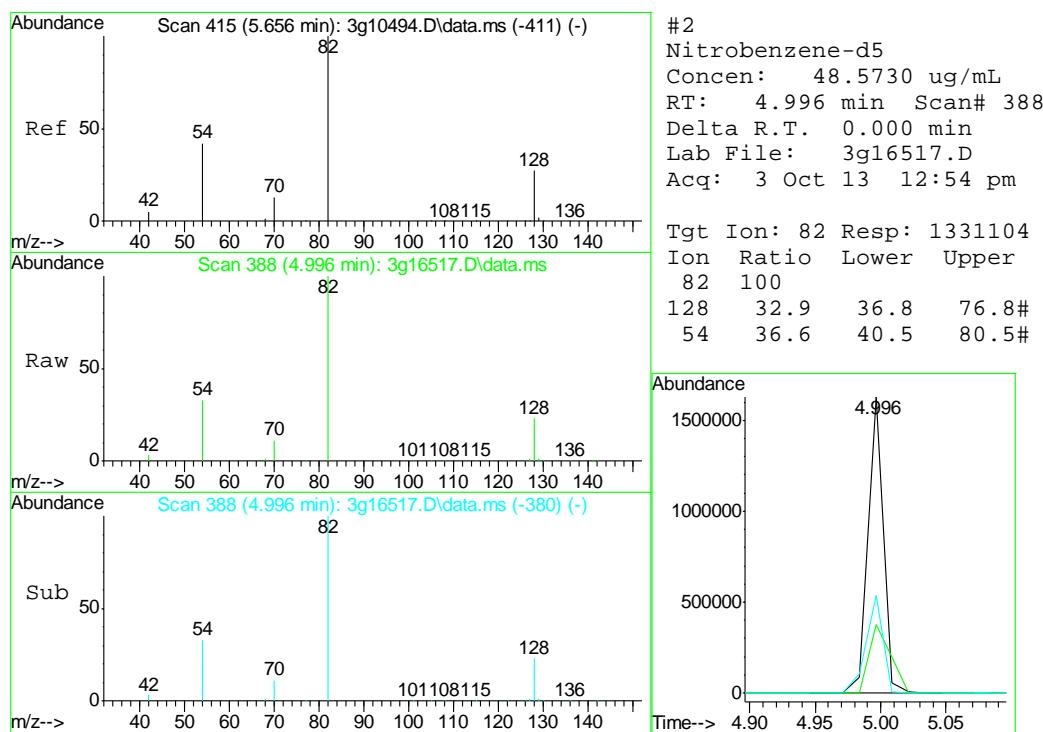
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 Quant Title : PAHSIM BASE
 QLast Update : Tue Sep 24 08:29:29 2013
 Response via : Initial Calibration

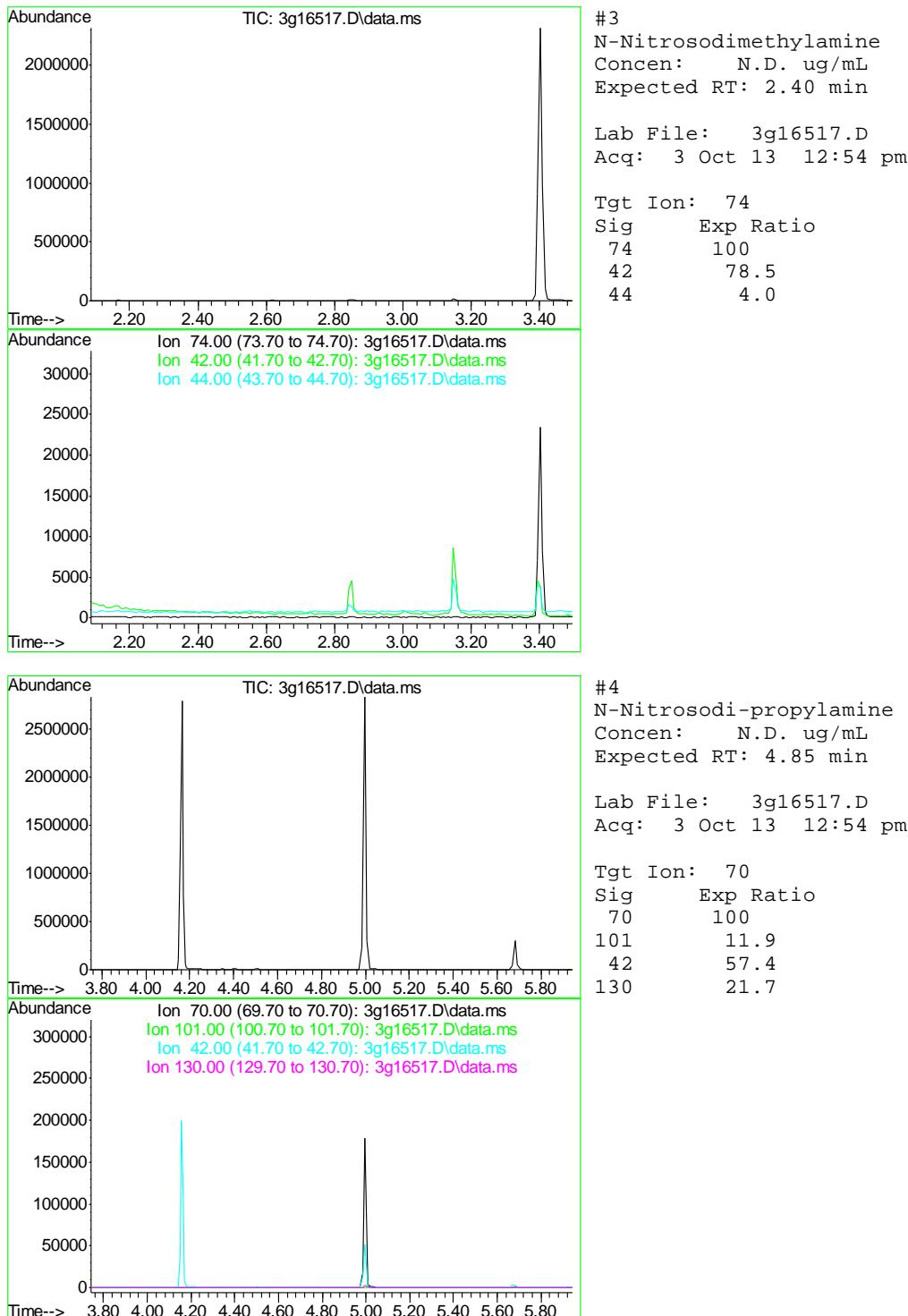


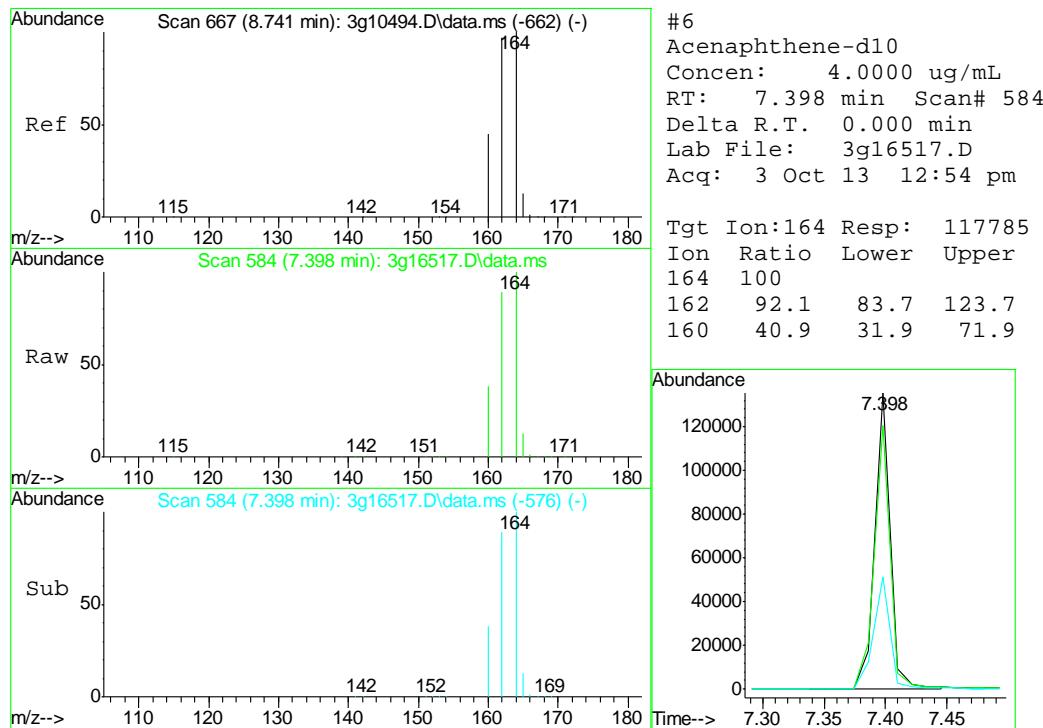
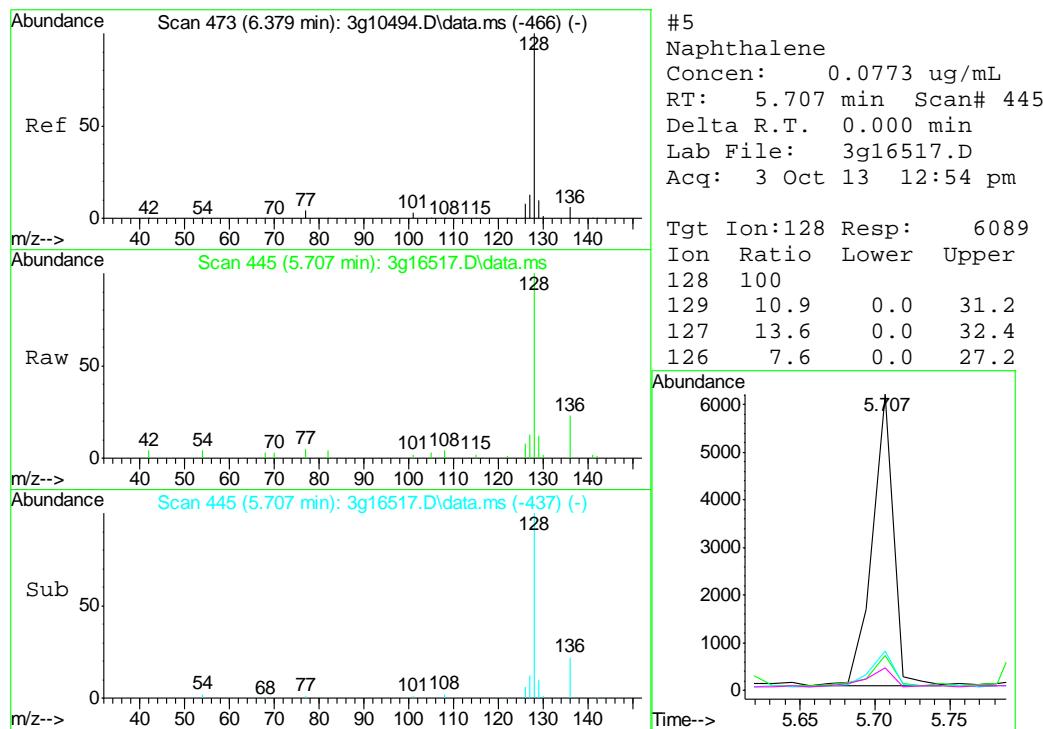


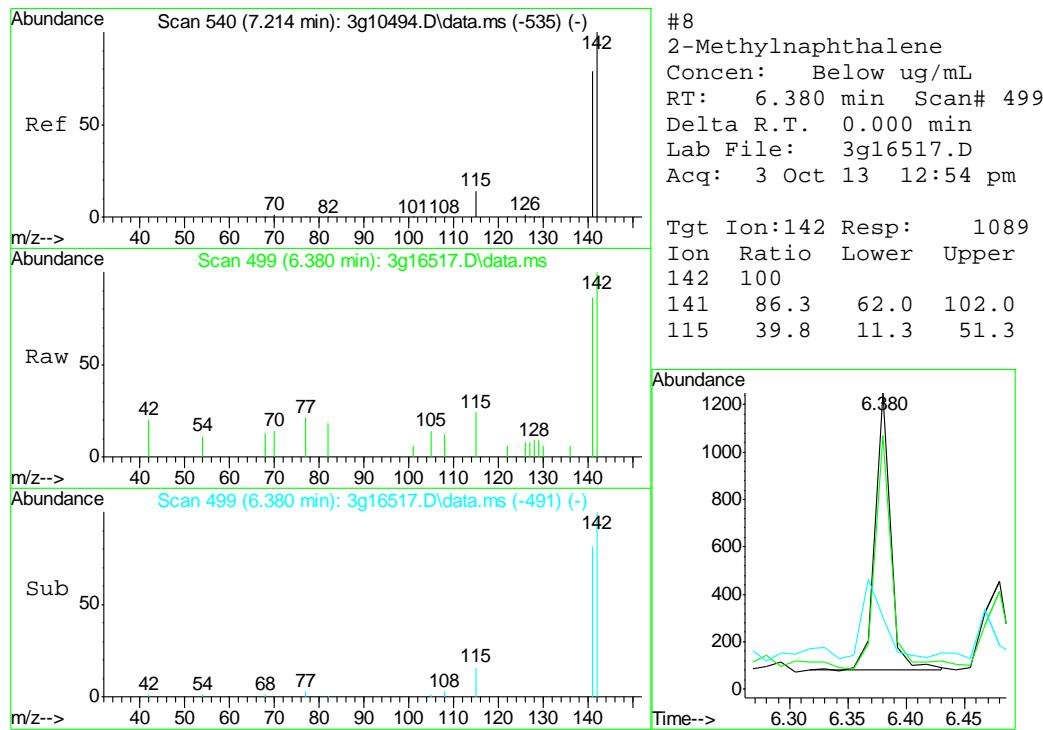
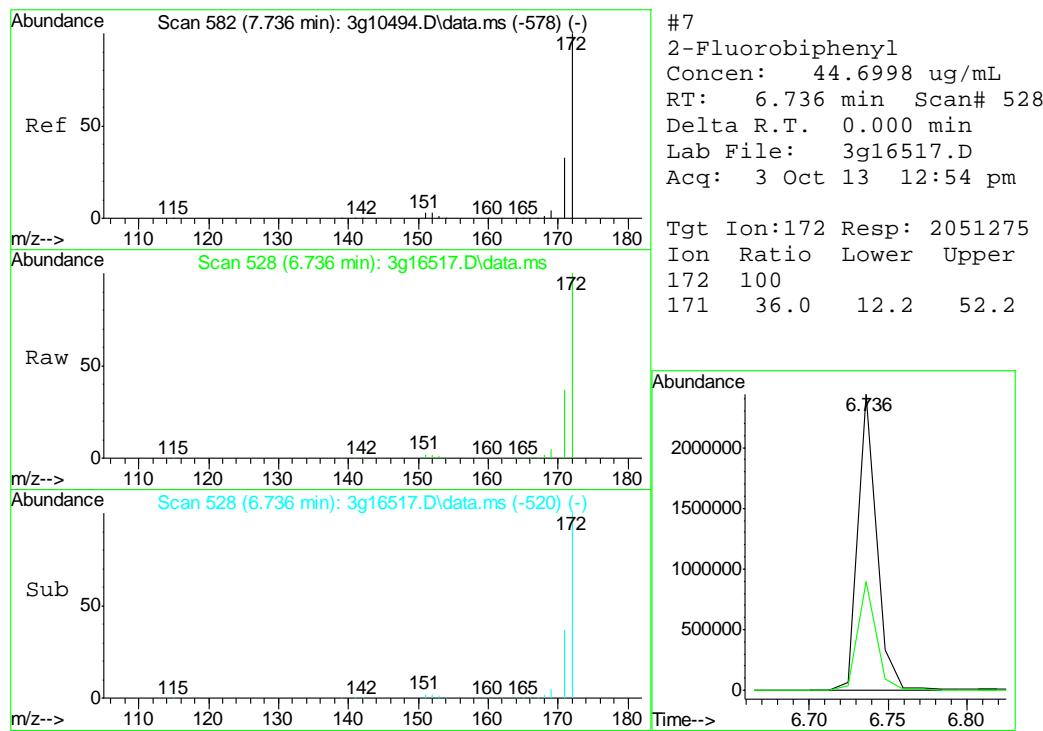
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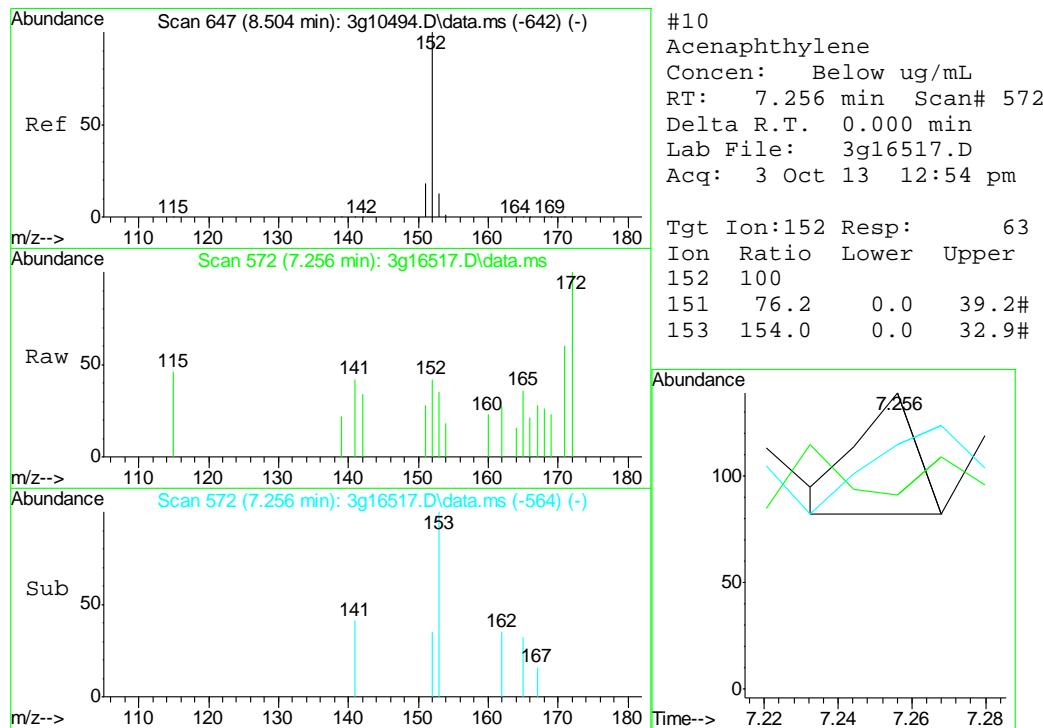
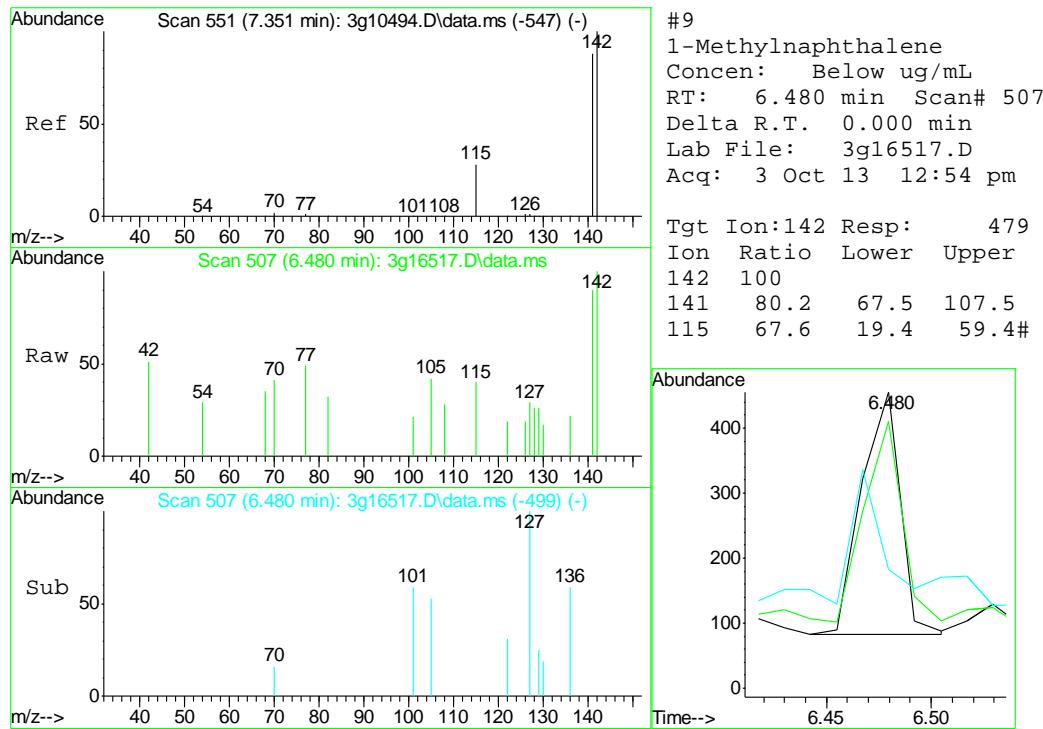
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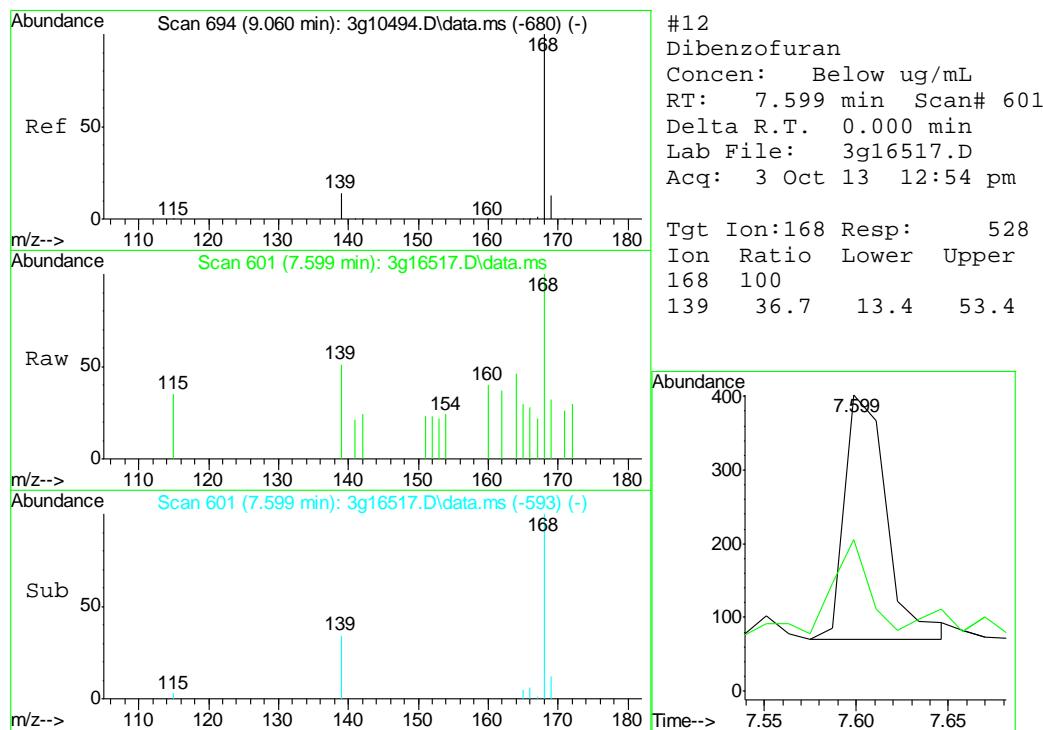
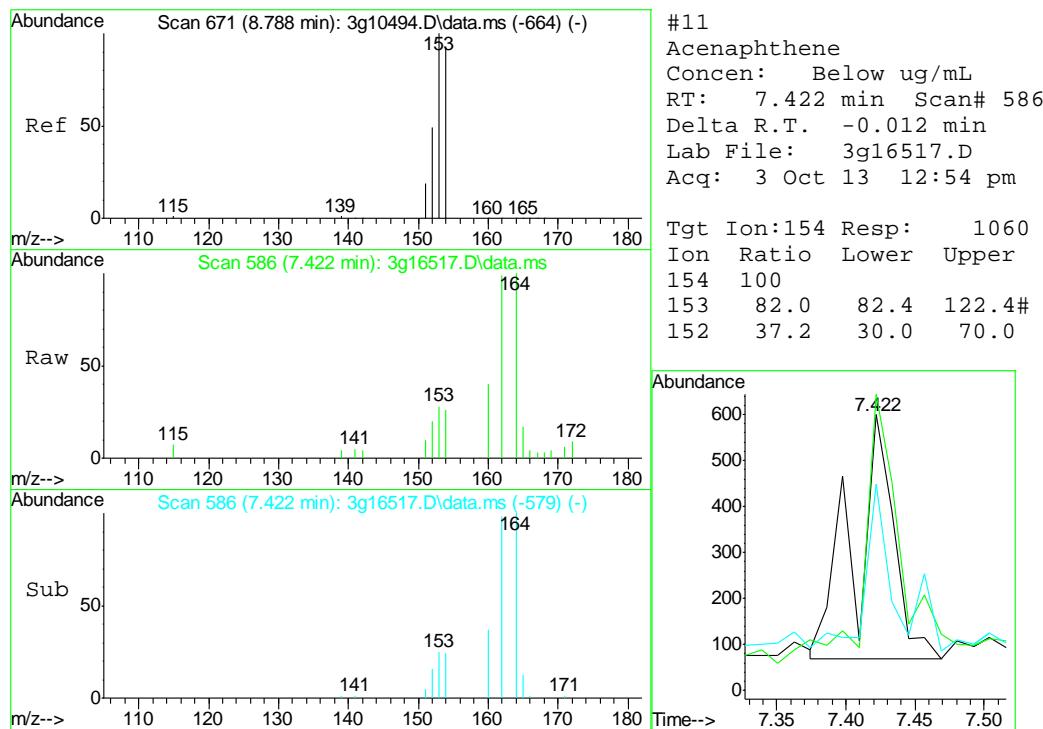


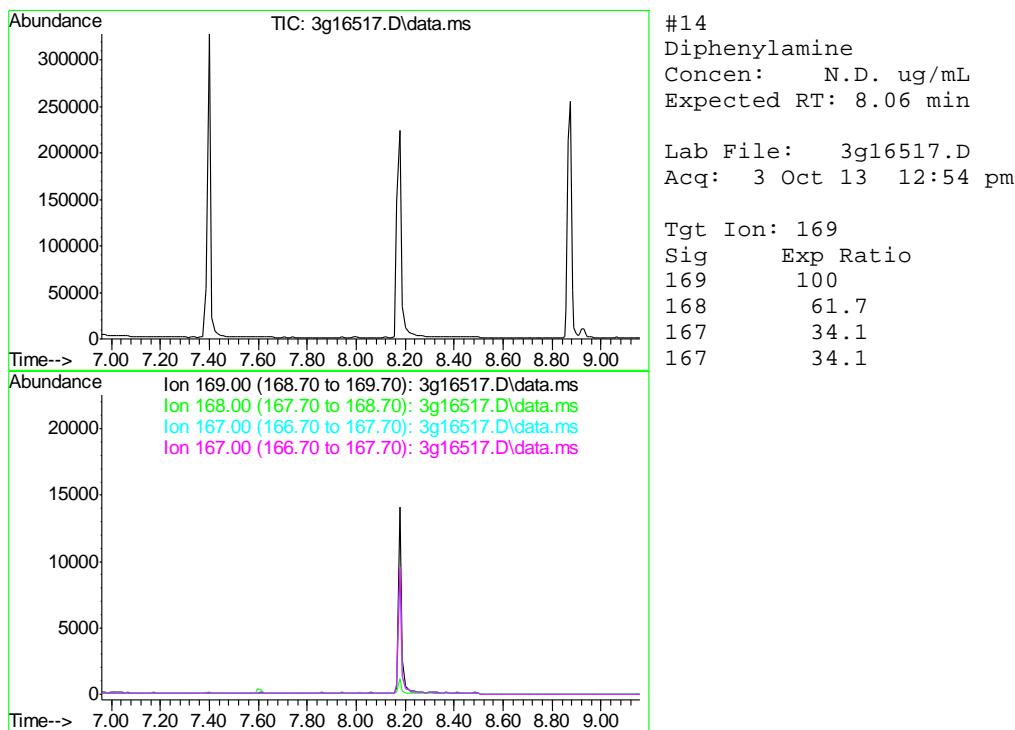
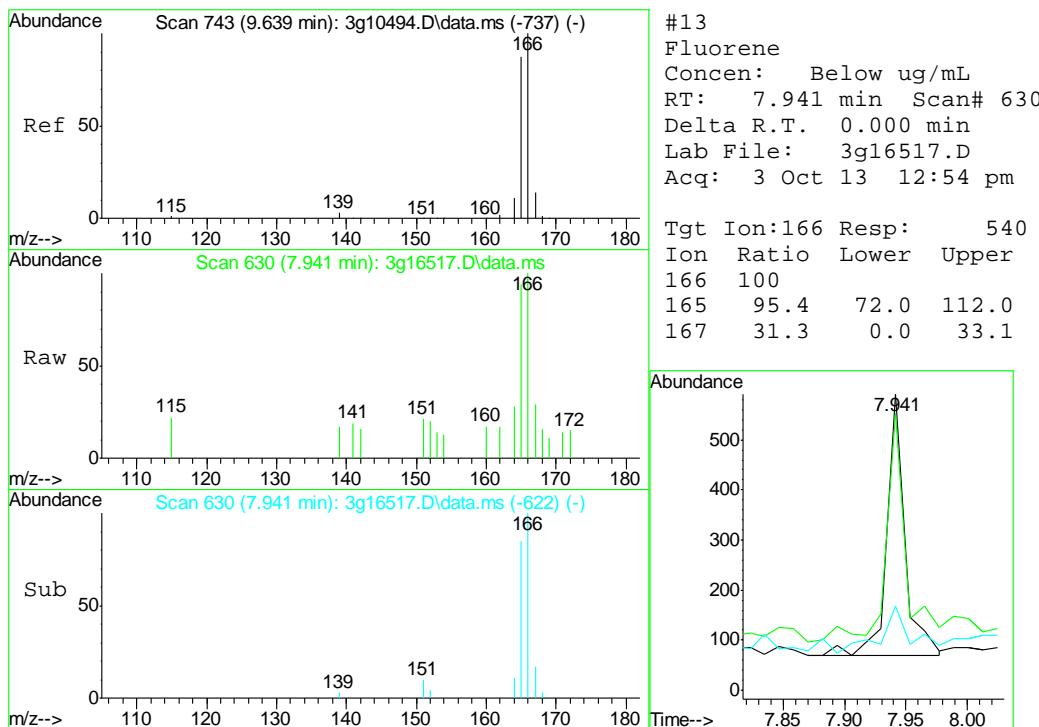


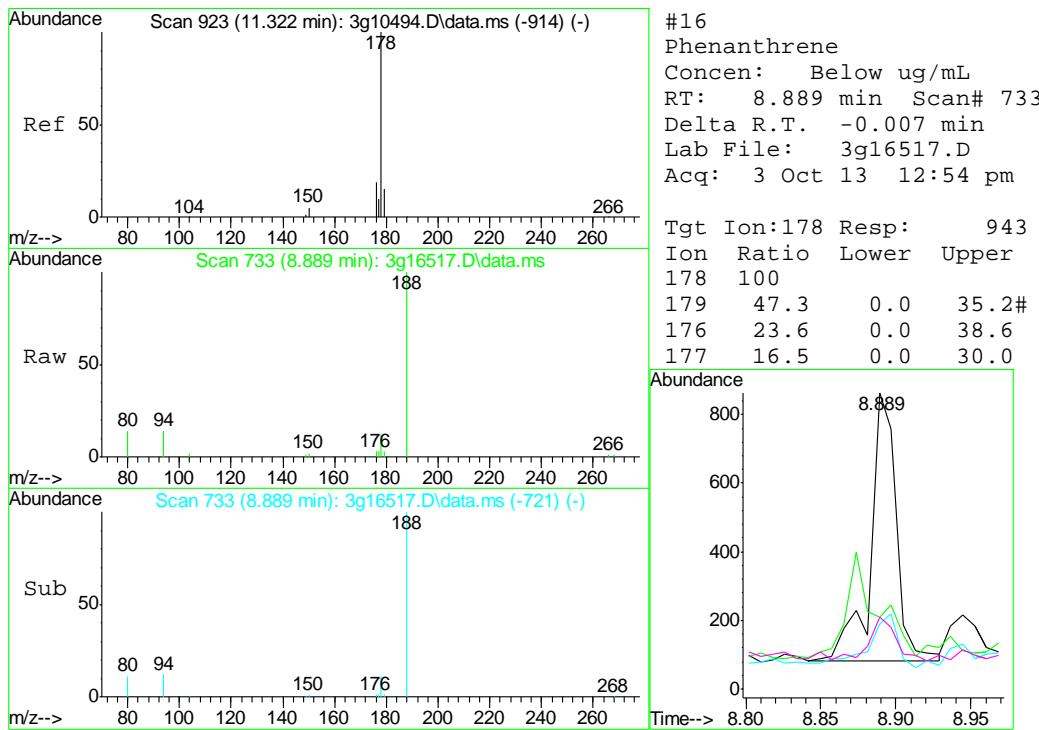
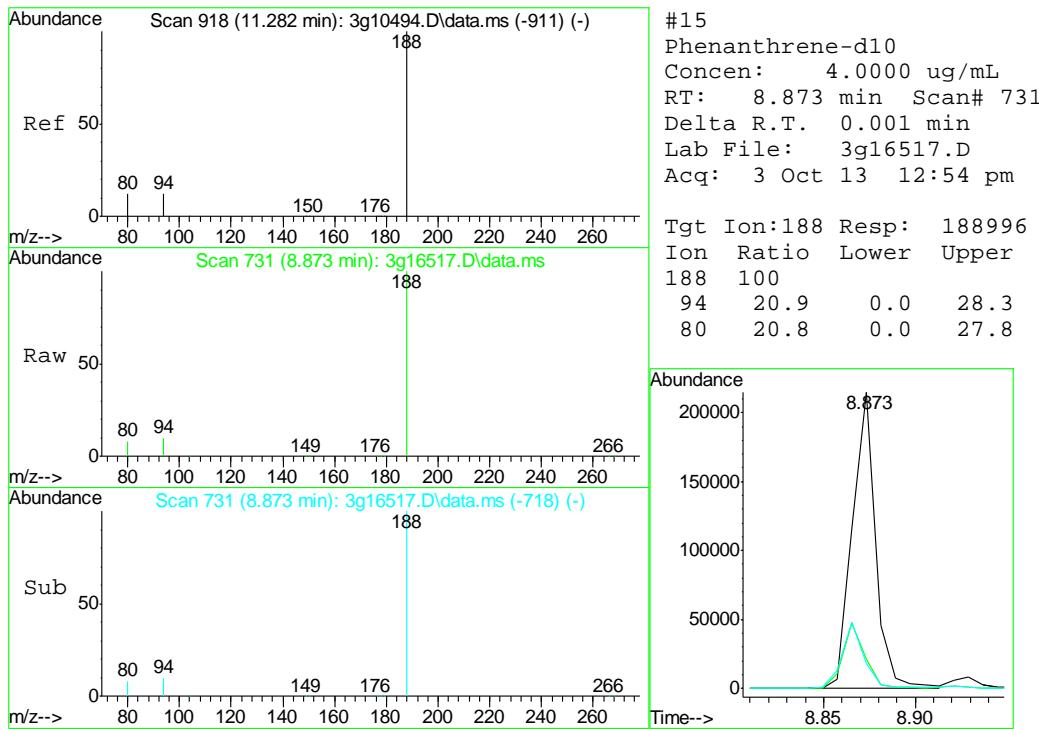


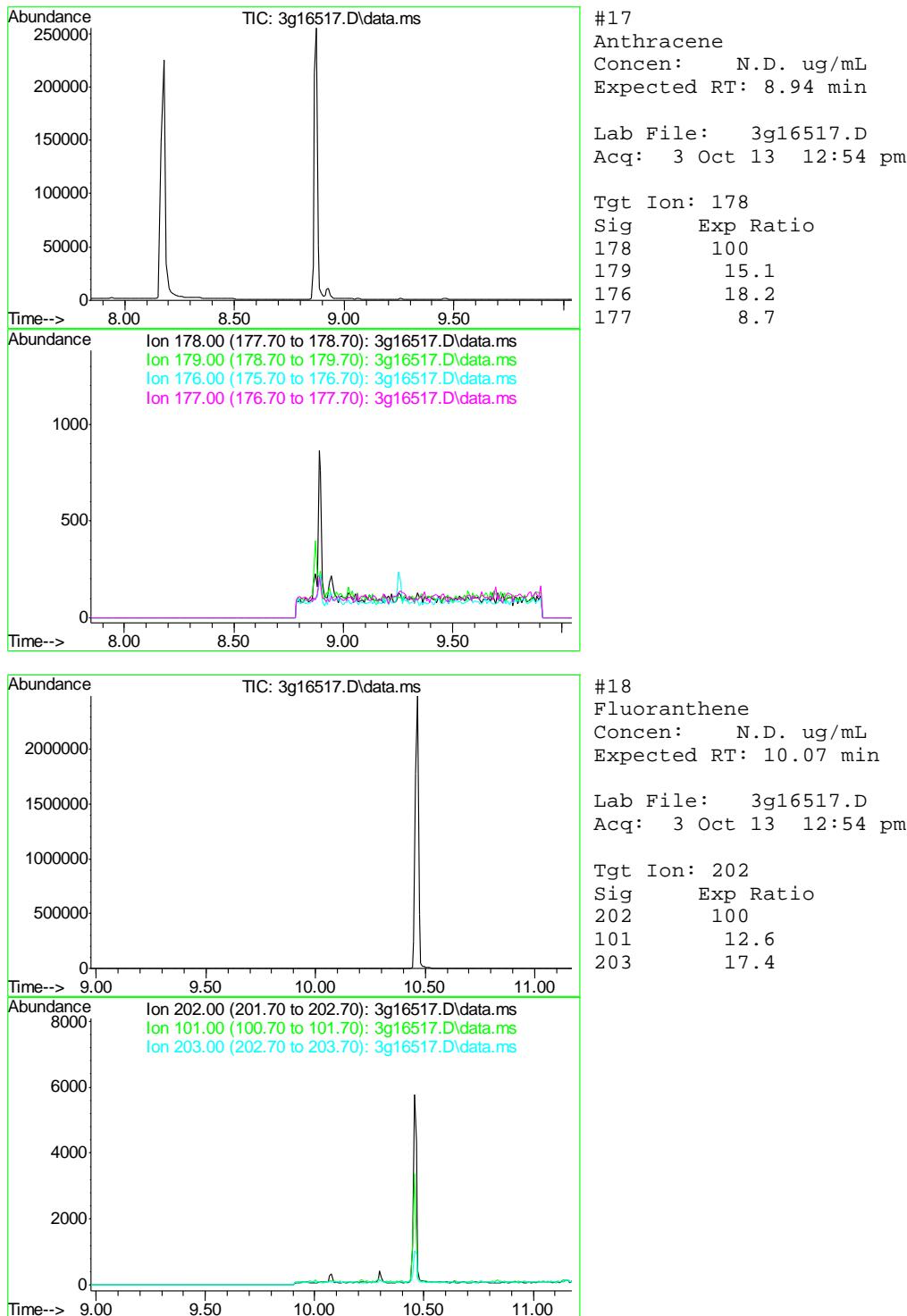


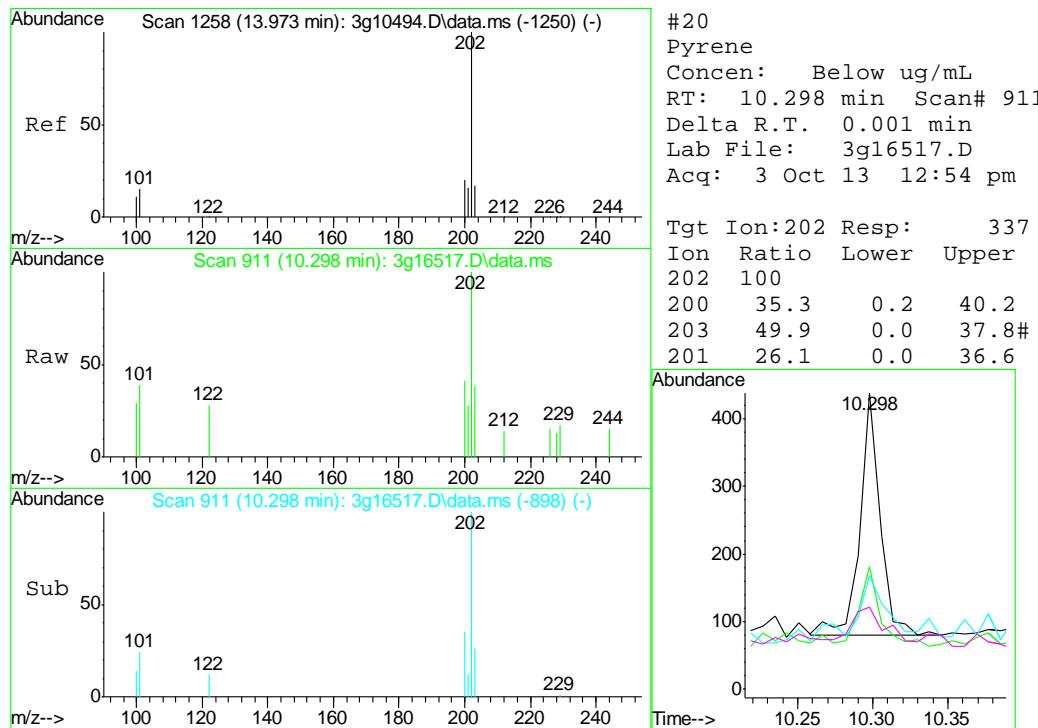
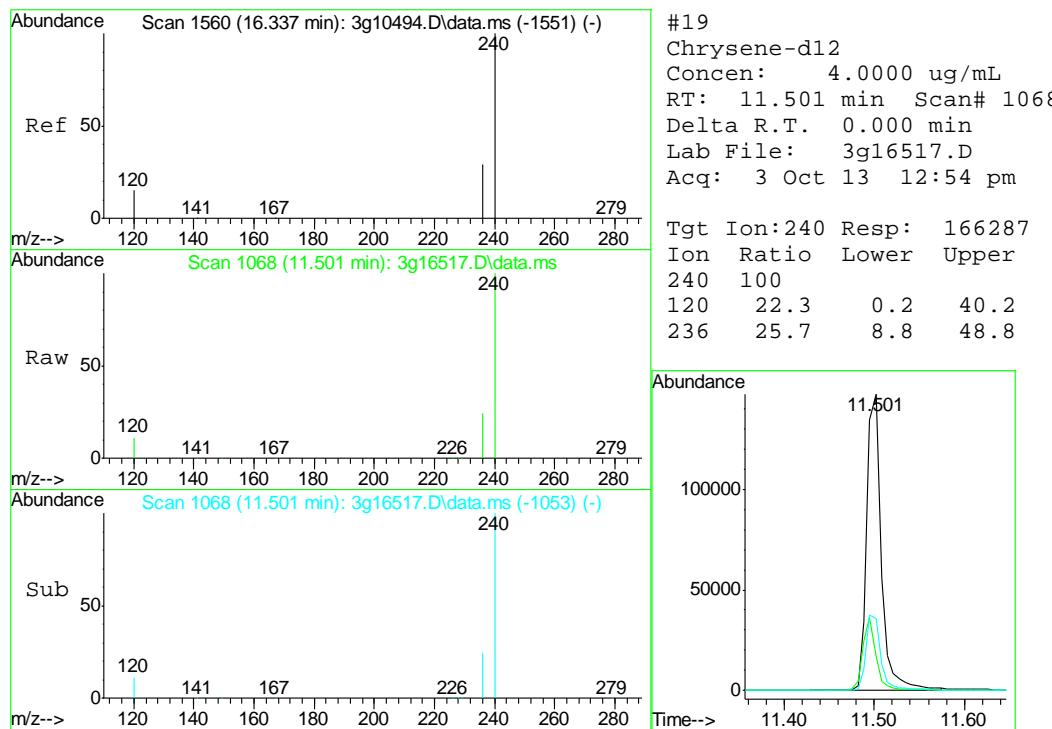


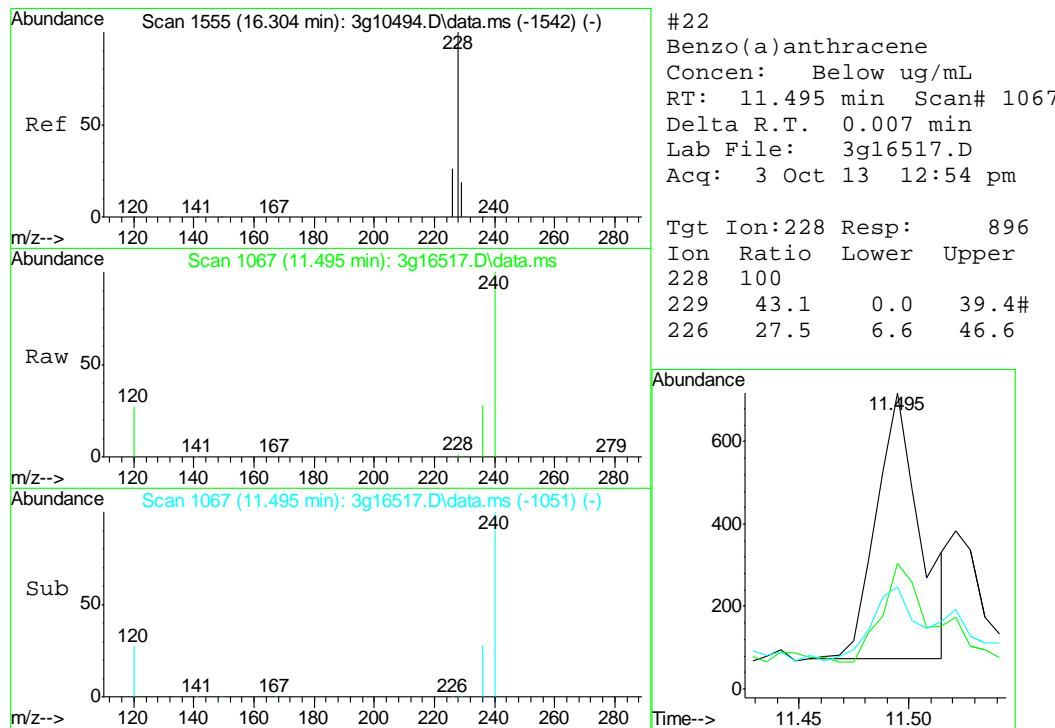
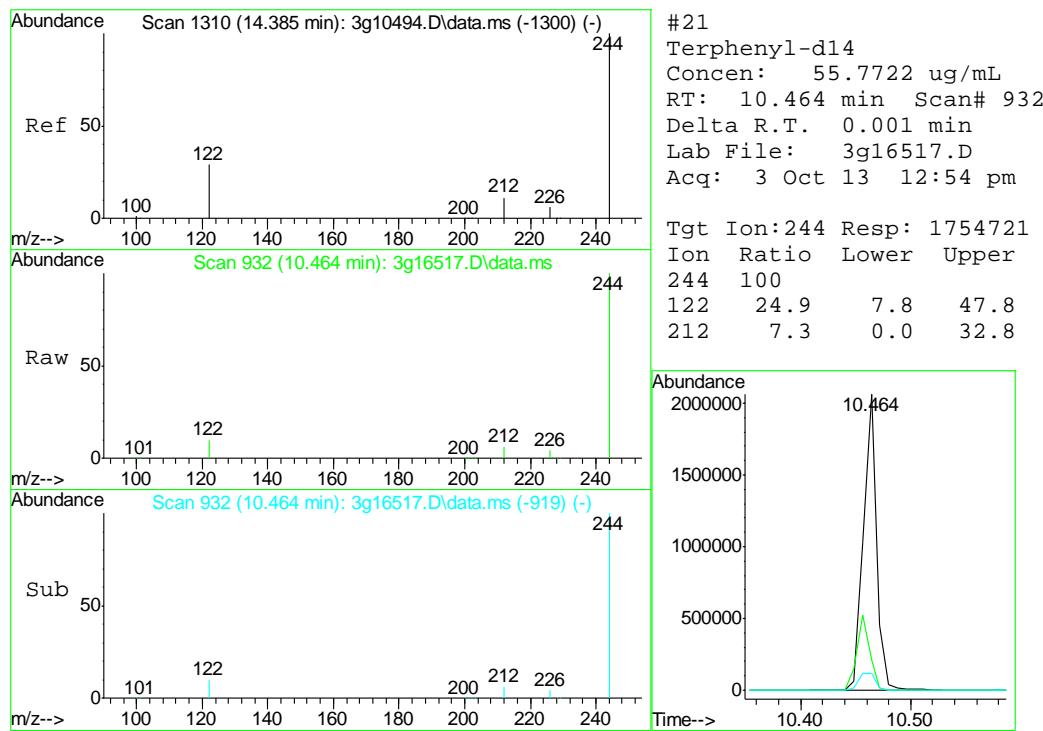


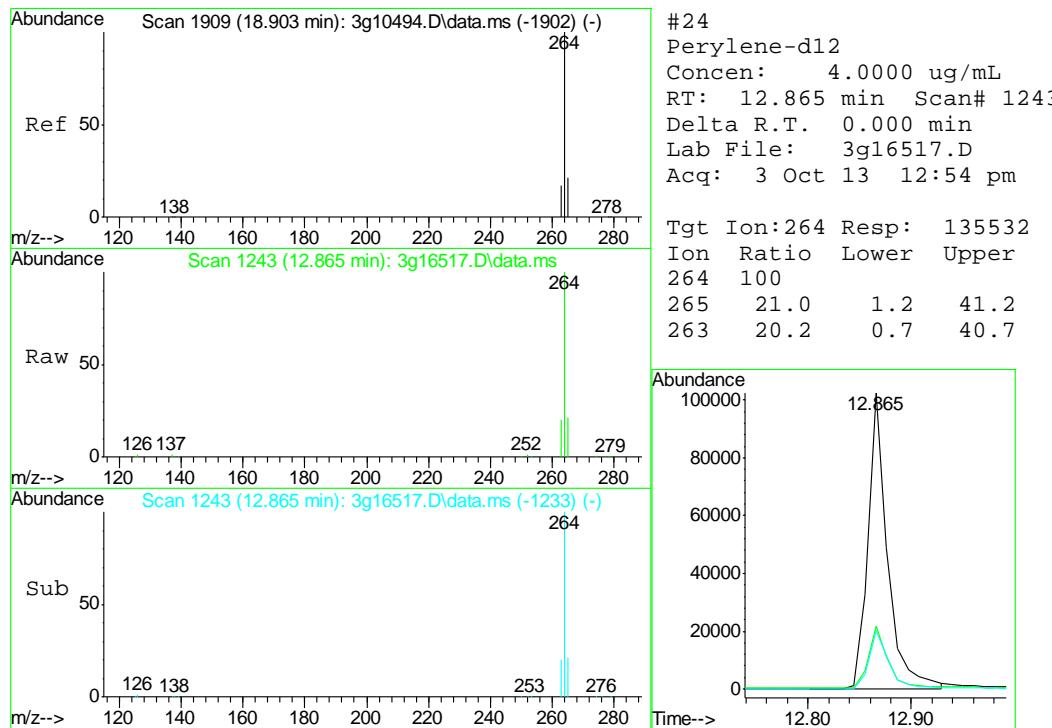
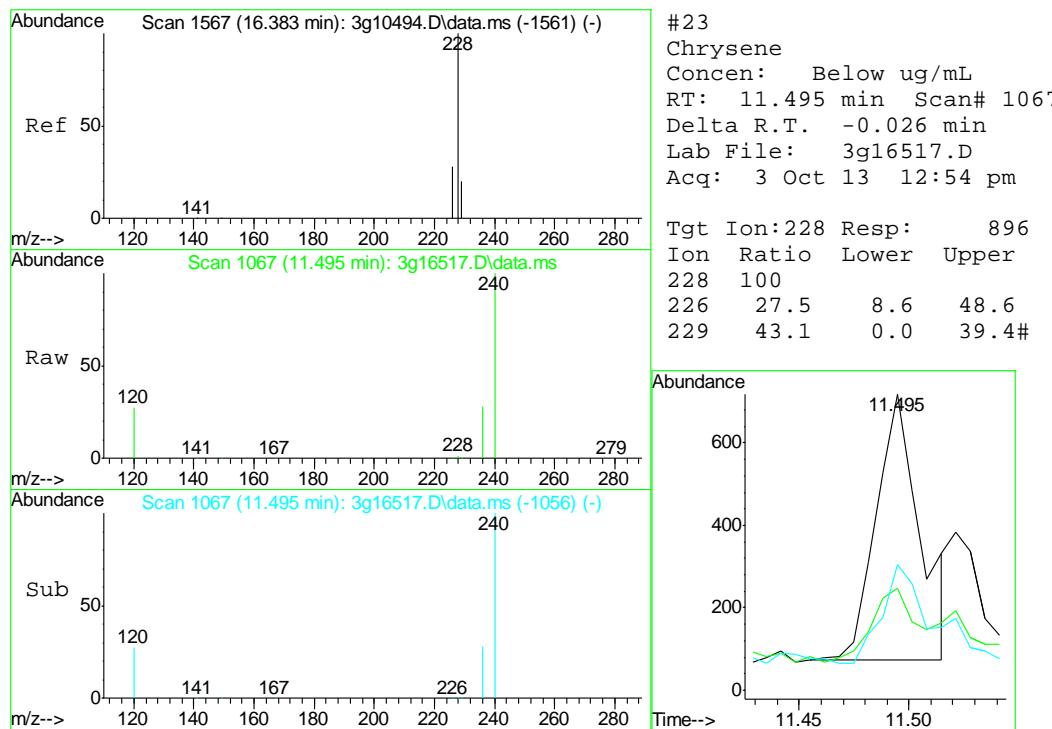


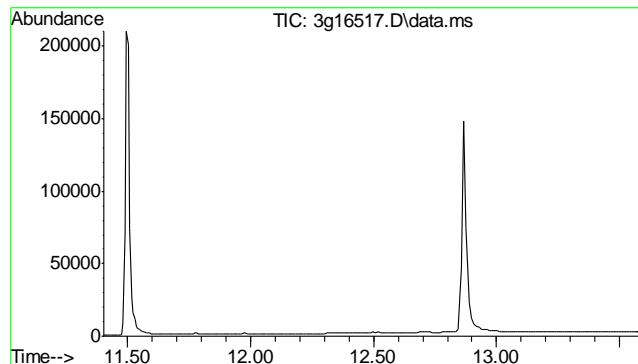








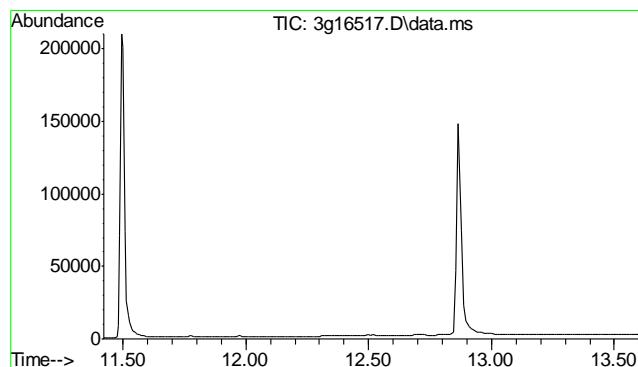
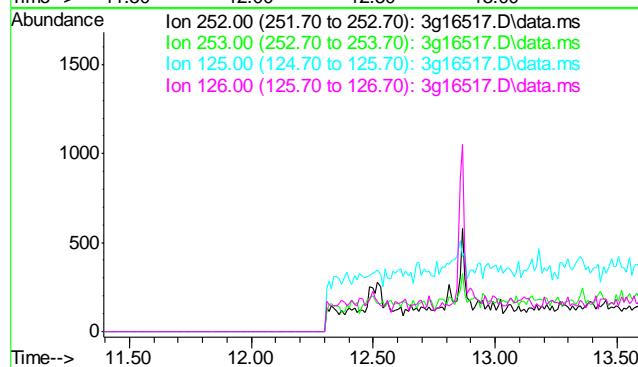




#25
Benzo(b)fluoranthene
Concen: N.D. ug/mL
Expected RT: 12.50 min

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

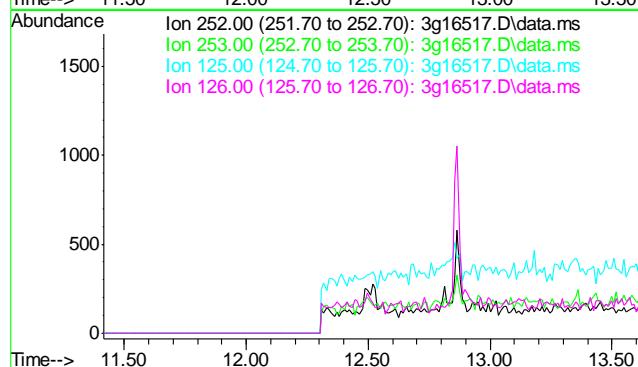
Tgt Ion: 252
Sig Exp Ratio
252 100
253 51.5
125 13.2
126 46.9



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

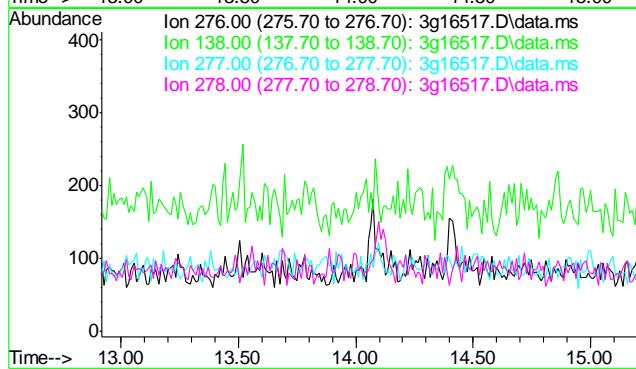
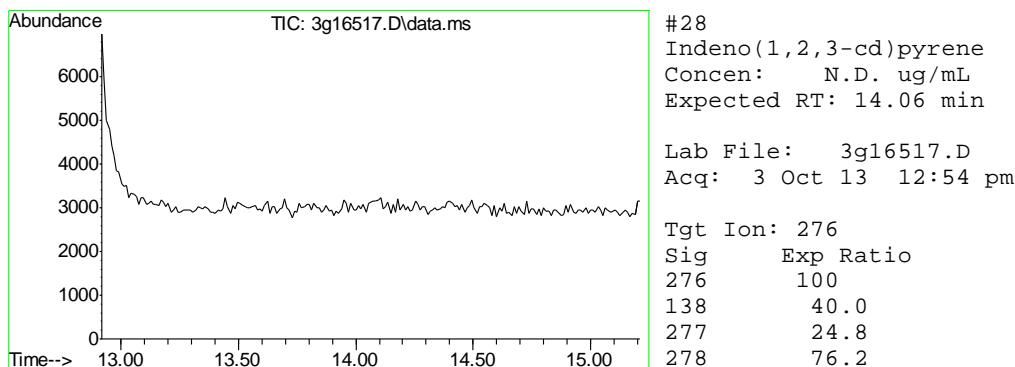
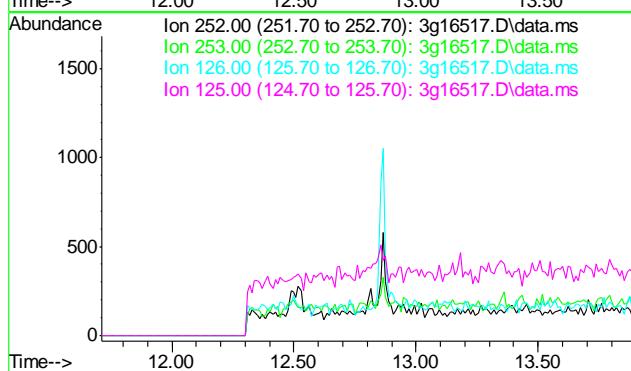
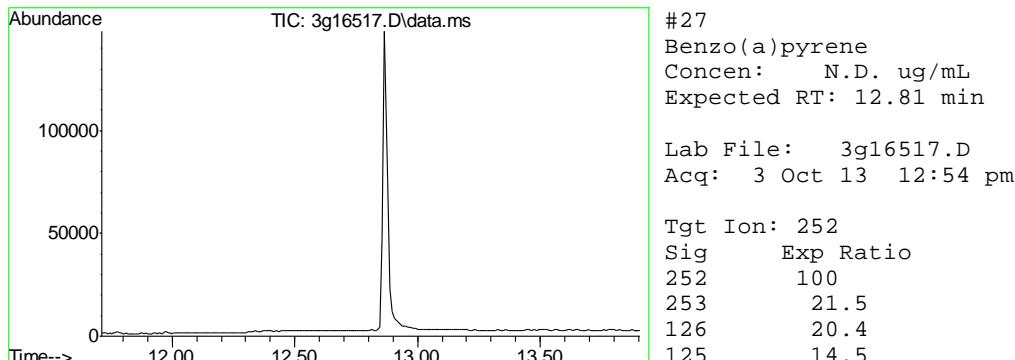
Lab File: 3g16517.D
Acq: 3 Oct 13 12:54 pm

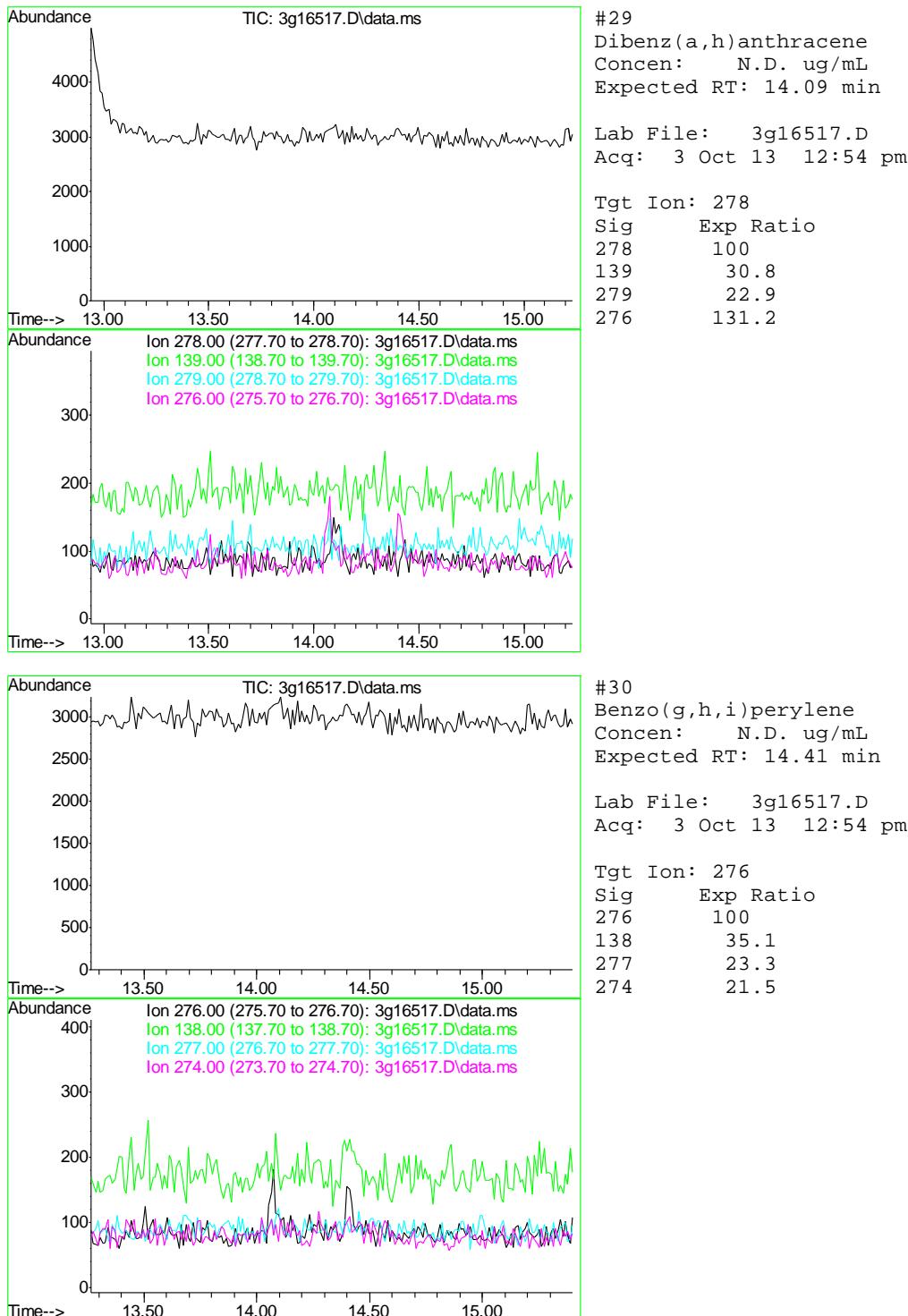
Tgt Ion: 252
Sig Exp Ratio
252 100
253 37.3
125 9.6
126 34.1



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: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1231-MB	GB22388.D	1	10/02/13	EV	n/a	n/a	GGB1231

The QC reported here applies to the following samples:

Method: SW846 8015B

D51203-1

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: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1231-BS	GB22387.D	1	10/02/13	EV	n/a	n/a	GGB1231

The QC reported here applies to the following samples:

Method: SW846 8015B

D51203-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	116	105	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	92%	60-140%

10.2.1
10

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D51200-1MS	GB22391.D	1	10/02/13	EV	n/a	n/a	GGB1231
D51200-1MSD	GB22392.D	1	10/02/13	EV	n/a	n/a	GGB1231
D51200-1	GB22390.D	1	10/02/13	EV	n/a	n/a	GGB1231

The QC reported here applies to the following samples:

Method: SW846 8015B

D51203-1

CAS No.	Compound	D51200-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		149	156	105	157	105	1	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D51200-1	Limits
120-82-1	1,2,4-Trichlorobenzene	93%	93%	86%	60-140%

* = Outside of Control Limits.

10.3.1
10



GC Volatiles

Raw Data

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson
10/07/13 13:45

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100213\GB22395.D\FID1A.CH Vial: 10
 Signal #2 : Y:\1\DATA\100213\GB22395.D\FID2B.CH
 Acq On : 2 Oct 2013 8:59 pm Operator: ELISEV
 Sample : D51203-1 Inst : GC/MS Ins
 Misc : GC3914,GGB1231,5.074,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 03 08:08:14 2013 Quant Results File: TB1125GB1125SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Oct 03 08:07:23 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	14.36	2420842	80.131 %	m
10) S 1,2,4-Trichlorobenzene (P)	14.36	11197034	84.794 %	m

Target Compounds

1) H TVH-Gasoline	7.28	4296220	0.061 mg/L
4) T Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T Benzene	4.14	104280	0.272 ug/L
6) T Toluene	7.66	107870	0.291 ug/L
7) T Ethylbenzene	0.00	0	N.D. ug/L d
8) T m,p-Xylene	10.46	167309	0.443 ug/L m
9) T o-Xylene	0.00	0	N.D. ug/L d
11) T Naphthalene	14.54	26409	0.153 ug/L m

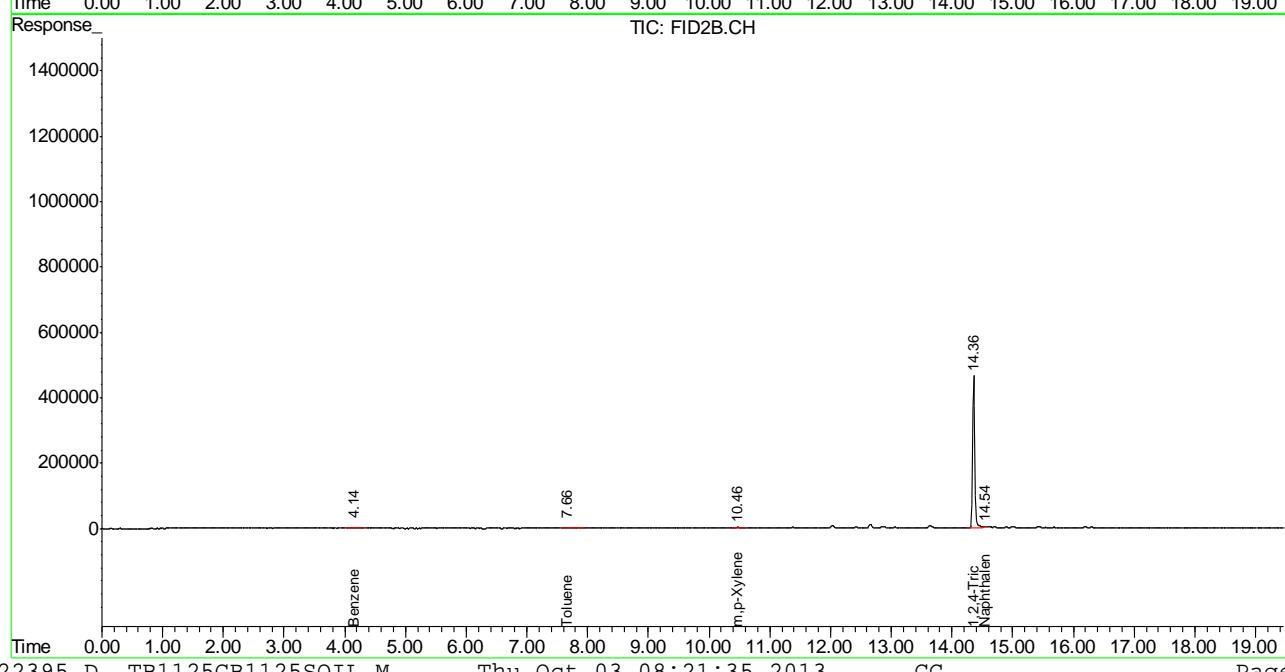
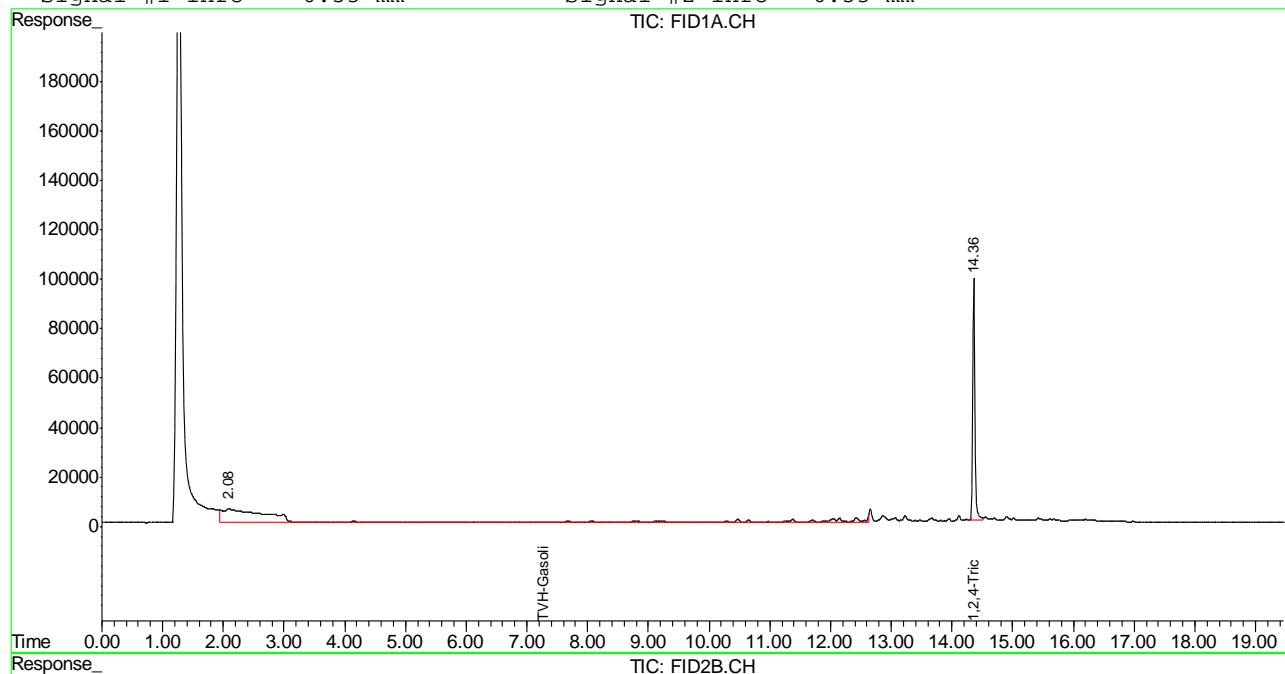
 (f)=RT Delta > 1/2 Window (m)=manual int.
 GB22395.D TB1125GB1125SOIL.M Thu Oct 03 08:21:35 2013 GC

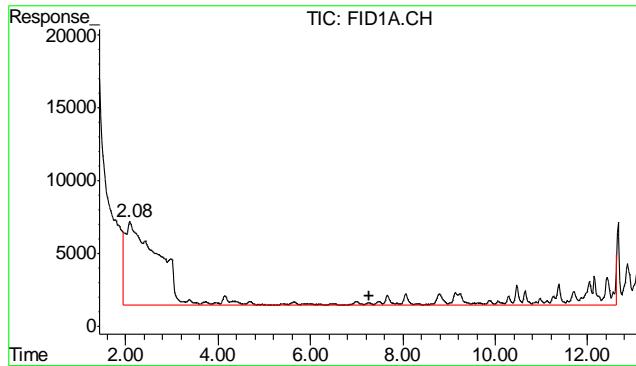
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100213\GB22395.D\FID1A.CH Vial: 10
 Signal #2 : Y:\1\DATA\100213\GB22395.D\FID2B.CH
 Acq On : 2 Oct 2013 8:59 pm Operator: ELISEV
 Sample : D51203-1 Inst : GC/MS Ins
 Misc : GC3914,GGB1231,5.074,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 3 8:25 2013 Quant Results File: TB1125GB1125SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Oct 03 08:07:23 2013
 Response via : Multiple Level Calibration
 DataAcq Meth : TVB4.M

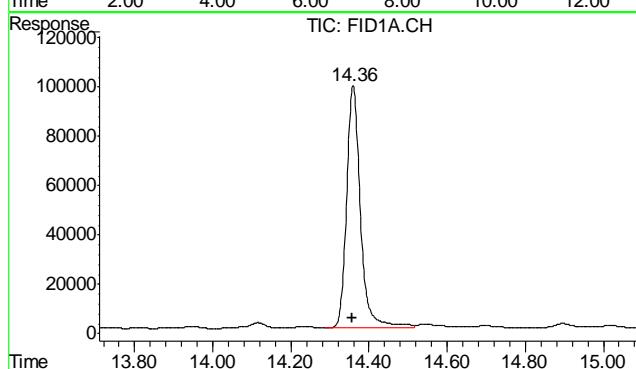
Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





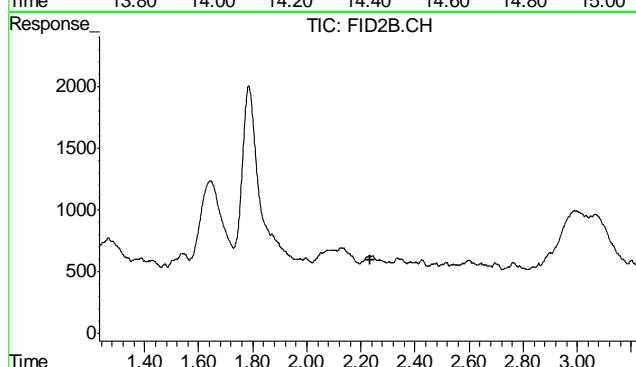
#1 TVH-Gasoline

R.T.: 7.280 min
 Delta R.T.: 0.000 min
 Response: 4296220
 Conc: 0.06 mg/L m



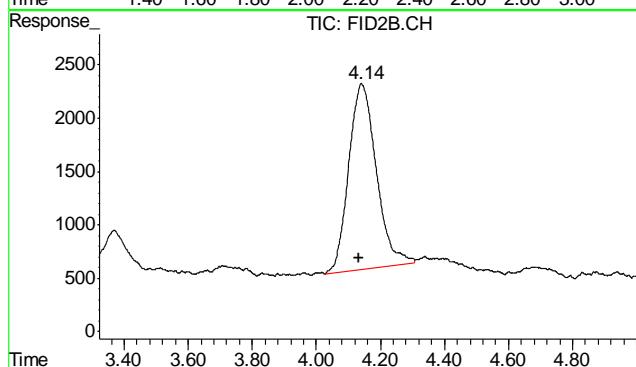
#2 1,2,4-Trichlorobenzene

R.T.: 14.359 min
 Delta R.T.: 0.002 min
 Response: 2420842
 Conc: 80.13 % m



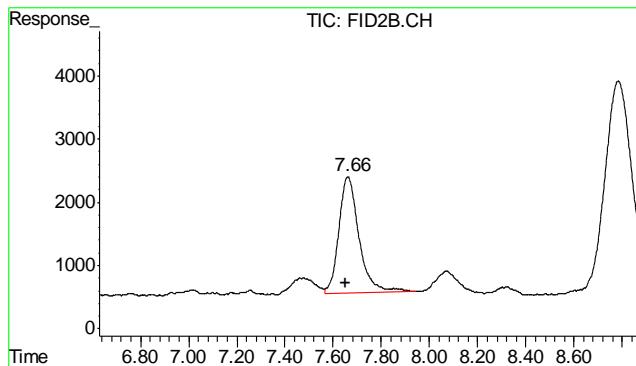
#4 Methyl-t-butyl-ether

R.T.: 0.000 min
 Exp R.T.: 2.233 min
 Response: 0
 Conc: N.D.

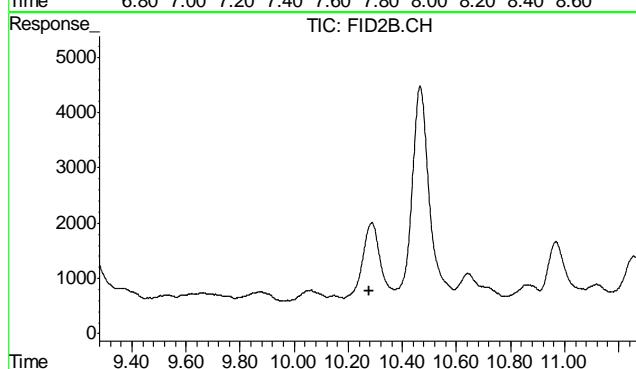


#5 Benzene

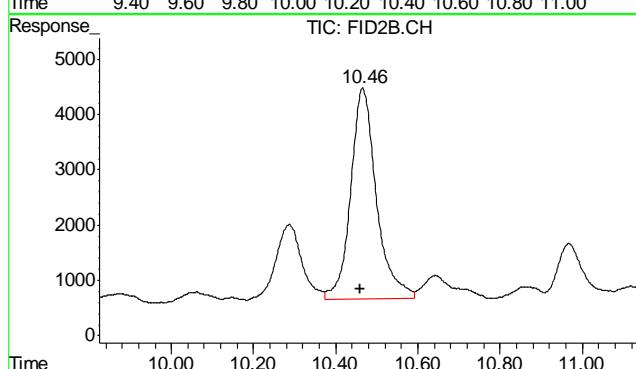
R.T.: 4.141 min
 Delta R.T.: 0.010 min
 Response: 104280
 Conc: 0.27 ug/L



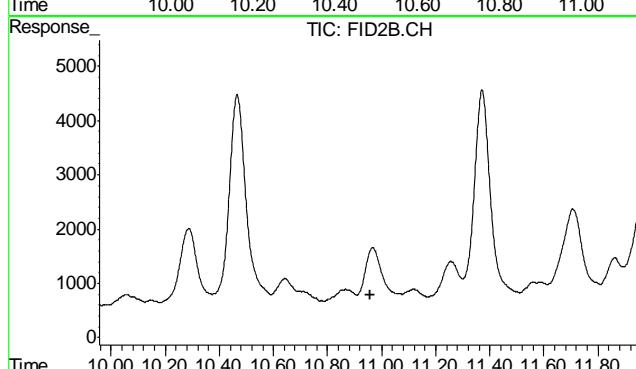
#6 Toluene
R.T.: 7.661 min
Delta R.T.: 0.009 min
Response: 107870
Conc: 0.29 ug/L



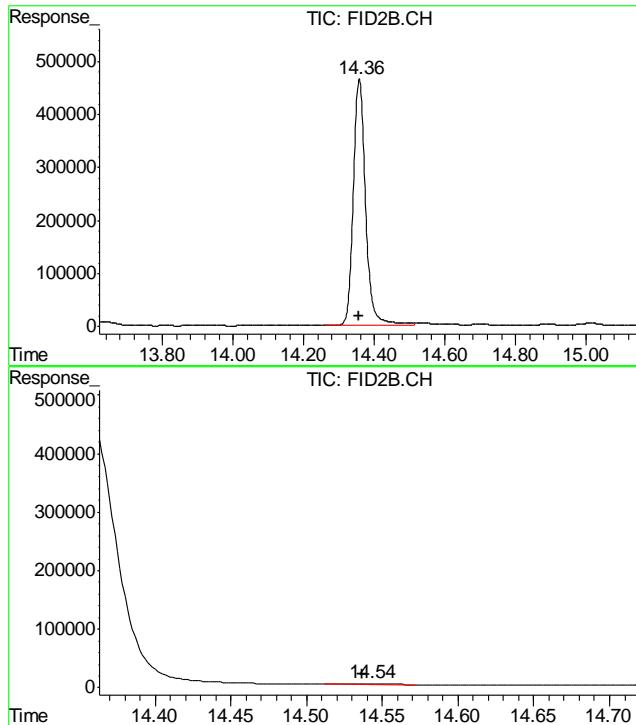
#7 Ethylbenzene
R.T.: 0.000 min
Exp R.T. : 10.279 min
Response: 0
Conc: N.D.



#8 m,p-Xylene
R.T.: 10.465 min
Delta R.T.: 0.005 min
Response: 167309
Conc: 0.44 ug/L m



#9 o-Xylene
R.T.: 0.000 min
Exp R.T. : 10.956 min
Response: 0
Conc: N.D.



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

R.T.: 14.356 min
Delta R.T.: 0.000 min
Response: 11197034
Conc: 84.79 % m

#11 Naphthalene

R.T.: 14.542 min
Delta R.T.: 0.005 min
Response: 26409
Conc: 0.15 ug/L m

11.1

**Manual Integrations
APPROVED
(compounds with "m" flag)**
Jennifer Laidlaw
10/03/13 09:15

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100213\GB22388.D\FID1A.CH Vial: 3
 Signal #2 : Y:\1\DATA\100213\GB22388.D\FID2B.CH
 Acq On : 2 Oct 2013 4:51 pm Operator: ELISEV
 Sample : MB, S Inst : GC/MS Ins
 Misc : GC3914,GGB1231,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 03 08:07:46 2013 Quant Results File: TB1125GB1125SOIL.RES

Quant Method : C:\MSDCHEM\1...TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Oct 03 08:07:23 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	14.36	2563808	84.863 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	11827632	89.570 %	m

Target Compounds

1) H	TVH-Gasoline	7.28	4113266	0.059 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	7.65	113265	0.306 ug/L
7) T	Ethylbenzene	0.00	0	N.D. ug/L d
8) T	m,p-Xylene	10.46	150042	0.397 ug/L m
9) T	o-Xylene	0.00	0	N.D. ug/L d
11) T	Naphthalene	14.54	260490	1.511 ug/L m

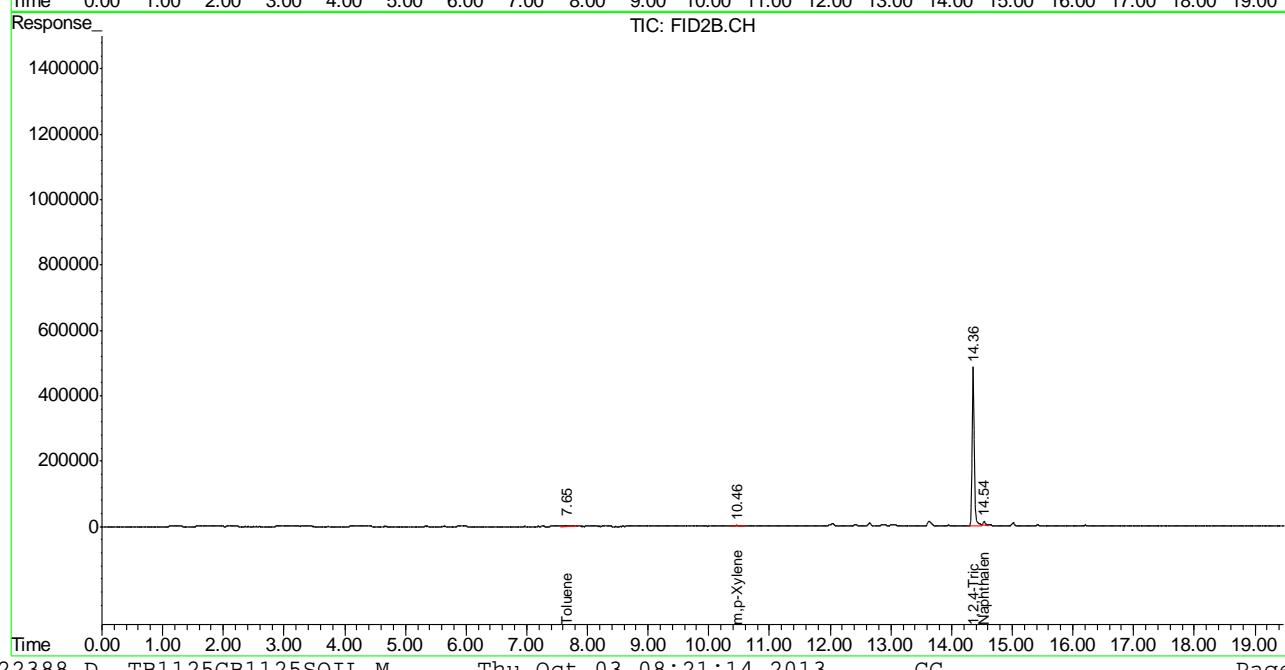
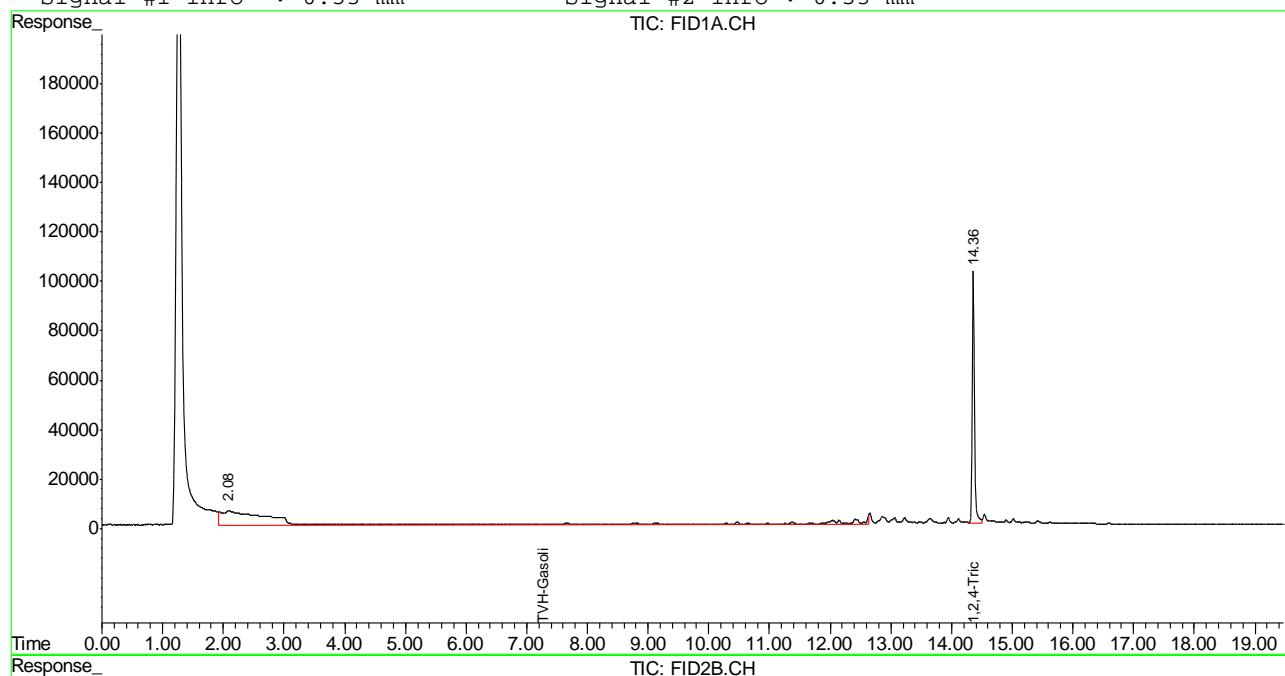
(f)=RT Delta > 1/2 Window (m)=manual int.
 GB22388.D TB1125GB1125SOIL.M Thu Oct 03 08:21:14 2013 GC

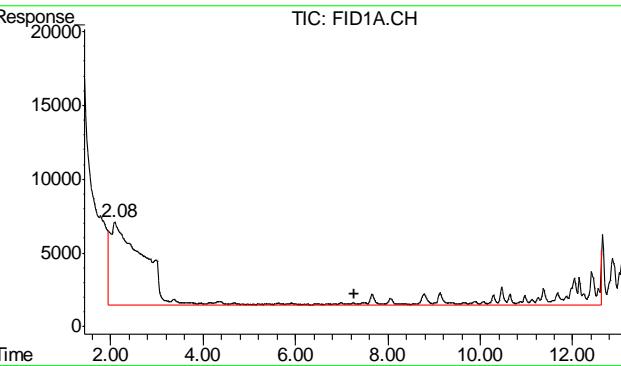
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\100213\GB22388.D\FID1A.CH Vial: 3
 Signal #2 : Y:\1\DATA\100213\GB22388.D\FID2B.CH
 Acq On : 2 Oct 2013 4:51 pm Operator: ELISEV
 Sample : MB, S Inst : GC/MS Ins
 Misc : GC3914, GGB1231, 5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Oct 3 8:17 2013 Quant Results File: TB1125GB1125SOIL.RES

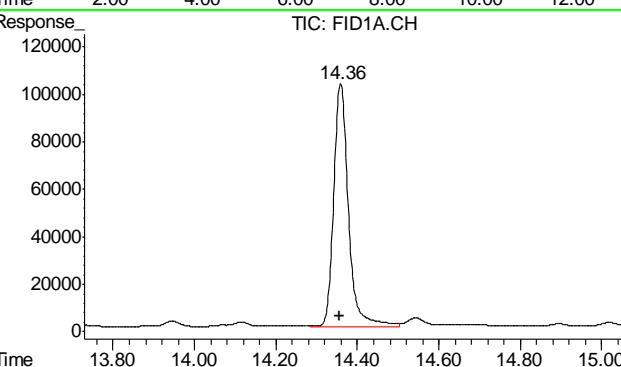
Quant Method : C:\MSDCHEM\1...\TB1125GB1125SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Oct 03 08:07:23 2013
 Response via : Multiple Level Calibration
 DataAcq Meth : TVB4.M

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

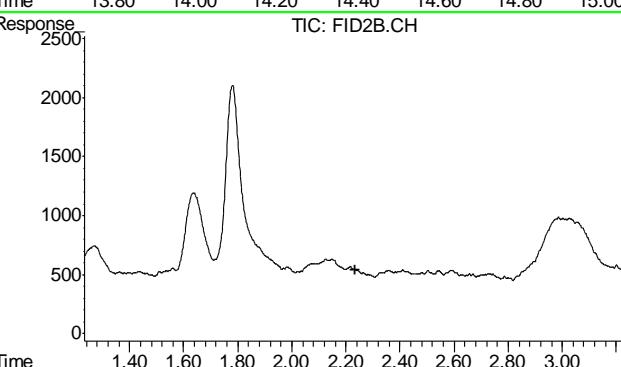




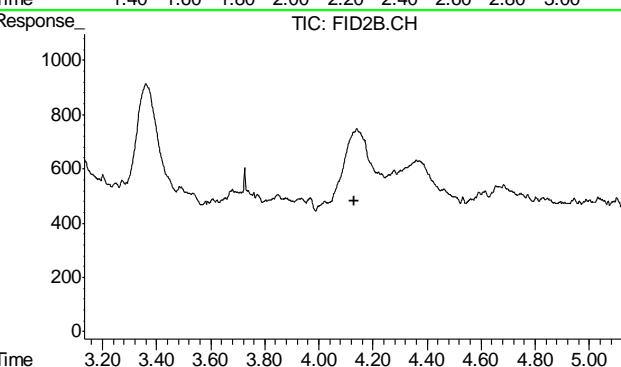
#1 TVH-Gasoline
R.T.: 7.280 min
Delta R.T.: 0.000 min
Response: 4113266
Conc: 0.06 mg/L m



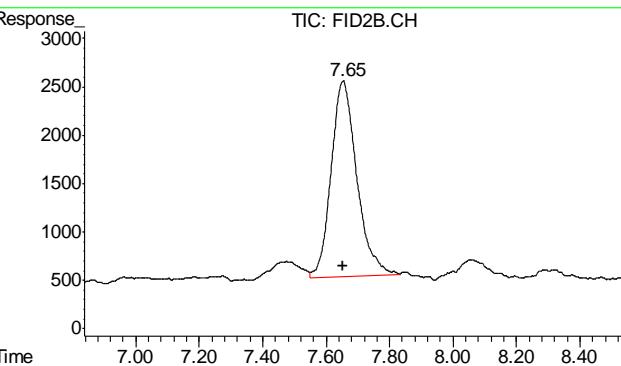
#2 1,2,4-Trichlorobenzene
R.T.: 14.358 min
Delta R.T.: 0.001 min
Response: 2563808
Conc: 84.86 % m



#4 Methyl-t-butyl-ether
R.T.: 0.000 min
Exp R.T. : 2.233 min
Response: 0
Conc: N.D.

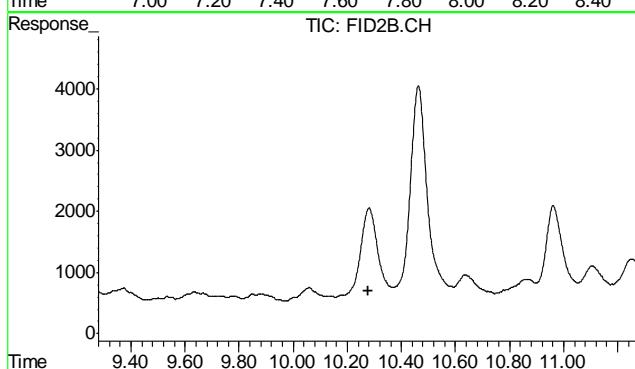


#5 Benzene
R.T.: 0.000 min
Exp R.T. : 4.132 min
Response: 0
Conc: N.D.



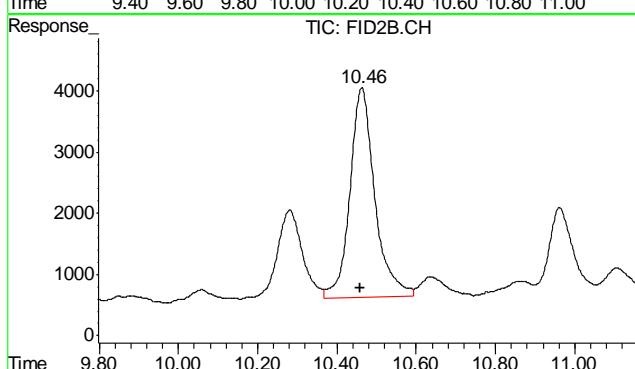
#6 Toluene

R.T.: 7.654 min
Delta R.T.: 0.002 min
Response: 113265
Conc: 0.31 ug/L



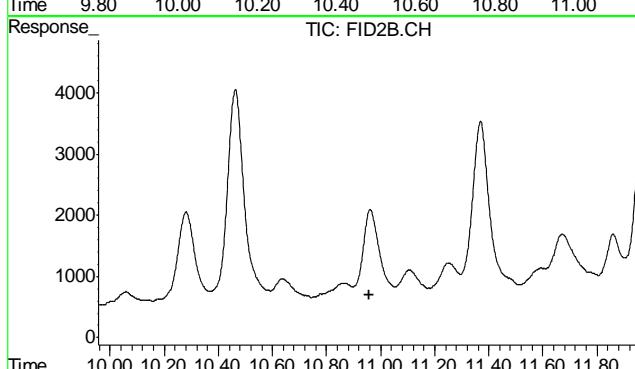
#7 Ethylbenzene

R.T.: 0.000 min
Exp R.T.: 10.279 min
Response: 0
Conc: N.D.



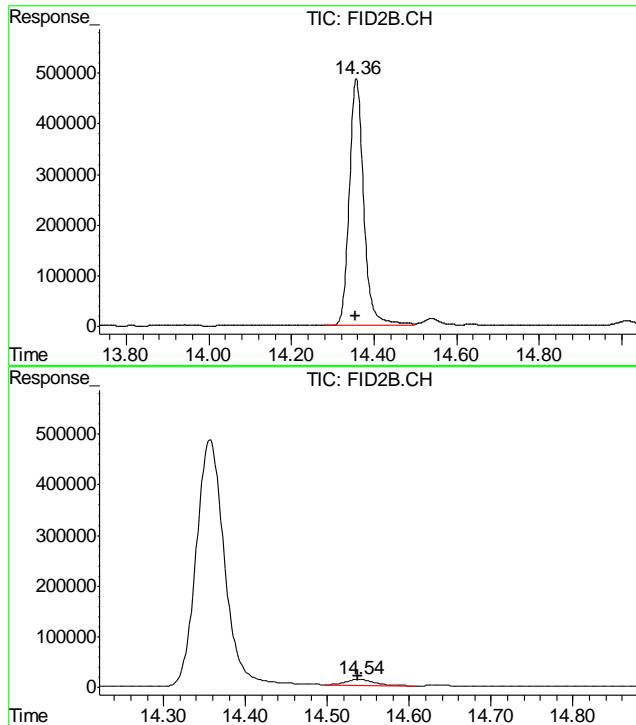
#8 m,p-Xylene

R.T.: 10.463 min
Delta R.T.: 0.003 min
Response: 150042
Conc: 0.40 ug/L m



#9 o-Xylene

R.T.: 0.000 min
Exp R.T.: 10.956 min
Response: 0
Conc: N.D.



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

R.T.: 14.356 min
Delta R.T.: 0.000 min
Response: 11827632
Conc: 89.57 % m

#11 Naphthalene

R.T.: 14.539 min
Delta R.T.: 0.002 min
Response: 260490
Conc: 1.51 ug/L m

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: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8682-MB	FH013720.D	1	10/04/13	TU	10/04/13	OP8682	GFH723

The QC reported here applies to the following samples:

Method: SW846-8015B

D51203-1

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	80% 20-130%

12.1.1

12

Blank Spike Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8682-BS	FH013722.D	1	10/04/13	TU	10/04/13	OP8682	GFH723

The QC reported here applies to the following samples:

Method: SW846-8015B

D51203-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	582	87	42-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	86%	20-130%

* = Outside of Control Limits.

12.2.1

12

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D51203

Account: XTOKWR XTO Energy

Project: FRU 197-31A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP8682-MS	FH013724.D	10	10/04/13	TU	10/04/13	OP8682	GFH723
OP8682-MSD	FH013726.D	10	10/04/13	TU	10/04/13	OP8682	GFH723
D51224-3	FH013732.D	10	10/04/13	TU	10/04/13	OP8682	GFH723

The QC reported here applies to the following samples:

Method: SW846-8015B

D51203-1

CAS No.	Compound	D51224-3		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	618		835	579	-5*	743	15*	25	20-150/30
CAS No.	Surrogate Recoveries	MS		MSD		D51224-3		Limits		
84-15-1	o-Terphenyl	59%		74%		78%		20-130%		

* = Outside of Control Limits.

12.3.1
12



GC Semi-volatiles

Raw Data

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH100413.SEC\
 Data File : FH013760.D
 Signal(s) : FID2B.ch
 Acq On : 5 Oct 2013 4:43 am
 Operator : TIMU
 Sample : D51203-1
 Misc : OP8682,GFH723,30.04,,,1,1
 ALS Vial : 74 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Oct 07 08:16:19 2013
 Quant Method : C:\msdchem\1\METHODS\DRD-GFH689R.M
 Quant Title : DRD-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	12.163	2262859533	1304.152	ug/ml
<hr/>				
Target Compounds				
2) H TPH-DRO (C10-C28)	9.818	997691352	709.303	ug/ml
<hr/>				

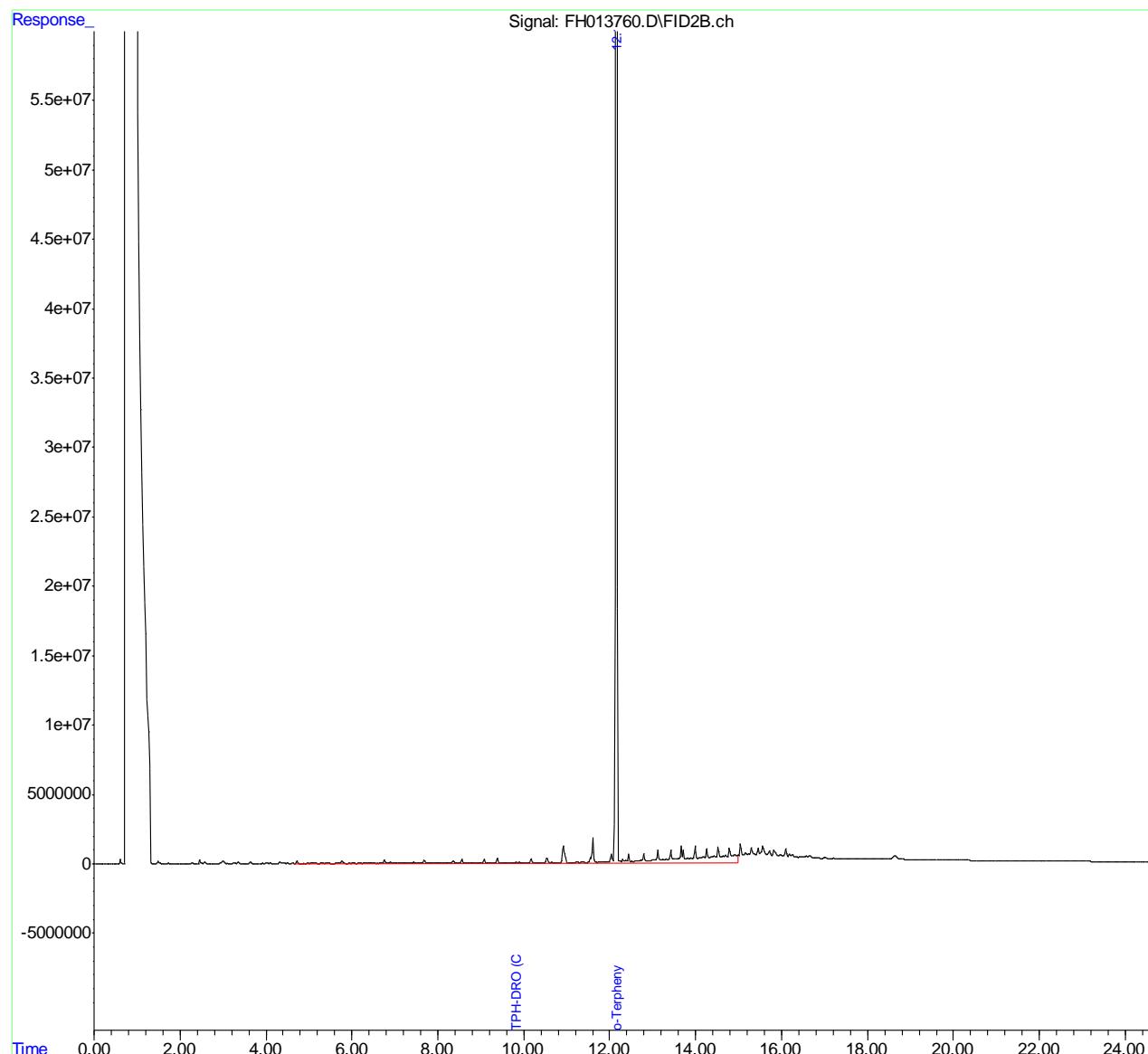
(f)=RT Delta > 1/2 Window (m)=manual int.

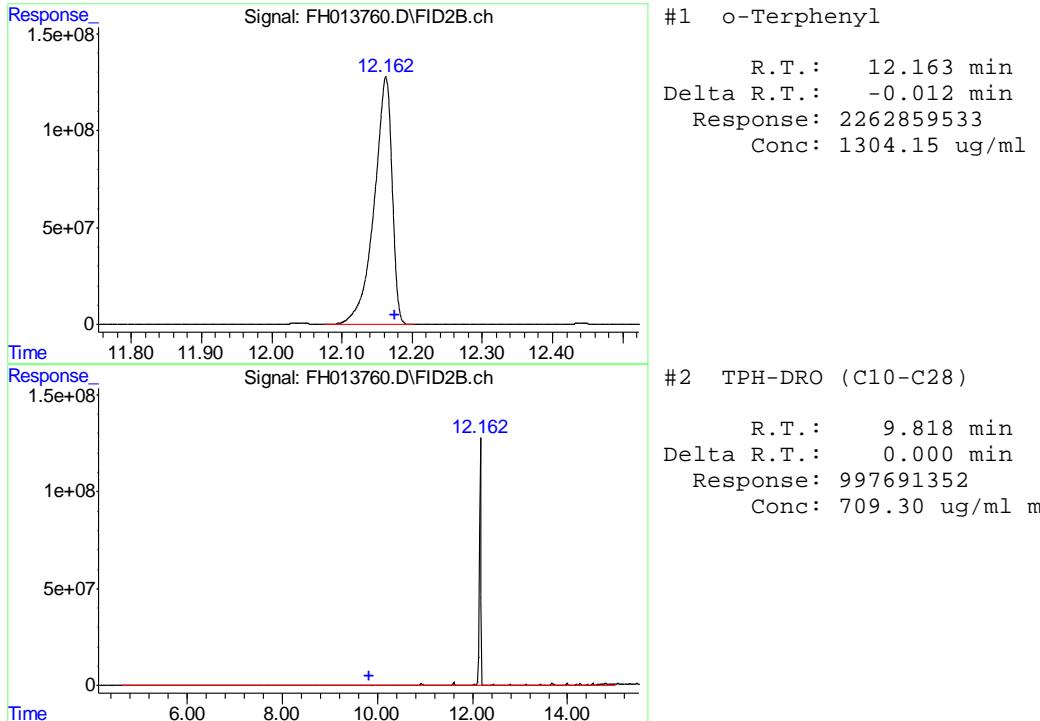
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH100413.SEC\
 Data File : FH013760.D
 Signal(s) : FID2B.ch
 Acq On : 5 Oct 2013 4:43 am
 Operator : TIMU
 Sample : D51203-1
 Misc : OP8682,GFH723,30.04,,,1,1
 ALS Vial : 74 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Oct 07 08:16:19 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 :





13.1.1

13

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH100413.SEC\
 Data File : FH013720.D
 Signal(s) : FID2B.ch
 Acq On : 4 Oct 2013 4:50 pm
 Operator : TIMU
 Sample : OP8682-MB
 Misc : OP8682,GFH723,30.00,,,1,1
 ALS Vial : 54 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Oct 07 08:29:11 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	12.171	2782035293	1603.368	ug/ml
<hr/>				
Target Compounds				
2) H TPH-DRO (C10-C28)	9.818	77975044	55.436	ug/ml
<hr/>				

(f)=RT Delta > 1/2 Window

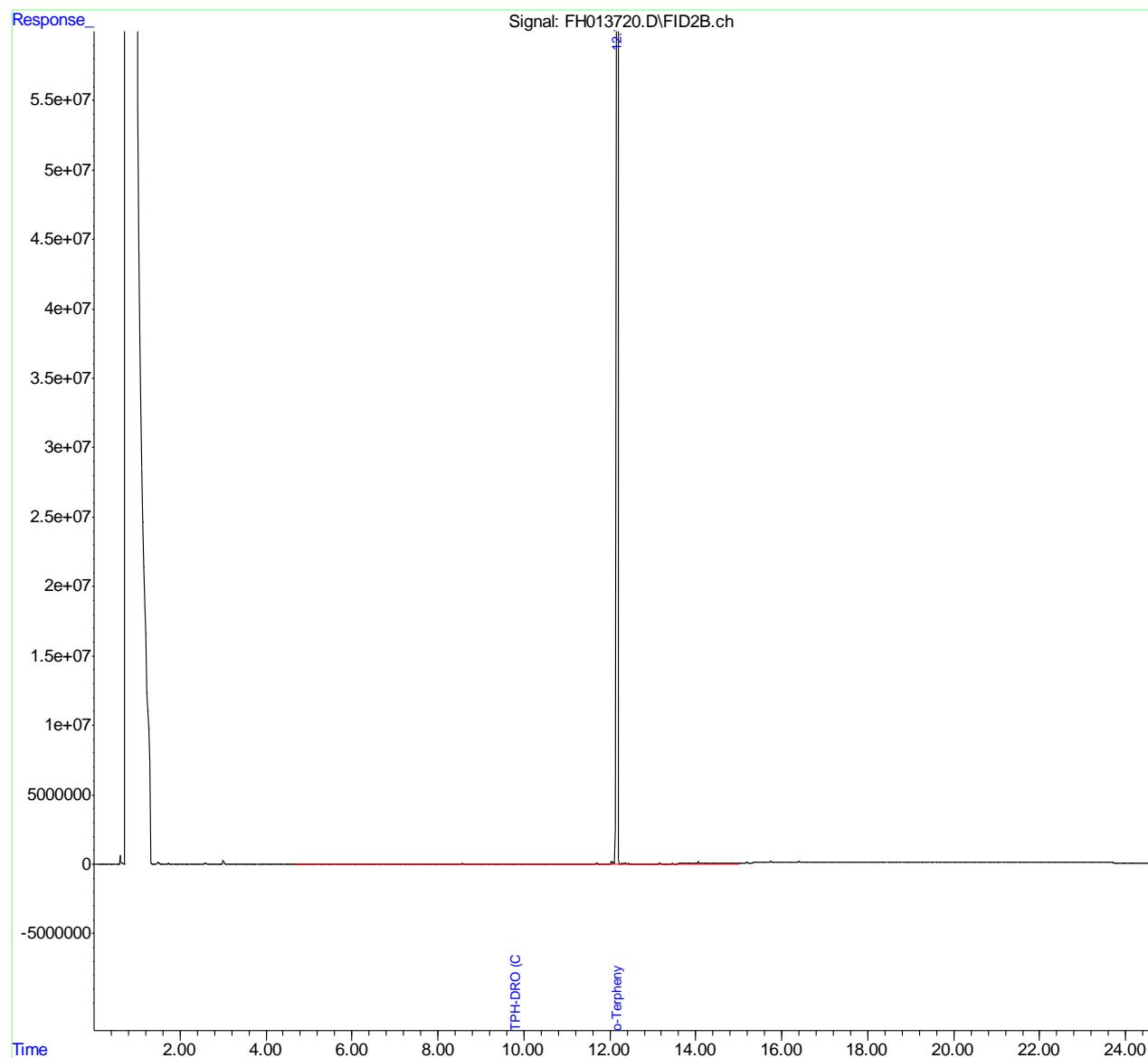
(m)=manual int.

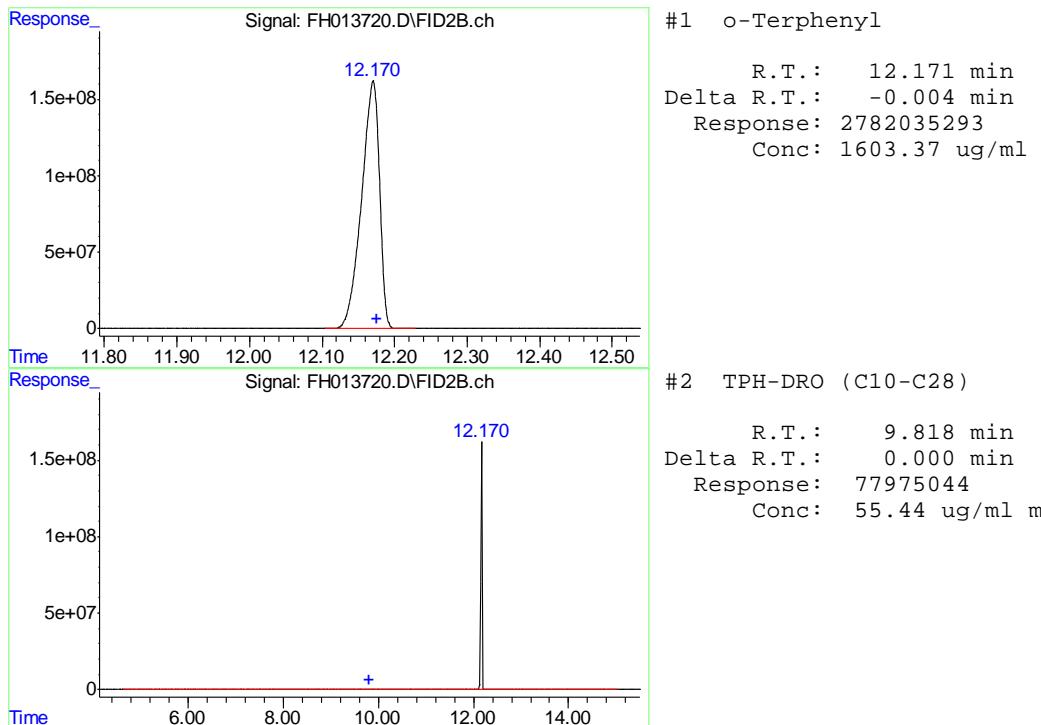
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH100413.SEC\
 Data File : FH013720.D
 Signal(s) : FID2B.ch
 Acq On : 4 Oct 2013 4:50 pm
 Operator : TIMU
 Sample : OP8682-MB
 Misc : OP8682,GFH723,30.00,,,1,1
 ALS Vial : 54 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Oct 07 08:29:11 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 :





13.2.1

13



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: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11290
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

10/03/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	.86	1.8		
Antimony	3.0	.21	.5		
Arsenic	2.5	.38	.63		
Barium	1.0	.02	.36	0.030	<1.0
Beryllium	1.0	.08	.06		
Boron	5.0	.08	.16		
Cadmium	1.0	.02	.28	0.010	<1.0
Calcium	40	.22	6.8		
Chromium	1.0	.03	.03	0.030	<1.0
Cobalt	0.50	.04	.039		
Copper	1.0	.12	.13	0.13	<1.0
Iron	7.0	.15	1.8		
Lead	5.0	.21	.25	-0.21	<5.0
Lithium	0.50	.04	.13		
Magnesium	20	.68	1.8		
Manganese	0.50	.001	.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	0.29	<5.0
Silicon	5.0	.47	1.1		
Silver	3.0	.03	.05	-0.050	<3.0
Sodium	40	.49	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	0.21	<3.0

Associated samples MP11290: D51203-1

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11290
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11290
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

10/03/13

Metal	D51202-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	225	467	235	102.2
Beryllium				
Boron	anr			
Cadmium	0.13	51.1	58.7	86.9
Calcium				
Chromium	42.0	97.2	58.7	94.9
Cobalt				
Copper	7.0	66.8	58.7	101.9
Iron				
Lead	8.1	110	117	86.8
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	14.2	63.3	58.7	85.5
Phosphorus				
Potassium				
Selenium	0.0	105	117	89.5
Silicon				
Silver	0.15	22.3	23.5	94.4
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	34.8	83.4	58.7	82.8

Associated samples MP11290: D51203-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11290
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11290
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 10/03/13

Metal	D51202-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	anr					
Barium	225	494	228	117.2	5.6	20
Beryllium						
Boron	anr					
Cadmium	0.13	48.3	57	84.6	5.6	20
Calcium						
Chromium	42.0	89.1	57	83.5	8.7	20
Cobalt						
Copper	7.0	64.7	57	101.3	3.2	20
Iron						
Lead	8.1	105	114	85.0	4.7	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	14.2	60.9	57	83.9	3.9	20
Phosphorus						
Potassium						
Selenium	0.0	100	114	87.8	4.9	20
Silicon						
Silver	0.15	21.0	22.8	91.5	6.0	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	34.8	79.8	57	79.0	4.4	20

Associated samples MP11290: D51203-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11290
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51203
 Account: XTOKRWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11290
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 10/03/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	196	200	98.0	80-120
Beryllium				
Boron	anr			
Cadmium	47.4	50	94.8	80-120
Calcium				
Chromium	50.5	50	101.0	80-120
Cobalt				
Copper	52.0	50	104.0	80-120
Iron				
Lead	97.5	100	97.5	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	48.3	50	96.6	80-120
Phosphorus				
Potassium				
Selenium	98.7	100	98.7	80-120
Silicon				
Silver	20.6	20	103.0	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	47.1	50	94.2	80-120

Associated samples MP11290: D51203-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11290
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51203
 Account: XTOKWR - XTO Energy
 Project: FRU 197-31A

QC Batch ID: MP11290
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: ug/l

Prep Date:

10/03/13

Metal	D51202-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	1960	2150	10.2*(a)	0-10
Beryllium				
Boron	anr			
Cadmium	1.10	0.00	100.0(b)	0-10
Calcium				
Chromium	358	403	12.6*(a)	0-10
Cobalt				
Copper	61.0	67.0	11.9 (b)	0-10
Iron				
Lead	91.7	59.5	14.3 (b)	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	112	131	16.1*(a)	0-10
Phosphorus				
Potassium				
Selenium	0.00	0.00	NC	0-10
Silicon				
Silver	1.30	1.50	15.4 (b)	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	321	357	19.0*(a)	0-10

Associated samples MP11290: D51203-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11290
Matrix Type: SOLID

Methods: SW846 6010C
Units: ug/l

Prep Date:

Metal

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11291
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date: 10/03/13

Metal	RL	IDL	MDL	MB raw	final
Arsenic	0.10	.0085	.024	0.0076	<0.10

Associated samples MP11291: D51203-1

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: D51203
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11291
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date: 10/03/13

Metal	D51202-1 Original MS	Spikelot ICPALL2	QC % Rec	QC Limits
Arsenic	7.0	123	117	98.8 75-125

Associated samples MP11291: D51203-1

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: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11291
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date:

10/03/13

Metal	D51202-1 Original	MSD	Spikelot ICPALL2	MSD % Rec	RPD	QC Limit
Arsenic	7.0	118	114	97.4	4.1	20

Associated samples MP11291: D51203-1

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: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11291
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date: 10/03/13

Metal	BSP Result	Spikelot ICPALL2	QC % Rec	Limits
Arsenic	96.8	100	96.8	80-120

Associated samples MP11291: D51203-1

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: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11291
Matrix Type: SOLID

Methods: SW846 6020A
Units: ug/l

Prep Date:

10/03/13

Metal	D51202-1	Original	SDL	5:25	%DIF	QC	Limits
Arsenic	60.0	60.9	1.6			0-10	

Associated samples MP11291: D51203-1

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51203
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11292
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	-0.00086	<0.10

Associated samples MP11292: D51203-1

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

14.3.1
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
Account: XTOKWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11292
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 10/04/13

Metal	D51148-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.037	0.45	0.389	106.2 75-125

Associated samples MP11292: D51203-1

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: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11292
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

10/04/13

Metal	D51148-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.037	0.42	0.364	105.3	6.9	20

Associated samples MP11292: D51203-1

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: D51203
Account: XTOKRWR - XTO Energy
Project: FRU 197-31A

QC Batch ID: MP11292
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 10/04/13

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.41	0.4	102.5	80-120

Associated samples MP11292: D51203-1

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

14.3.3
14

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51203
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: D51203-1A

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D51203
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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
 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 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: D51203-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
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

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
 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: D51203-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D51203
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

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51203
 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: D51203-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D51203
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: D51203
 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	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: D51203-1A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D51203
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: D51203
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	GP11117/GN22215	1.0	0.32	mg/kg	106	104	97.7	80-120%
Chromium, Hexavalent	GP11117/GN22215	1.0	0.36	mg/kg	106	104	97.9	80-120%
Specific Conductivity	GP11110/GN22203			umhos/cm	9979	9980	100.0	90-110%
pH	GN22154			su	8.00	8.03	100.4	99.3-100.7%

Associated Samples:
Batch GN22154: D51203-1
Batch GP11110: D51203-1
Batch GP11117: D51203-1
(*) Outside of QC limits

DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51203
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	GP11117/GN22215 GN22168	D51202-1 D51122-1	mg/kg mv	0.38 133	0.50 130	26.2(a) 2.3	0-20% 0-20%

Associated Samples:

Batch GN22168: D51203-1

Batch GP11117: D51203-1

(*) Outside of QC limits

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

MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51203
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	GP11117/GN22215	D51202-1	mg/kg	0.38	40	35.8	89.5	75-125%

Associated Samples:

Batch GP11117: D51203-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D51203
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	GP11117/GN22215	D51202-1	mg/kg	0.38	40	35.4	1.2	20%

Associated Samples:

Batch GP11117: D51203-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits