



01/17/13



Technical Report for

XTO Energy

PCU 296-5A

1210-04

Accutest Job Number: D42511

Sampling Date: 01/07/13

Report to:

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Total number of pages in report: 150



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, appearing to read "H. Madadian".

Brad Madadian
Laboratory Director

Client Service contact: Renea Jackson 303-425-6021

Certifications: CO, ID, NE, NM, ND (R-027) (PW), UT (NELAP CO00049), TX (T104704511-12-1)

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Sample Summary

XTO Energy

Job No: D42511

PCU 296-5A

Project No: 1210-04

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D42511-1	01/07/13	13:20 DS	01/10/13	SO	Soil	CUT 1 SUBLINER COMP
D42511-1A	01/07/13	13:20 DS	01/10/13	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 D42511

Site: PCU 296-5A

Report Date 1/17/2013 1:09:39 PM

On 01/10/2013, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 4.1 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D42511 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: V3V1327
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42512-1MS, D42512-1MSD were used as the QC samples indicated.

Extractables by GCMS By Method SW846 8270C BY SIM

Matrix SO	Batch ID: OP7223
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42510-1MS, D42510-1MSD were used as the QC samples indicated.
- The matrix spike (MS) recovery(s) of Naphthalene are outside control limits. Outside control limits due to high level in sample relative to spike amount.
- The RPD(s) for the MS and MSD recoveries of Naphthalene are outside control limits for sample OP7223-MSD. Variability of recovery may be due to sample matrix/homogeneity.

Volatiles by GC By Method SW846 8015B

Matrix SO	Batch ID: GGB1044
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42435-1MS, D42435-1MSD were used as the QC samples indicated.

Extractables by GC By Method SW846-8015B

Matrix SO	Batch ID: OP7222
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D42509-2MS, D42509-2MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

Metals By Method SW846 6010C

Matrix AQ

Batch ID: MP9237

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42427-2MS, D42427-2MSD, D42427-2SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Calcium, Sodium are outside control limits. Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.
- The matrix spike duplicate (MSD) recovery(s) of Calcium, Sodium are outside control limits. Probable cause due to matrix interference.
- The serial dilution RPD(s) for Magnesium, Calcium, Sodium are outside control limits for sample MP9237-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP9237-SD1 for Calcium: Serial dilution indicates possible matrix interference.
- MP9237-SD1 for Sodium: Serial dilution indicates possible matrix interference.

Matrix SO

Batch ID: MP9242

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42510-1MS, D42510-1MSD, D42510-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The serial dilution RPD(s) for Chromium, Lead, Zinc are outside control limits for sample MP9242-SD1. Probable cause due to sample homogeneity.
- MP9242-SD1 for Zinc, Chromium, Lead: Serial dilution indicates possible matrix interference.

Metals By Method SW846 6020A

Matrix SO

Batch ID: MP9243

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42510-1MS, D42510-1SDL, D42510-1MSD were used as the QC samples for the metals analysis.

Metals By Method SW846 7471B

Matrix SO

Batch ID: MP9244

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42445-1MS, D42445-1MSD were used as the QC samples for the metals analysis.

Wet Chemistry By Method ASTM D1498-76M

Matrix SO

Batch ID: GN18410

- Sample(s) D42511-1DUP were used as the QC samples for the Redox Potential Vs H₂ analysis.

Wet Chemistry By Method SM 2510B-2011 MOD

Matrix SO

Batch ID: GP9098

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

Wet Chemistry By Method SM19 2540B M

Matrix SO

Batch ID: GN18390

- The data for SM19 2540B M meets quality control requirements.

Wet Chemistry By Method SW846 3060A/7196A

Matrix SO

Batch ID: GP9086

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42556-1DUP, D42556-1MS, D42556-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

Wet Chemistry By Method SW846 3060A/7196A M

Matrix SO

Batch ID: R15683

- The data for SW846 3060A/7196A M meets quality control requirements.
- D42511-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Wet Chemistry By Method SW846 9045D

Matrix SO

Batch ID: GN18407

- The following samples were run outside of holding time for method SW846 9045D: D42511-1

Wet Chemistry By Method USDA HANDBOOK 60

Matrix SO

Batch ID: MP9237

- D42511-1A 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: D42511
Account: XTO Energy
Project: PCU 296-5A
Collected: 01/07/13

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

D42511-1 CUT 1 SUBLINER COMP

Ethylbenzene	0.0319 J	0.15	0.028	mg/kg	SW846 8260B
Naphthalene	0.0222	0.014	0.013	mg/kg	SW846 8270C BY SIM
TPH-DRO (C10-C28)	66.6	8.2	4.9	mg/kg	SW846-8015B
Arsenic	8.1	0.12		mg/kg	SW846 6020A
Barium	358	1.2		mg/kg	SW846 6010C
Chromium	56.8	1.2		mg/kg	SW846 6010C
Copper	10.4	1.2		mg/kg	SW846 6010C
Lead	9.6	6.2		mg/kg	SW846 6010C
Nickel	19.7	3.7		mg/kg	SW846 6010C
Zinc	40.0	3.7		mg/kg	SW846 6010C
Specific Conductivity	1560	1.0		umhos/cm	SM 2510B-2011 MOD
Chromium, Trivalent ^a	56.8	2.2		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	137			mv	ASTM D1498-76M
pH	11.08			su	SW846 9045D

D42511-1A CUT 1 SUBLINER COMP

Calcium	23.4	2.0	mg/l	SW846 6010C
Magnesium	3.13	1.0	mg/l	SW846 6010C
Sodium	275	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	14.2		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:** D42511-1**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** PCU 296-5A**Date Sampled:** 01/07/13**Date Received:** 01/10/13**Percent Solids:** 81.0

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3V22567.D	1	01/12/13	BD	n/a	n/a	V3V1327
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.03 g	5.0 ml	100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	0.073	0.037	mg/kg	
108-88-3	Toluene	ND	0.15	0.073	mg/kg	
100-41-4	Ethylbenzene	0.0319	0.15	0.028	mg/kg	J
1330-20-7	Xylene (total)	ND	0.29	0.15	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	86%		64-130%
460-00-4	4-Bromofluorobenzene	105%		62-131%
17060-07-0	1,2-Dichloroethane-D4	87%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	01/07/13
Lab Sample ID:	D42511-1	Date Received:	01/10/13
Matrix:	SO - Soil	Percent Solids:	81.0
Method:	SW846 8270C BY SIM	SW846 3546	
Project:	PCU 296-5A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G12978.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
Run #2							

	Initial Weight	Final Volume
Run #1	30.0 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	0.010	0.0053	mg/kg	
120-12-7	Anthracene	ND	0.010	0.0053	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.010	0.0053	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.010	0.0053	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.010	0.0053	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.010	0.0053	mg/kg	
218-01-9	Chrysene	ND	0.010	0.0053	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.010	0.0053	mg/kg	
206-44-0	Fluoranthene	ND	0.010	0.0053	mg/kg	
86-73-7	Fluorene	ND	0.010	0.0053	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.010	0.0053	mg/kg	
91-20-3	Naphthalene	0.0222	0.014	0.013	mg/kg	
129-00-0	Pyrene	ND	0.010	0.0053	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	78%		10-159%
321-60-8	2-Fluorobiphenyl	78%		19-131%
1718-51-0	Terphenyl-d14	96%		18-150%

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

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Report of Analysis

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Client Sample ID: CUT 1 SUBLINER COMP**Lab Sample ID:** D42511-1**Date Sampled:** 01/07/13**Matrix:** SO - Soil**Date Received:** 01/10/13**Method:** SW846 8015B**Percent Solids:** 81.0**Project:** PCU 296-5A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB19112.D	1	01/11/13	SK	n/a	n/a	GGB1044
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	15	7.3	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	90%		60-140%		

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Report of Analysis

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Client Sample ID: CUT 1 SUBLINER COMP**Lab Sample ID:** D42511-1**Date Sampled:** 01/07/13**Matrix:** SO - Soil**Date Received:** 01/10/13**Method:** SW846-8015B SW846 3546**Percent Solids:** 81.0**Project:** PCU 296-5A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FH008578.D	1	01/15/13	AV	01/14/13	OP7222	GFH472
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)	66.6	8.2	4.9	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	67%		35-130%		

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

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Report of Analysis

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Client Sample ID: CUT 1 SUBLINER COMP**Lab Sample ID:** D42511-1**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/07/13**Date Received:** 01/10/13**Percent Solids:** 81.0**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	8.1	0.12	mg/kg	5	01/14/13	01/16/13 JB	SW846 6020A ³	SW846 3050B ⁵
Barium	358	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Cadmium	< 1.2	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Chromium	56.8	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Copper	10.4	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Lead	9.6	6.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Mercury	< 0.10	0.10	mg/kg	1	01/15/13	01/15/13 JB	SW846 7471B ²	SW846 7471B ⁶
Nickel	19.7	3.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Selenium	< 6.2	6.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Silver	< 3.7	3.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴
Zinc	40.0	3.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C ¹	SW846 3050B ⁴

- (1) Instrument QC Batch: MA3169
- (2) Instrument QC Batch: MA3177
- (3) Instrument QC Batch: MA3182
- (4) Prep QC Batch: MP9242
- (5) Prep QC Batch: MP9243
- (6) Prep QC Batch: MP9244

RL = Reporting Limit

Report of Analysis

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Client Sample ID: CUT 1 SUBLINER COMP**Lab Sample ID:** D42511-1**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/07/13**Date Received:** 01/10/13**Percent Solids:** 81.0**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	1560	1.0	umhos/cm	1	01/15/13	KB	SM 2510B-2011 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	01/14/13	KB	SW846 3060A/7196A
Chromium, Trivalent ^a	56.8	2.2	mg/kg	1	01/14/13 17:18	JB	SW846 3060A/7196A M
Redox Potential Vs H2	137		mv	1	01/11/13	CT	ASTM D1498-76M
Solids, Percent	81		%	1	01/11/13	SWT	SM19 2540B M
pH	11.08		su	1	01/11/13 13:30	CT	SW846 9045D

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

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	01/07/13
Lab Sample ID:	D42511-1A	Date Received:	01/10/13
Matrix:	SO - Soil	Percent Solids:	81.0
Project:	PCU 296-5A		

SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	23.4	2.0	mg/l	1	01/11/13	01/11/13 JB	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	3.13	1.0	mg/l	1	01/11/13	01/11/13 JB	SW846 6010C ¹	SW846 3010A/M ²
Sodium	275	2.0	mg/l	1	01/11/13	01/11/13 JB	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA3166

(2) Prep QC Batch: MP9237

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 1 SUBLINER COMP	Date Sampled:	01/07/13
Lab Sample ID:	D42511-1A	Date Received:	01/10/13
Matrix:	SO - Soil	Percent Solids:	81.0
Project:	PCU 296-5A		

General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	14.2		ratio	1	01/11/13 17:52	JB	USDA HANDBOOK 60

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

RL = Reporting Limit



Misc. Forms

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Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody



CHAIN OF CUSTODY

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

FED-EX Tracking #	Bottle Order Control #
	D42511

Client / Reporting Information		Project Information		Requested Analyses (see TEST CODE sheet)										Matrix Codes		
Company Name KRW Consulting	Project Name: XTO PCU296-5A	Street Address 8000 West 14th Street; Suite 200	Street	Billing Information (if different from Report to)										DW - Drinking Water		
City Lakewood, CO 80214	City	State	Company Name XTO Energy											GW - Ground Water		
Project Contact Dwayne Knudson	Project # 1210-04		Street Address 21459 CR 5											WW - Water		
Phone # 970-488-1098	Client Purchase Order #		City Rifle, CO 81650											SW - Surface Water		
Sampler(s) Name(s) DAVID STADERS	Project Manager Joe Hess		Attention: Jessica Dooling											SO - Soil		
				Number of preserved Bottles	HCl	NH3	HNO3	H2SO4	NONE	D Water	MEOH	EDD	Butylate	SL - Sludge		
Accutest Sample #	Field ID / Point of Collection CUT 1 SUBLINER COMP	MEOH/ED Vial #	Date 1-7-13	Time 13:20	Sampled by	# of bottles 50									SED - Sediment	
															Cl - Oil	
															LIO - Other Liquid	
															AIR - Air	
															SOL - Other Solid	
															WP - Wipe	
															FB - Field Blank	
															RB - Rinse Blank	
															TB - Trip Blank	
															LAB USE ONLY	
															01	
															DRY 10	
Turnaround Time (Business days)		Data Deliverable Information										Comments / Special Instructions				
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 6 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"/>		Approved By (Accutest PM): / Date: <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 Commercial "A" = Results Only Commercial "B" = Results + QC Summary Commercial BN = Results/QC/Narrative (+ = chromatograms)										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: 1	Date/Time: 1/16/13 16:00	Received By: 1	Relinquished By: 2	Date Time:	Received By:	Relinquished By: 4	Date/Time: 1/16/13 16:00	Received By: 4	Relinquished By: 5	Received By: 5	Custody Seal # 1000	<input checked="" type="checkbox"/> Intact	<input type="checkbox"/> Not Intact	Preserved where applicable	Office Temp. EN	Cooler Temp. 410
Relinquished by Sampler: 3	Date/Time:	Received By: 3	Relinquished By: 4	Date/Time:	Received By: 4	Relinquished By: 5	Date/Time:	Received By: 5	Relinquished By: 5	Received By: 5						

D42511: Chain of Custody

Page 1 of 2



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D42511

Client: KRW

Immediate Client Services Action Required: No

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

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO PCU 296-5A

Airbill #'s: HDCO

Cooler SecurityY 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 TemperatureY or N

1. Temp criteria achieved:
2. Cooler temp verification: Infared gun
3. Cooler media: Ice (bag)

Quality Control PreservationY or N

N/A

1. Trip Blank present / cooler:
2. Trip Blank listed on COC:
3. Samples preserved properly:
4. VOCs headspace free:

Sample Integrity - DocumentationY or N

1. Sample labels present on bottles:
2. Container labeling complete:
3. Sample container label / COC agree:

Sample Integrity - ConditionY or N

1. Sample recvd within HT:
2. All containers accounted for:
3. Condition of sample: Intact

Sample Integrity - InstructionsY or N

N/A

1. Analysis requested is clear:
2. Bottles received for unspecified tests:
3. Sufficient volume rec'd for analysis:
4. Compositing instructions clear:
5. Filtering instructions clear:

Comments

Accutest Laboratories
V:(303) 425-60214036 Youngfield Street
F: (303) 425-6854Wheat Ridge, CO
www.accutest.com

5.1

5

D42511: Chain of Custody**Page 2 of 2**



GC/MS Volatiles

QC Data Summaries

Includes the following where applicable:

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



Method Blank Summary

Page 1 of 1

Job Number: D42511
Account: XTOKRWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1327-MB	3V22560.D	1	01/12/13	BD	n/a	n/a	V3V1327

The QC reported here applies to the following samples:

Method: SW846 8260B

D42511-1

6.1

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	89%
460-00-4	4-Bromofluorobenzene	97%
17060-07-0	1,2-Dichloroethane-D4	98%

Blank Spike Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1327-BS	3V22561.D	1	01/12/13	BD	n/a	n/a	V3V1327

The QC reported here applies to the following samples:

Method: SW846 8260B

D42511-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	47.4	95	70-130
100-41-4	Ethylbenzene	50	48.7	97	70-130
108-88-3	Toluene	50	47.3	95	70-130
1330-20-7	Xylene (total)	150	149	99	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	91%	64-130%
460-00-4	4-Bromofluorobenzene	109%	62-131%
17060-07-0	1,2-Dichloroethane-D4	88%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D42512-1MS	3V22564.D	1	01/12/13	BD	n/a	n/a	V3V1327
D42512-1MSD	3V22565.D	1	01/12/13	BD	n/a	n/a	V3V1327
D42512-1	3V22563.D	1	01/12/13	BD	n/a	n/a	V3V1327

The QC reported here applies to the following samples:

Method: SW846 8260B

D42511-1

CAS No.	Compound	D42512-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	72.3		3580	3400	93	3340	91	2	64-139/30
100-41-4	Ethylbenzene	35.0	J	3580	3530	98	3500	97	1	68-136/30
108-88-3	Toluene	138	J	3580	3330	89	3300	88	1	60-130/30
1330-20-7	Xylene (total)	212	J	10700	10800	99	10800	99	0	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D42512-1	Limits
2037-26-5	Toluene-D8	86%	87%	84%	64-130%
460-00-4	4-Bromofluorobenzene	113%	112%	104%	62-131%
17060-07-0	1,2-Dichloroethane-D4	90%	86%	90%	70-130%

* = Outside of Control Limits.

6.3.1
6



GC/MS Volatiles

Raw Data

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3011113.S\
 Data File : 3V22567.D
 Acq On : 12 Jan 2013 5:39 am
 Operator : BRETD
 Sample : D42511-1
 Misc : MS5218,V3V1327,5.034,,100,5,1
 ALS Vial : 34 Sample Multiplier: 1

Quant Time: Jan 15 08:23:28 2013
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M
 Quant Title : 8260
 QLast Update : Thu Jan 03 11:40:16 2013
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.861	168	354648	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.656	114	545471	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.296	117	612959	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.288	152	363792	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.252	102	34366	43.30	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	86.60%	
61) Toluene-d8	14.052	98	632427	42.85	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	85.70%	
69) 4-Bromofluorobenzene	16.246	95	332232	52.26	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	104.52%	

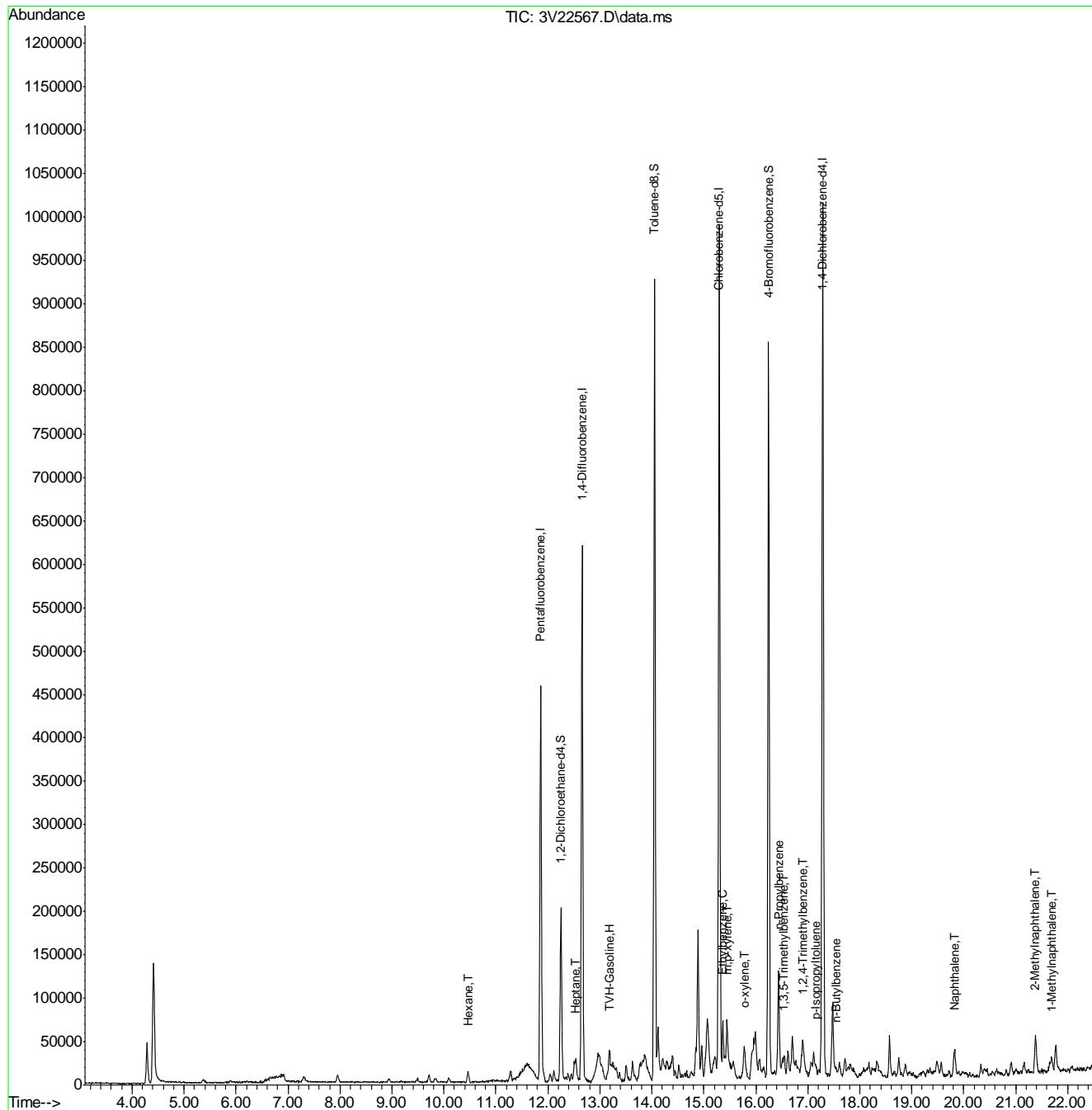
Target Compounds						Qvalue
1) TVH-Gasoline	13.200	TIC	3686969m	125.64	ug/l	
41) Hexane	10.462	57	6287	0.98	ug/l	100
43) Heptane	12.535	43	11700	1.60	ug/l	93
66) Ethylbenzene	15.367	91	9736	0.44	ug/l	89
72) m,p-xylene	15.447	106	15298	1.63	ug/l	91
73) o-xylene	15.797	106	2967	0.32	ug/l	89
77) n-Propylbenzene	16.426	91	6382	0.25	ug/l	87
80) 1,3,5-Trimethylbenzene	16.544	105	3948m	0.21	ug/l	
82) 1,2,4-Trimethylbenzene	16.894	105	8722	0.46	ug/l	89
86) p-Isopropyltoluene	17.157	119	7236	0.34	ug/l	89
88) n-Butylbenzene	17.545	91	4046	0.23	ug/l	85
91) Naphthalene	19.842	128	16004	1.84	ug/l	100
94) 2-Methylnaphthalene	21.378	142	32084	3.80	ug/l	96
95) 1-Methylnaphthalene	21.689	142	12578	1.57	ug/l	97

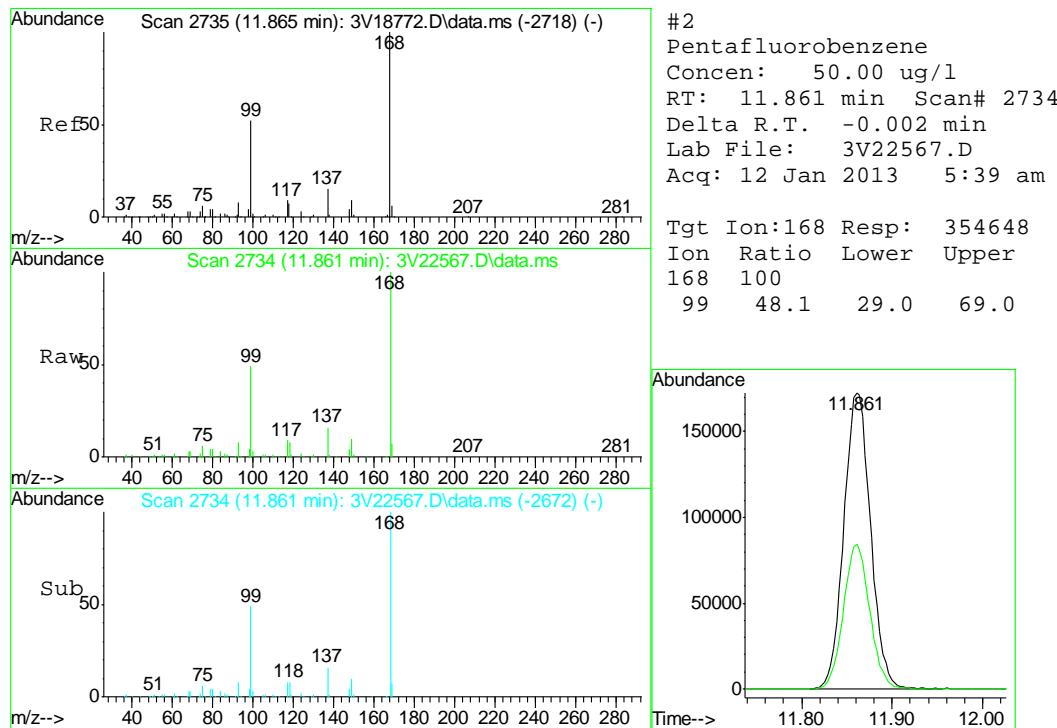
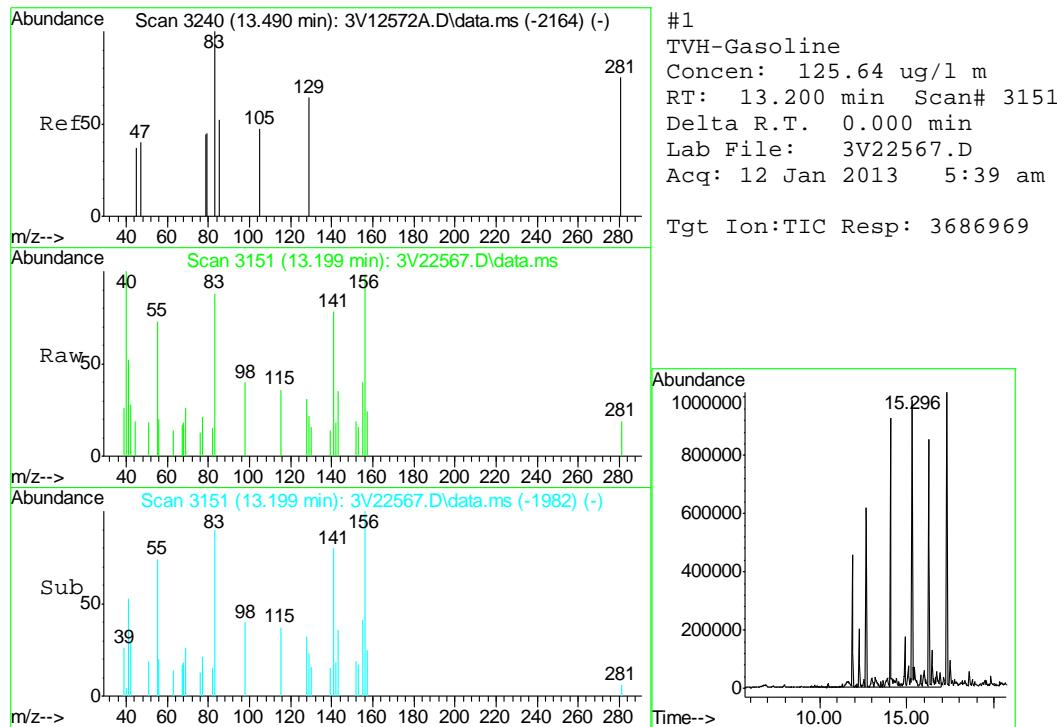
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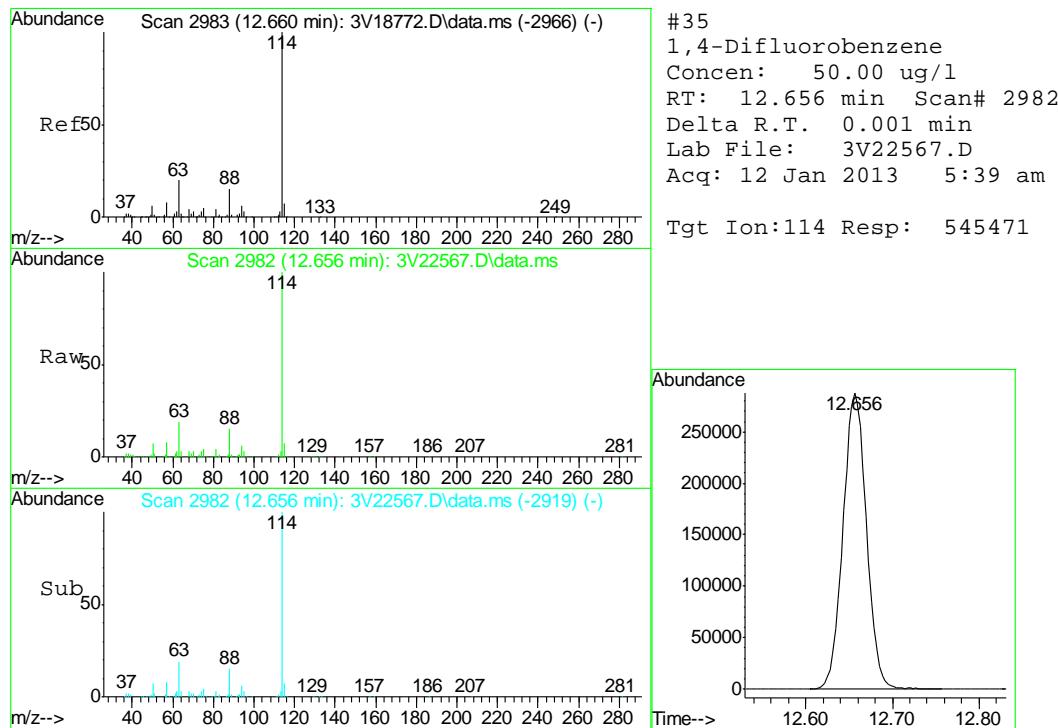
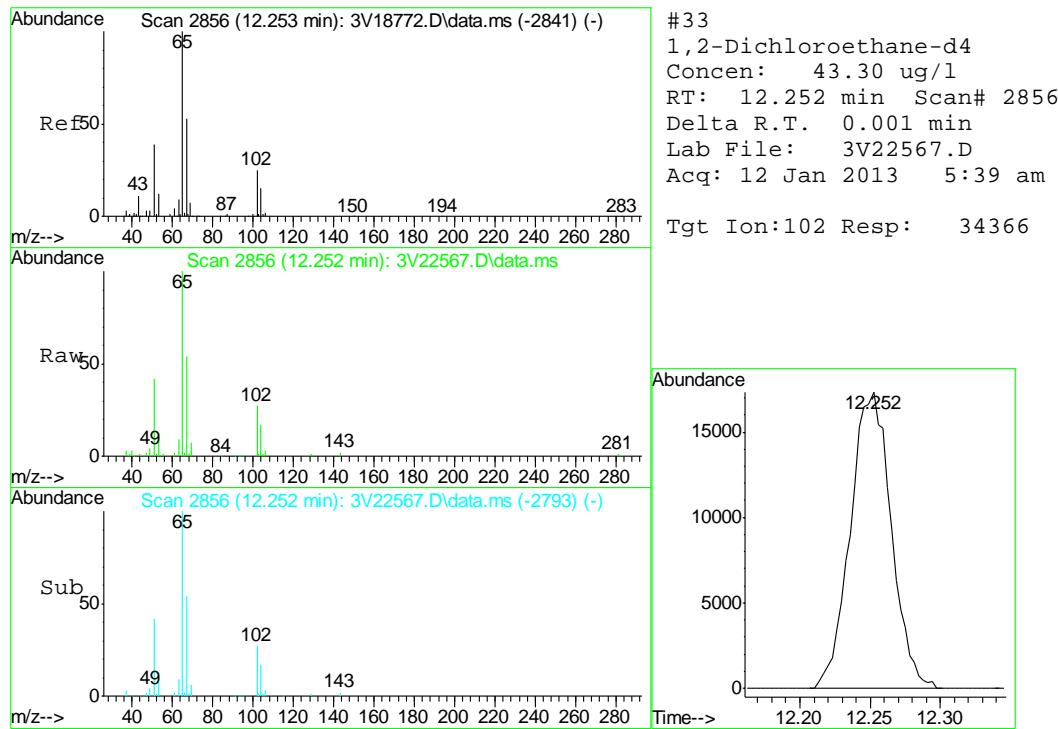
Quantitation Report (QT Reviewed)

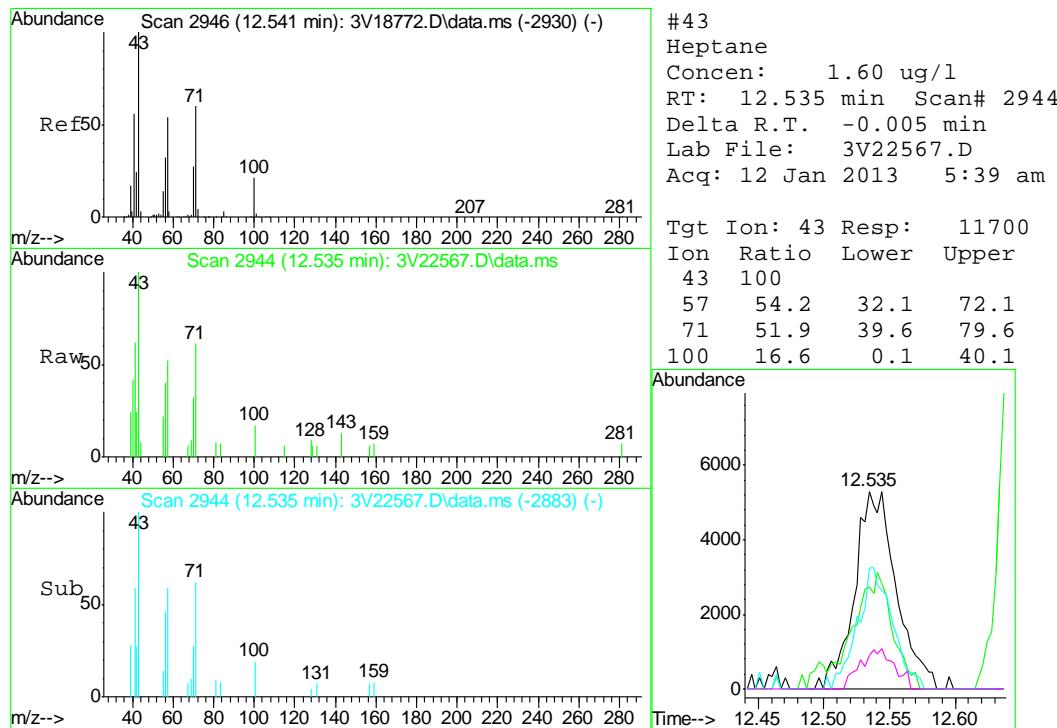
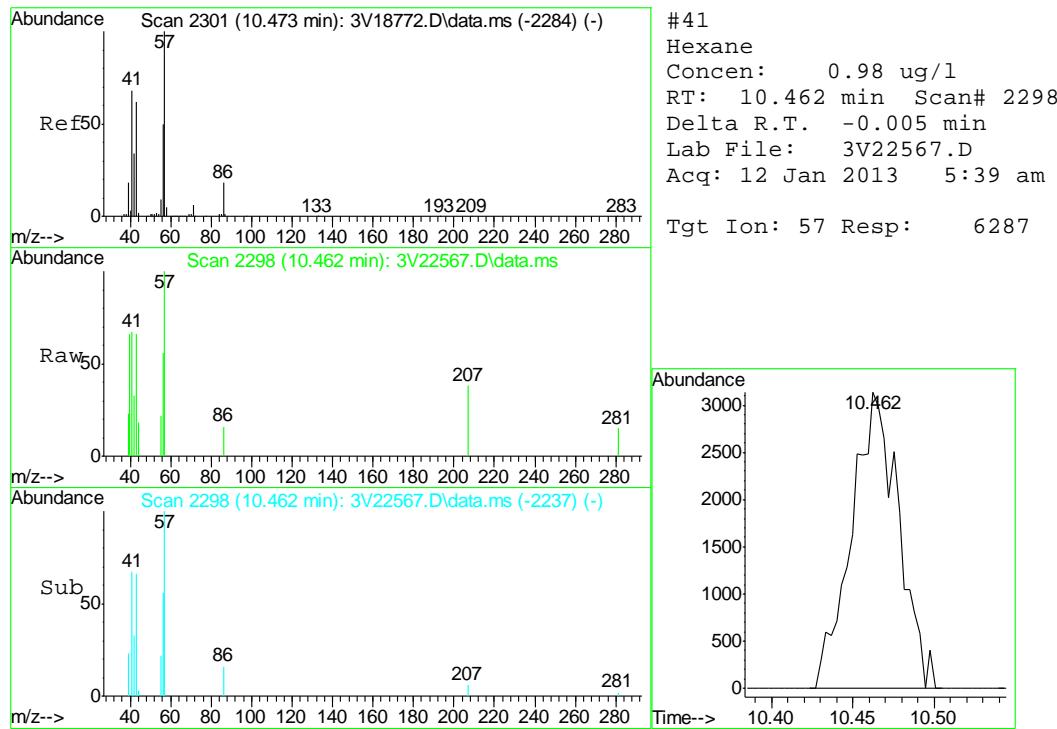
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 Data File : 3V22567.D
 Acq On : 12 Jan 2013 5:39 am
 Operator : BRETD
 Sample : D42511-1
 Misc : MS5218,V3V1327,5.034,,100,5,1
 ALS Vial : 34 Sample Multiplier: 1

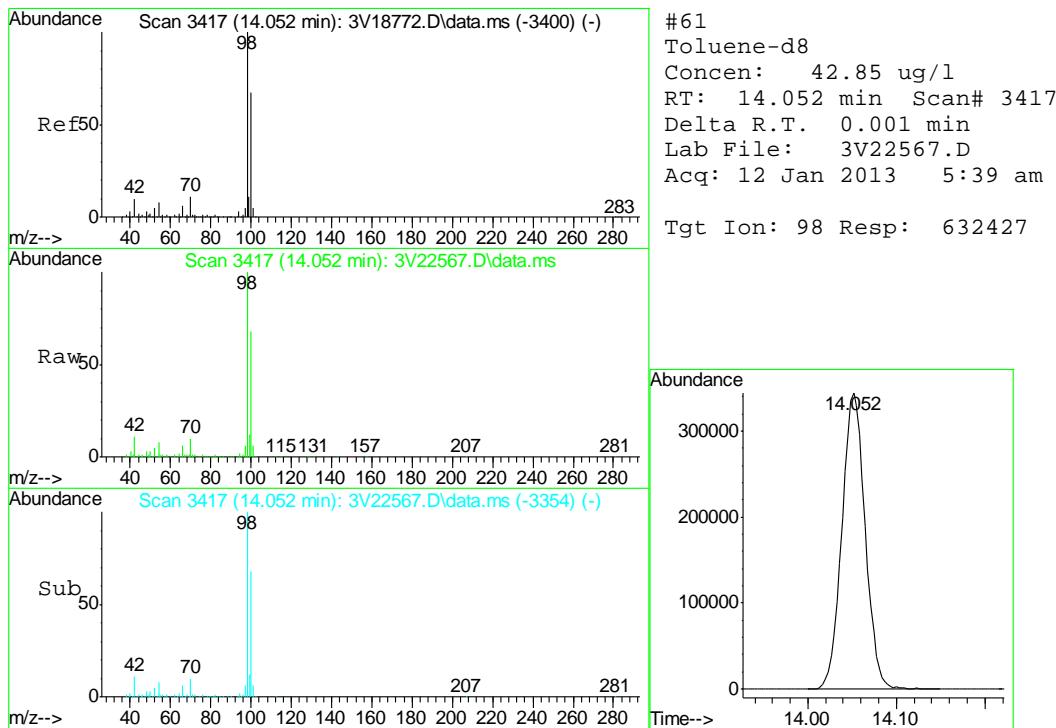
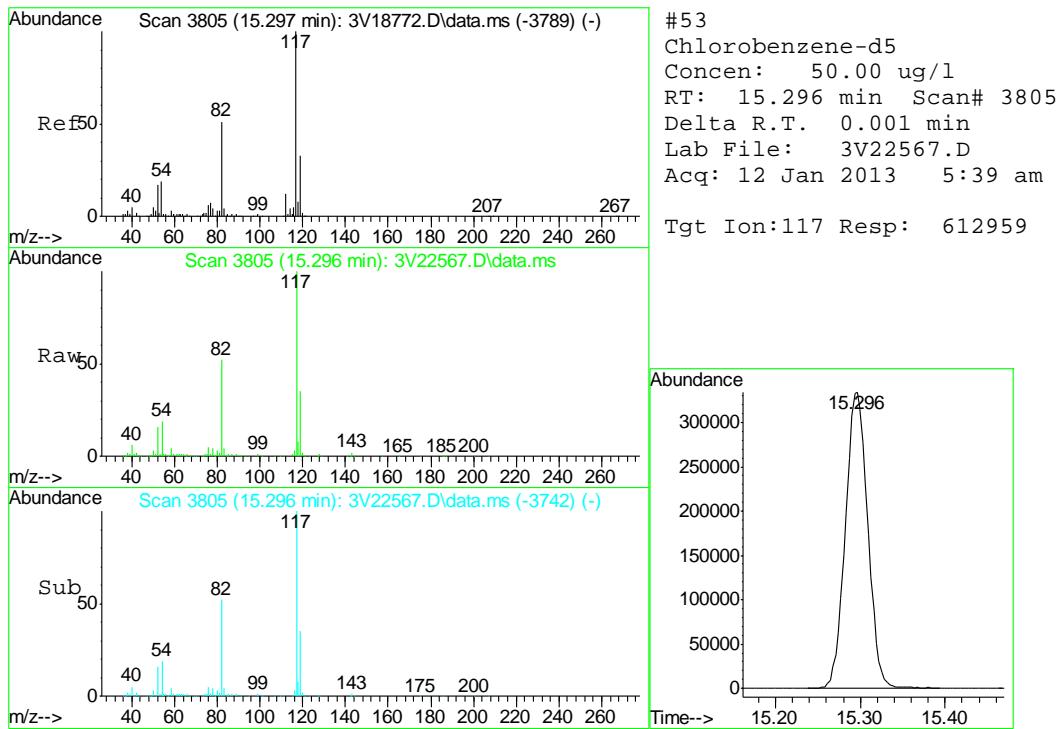
Quant Time: Jan 15 08:23:28 2013
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 Quant Title : 8260
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 Response via : Initial Calibration

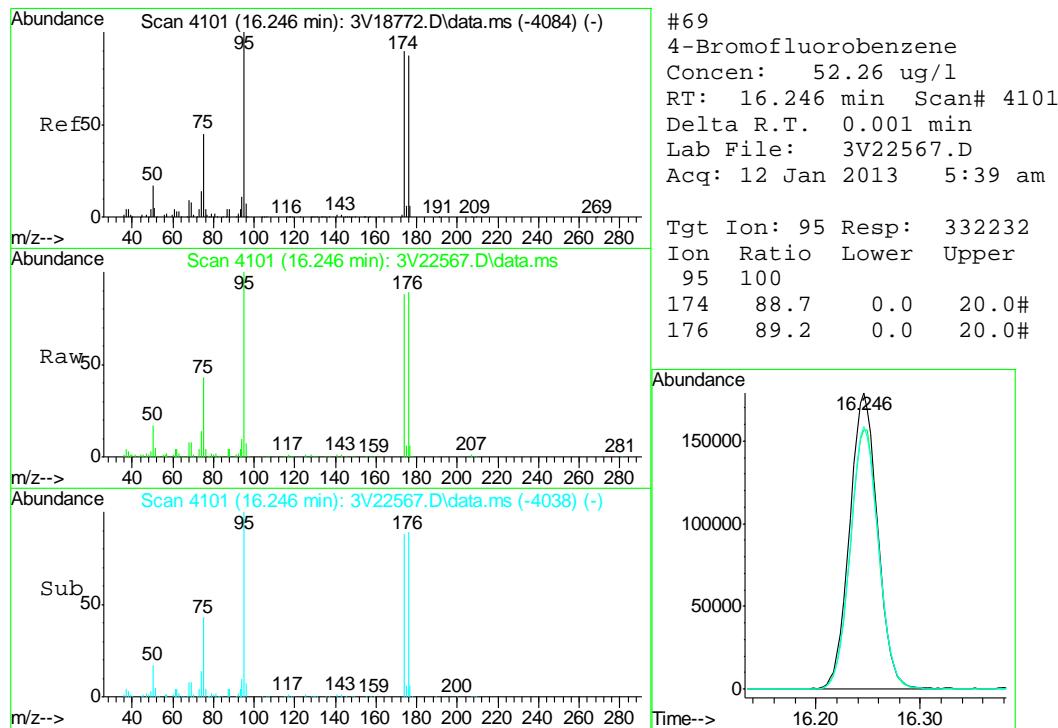
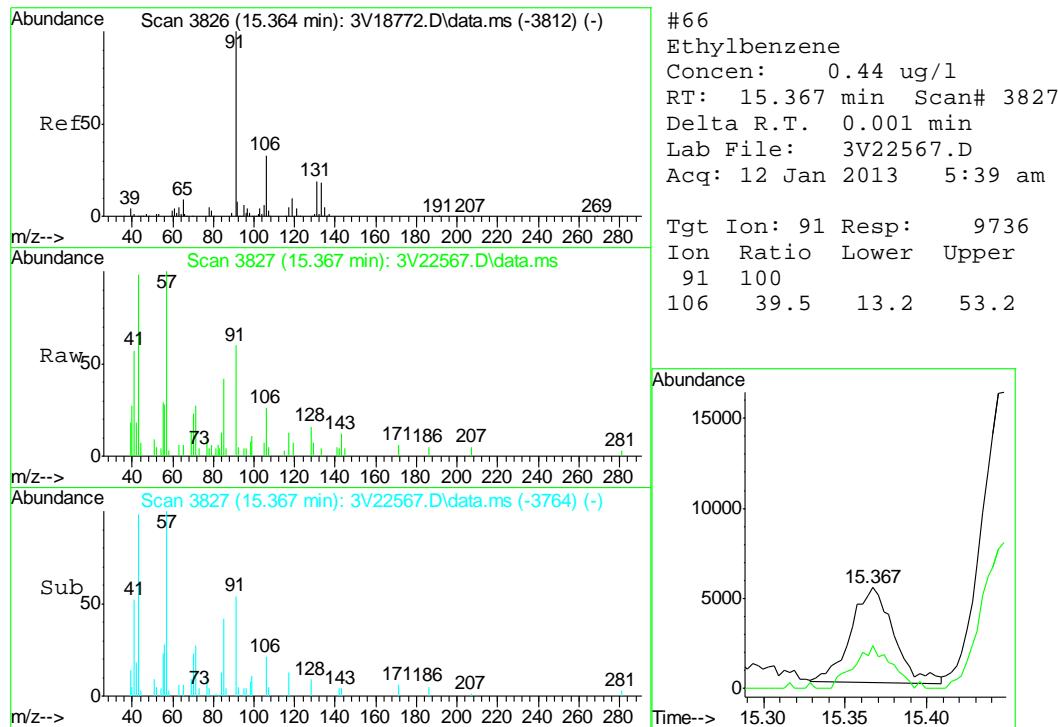


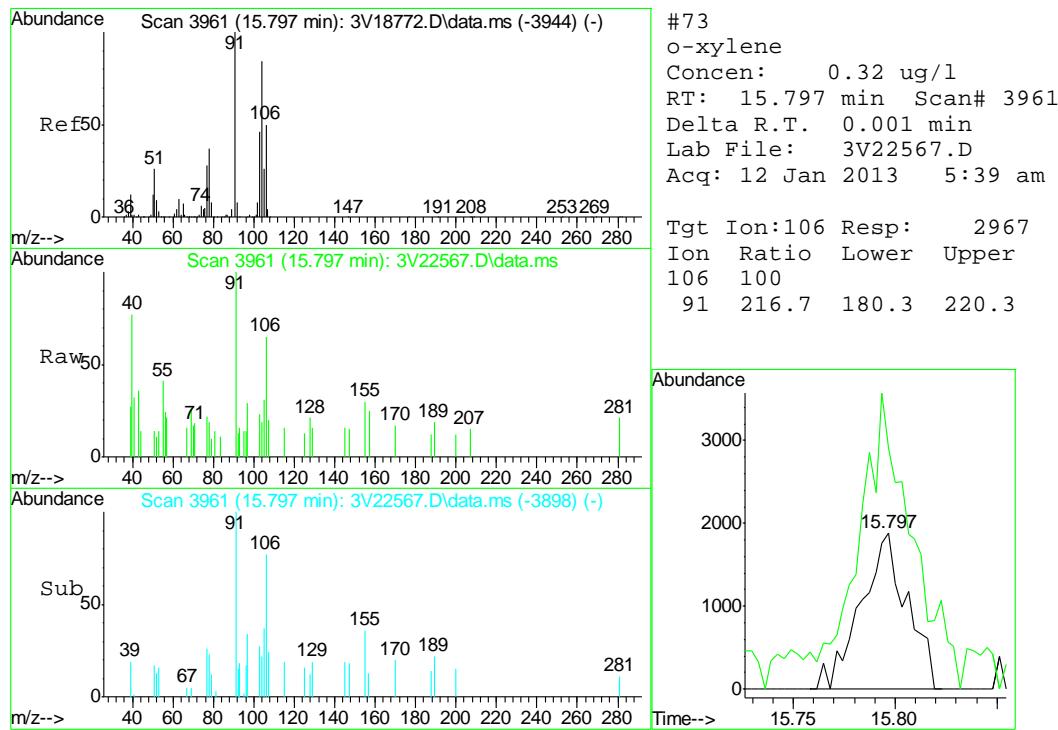
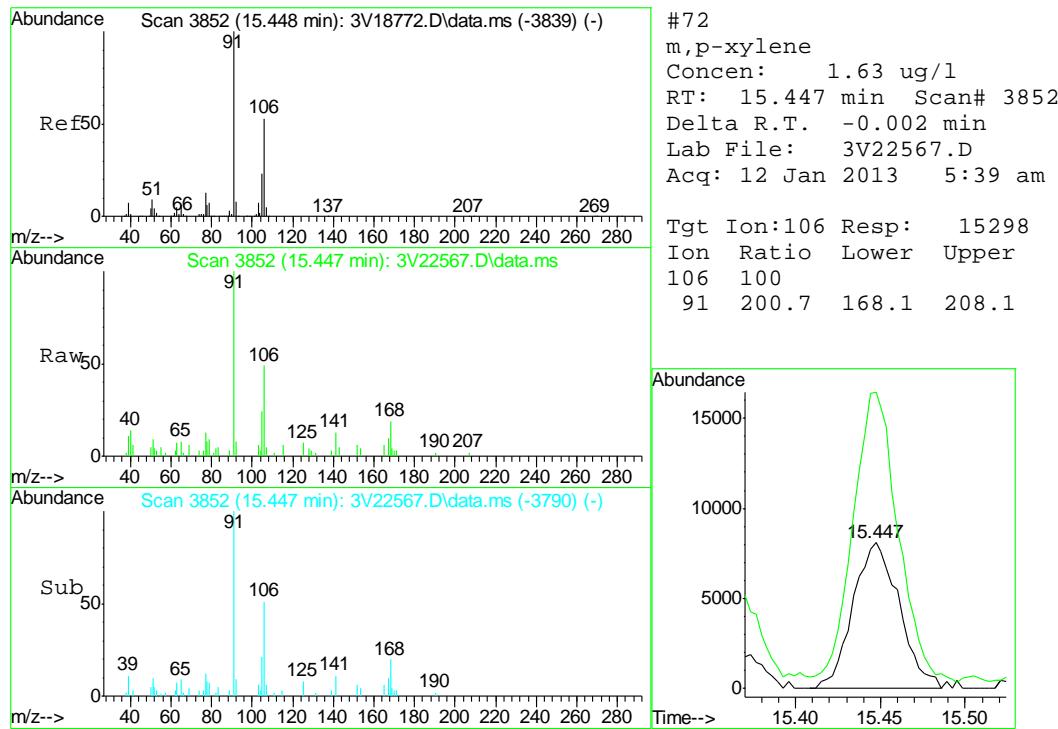


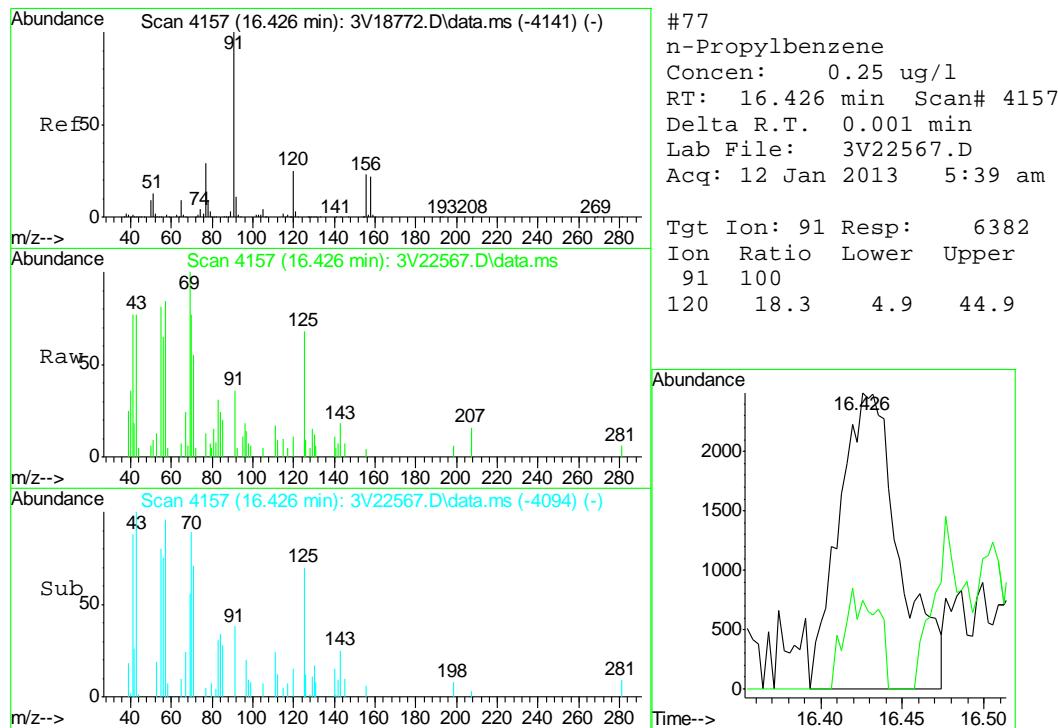
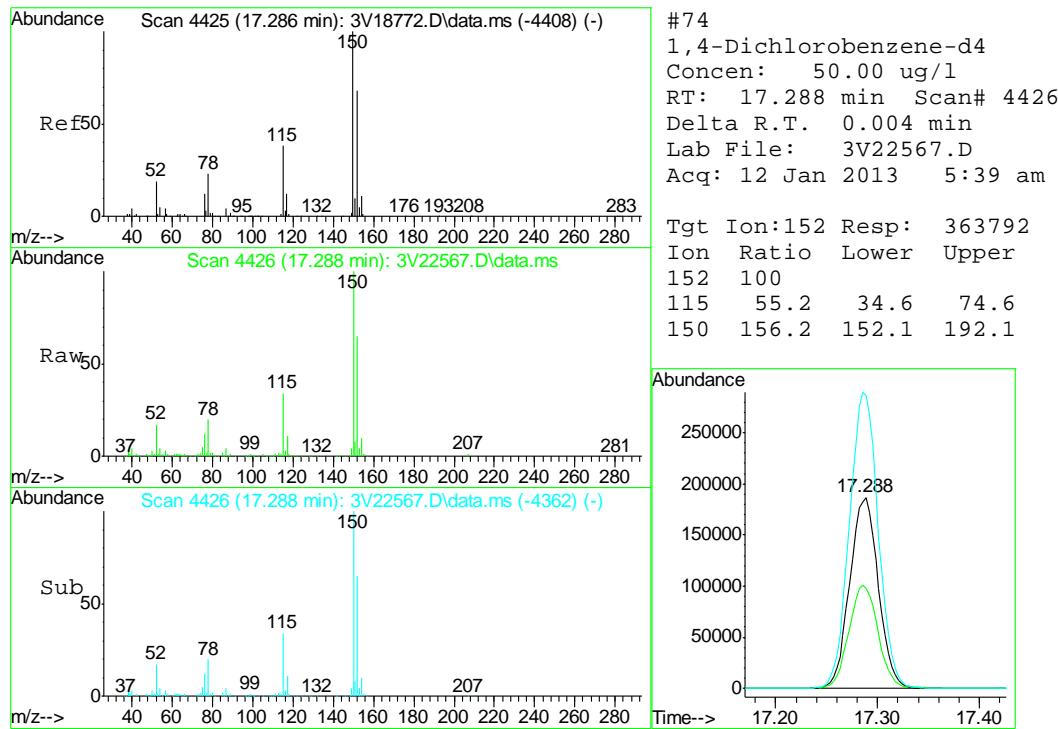


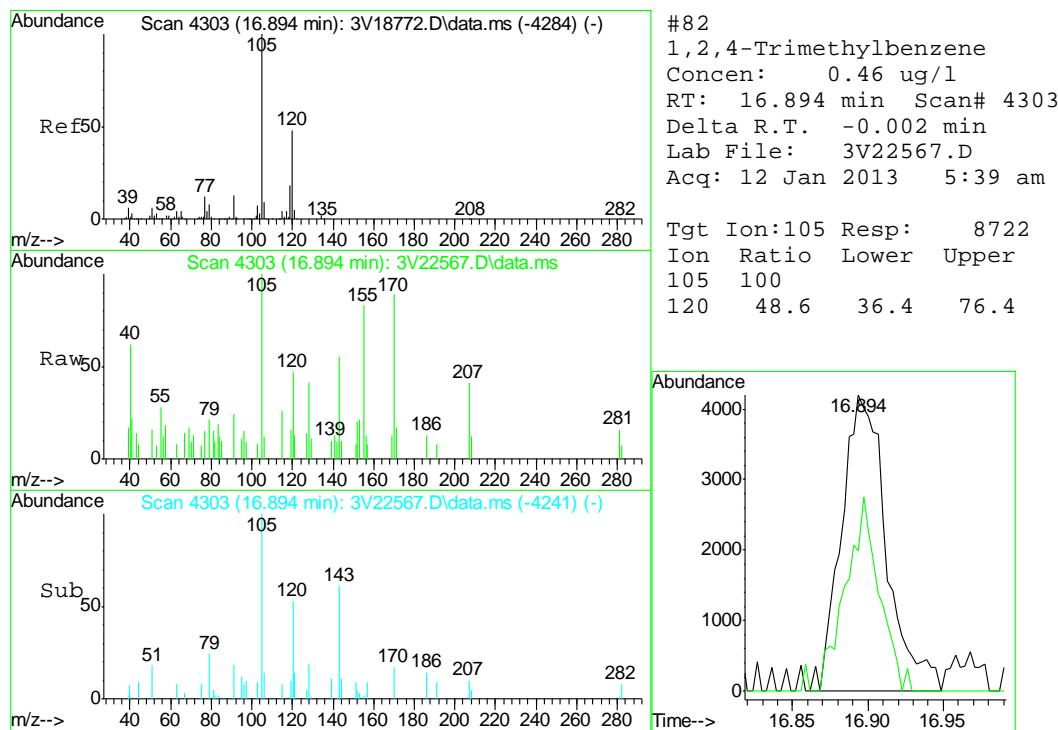
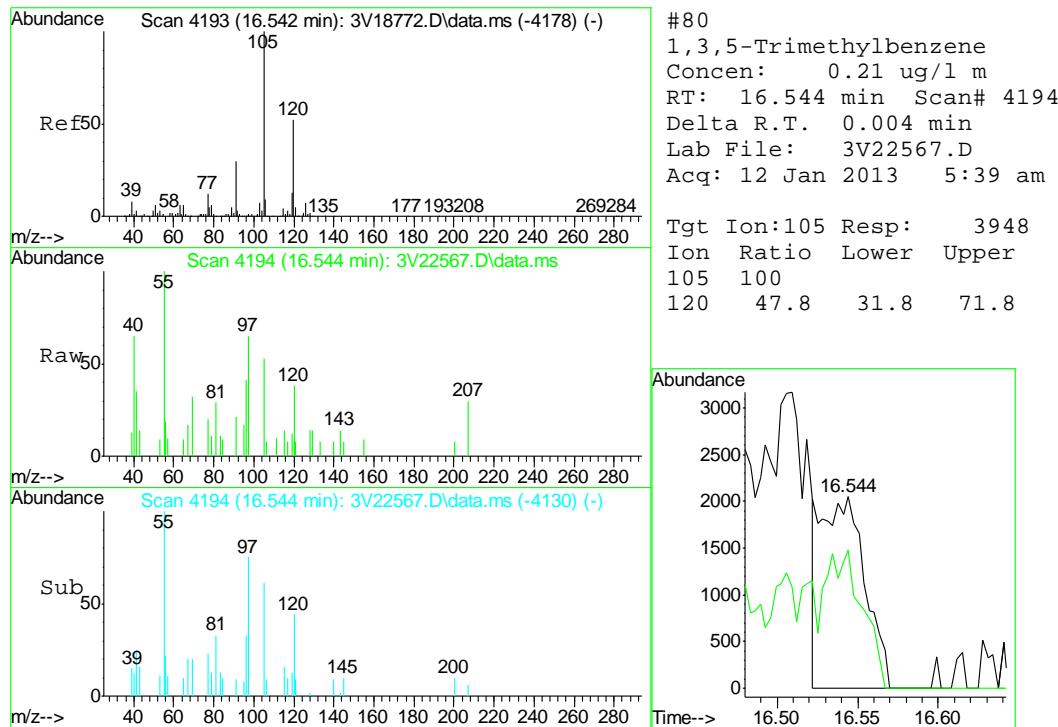


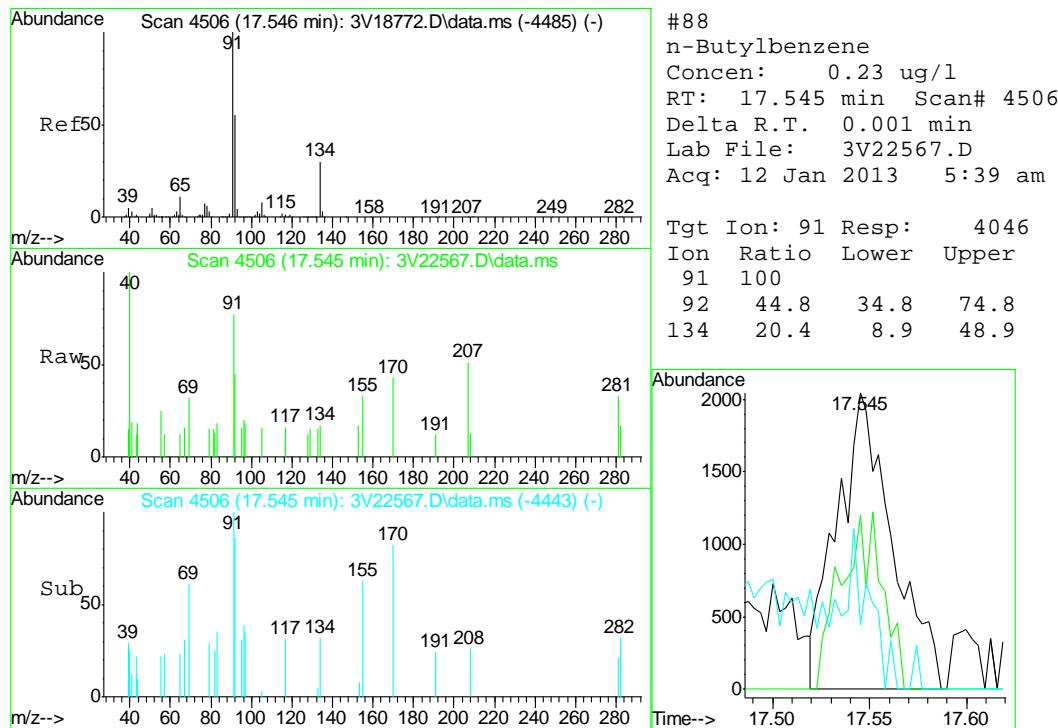
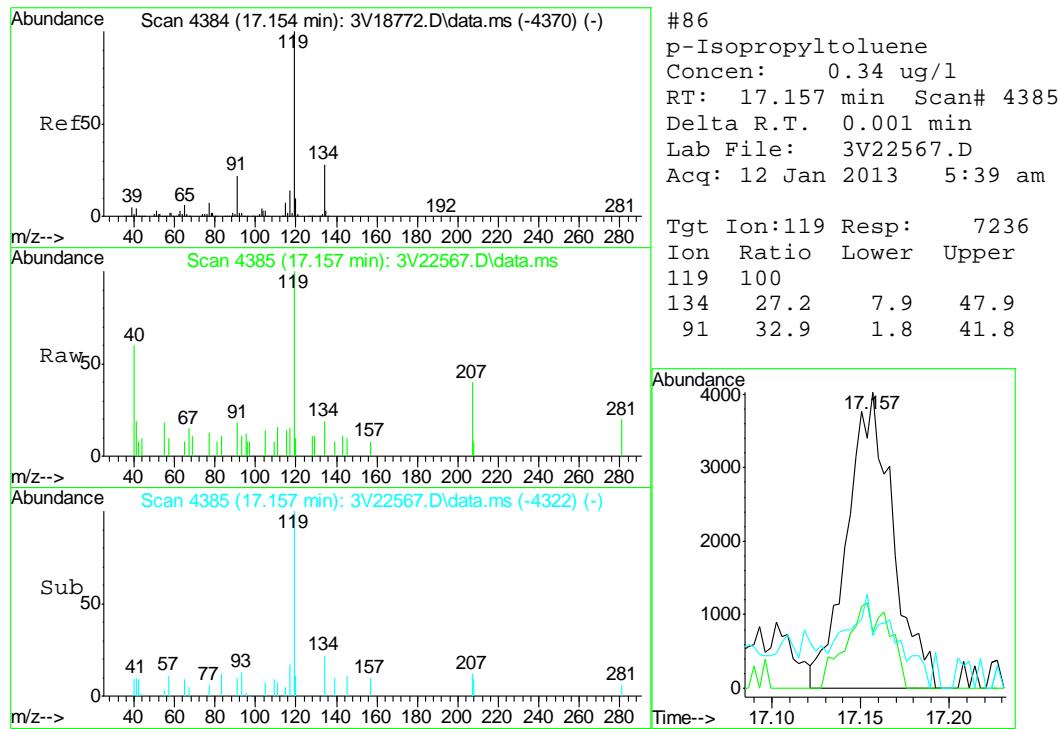


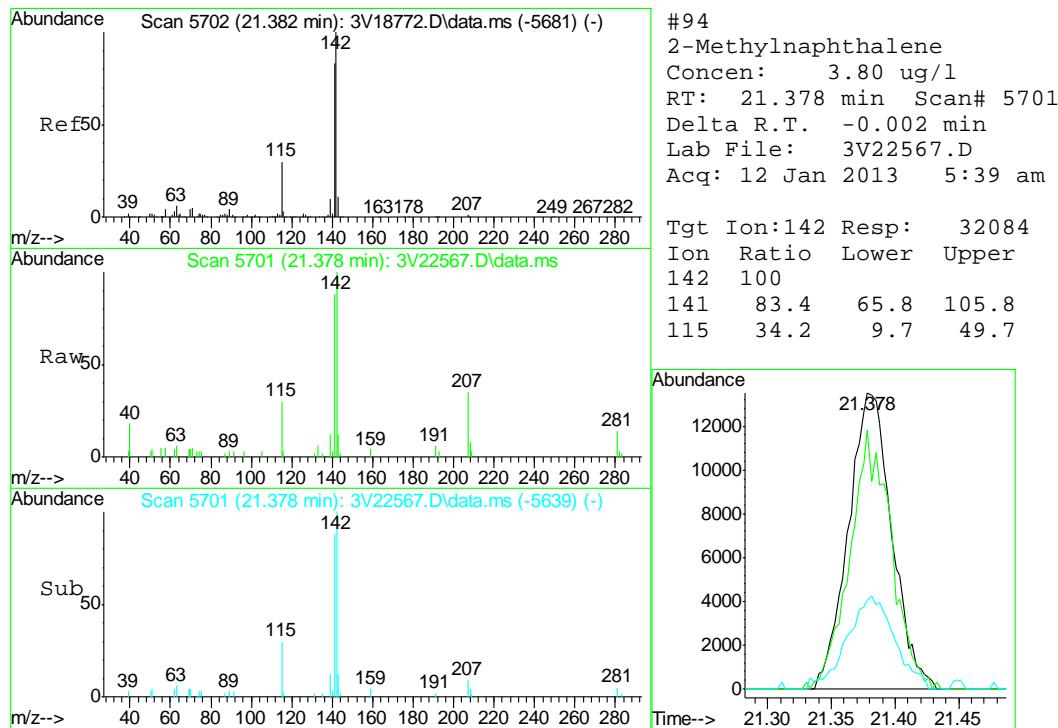
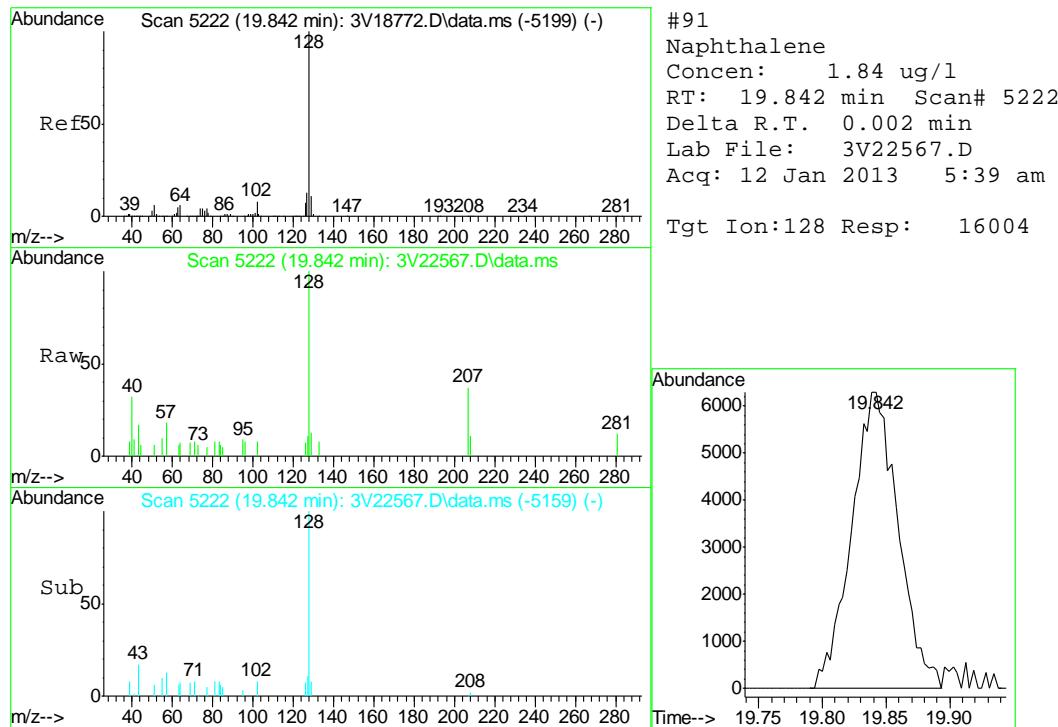


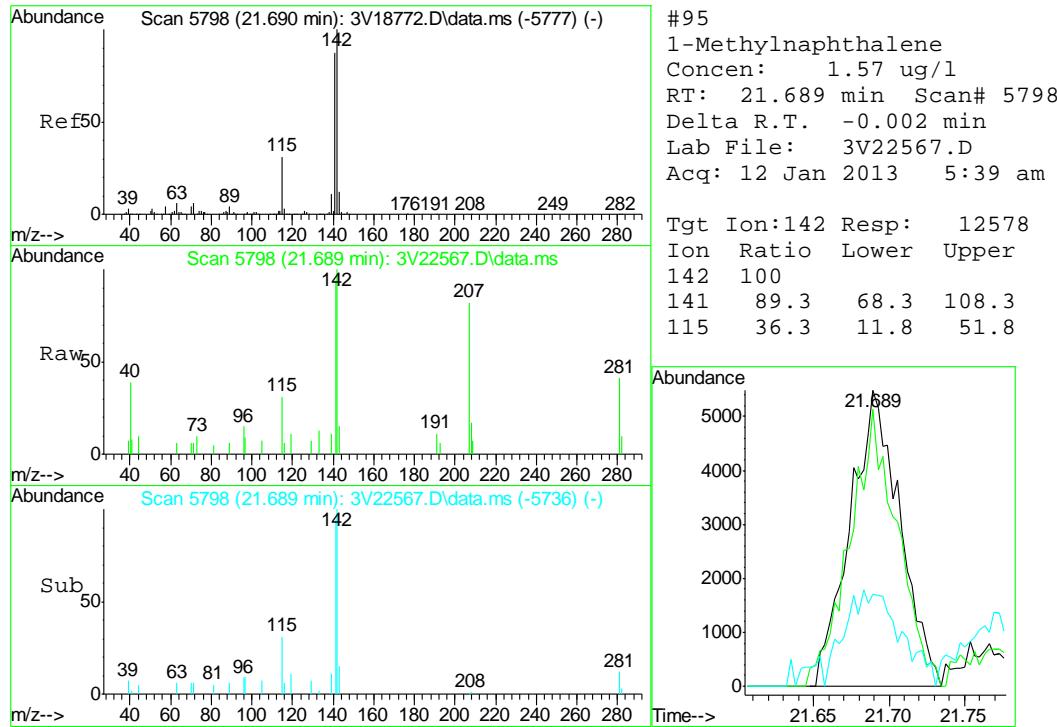












Quantitation Report (QT Reviewed)

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 Data File : 3V22560.D
 Acq On : 12 Jan 2013 2:00 am
 Operator : BRETD
 Sample : MB
 Misc : MS5218,V3V1327,5.00,,100,5,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jan 15 08:11:29 2013
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M
 Quant Title : 8260
 QLast Update : Thu Jan 03 11:40:16 2013
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.863	168	308471	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.659	114	500082	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.295	117	525362	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.287	152	304372	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.251	102	33686	48.79	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	97.58%	
61) Toluene-d8	14.054	98	562082	44.44	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	88.88%	
69) 4-Bromofluorobenzene	16.245	95	265181	48.67	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	97.34%	

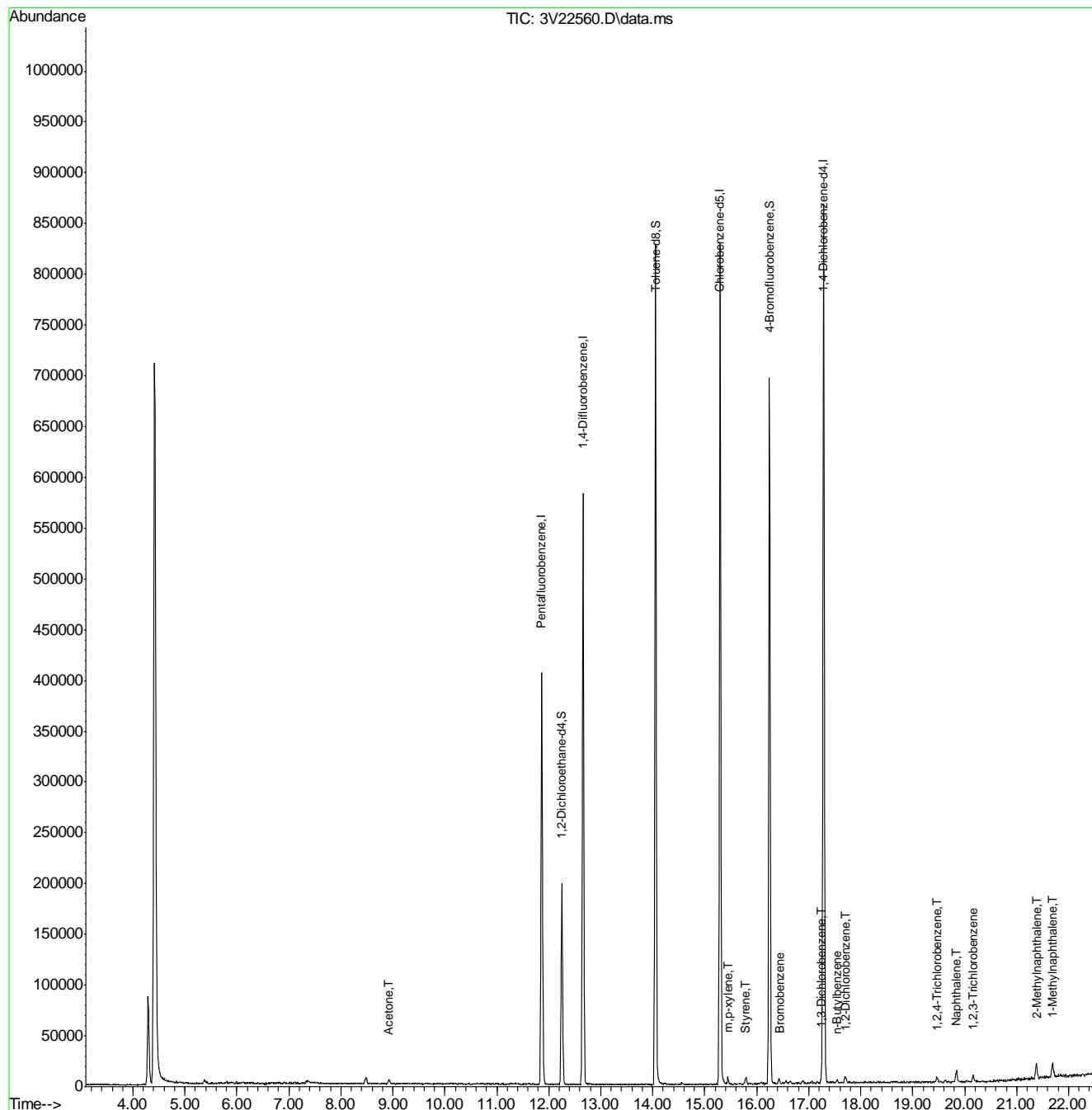
Target Compounds					Qvalue
15) Acetone	8.925	43	6973	0.73	ug/l 100
70) Bromobenzene	16.441	156	1174	0.21	ug/l 90
71) Styrene	15.796	104	1960	0.56	ug/l 86
72) m,p-xylene	15.453	106	2503	0.31	ug/l 90
84) 1,3-Dichlorobenzene	17.239	146	3019	0.28	ug/l 96
87) 1,2-Dichlorobenzene	17.704	146	3319	0.32	ug/l 96
88) n-Butylbenzene	17.541	91	3129	0.22	ug/l 88
90) 1,2,4-Trichlorobenzene	19.462	180	3118	0.48	ug/l 92
91) Naphthalene	19.838	128	16063	2.00	ug/l 100
93) 1,2,3-Trichlorobenzene	20.168	180	3642	0.57	ug/l 94
94) 2-Methylnaphthalene	21.384	142	13326	1.89	ug/l 97
95) 1-Methylnaphthalene	21.692	142	11465	1.71	ug/l 98

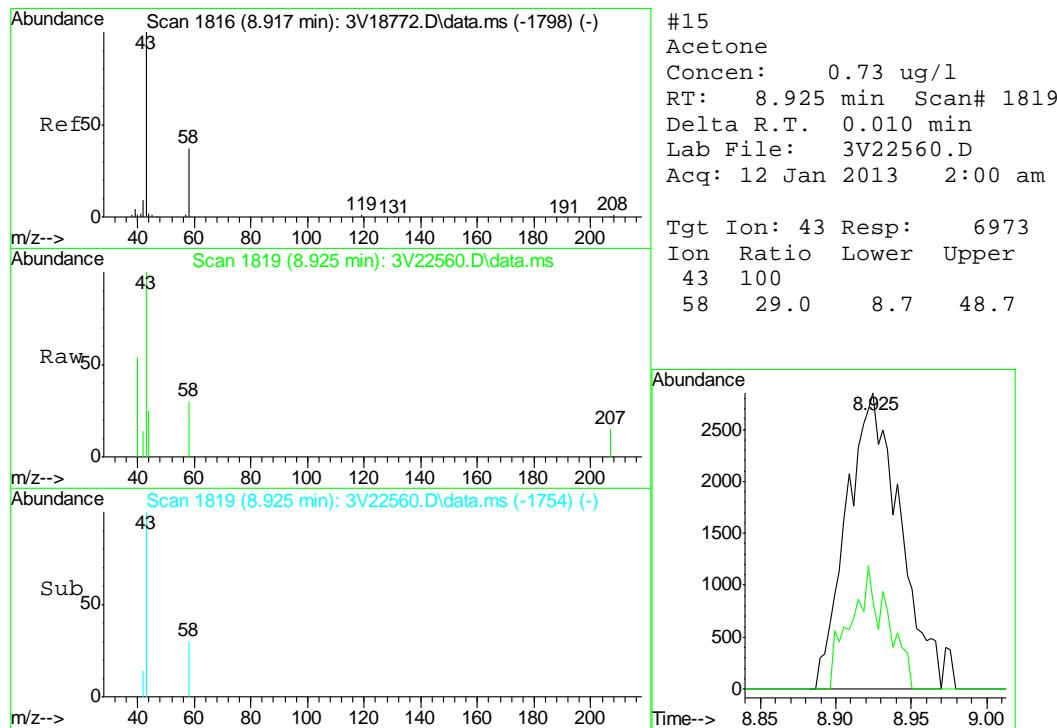
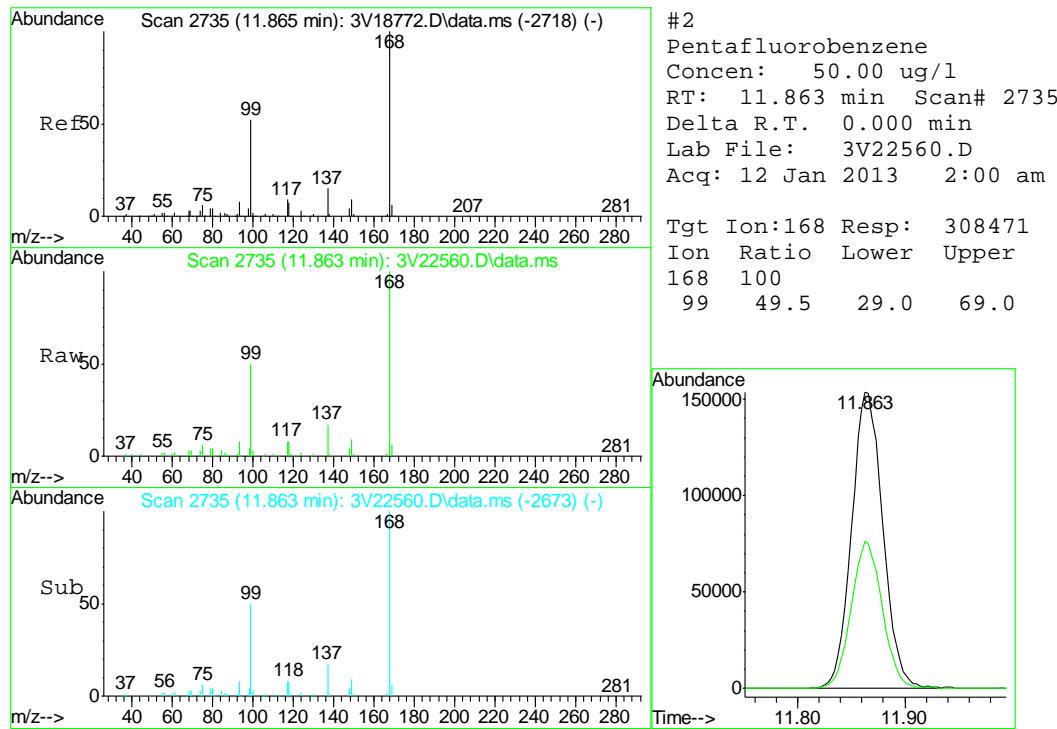
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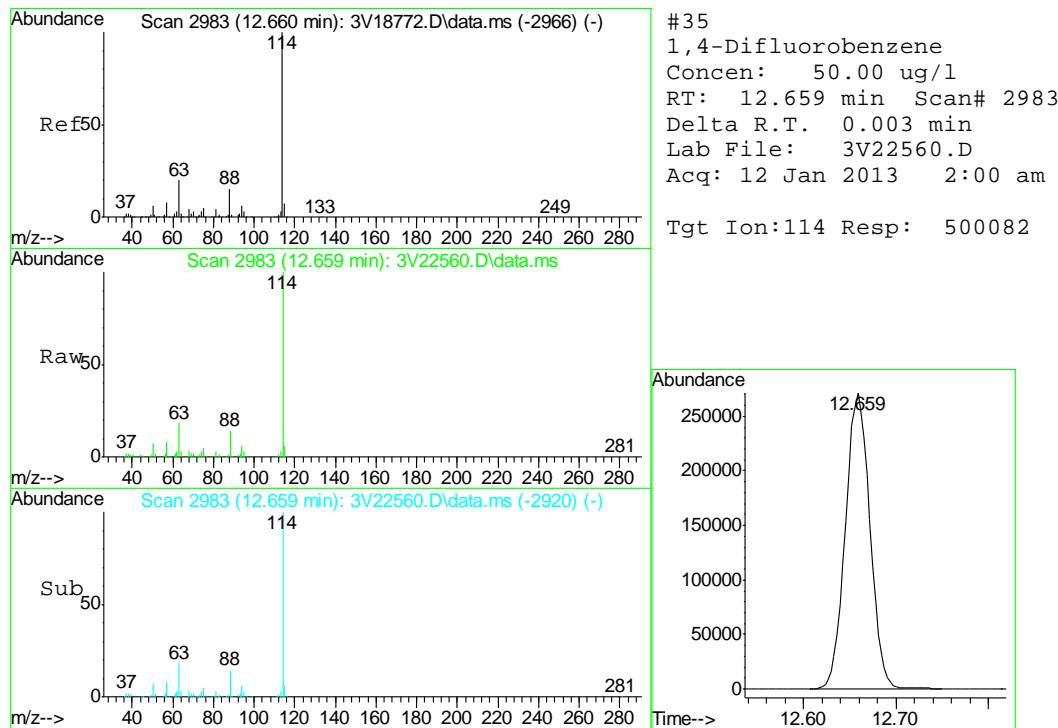
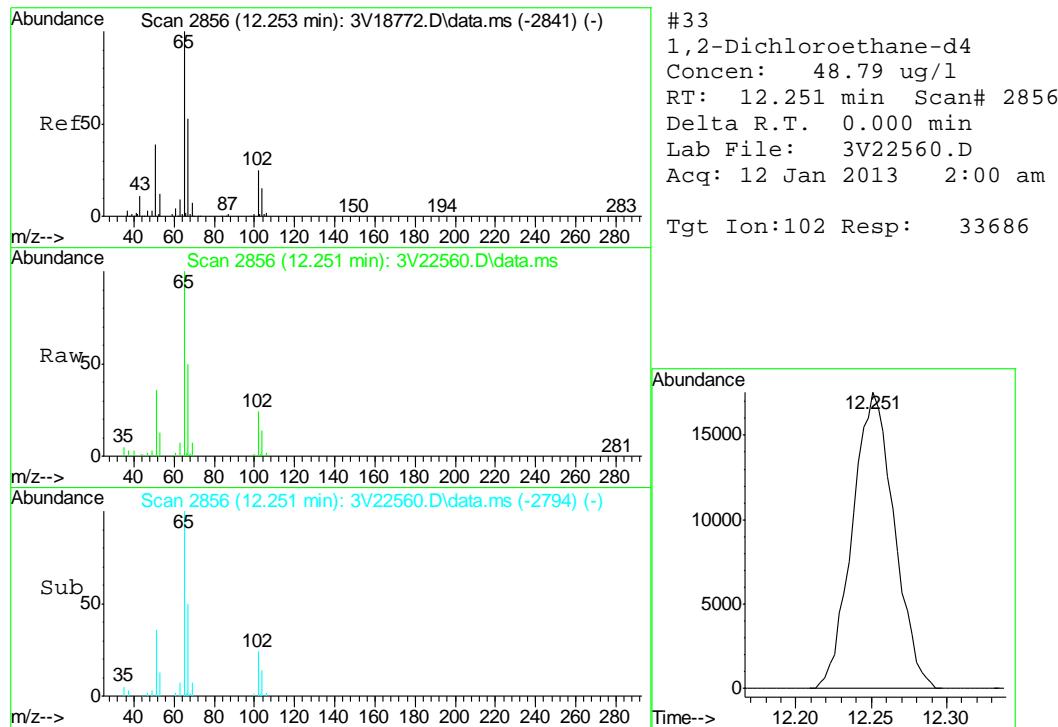
Quantitation Report (QT Reviewed)

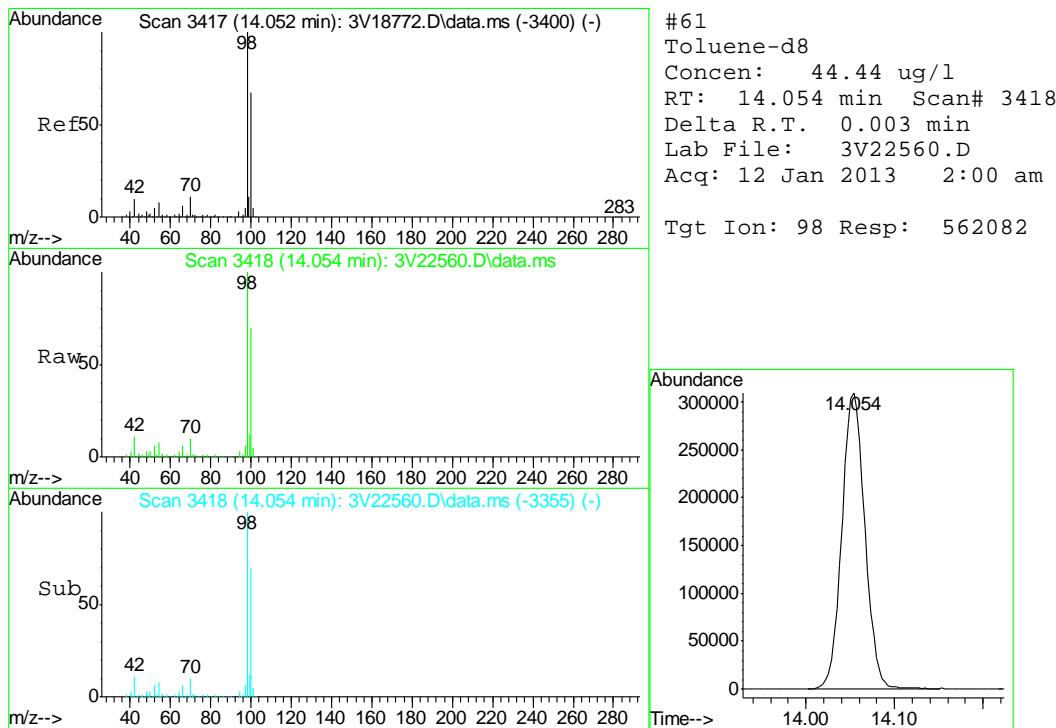
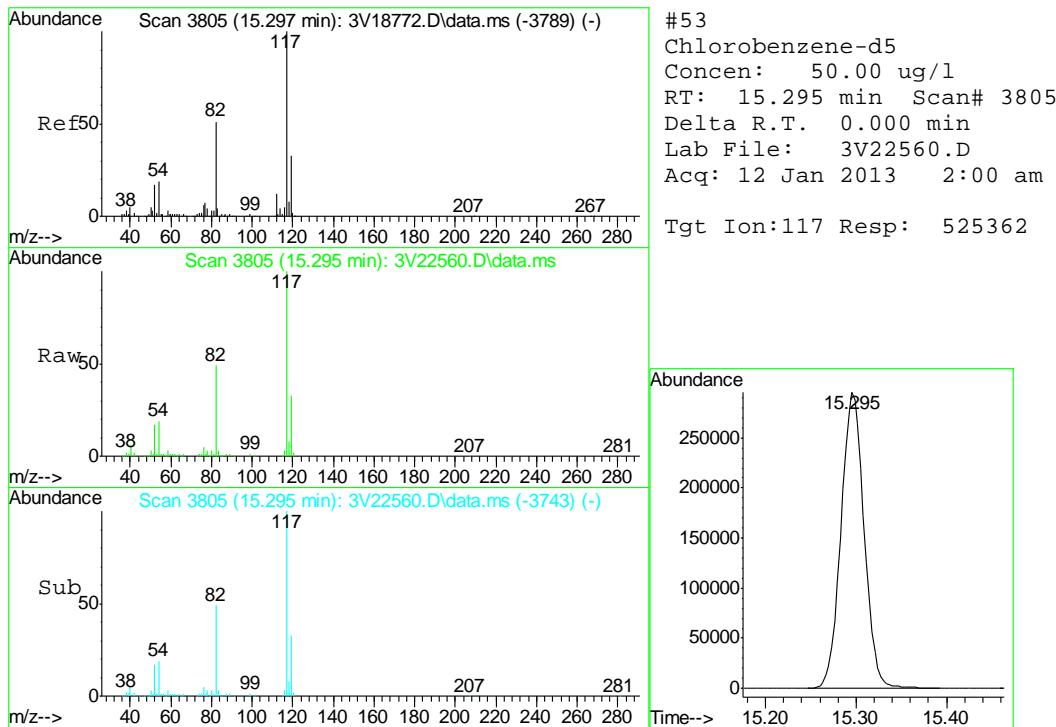
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 Acq On : 12 Jan 2013 2:00 am
 Operator : BRETD
 Sample : MB
 Misc : MS5218,V3V1327,5.00,,100,5,1
 ALS Vial : 27 Sample Multiplier: 1

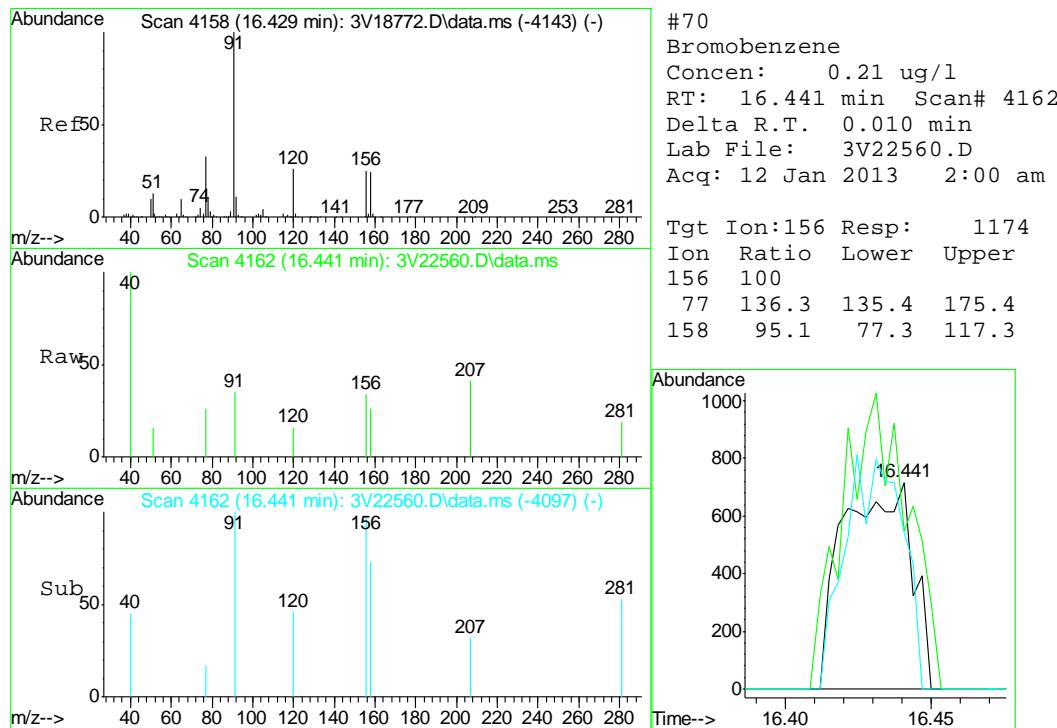
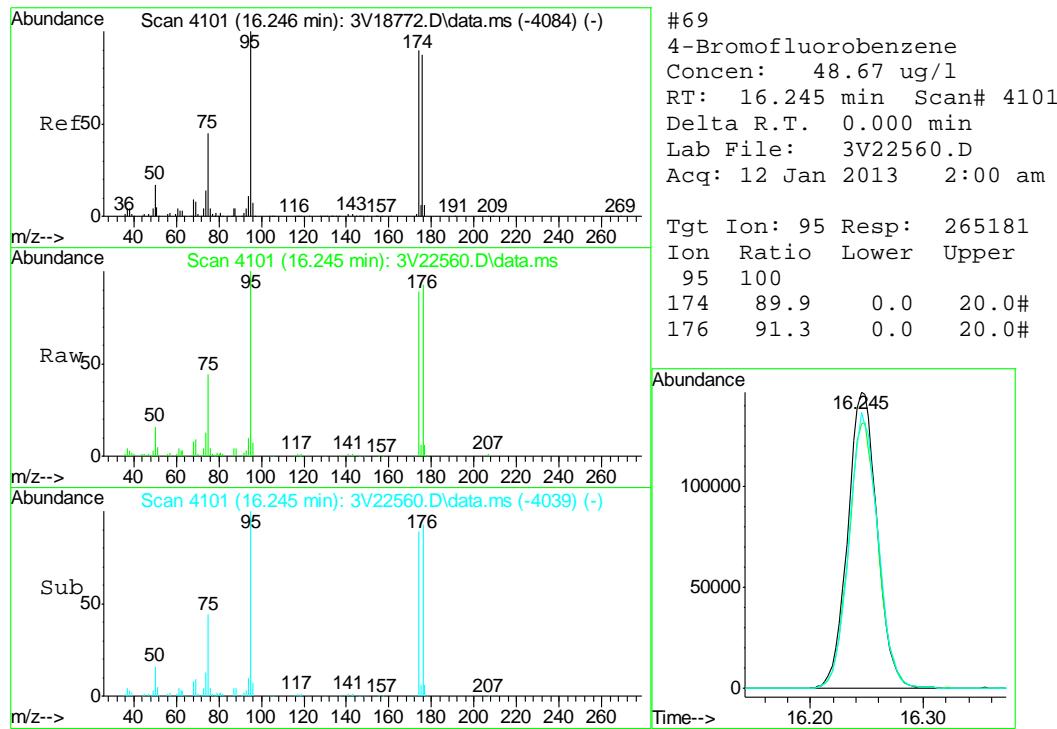
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 Response via : Initial Calibration

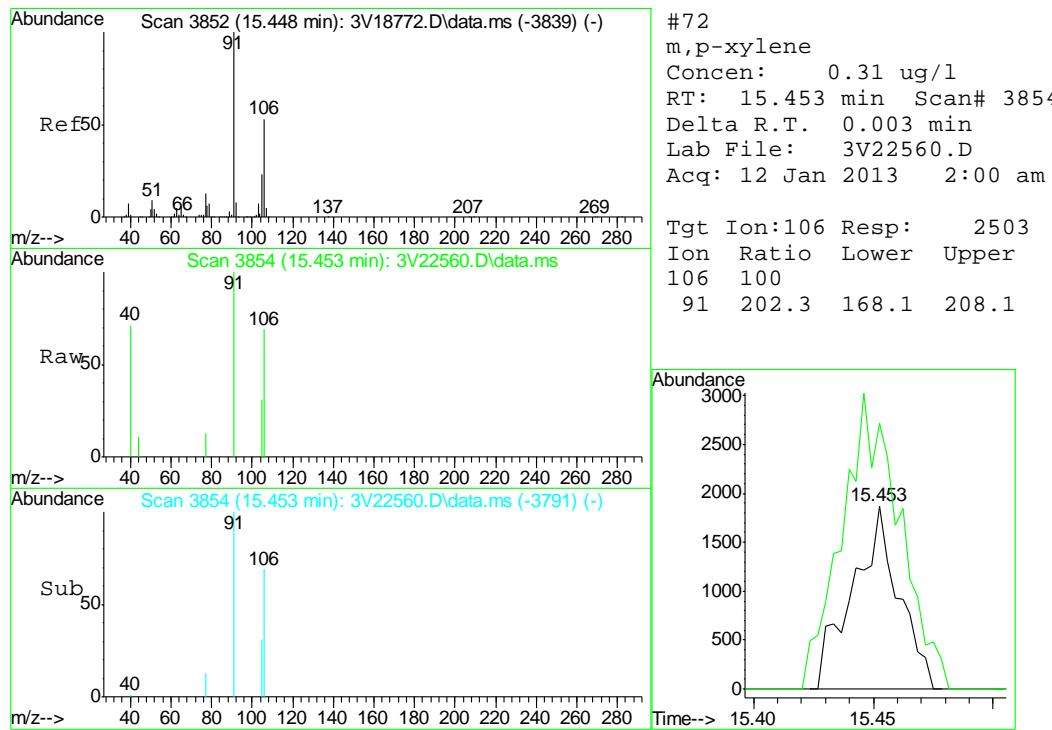
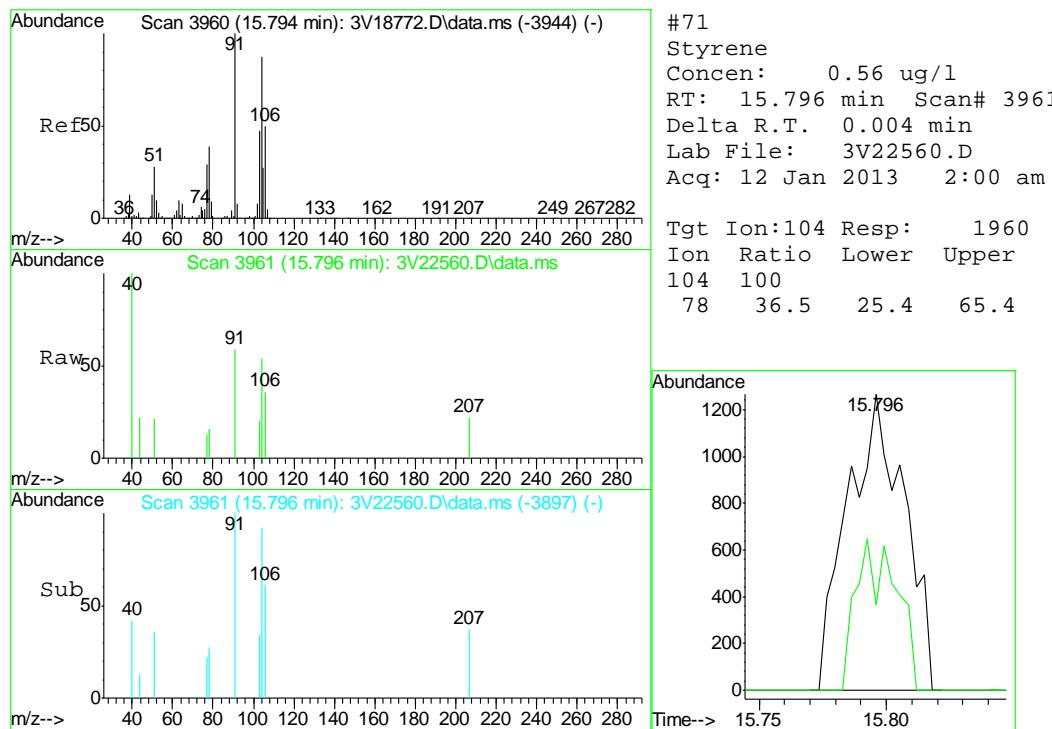


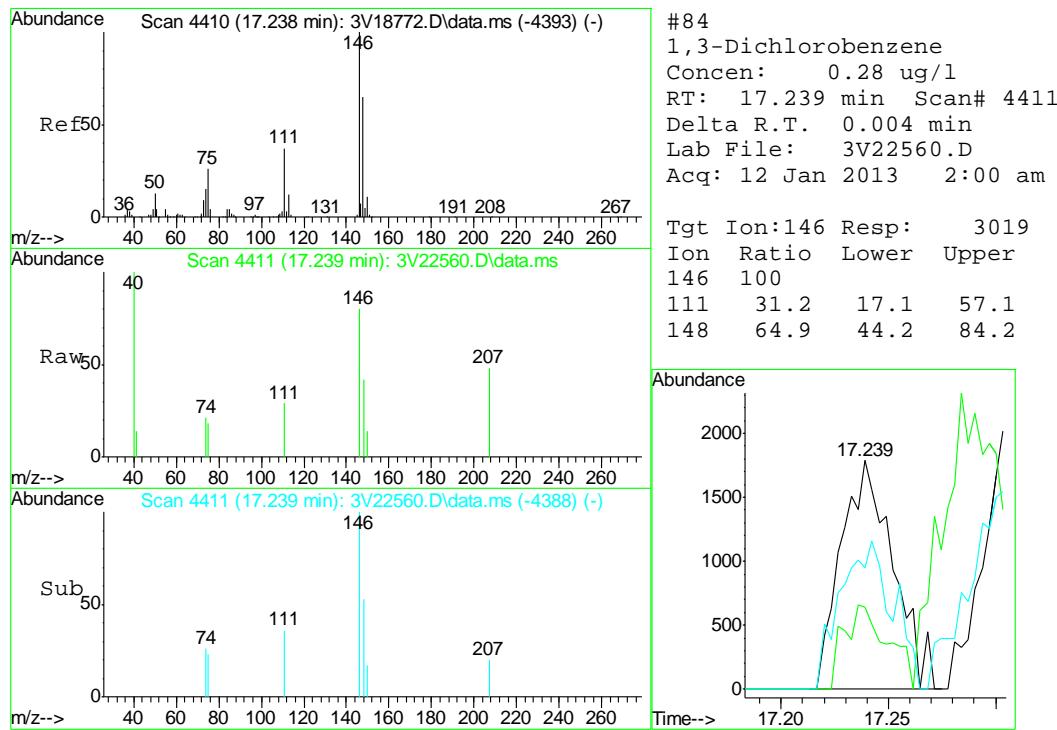
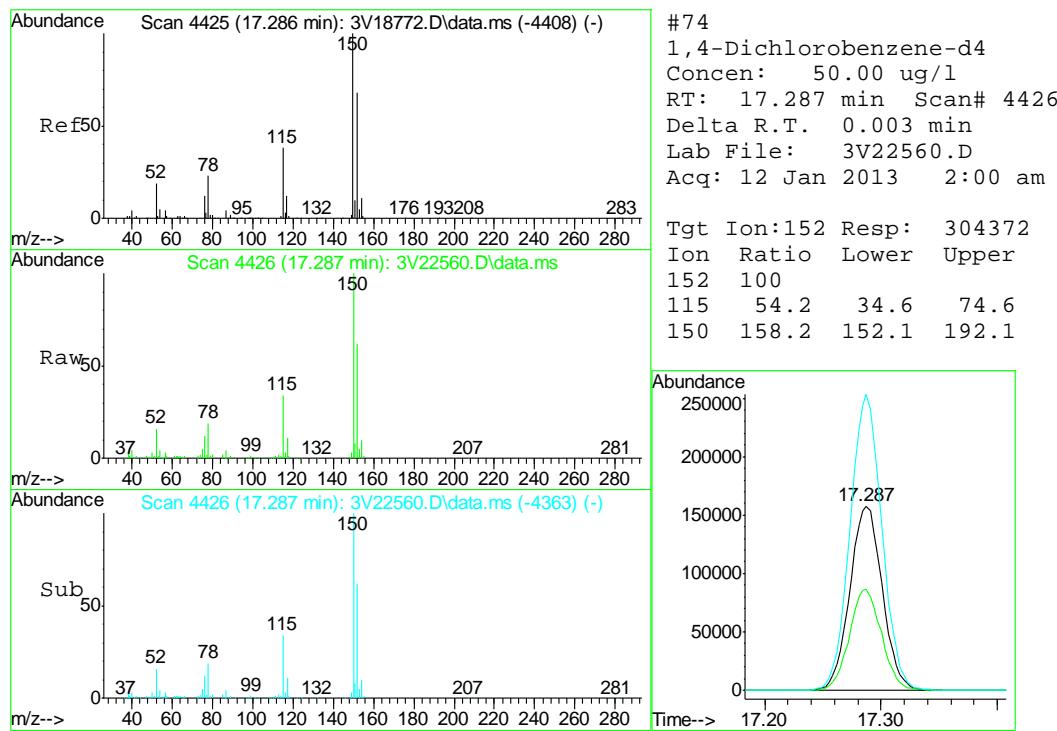


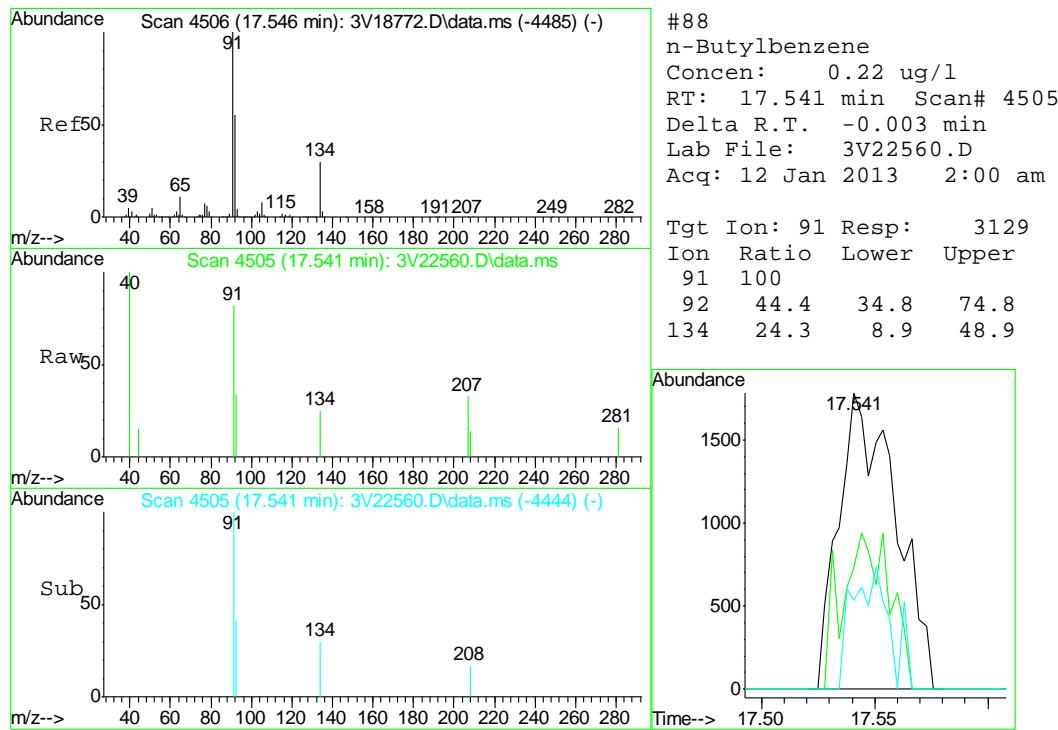
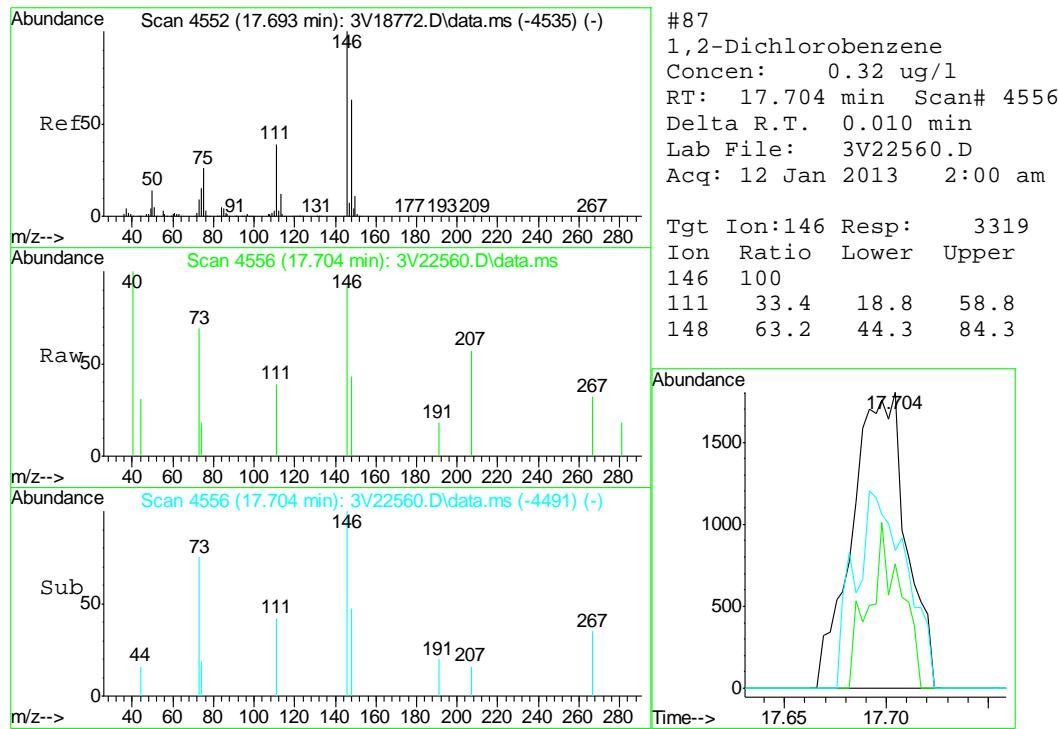


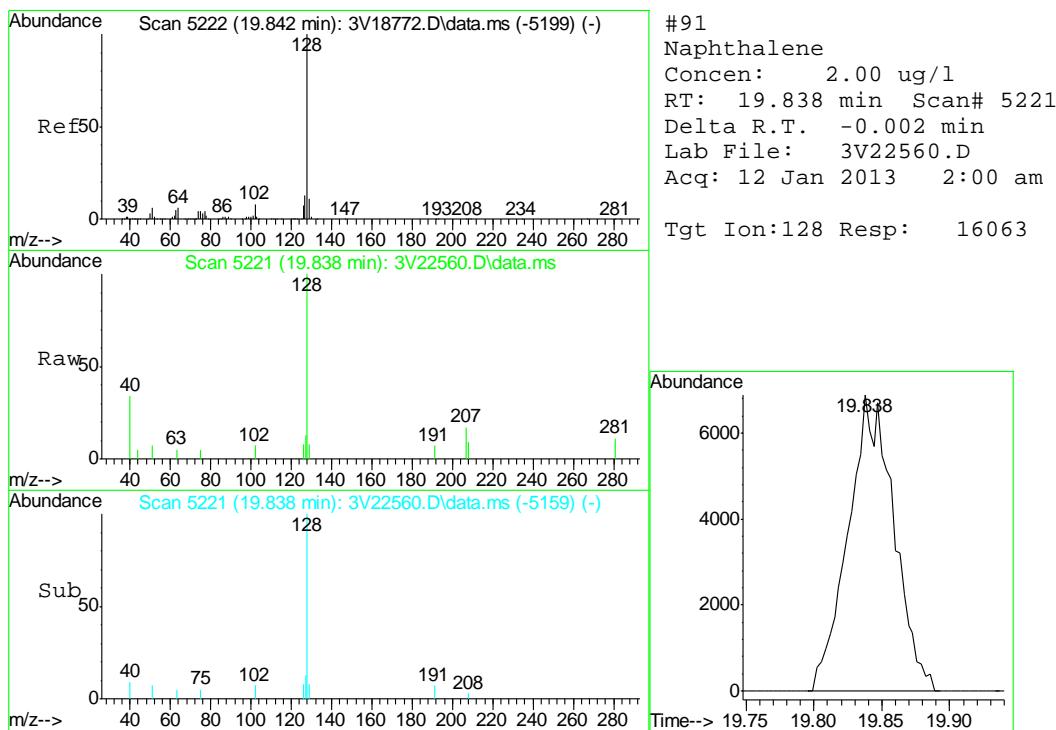
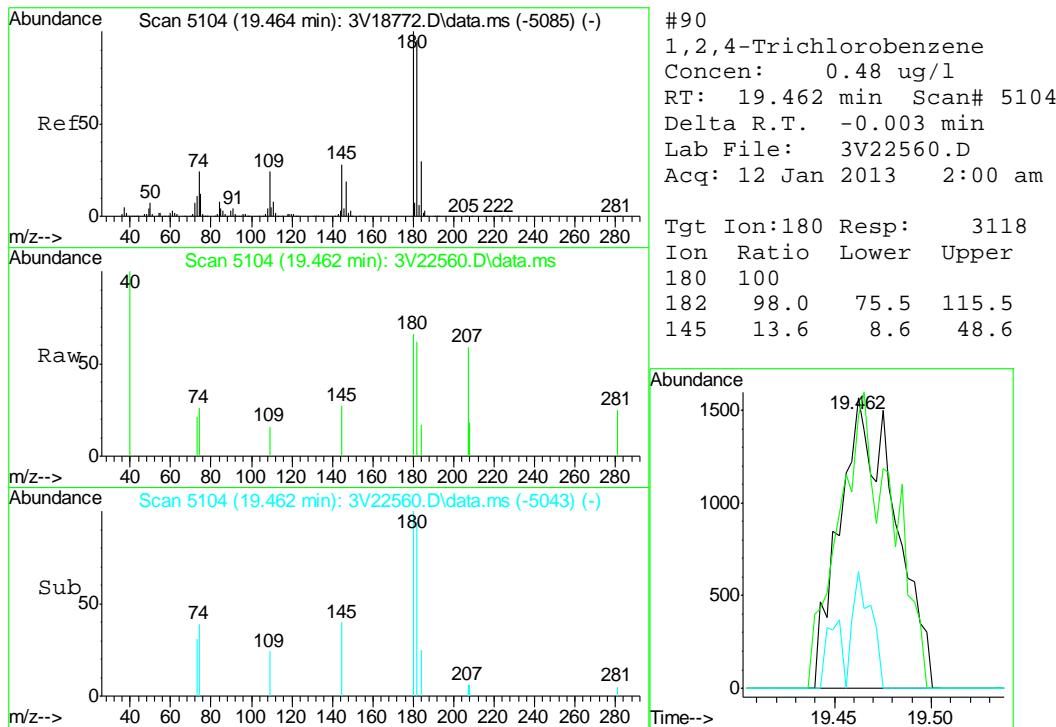


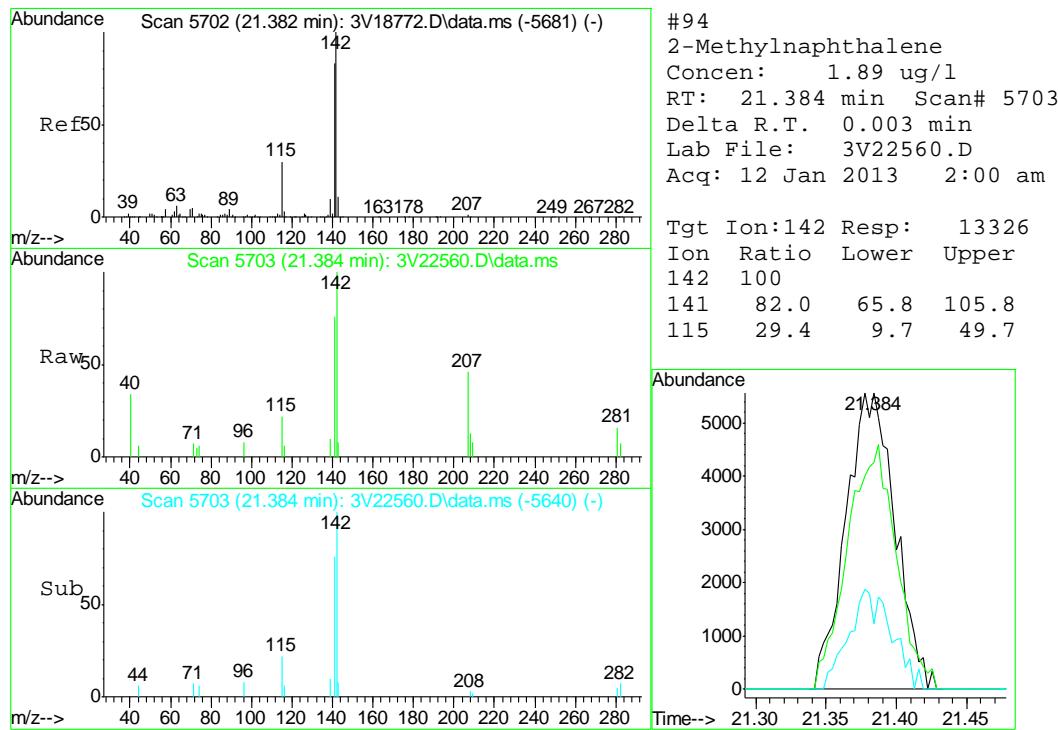
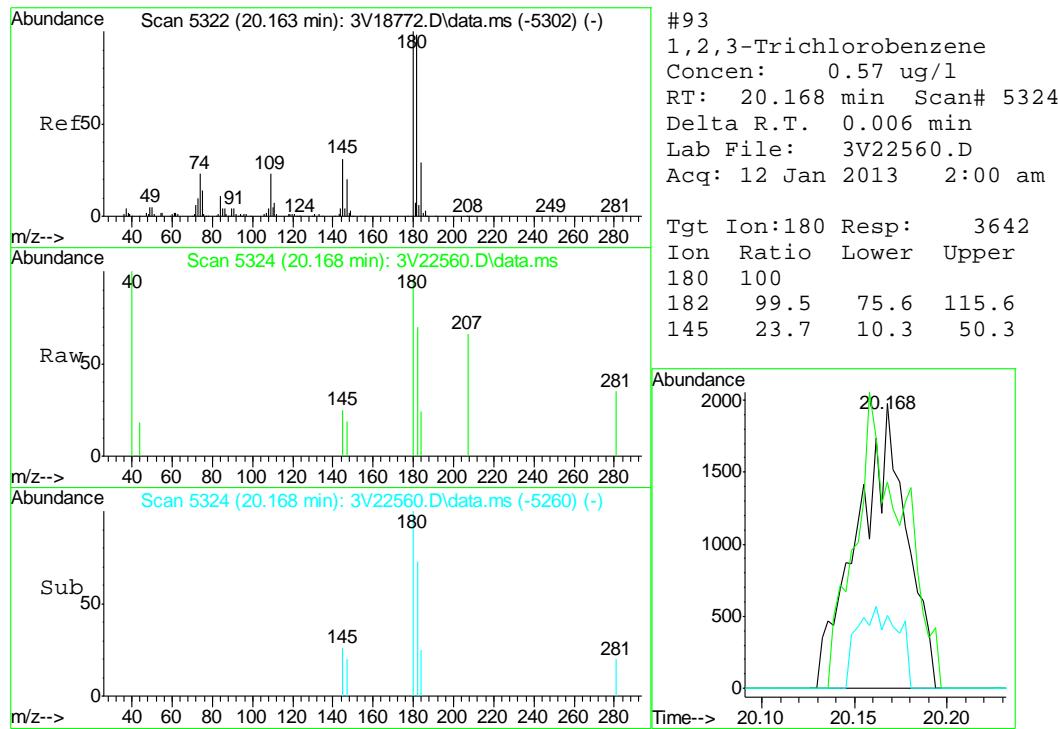


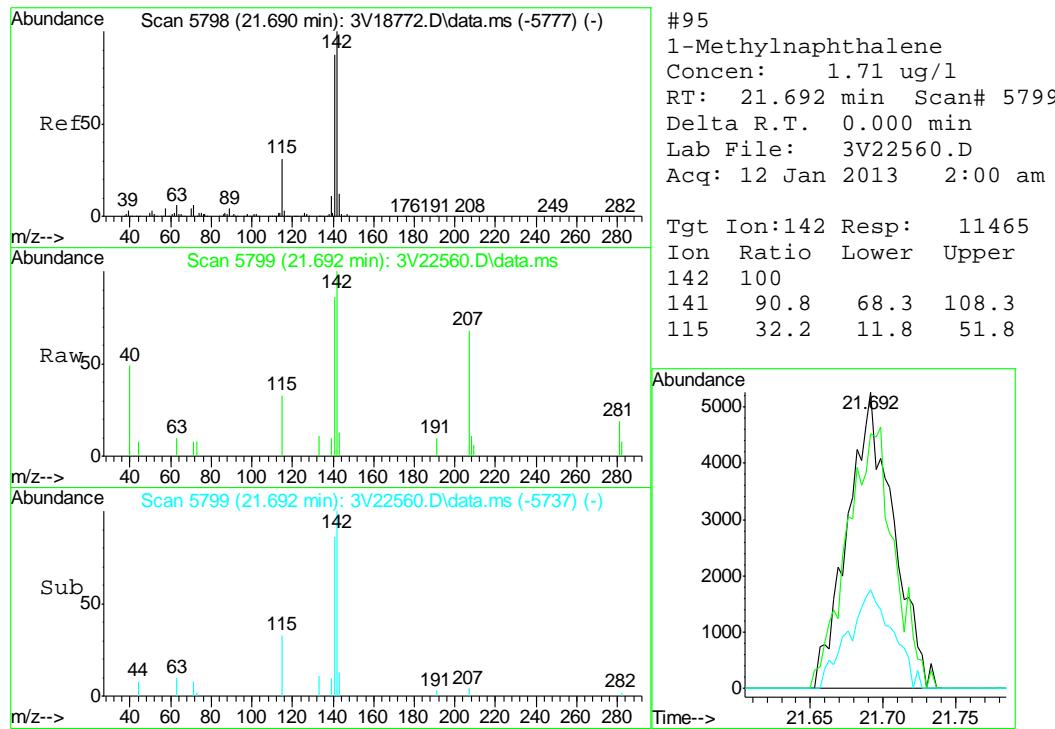














GC/MS Semi-volatiles

QC Data Summaries

∞

Includes the following where applicable:

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

Method Blank Summary

Job Number: D42511
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7223-MB	3G12973.D	1	01/15/13	DC	01/14/13	OP7223	E3G621

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42511-1

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	4.3	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	81% 10-159%
321-60-8	2-Fluorobiphenyl	85% 19-131%
1718-51-0	Terphenyl-d14	105% 18-150%

Blank Spike Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7223-BS	3G12974.D	1	01/15/13	DC	01/14/13	OP7223	E3G621

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42511-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	87.0	104	68-130
120-12-7	Anthracene	83.3	75.1	90	67-130
56-55-3	Benzo(a)anthracene	83.3	79.3	95	65-130
205-99-2	Benzo(b)fluoranthene	83.3	79.2	95	44-130
207-08-9	Benzo(k)fluoranthene	83.3	65.4	78	56-131
50-32-8	Benzo(a)pyrene	83.3	70.8	85	62-130
218-01-9	Chrysene	83.3	70.3	84	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	79.8	96	55-130
206-44-0	Fluoranthene	83.3	76.4	92	70-130
86-73-7	Fluorene	83.3	80.9	97	70-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	77.2	93	56-130
91-20-3	Naphthalene	83.3	71.1	85	70-130
129-00-0	Pyrene	83.3	74.5	89	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	103%	10-159%
321-60-8	2-Fluorobiphenyl	97%	19-131%
1718-51-0	Terphenyl-d14	107%	18-150%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7223-MS	3G12984.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
OP7223-MSD	3G12985.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
D42510-1	3G12983.D	1	01/15/13	DC	01/14/13	OP7223	E3G621

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42511-1

CAS No.	Compound	D42510-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND		98.3	76.7	78	76.8	78	0	25-151/30
120-12-7	Anthracene	ND		98.3	93.2	95	95.1	97	2	39-159/30
56-55-3	Benzo(a)anthracene	ND		98.3	107	109	109	111	2	39-168/30
205-99-2	Benzo(b)fluoranthene	ND		98.3	109	111	110	112	1	24-163/30
207-08-9	Benzo(k)fluoranthene	ND		98.3	70.7	72	74.9	76	6	10-188/30
50-32-8	Benzo(a)pyrene	ND		98.3	77.1	78	79.0	80	2	32-144/30
218-01-9	Chrysene	25.4		98.3	97.1	73	101	77	4	43-150/30
53-70-3	Dibenzo(a,h)anthracene	ND		98.3	76.5	78	79.8	81	4	21-152/30
206-44-0	Fluoranthene	ND		98.3	101	103	102	104	1	36-157/30
86-73-7	Fluorene	198		98.3	260	63	249	52	4	10-182/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND		98.3	73.2	74	76.2	77	4	20-154/30
91-20-3	Naphthalene	1070		98.3	1400	336* a	909	-164* a	43* b	10-163/30
129-00-0	Pyrene	31.3		98.3	126	96	127	97	1	25-180/30

CAS No.	Surrogate Recoveries	MS	MSD	D42510-1	Limits
4165-60-0	Nitrobenzene-d5	141%	110%	141%	10-159%
321-60-8	2-Fluorobiphenyl	63%	56%	61%	19-131%
1718-51-0	Terphenyl-d14	98%	96%	91%	18-150%

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

(b) Variability of recovery may be due to sample matrix/homogeneity.

* = Outside of Control Limits.

8.3.1
8



GC/MS Semi-volatiles

Raw Data

Quantitation Report (QT Reviewed)

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson
01/16/13 11:51

Data Path : C:\msdchem\1\DATA\011513\
Data File : 3g12978.D
Acq On : 15 Jan 2013 12:18 pm
Operator : DONC
Sample : D42511-1
Misc : OP7223,E3G621,30.01,,,1,1
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 15 13:56:31 2013
Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M
Quant Title : PAHSIM BASE
QLast Update : Thu Jan 10 14:18:35 2013
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.621	136	148616	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.326	164	84132	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.812	188	143655	4.0000	ug/mL	0.00
19) Chrysene-d12	11.443	240	104141	4.0000	ug/mL	0.00
24) Perylene-d12	12.810	264	86662	4.0000	ug/mL	0.00

System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	523310	39.1476	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	78.30%
7) 2-Fluorobiphenyl	6.664	172	1261991	38.8263	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	77.66%
21) Terphenyl-d14	10.402	244	682147	48.1388	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	96.28%

Target Compounds

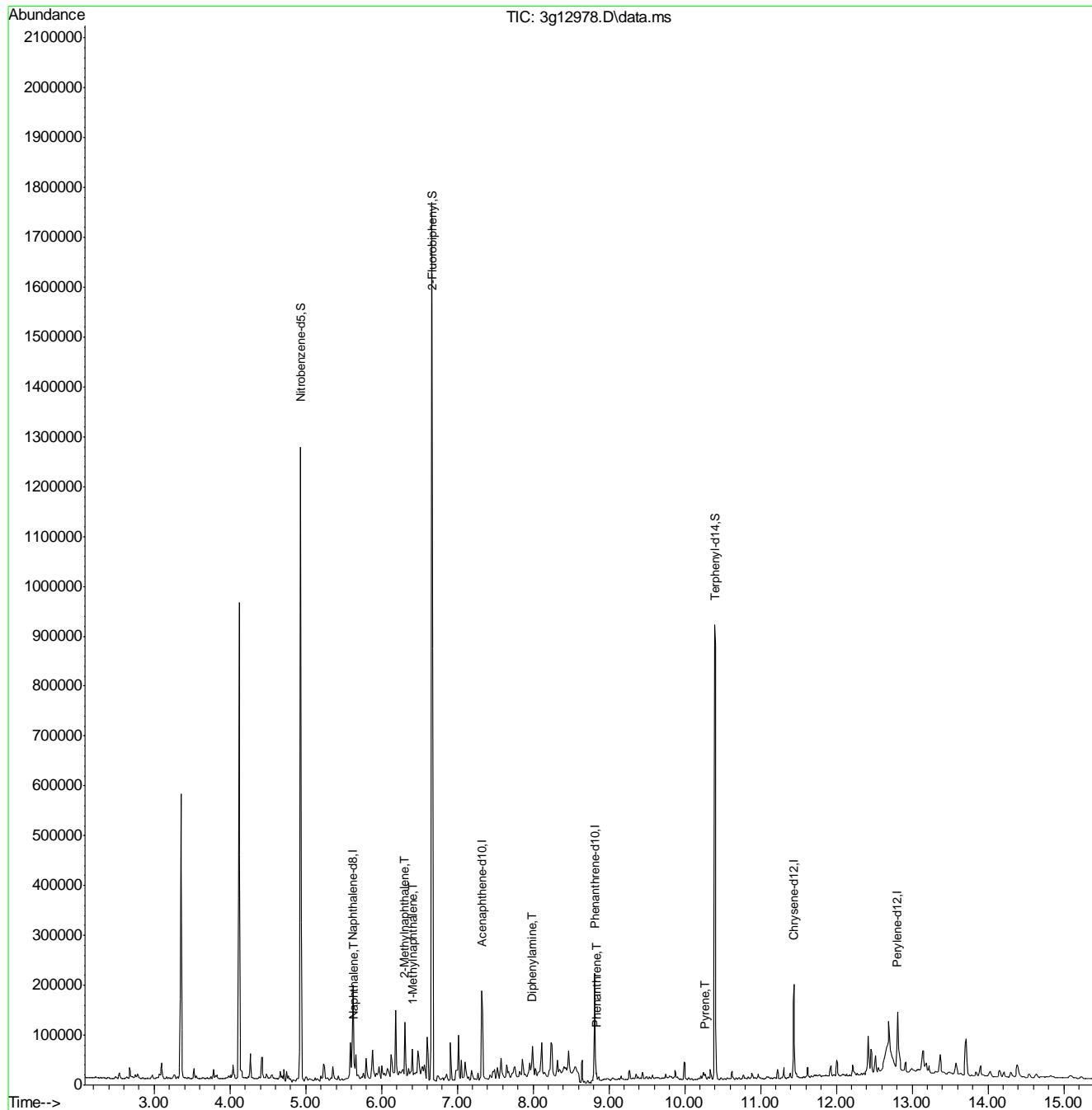
				Qvalue
3) N-Nitrosodimethylamine	2.342	74	74	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.633	128	23611	0.5400 ug/mL 76
8) 2-Methylnaphthalene	6.306	142	43243	1.5995 ug/mL 92
9) 1-Methylnaphthalene	6.406	142	19288	0.8157 ug/mL 88
10) Acenaphthylene	7.184	152	1403	N.D.
11) Acenaphthene	0.000	154	0	N.D. d
12) Dibenzofuran	0.000	168	0	N.D. d
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	7.987	169	31626m	1.1980 ug/mL
16) Phenanthrene	8.835	178	14221	0.2562 ug/mL# 61
17) Anthracene	8.915	178	1398	N.D.
18) Fluoranthene	10.015	202	1072	N.D.
20) Pyrene	10.268	202	6185	0.1112 ug/mL# 59
22) Benzo(a)anthracene	11.463	228	2101	N.D.
23) Chrysene	11.463	228	2129	N.D.
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d
26) Benzo(k)fluoranthene	12.452	252	2364	N.D.
27) Benzo(a)pyrene	0.000	252	0	N.D. d
28) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D. d
29) Dibenz(a,h)anthracene	0.000	278	0	N.D. d
30) Benzo(g,h,i)perylene	14.408	276	378	N.D.

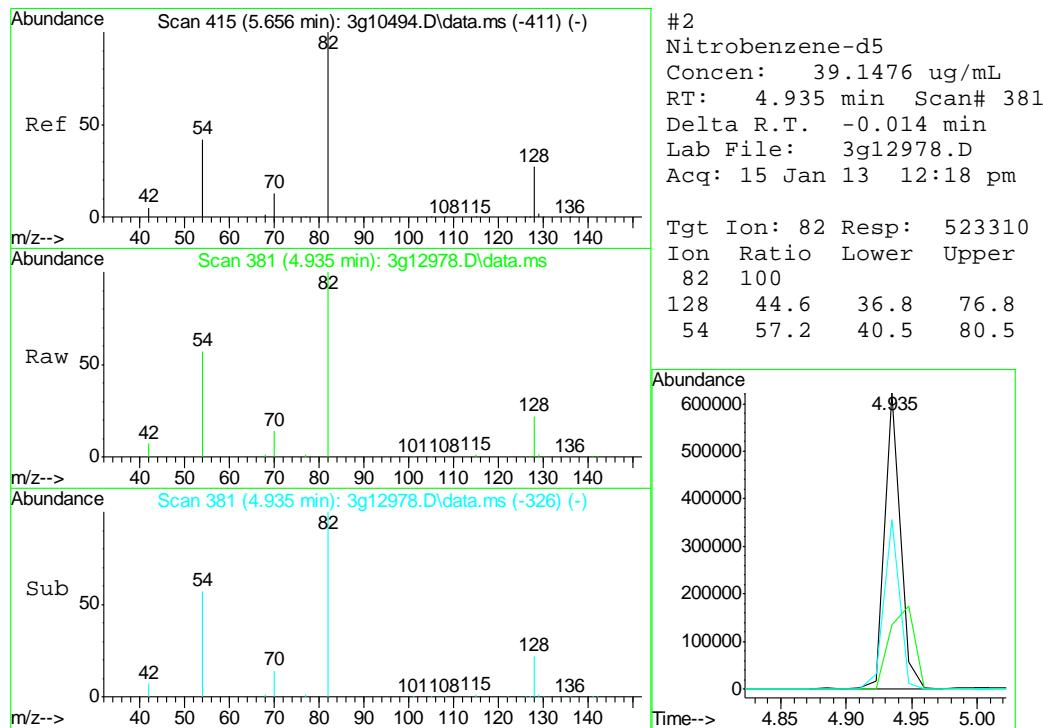
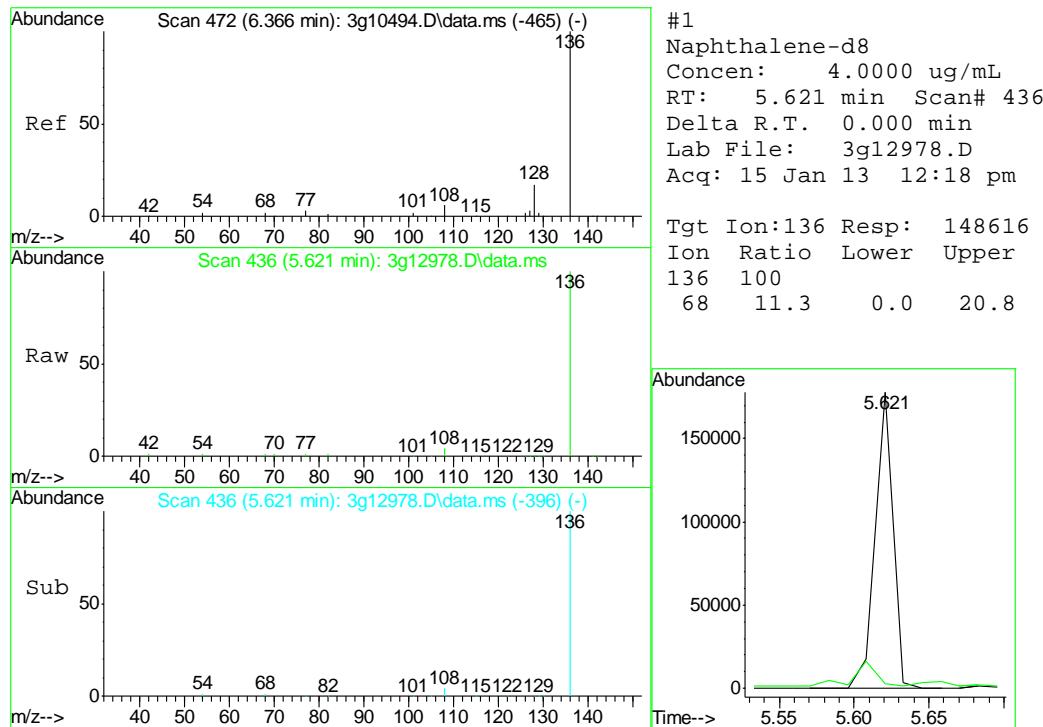
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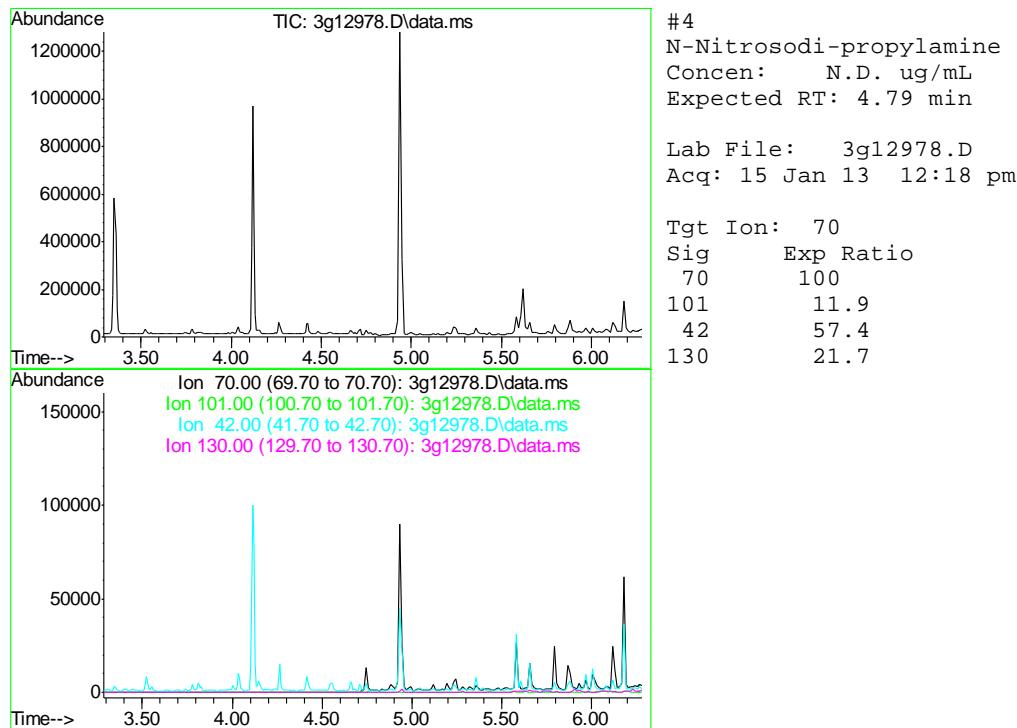
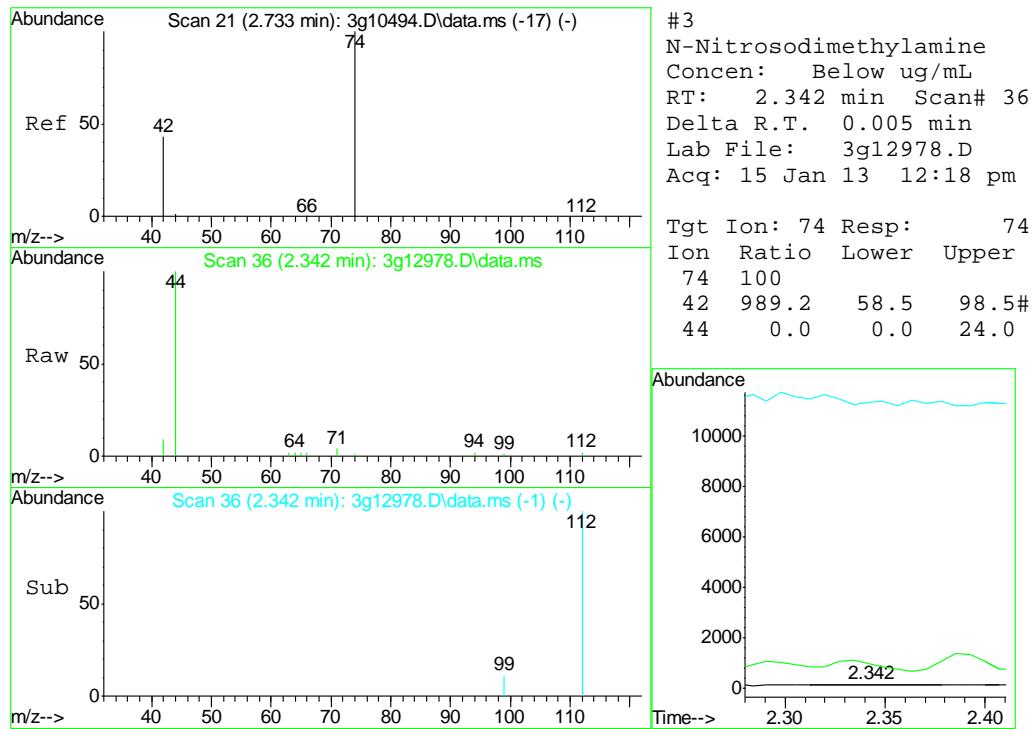
Quantitation Report (QT Reviewed)

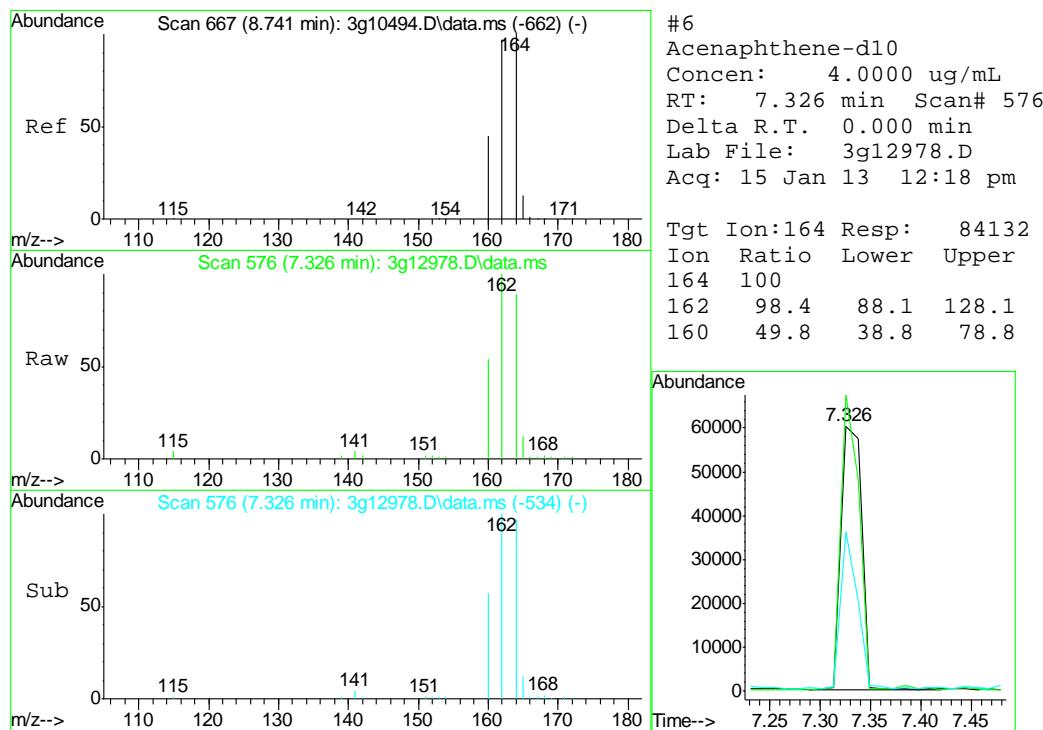
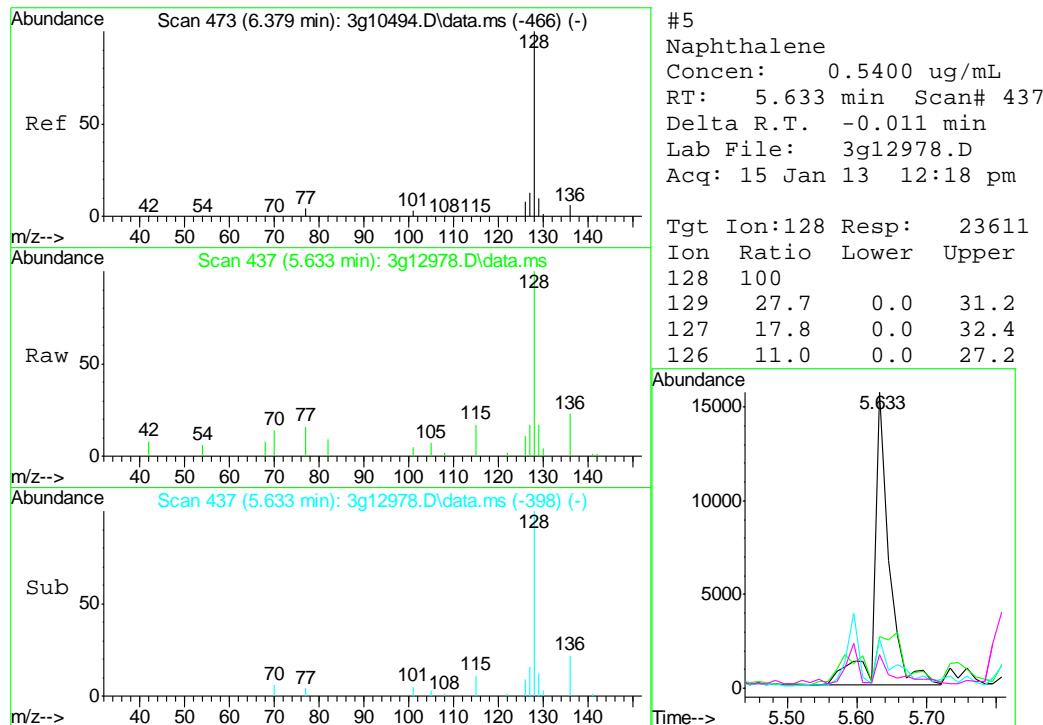
Data Path : C:\msdchem\1\DATA\011513\
 Data File : 3g12978.D
 Acq On : 15 Jan 2013 12:18 pm
 Operator : DONC
 Sample : D42511-1
 Misc : OP7223,E3G621,30.01,,,1,1
 ALS Vial : 9 Sample Multiplier: 1

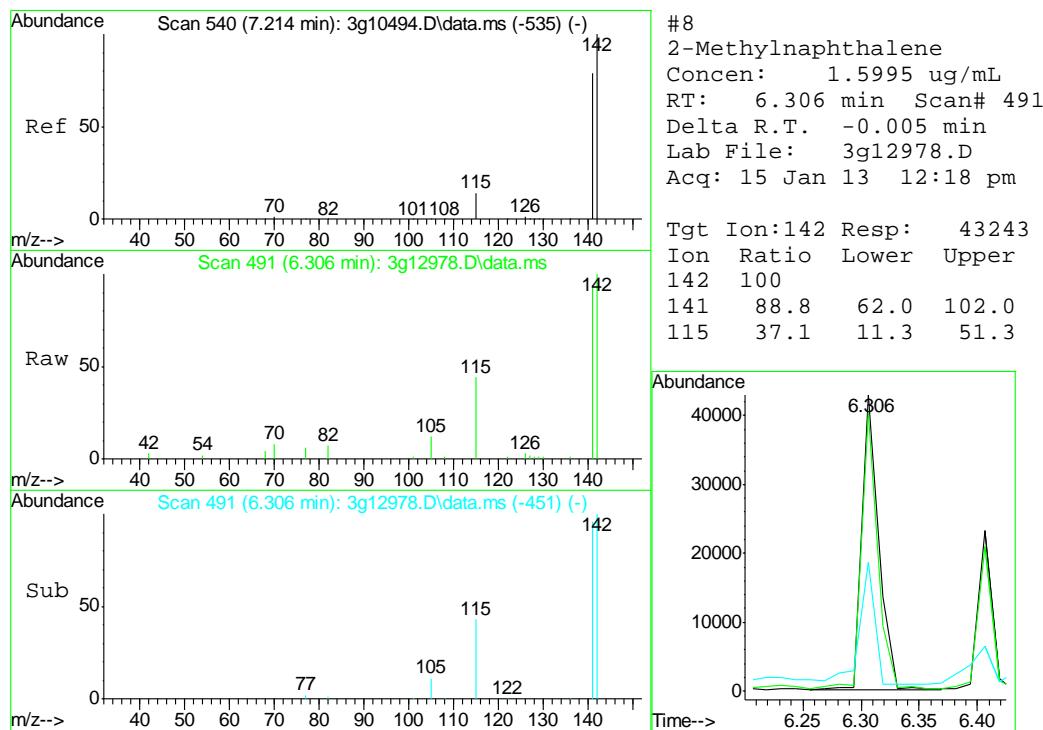
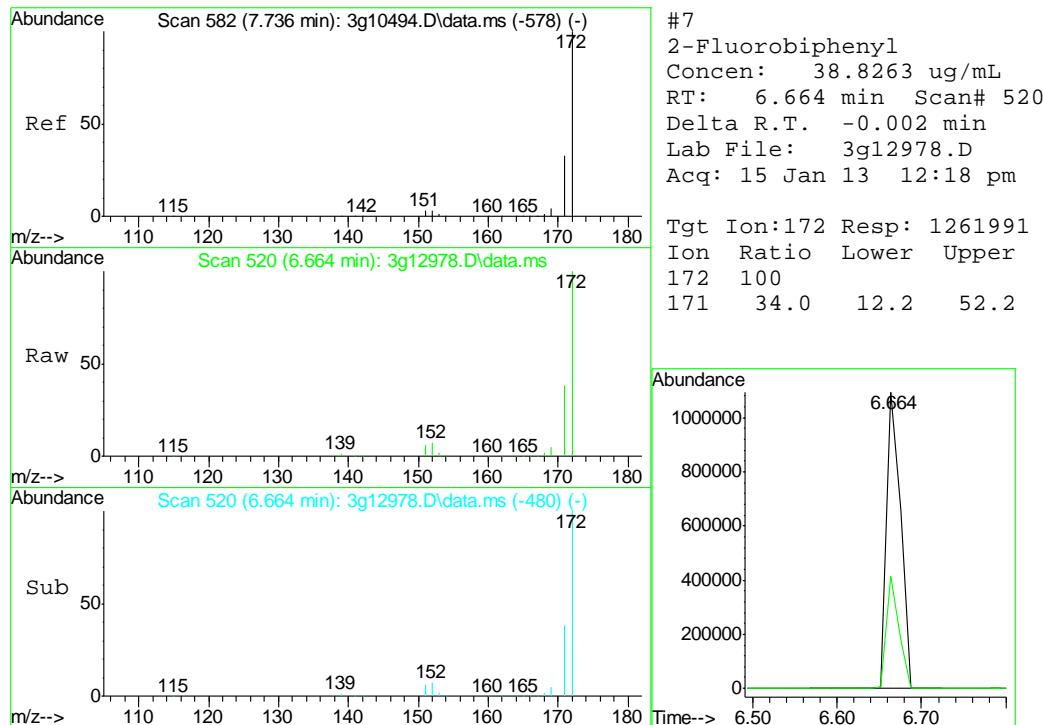
Quant Time: Jan 15 13:56:31 2013
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Jan 10 14:18:35 2013
 Response via : Initial Calibration

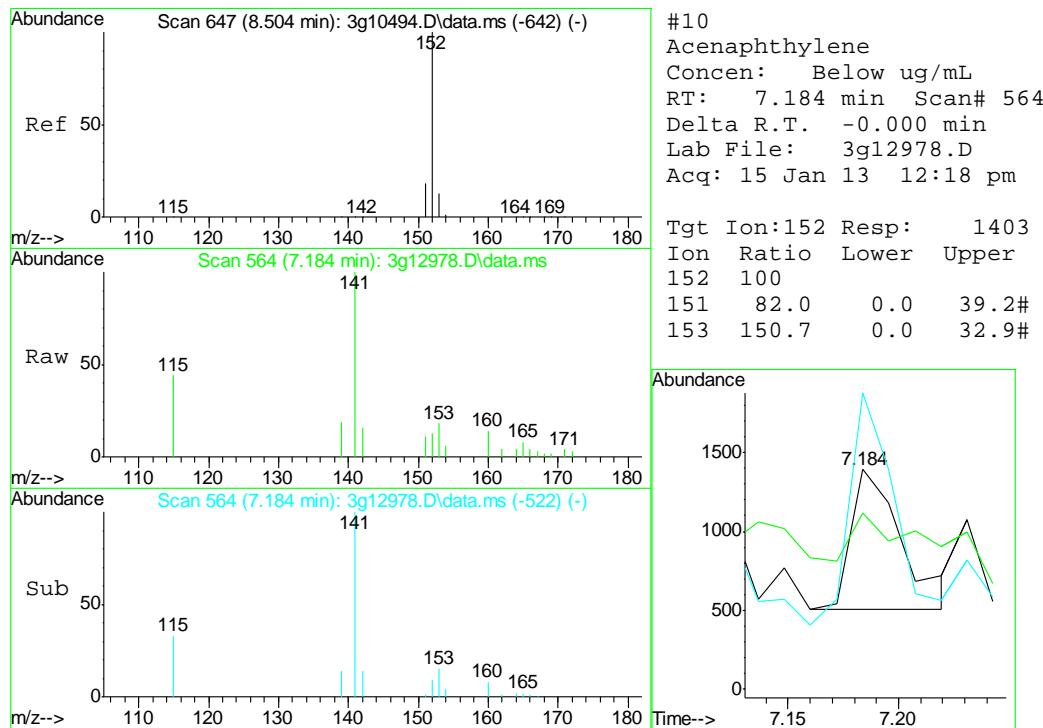
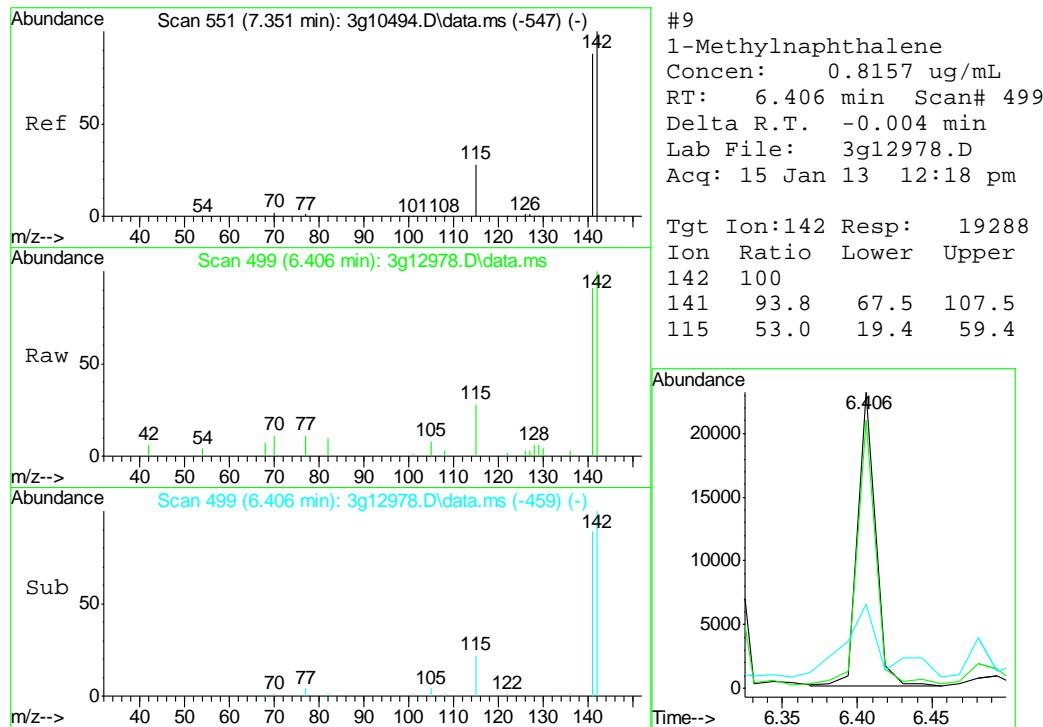


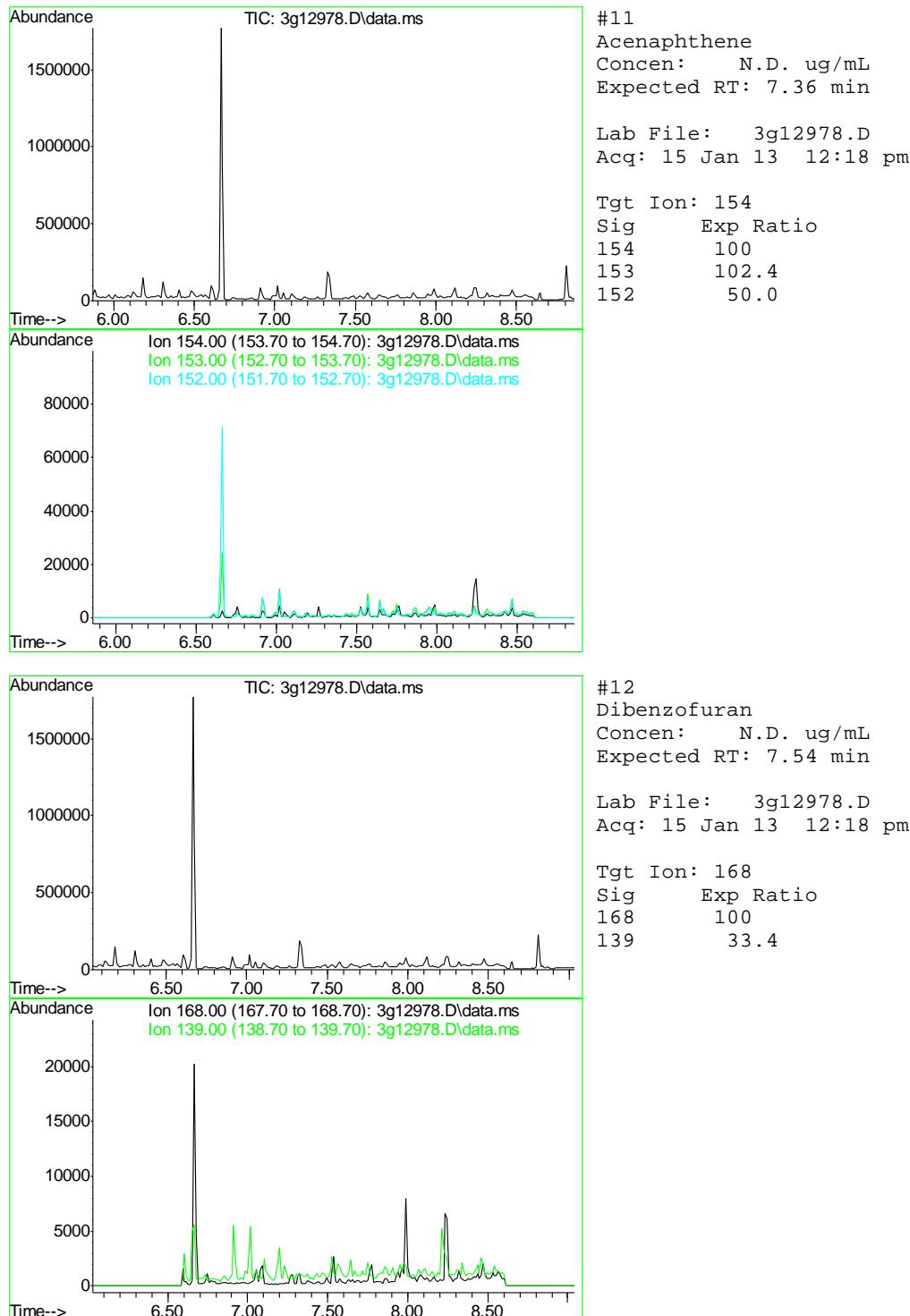


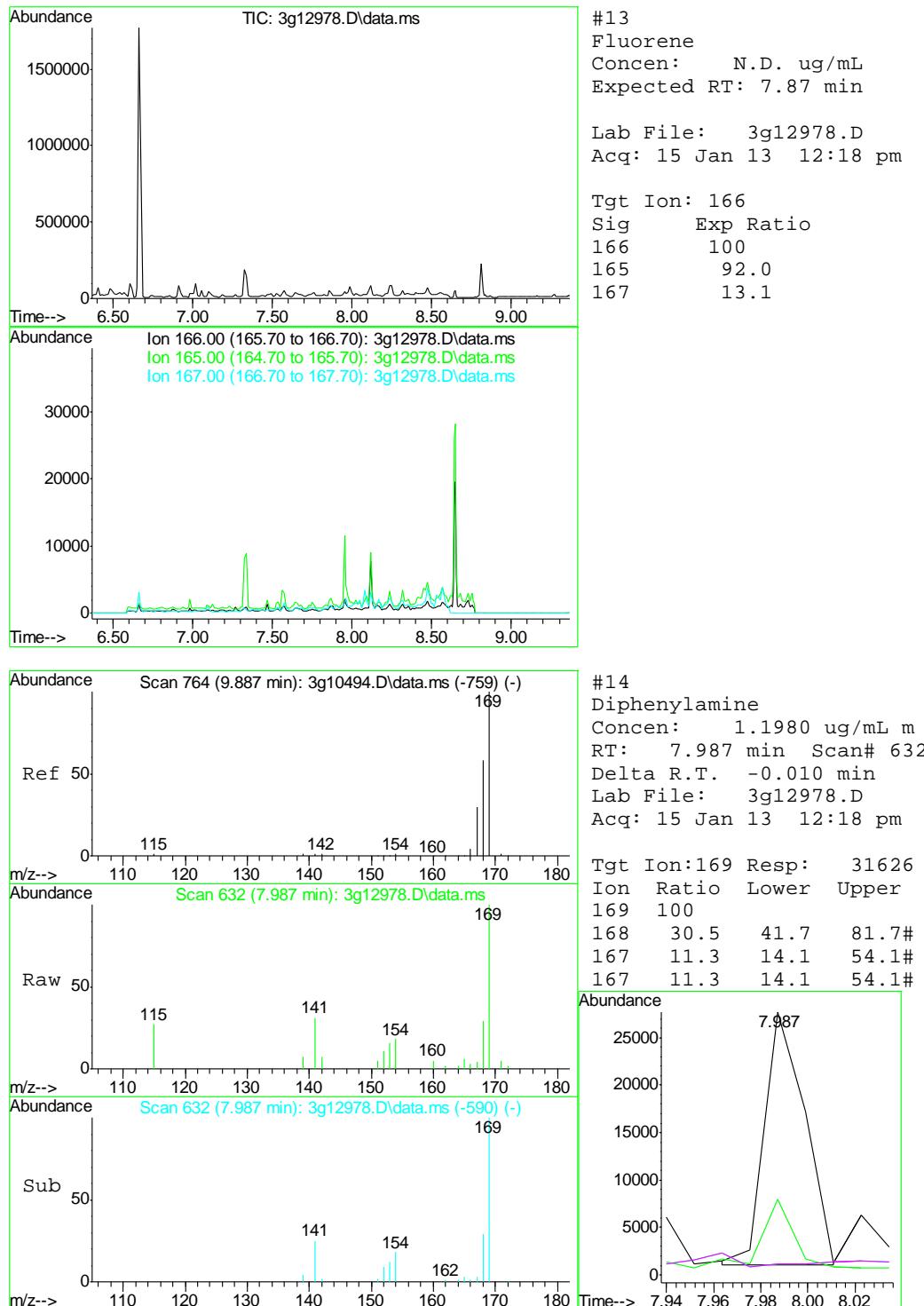


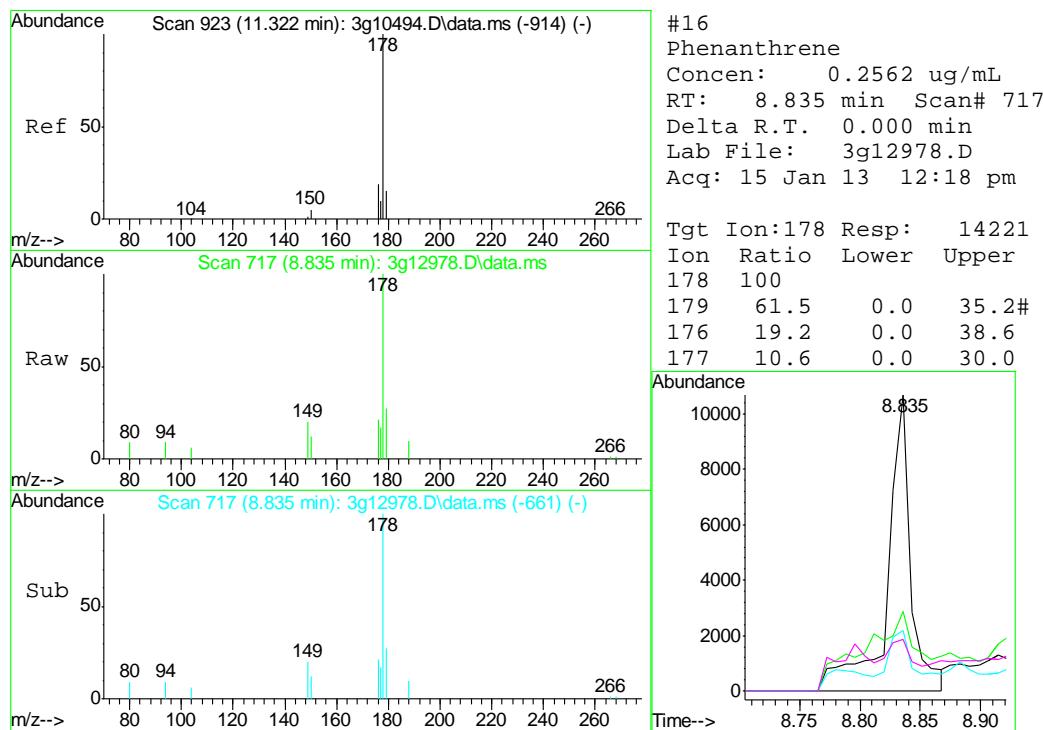
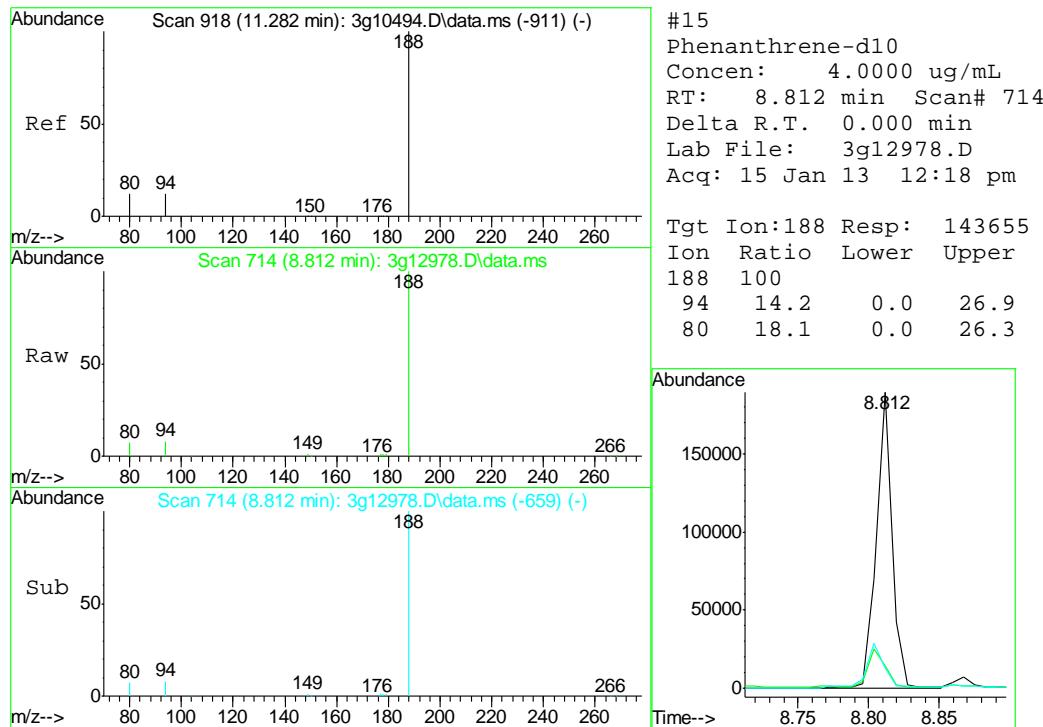


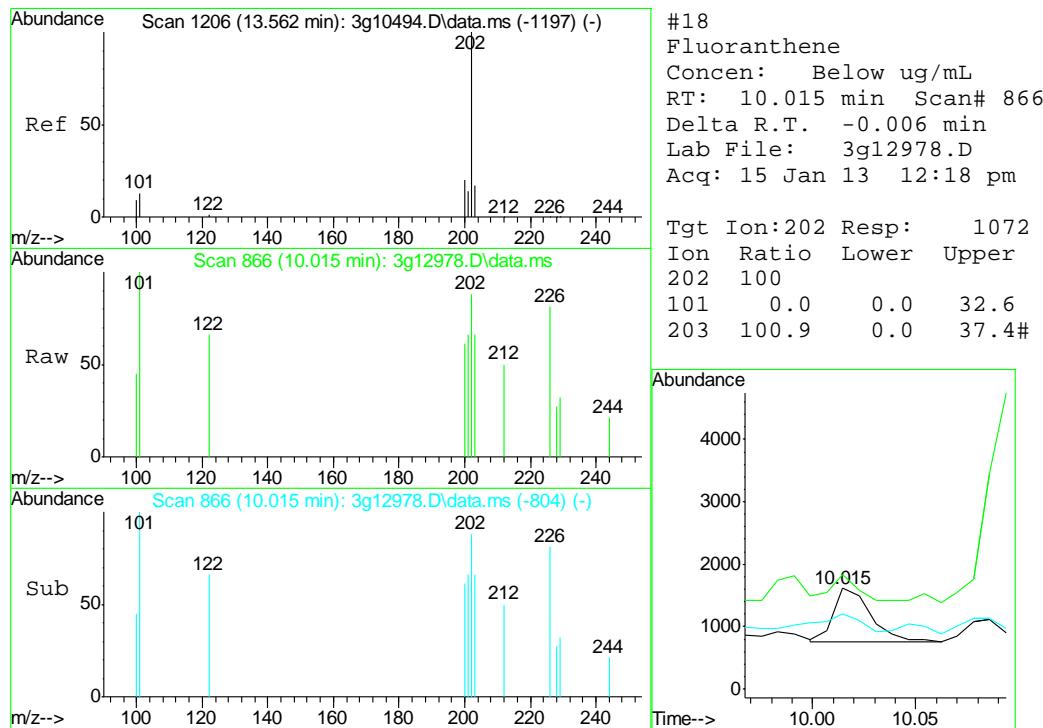
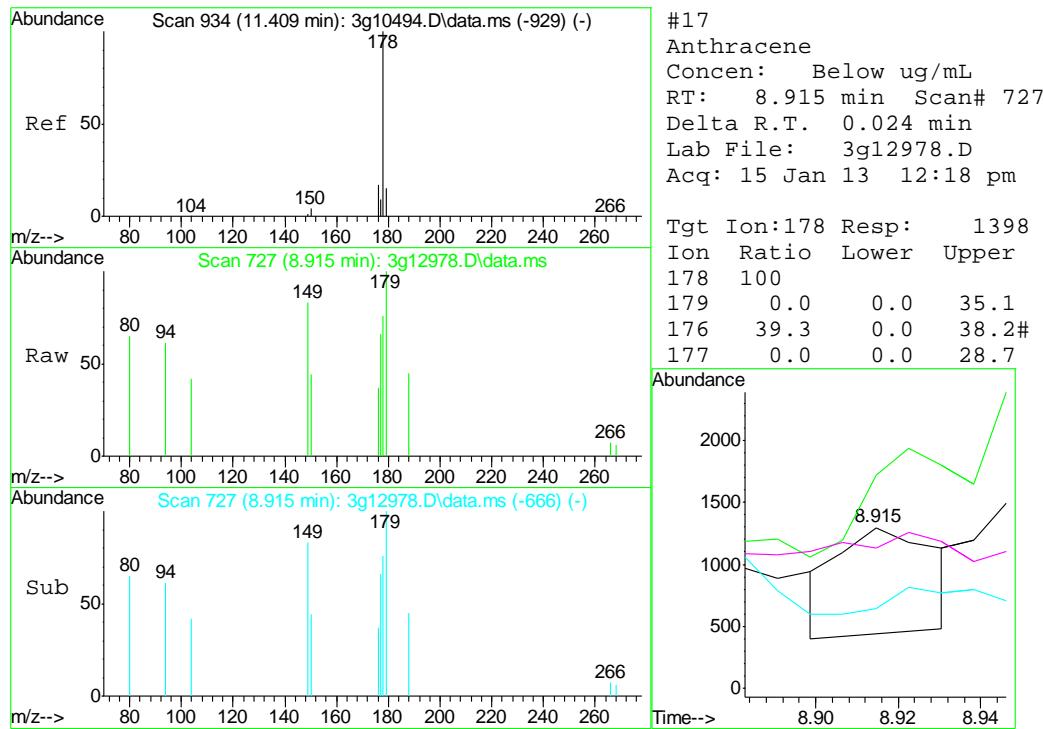


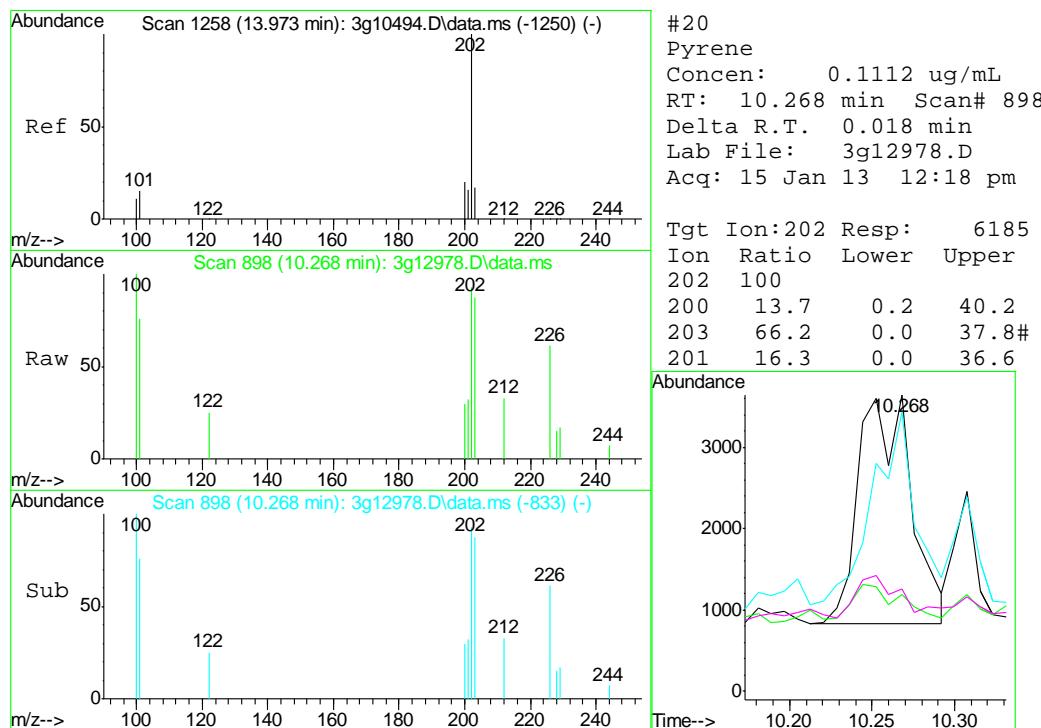
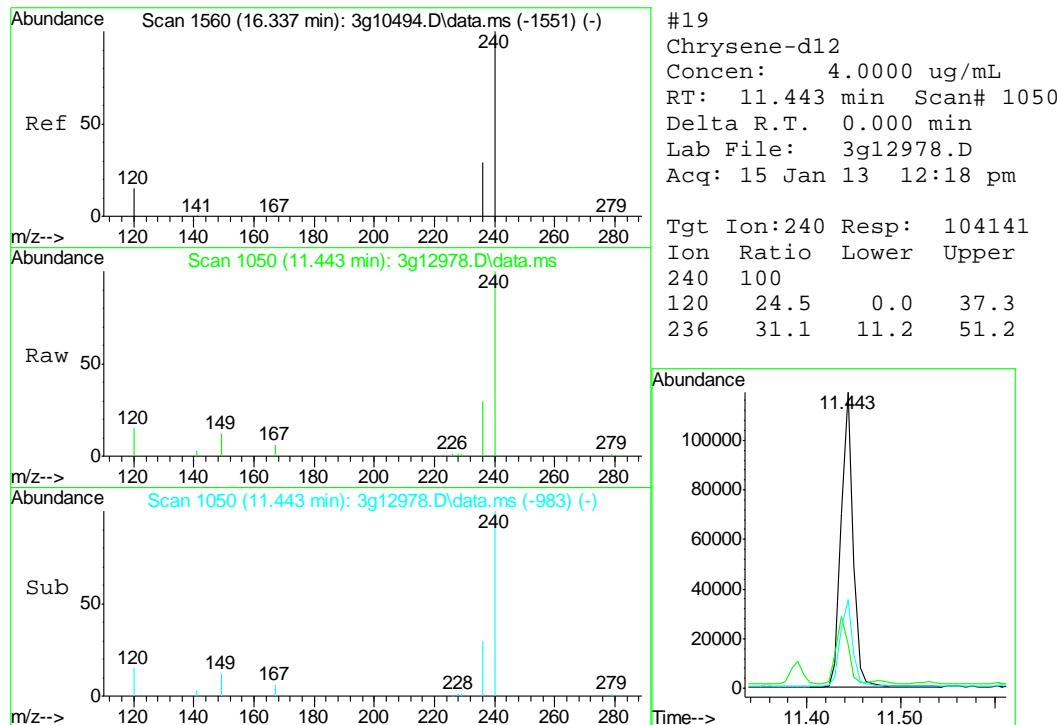


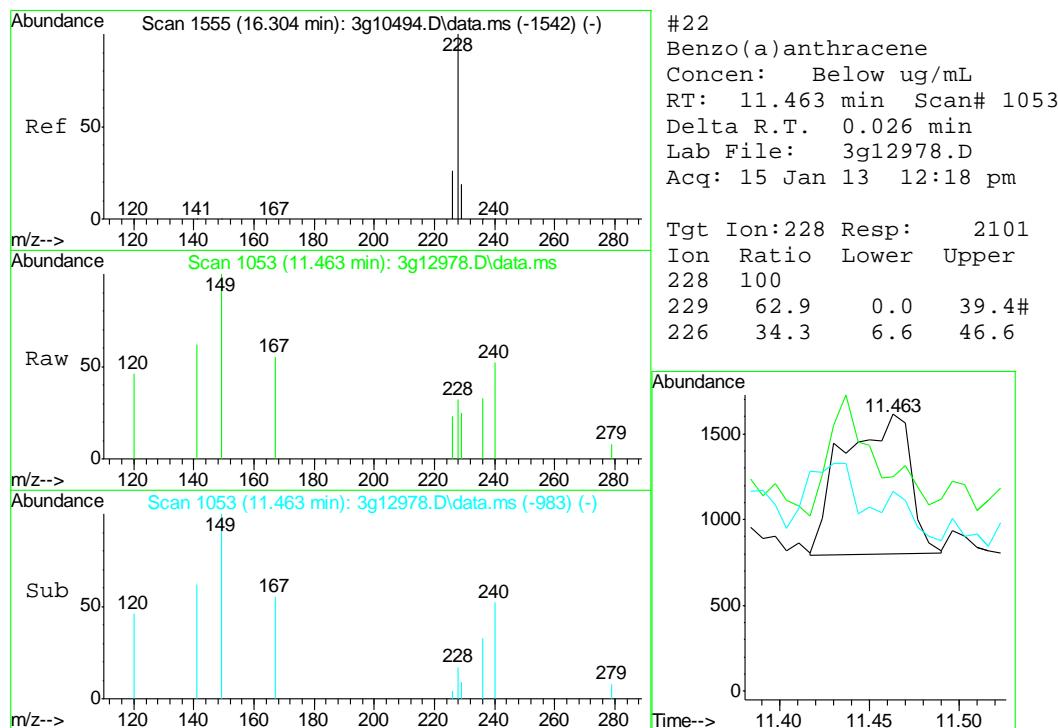
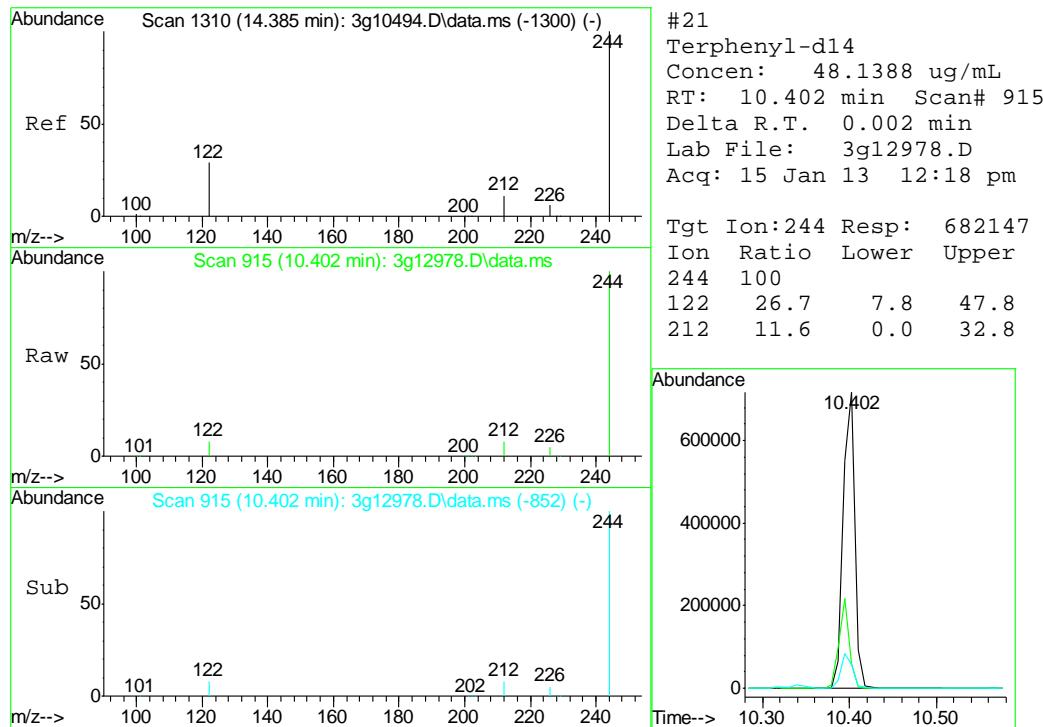


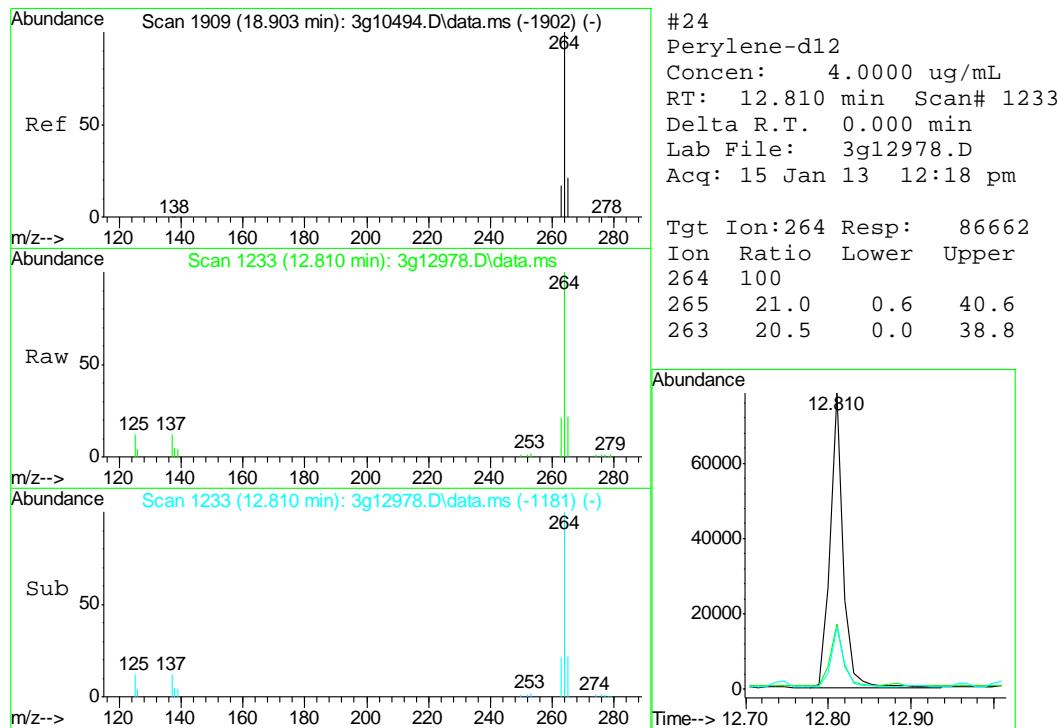
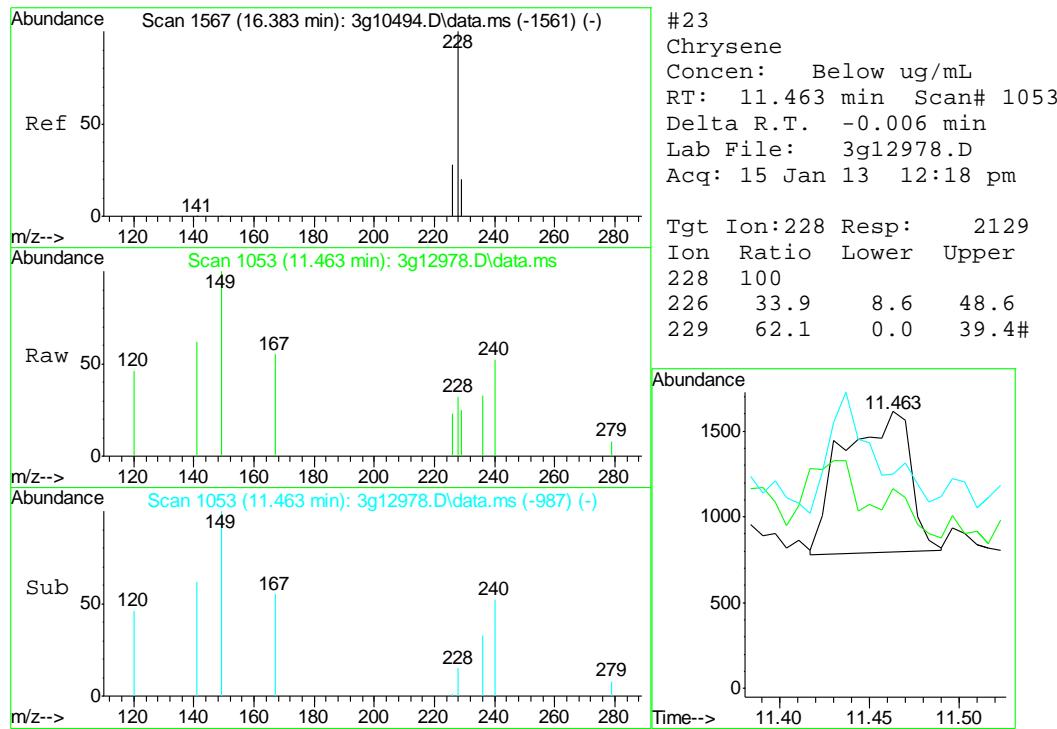


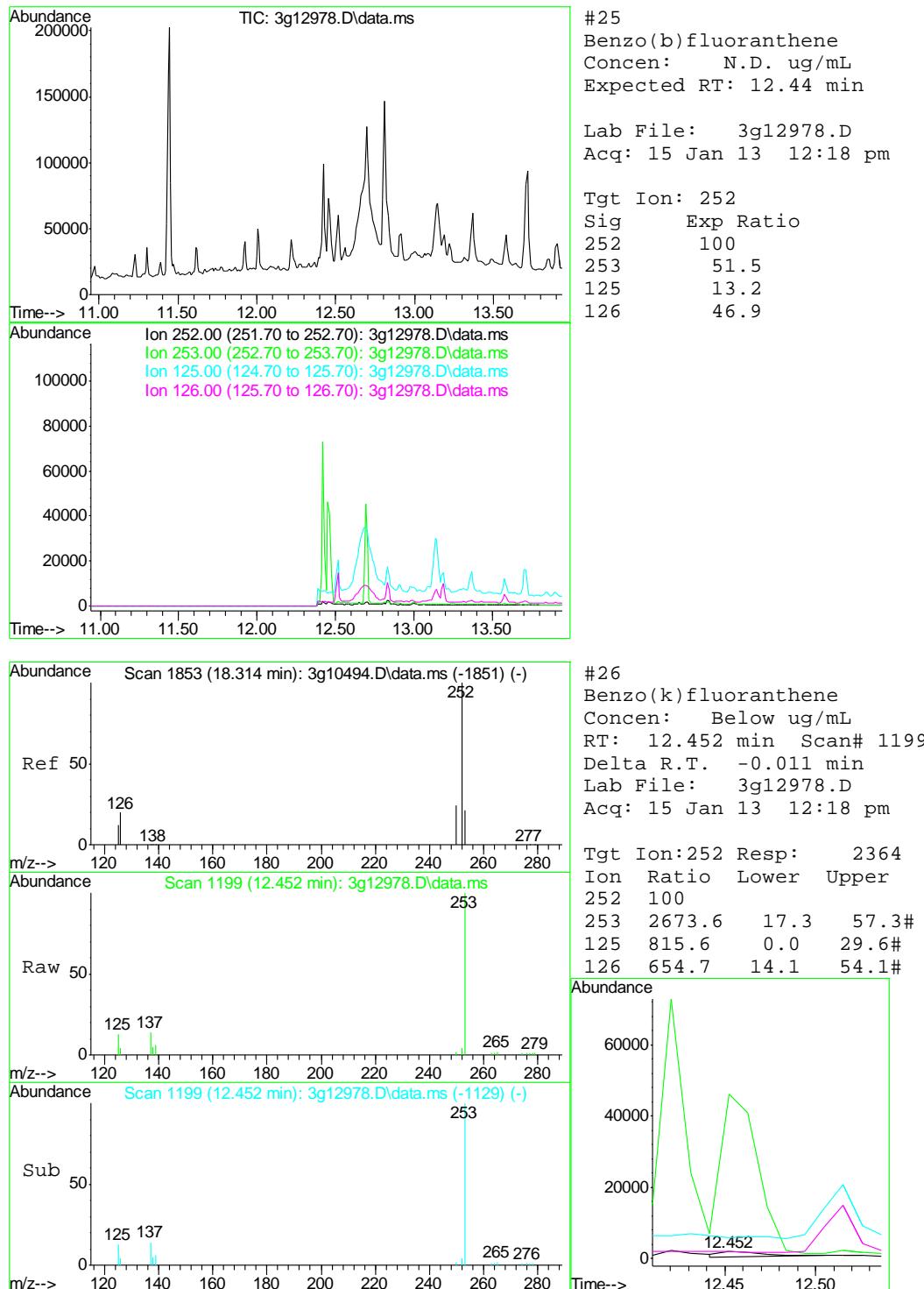


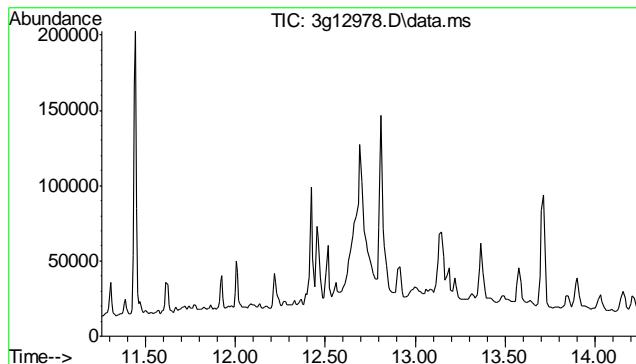




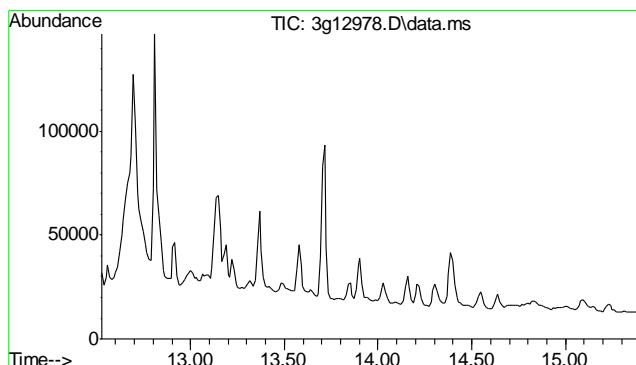
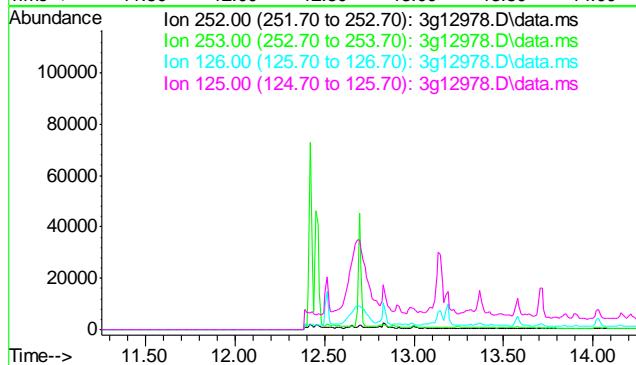




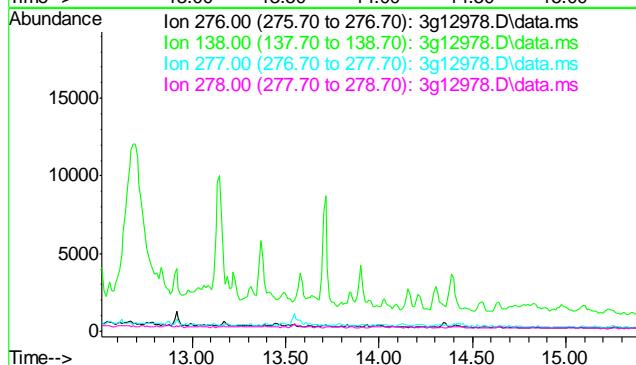


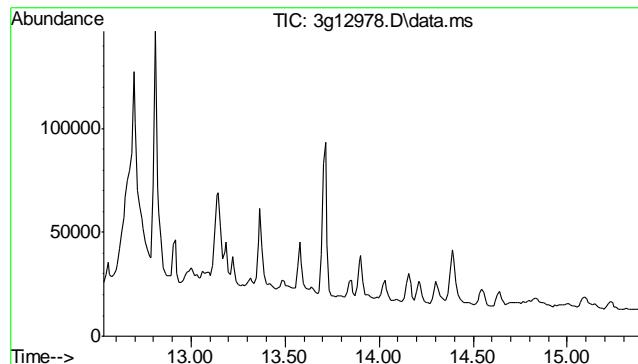


#27
 Benzo(a)pyrene
 Concen: N.D. ug/mL
 Expected RT: 12.76 min
 Lab File: 3g12978.D
 Acq: 15 Jan 13 12:18 pm
 Tgt Ion: 252
 Sig Exp Ratio
 252 100
 253 21.5
 126 20.4
 125 14.5

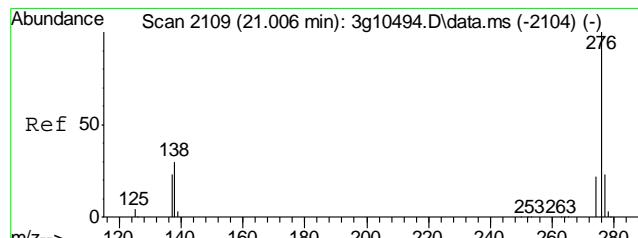
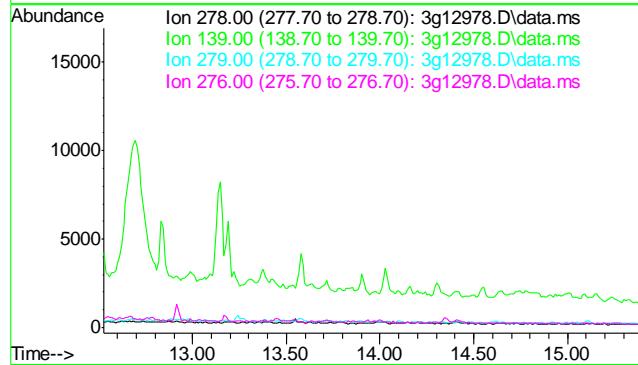


#28
 Indeno(1,2,3-cd)pyrene
 Concen: N.D. ug/mL
 Expected RT: 14.02 min
 Lab File: 3g12978.D
 Acq: 15 Jan 13 12:18 pm
 Tgt Ion: 276
 Sig Exp Ratio
 276 100
 138 40.0
 277 24.8
 278 76.2

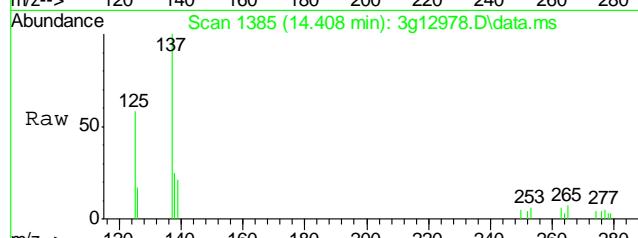




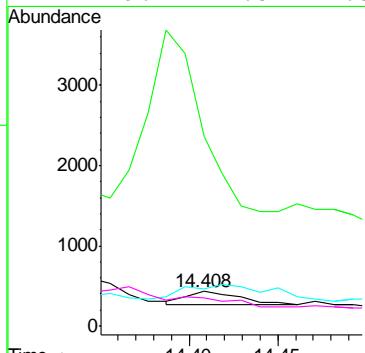
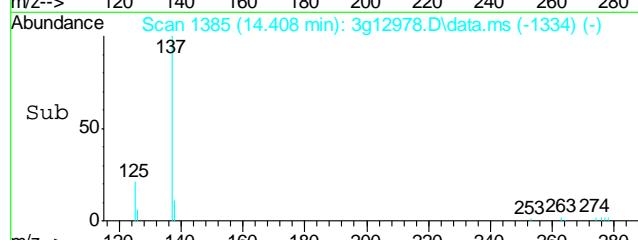
#29
Dibenz(a,h)anthracene
Concen: N.D. ug/mL
Expected RT: 14.03 min
Lab File: 3g12978.D
Acq: 15 Jan 13 12:18 pm
Tgt Ion: 278
Sig Exp Ratio
278 100
139 30.8
279 22.9
276 131.2



#30
Benzo(g,h,i)perylene
Concen: Below ug/mL
RT: 14.408 min Scan# 1385
Delta R.T. 0.034 min
Lab File: 3g12978.D
Acq: 15 Jan 13 12:18 pm



Tgt Ion: 276 Resp: 378
Ion Ratio Lower Upper
276 100
138 1253.7 15.1 55.1#
277 179.9 3.3 43.3#
274 229.1 1.5 41.5#



Quantitation Report (QT Reviewed)

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson
01/16/13 11:51

Data Path : C:\msdchem\1\DATA\011513\
Data File : 3g12973.D
Acq On : 15 Jan 2013 10:14 am
Operator : DONC
Sample : OP7223-MB
Misc : OP7223,E3G621,30.00,,,1,1
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jan 15 13:46:44 2013
Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M
Quant Title : PAHSIM BASE
QLast Update : Thu Jan 10 14:18:35 2013
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.621	136	122903	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.337	164	71901	4.0000	ug/mL	0.01
15) Phenanthrene-d10	8.812	188	128999	4.0000	ug/mL	0.00
19) Chrysene-d12	11.443	240	101544	4.0000	ug/mL	0.00
24) Perylene-d12	12.810	264	84293	4.0000	ug/mL	0.00

System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	445750	40.3218	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	80.64%
7) 2-Fluorobiphenyl	6.676	172	1170373	42.6075	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	85.22%
21) Terphenyl-d14	10.402	244	725496	52.5073	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	105.02%

Target Compounds Qvalue

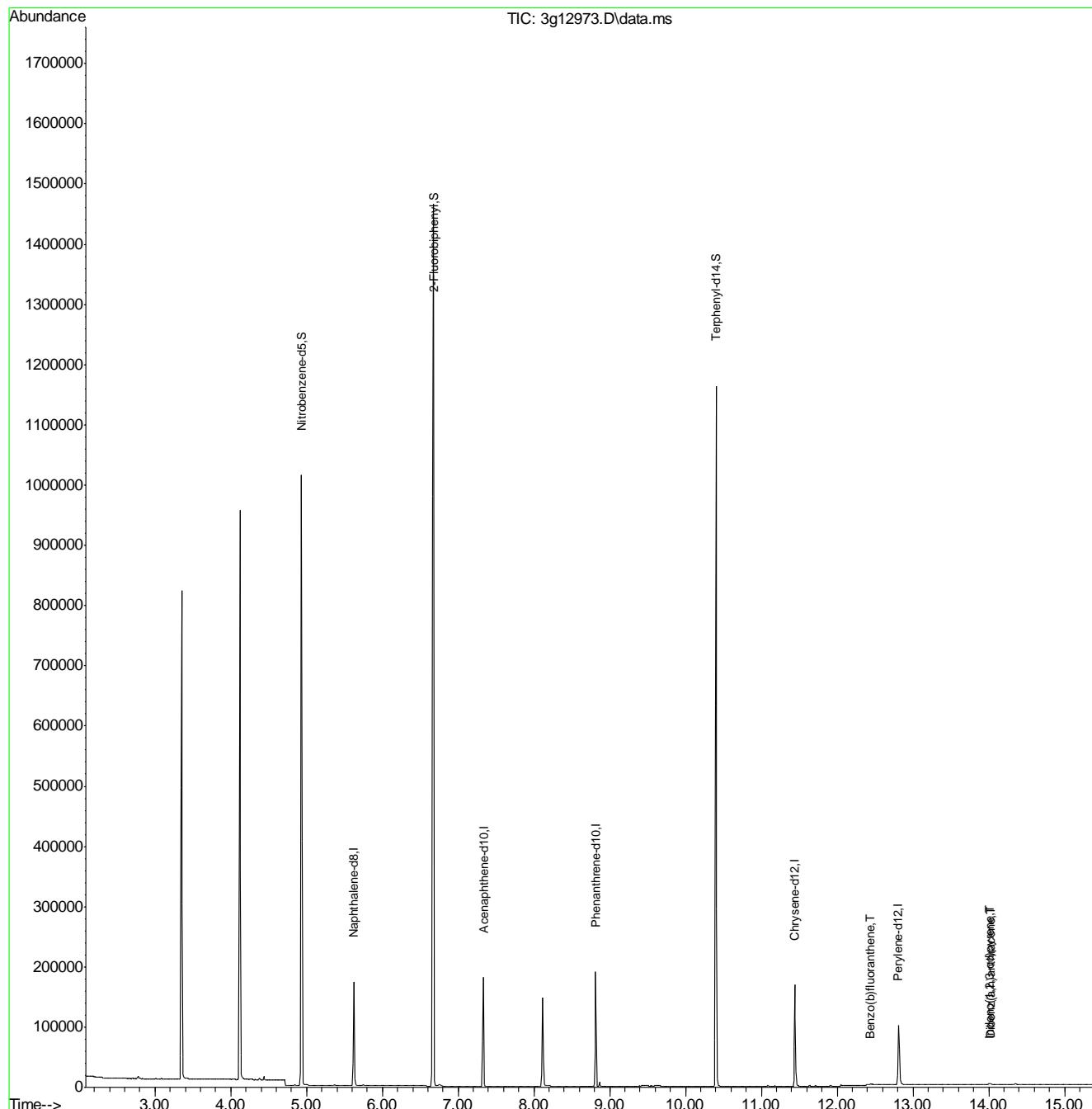
3) N-Nitrosodimethylamine	2.356	74	24	N.D.		
4) N-Nitrosodi-propylamine	0.000	70	0	N.D.	d	
5) Naphthalene	5.646	128	183	Below Cal	# 48	
8) 2-Methylnaphthalene	6.319	142	186	N.D.		
9) 1-Methylnaphthalene	6.394	142	78	N.D.		
10) Acenaphthylene	7.396	152	202	N.D.		
11) Acenaphthene	7.385	154	61	N.D.		
12) Dibenzofuran	7.857	168	73	N.D.		
13) Fluorene	0.000	166	0	N.D.	d	
14) Diphenylamine	0.000	169	0	N.D.	d	
16) Phenanthrene	8.812	178	152	N.D.		
17) Anthracene	8.891	178	39	N.D.		
18) Fluoranthene	10.015	202	205	N.D.		
20) Pyrene	10.244	202	152	N.D.		
22) Benzo(a)anthracene	11.437	228	1041	N.D.		
23) Chrysene	11.470	228	877	N.D.		
25) Benzo(b)fluoranthene	12.442	252	1135m	0.0649	ug/mL	
26) Benzo(k)fluoranthene	12.463	252	1206	N.D.		
27) Benzo(a)pyrene	0.000	252	0	N.D.	d	
28) Indeno(1,2,3-cd)pyrene	14.009	276	1461	0.0777	ug/mL	89
29) Dibenz(a,h)anthracene	14.019	278	1174	0.0897	ug/mL	92
30) Benzo(g,h,i)perylene	14.356	276	1418	N.D.		

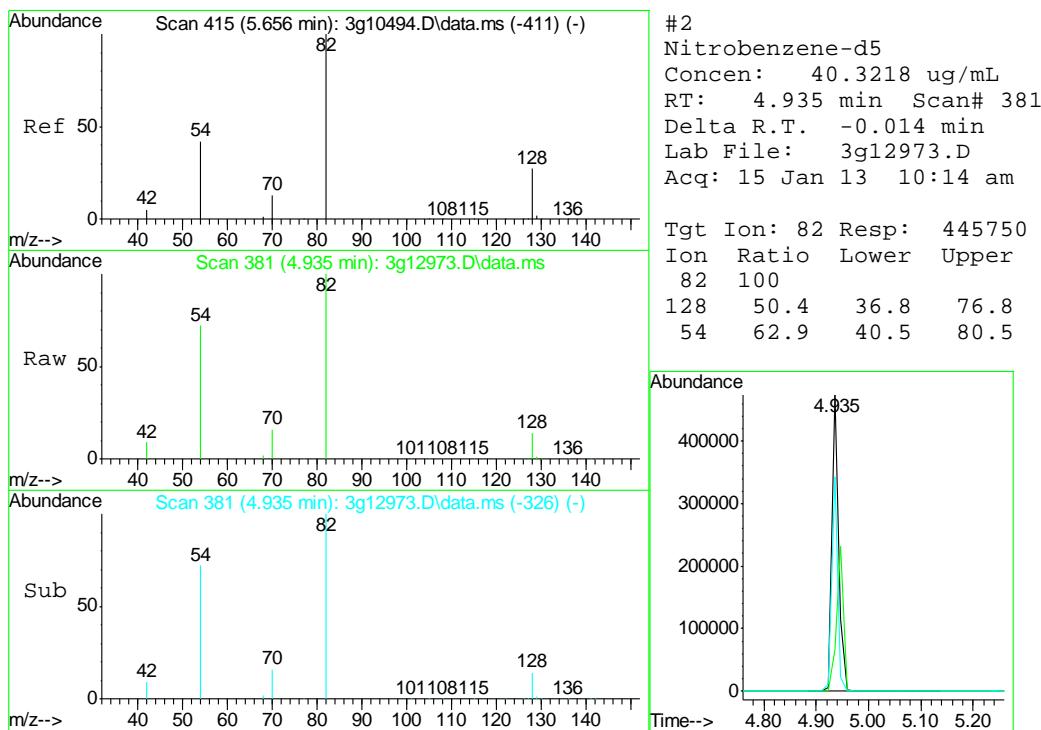
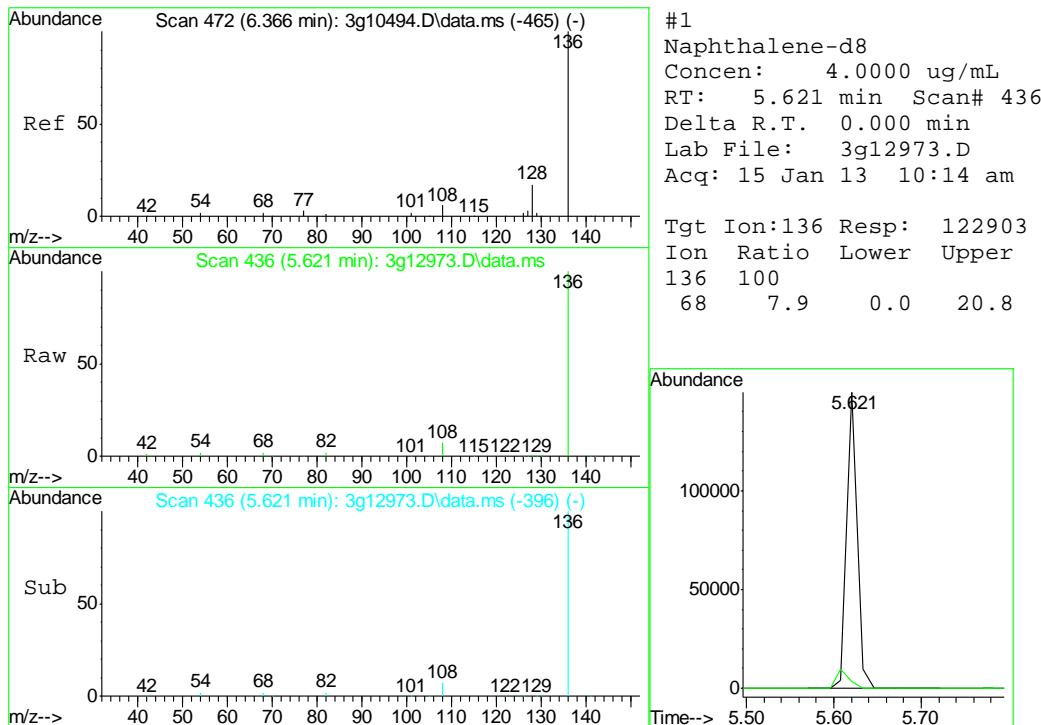
(#) = qualifier out of range (m) = manual integration (+) = signals summed

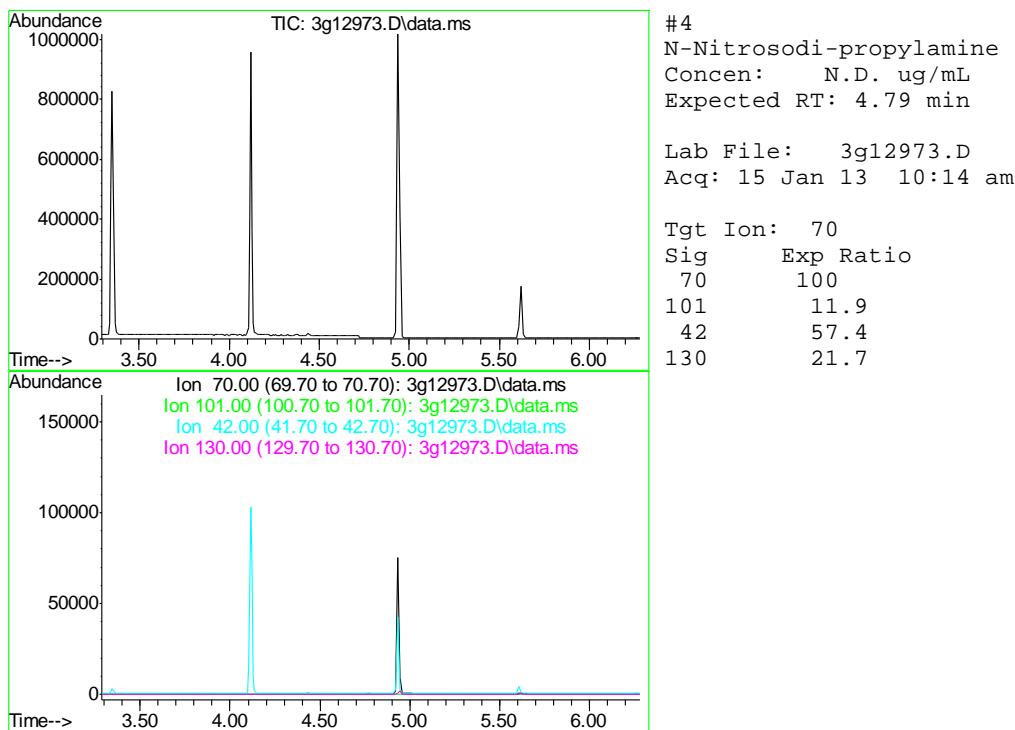
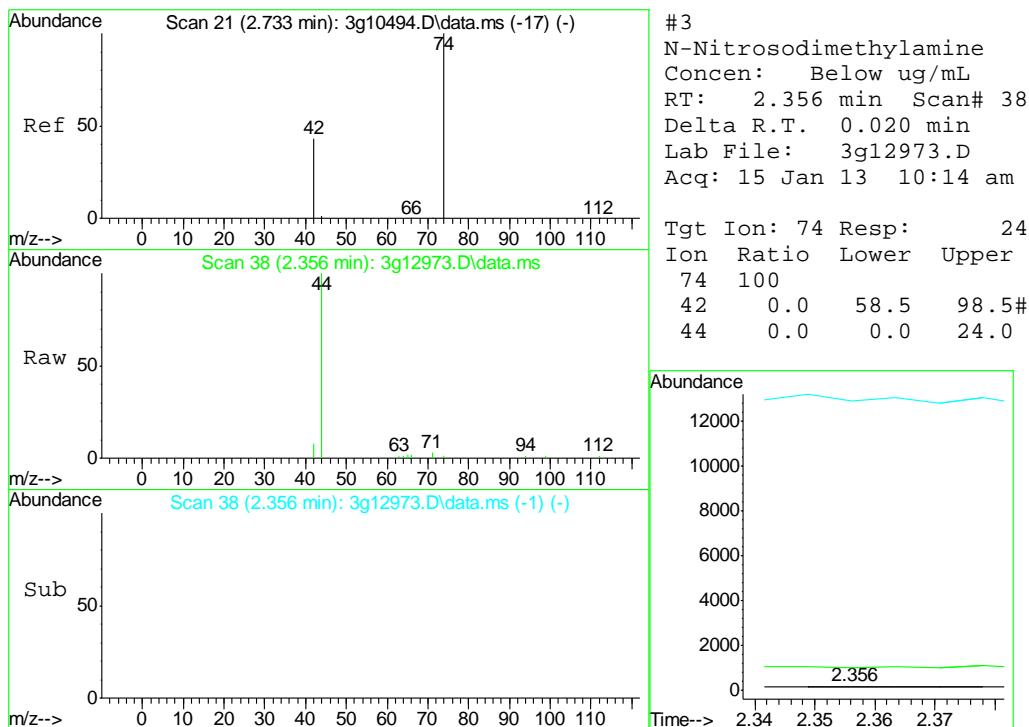
Quantitation Report (QT Reviewed)

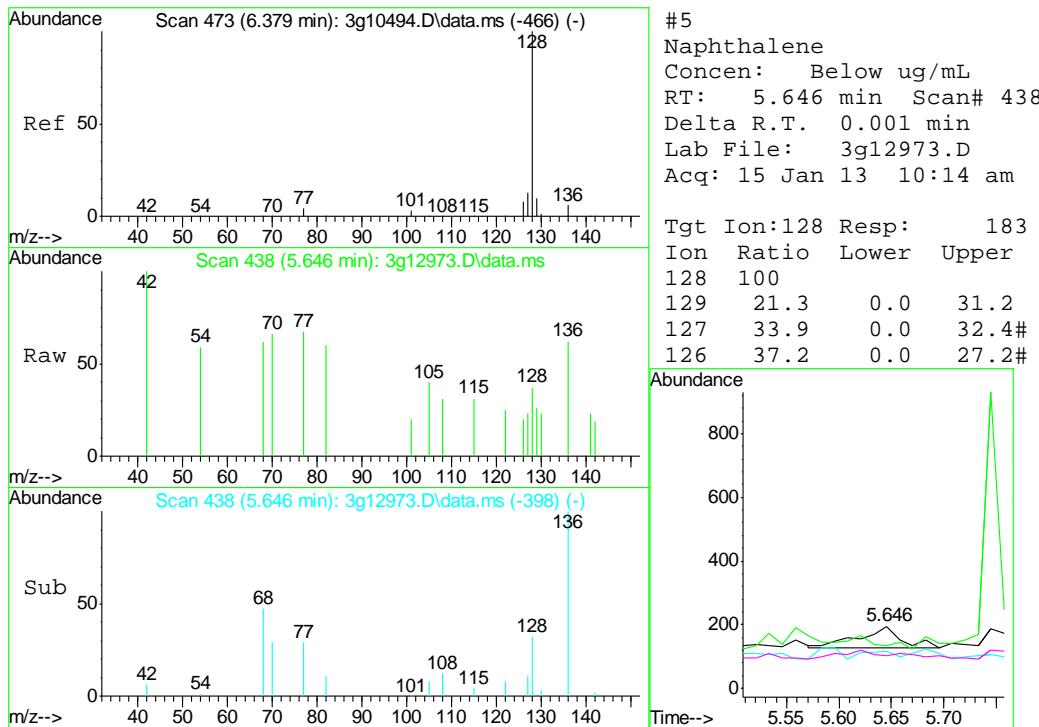
Data Path : C:\msdchem\1\DATA\011513\
 Data File : 3g12973.D
 Acq On : 15 Jan 2013 10:14 am
 Operator : DONC
 Sample : OP7223-MB
 Misc : OP7223,E3G621,30.00,,,1,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jan 15 13:46:44 2013
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Jan 10 14:18:35 2013
 Response via : Initial Calibration



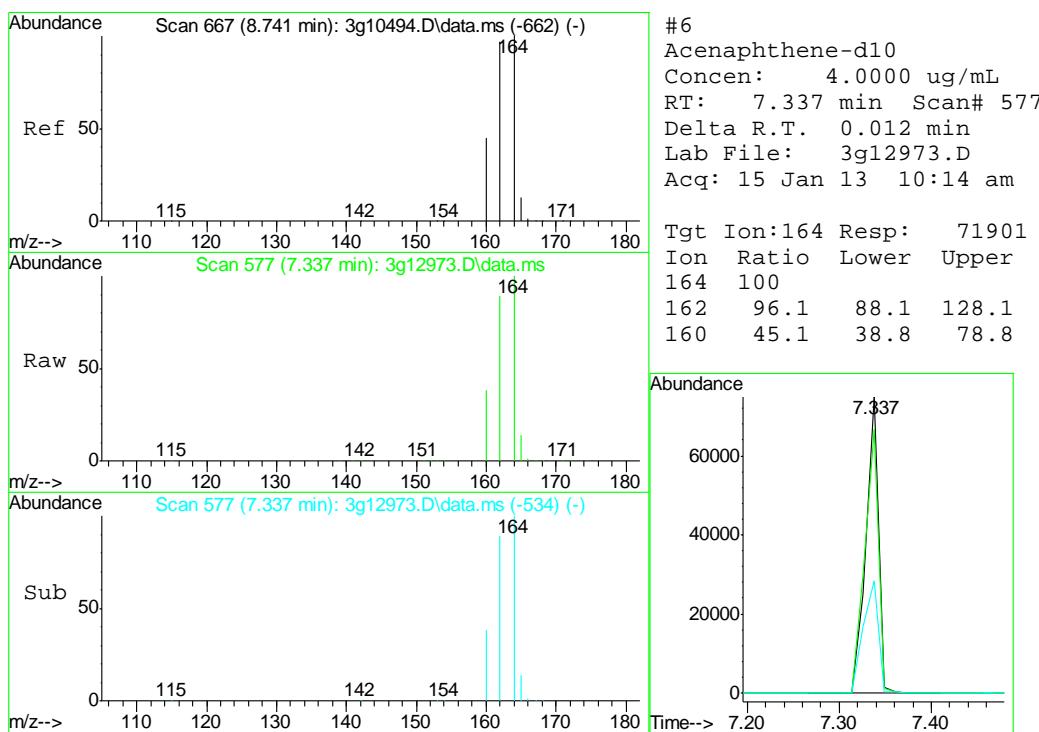


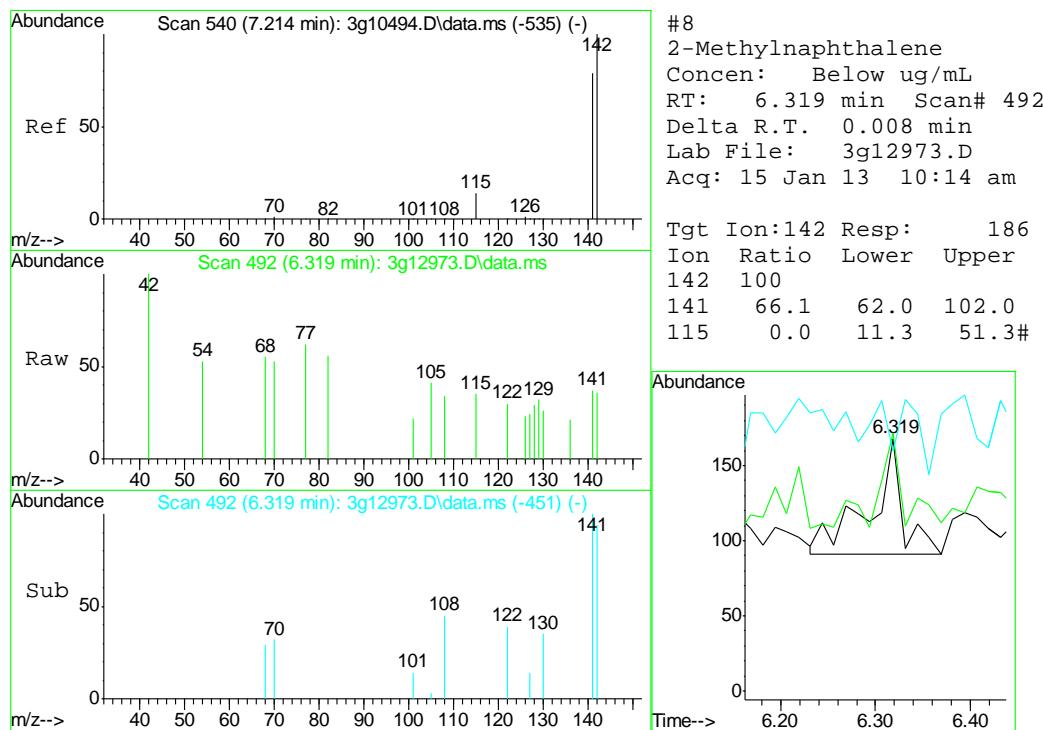
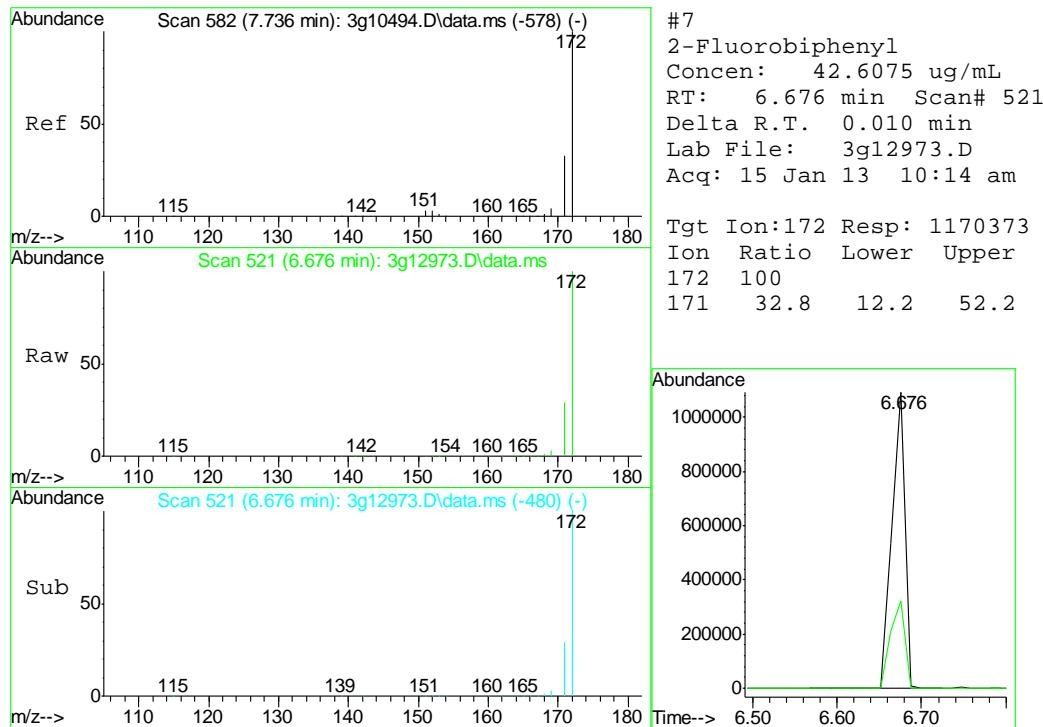


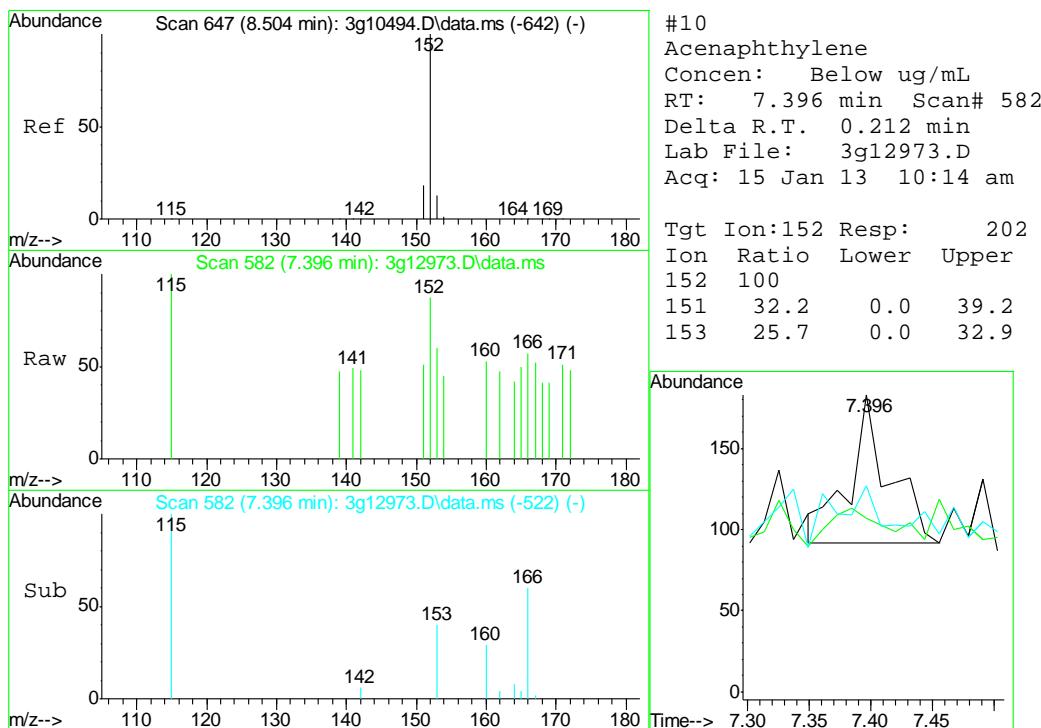
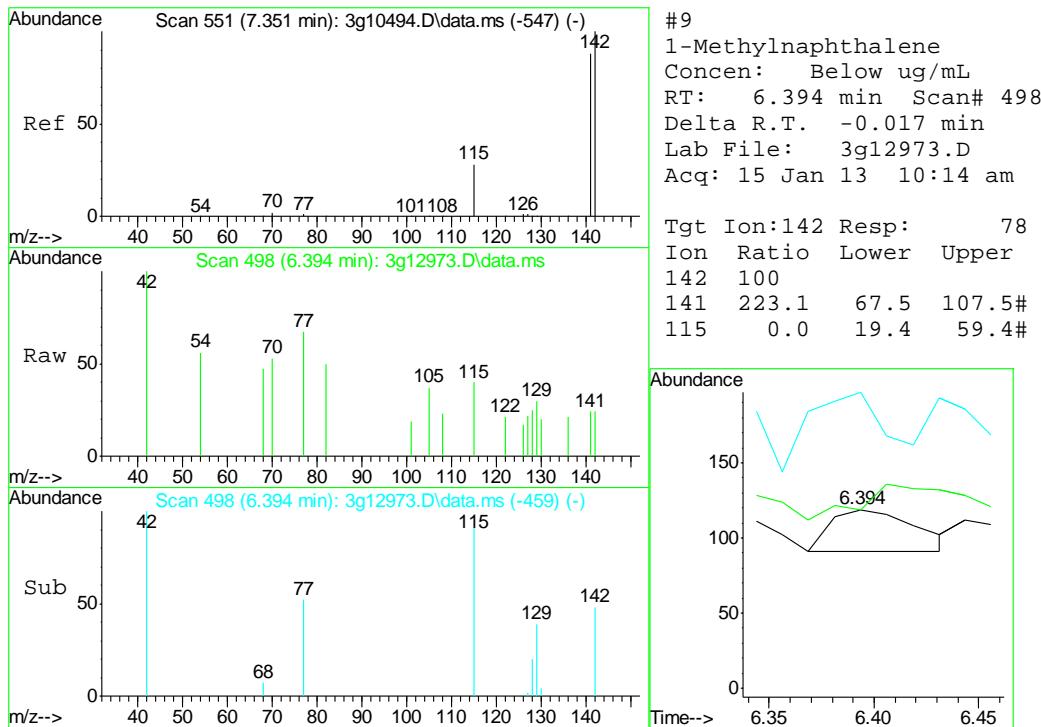


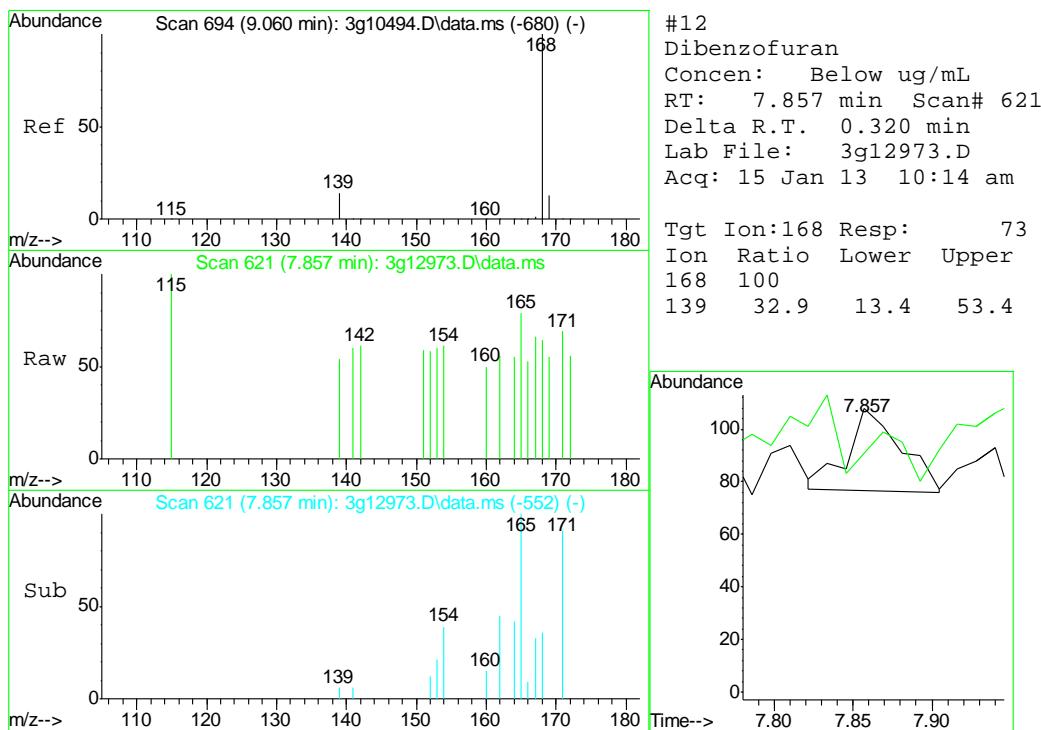
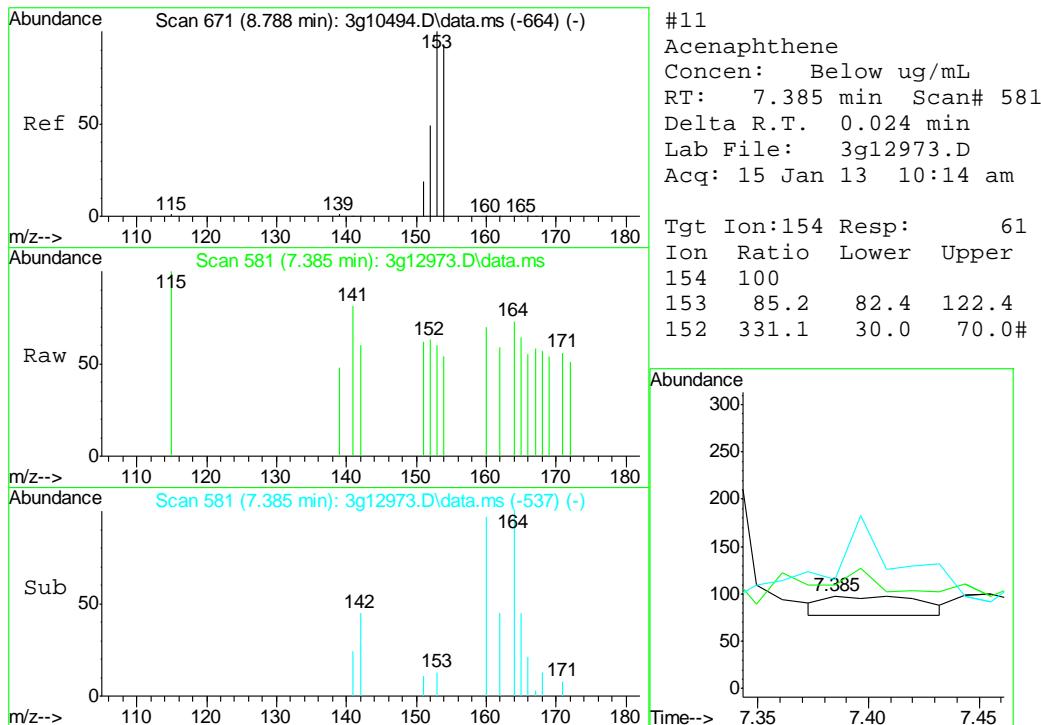
9.2.1

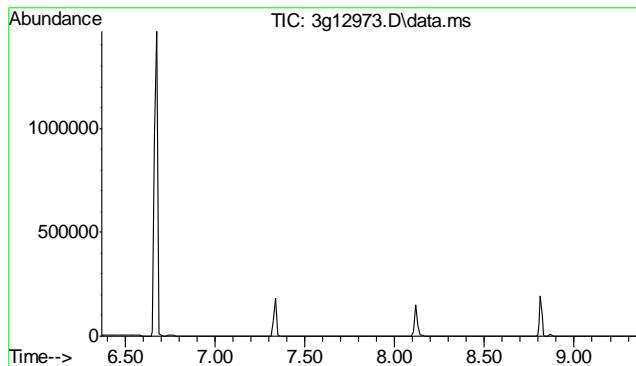
9







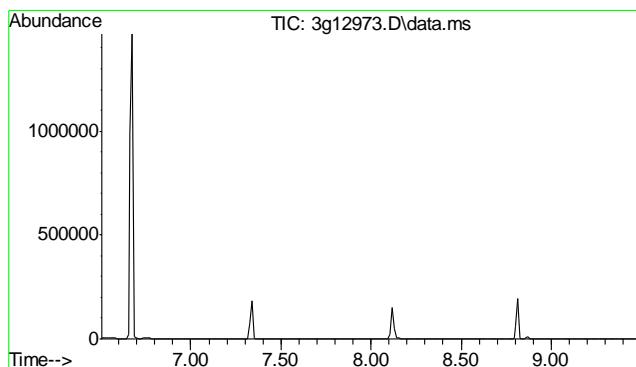
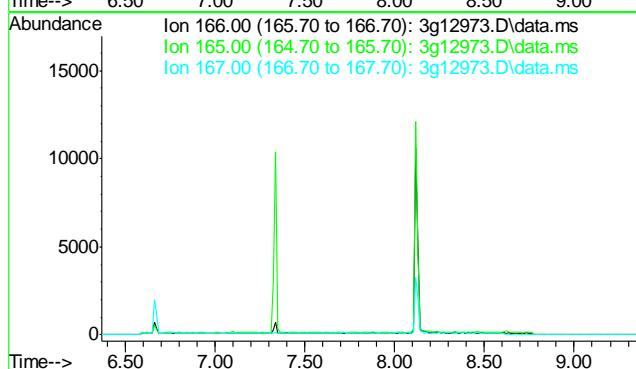




#13
Fluorene
Concen: N.D. ug/mL
Expected RT: 7.87 min

Lab File: 3g12973.D
Acq: 15 Jan 13 10:14 am

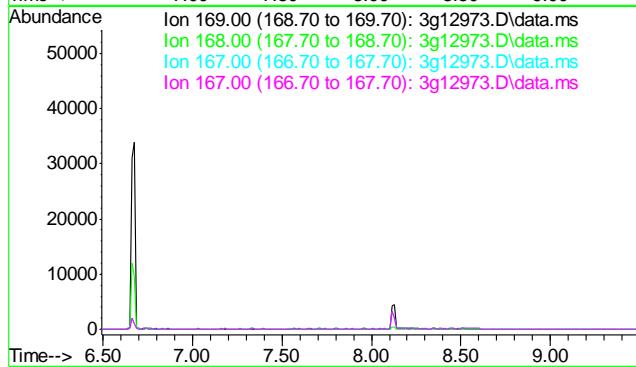
Tgt Ion: 166
Sig Exp Ratio
166 100
165 92.0
167 13.1



#14
Diphenylamine
Concen: N.D. ug/mL
Expected RT: 8.00 min

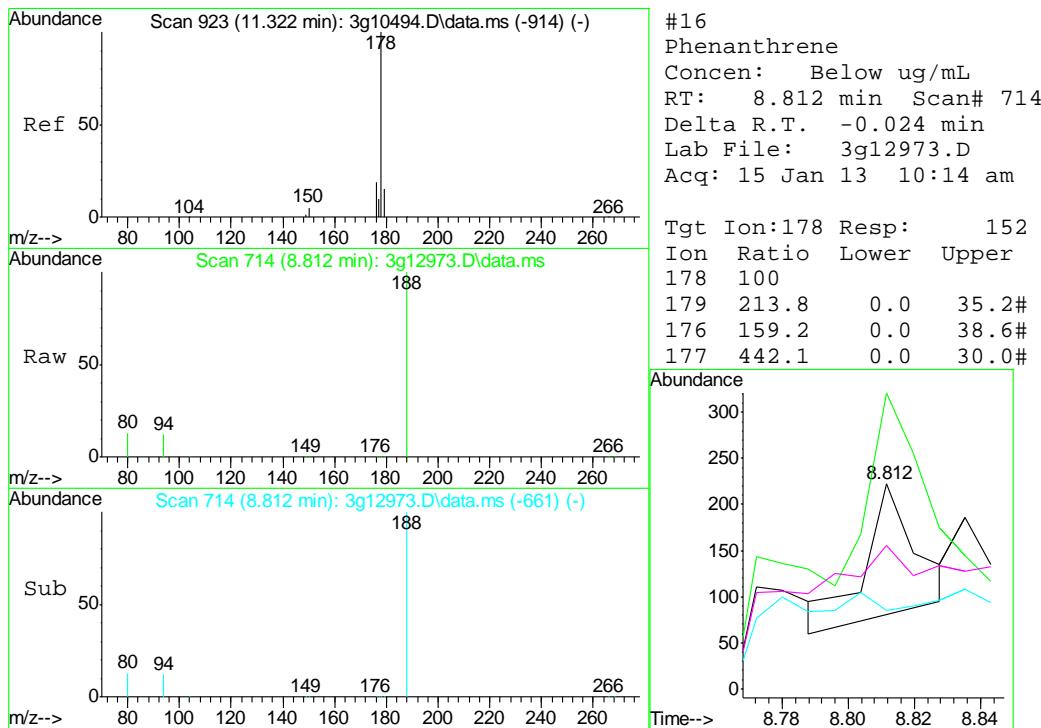
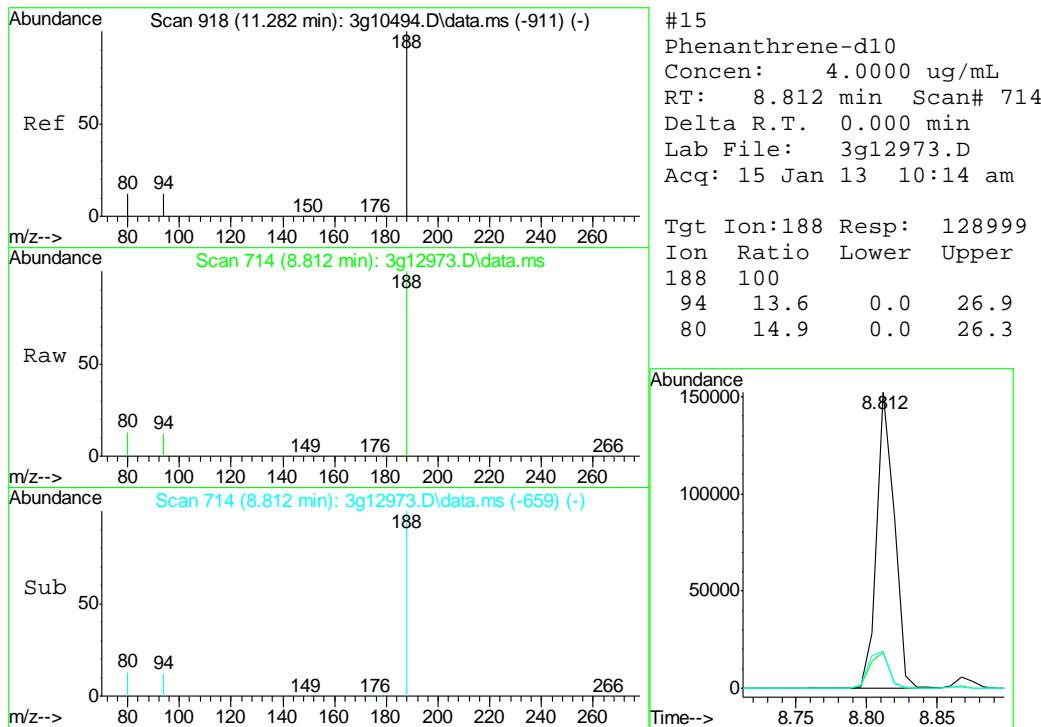
Lab File: 3g12973.D
Acq: 15 Jan 13 10:14 am

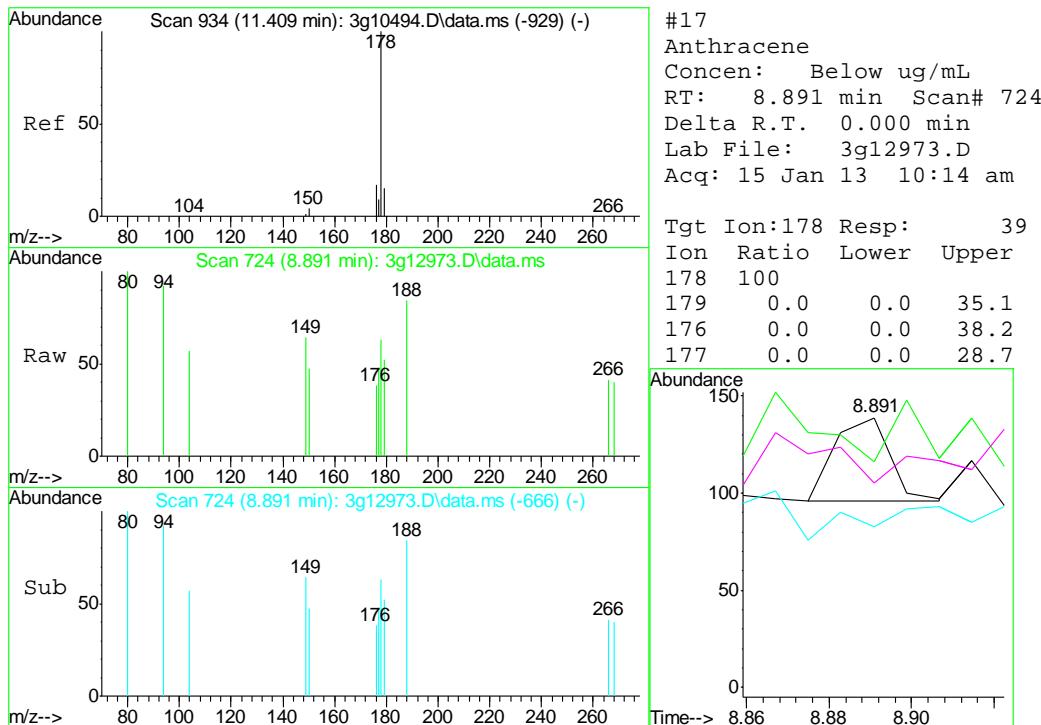
Tgt Ion: 169
Sig Exp Ratio
169 100
168 61.7
167 34.1
167 34.1



9.2.1

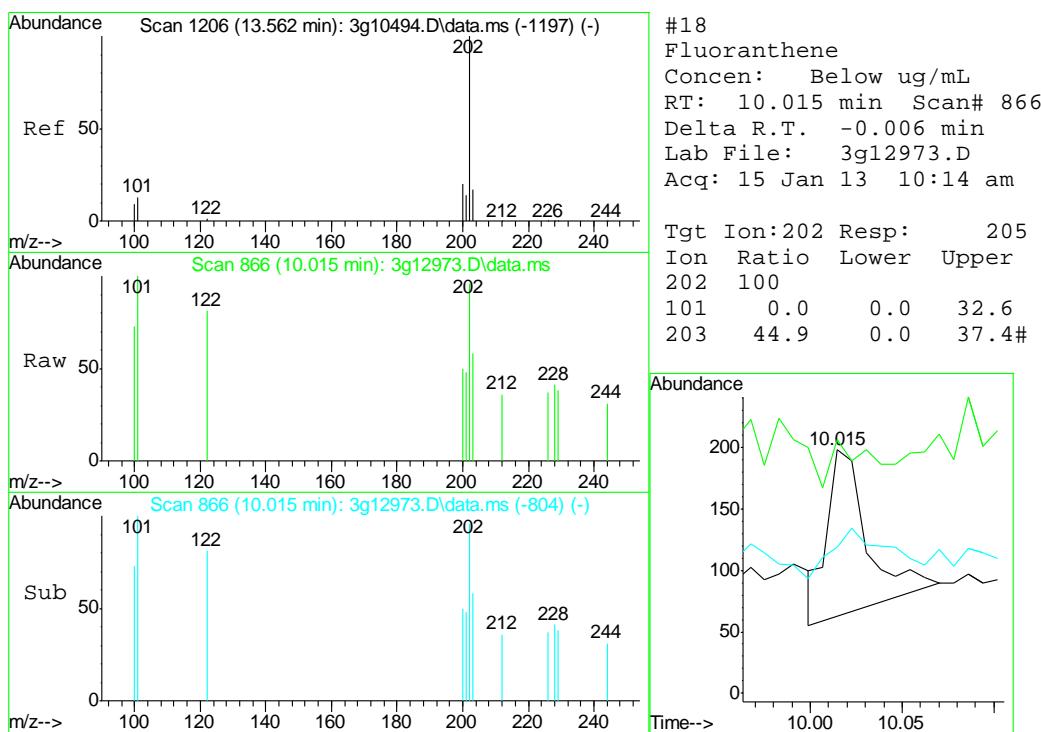
9

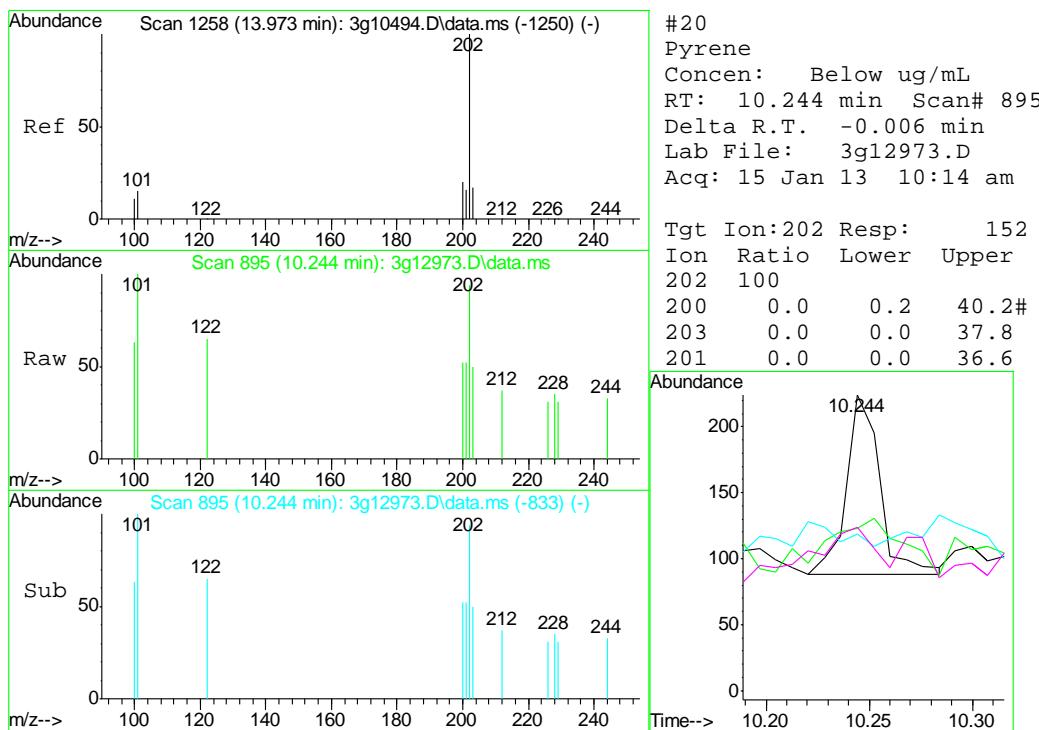
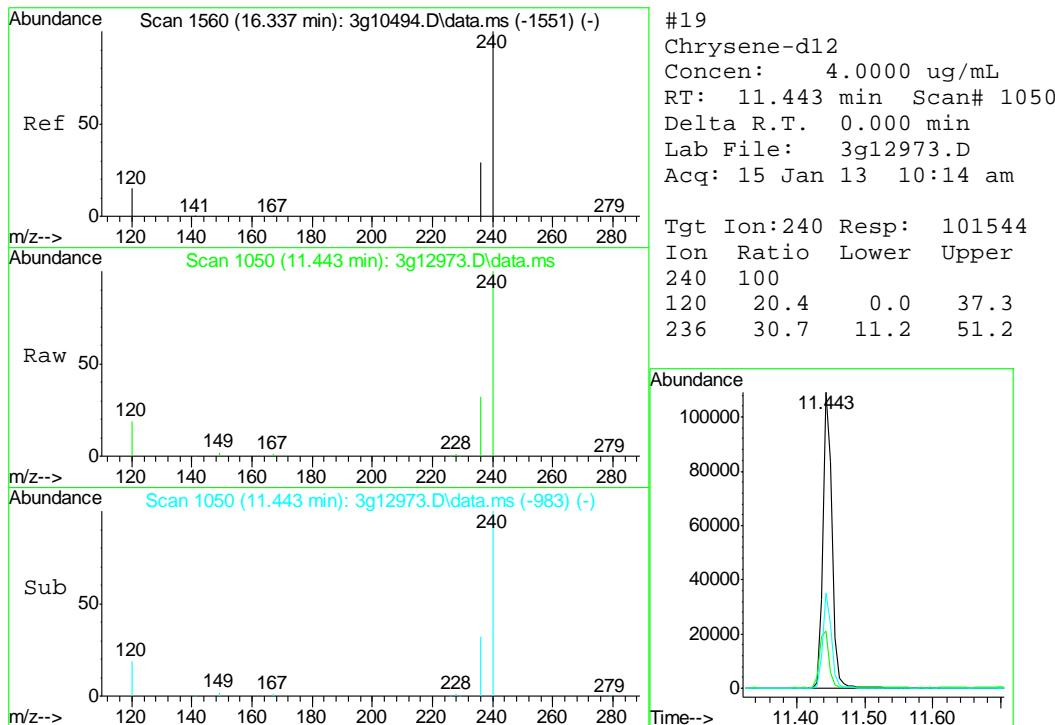


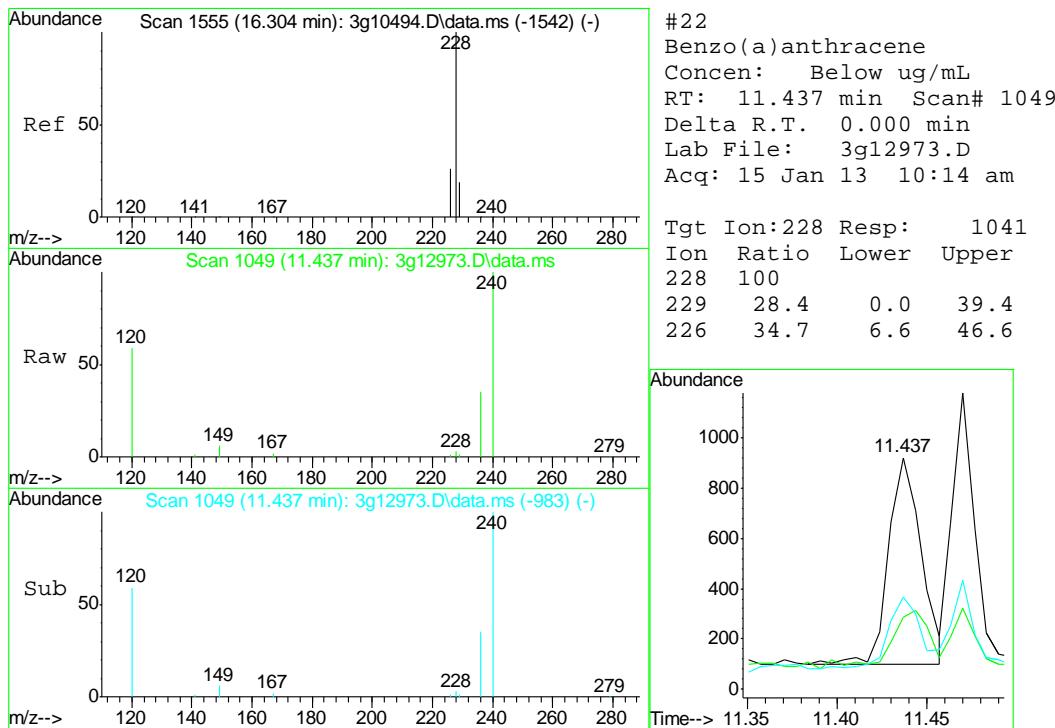
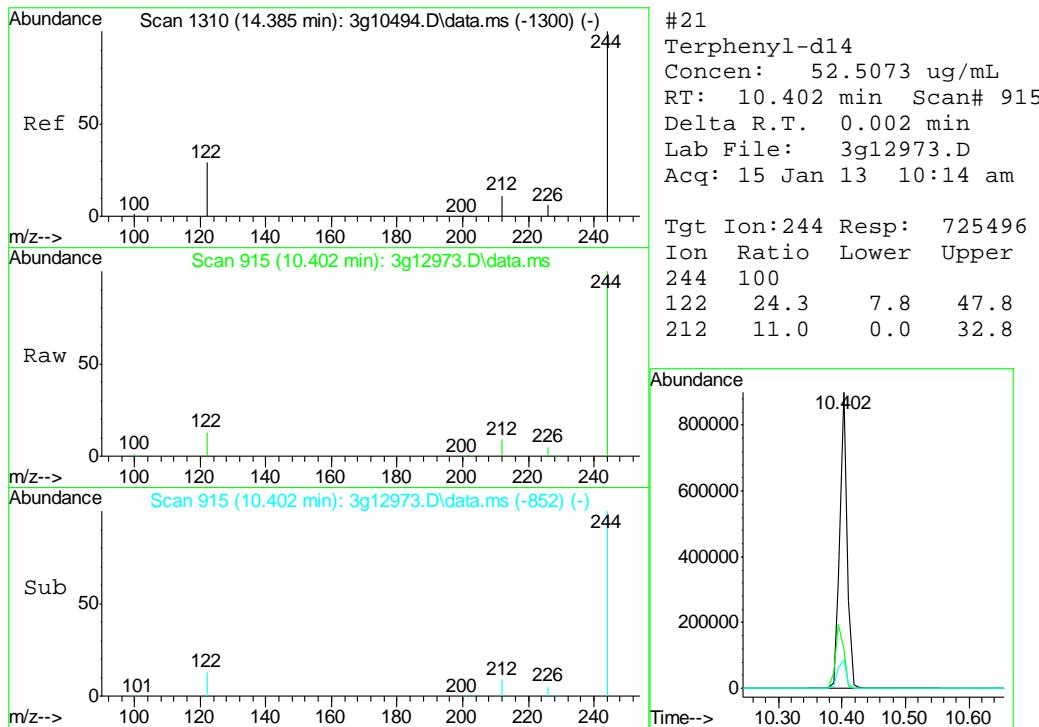


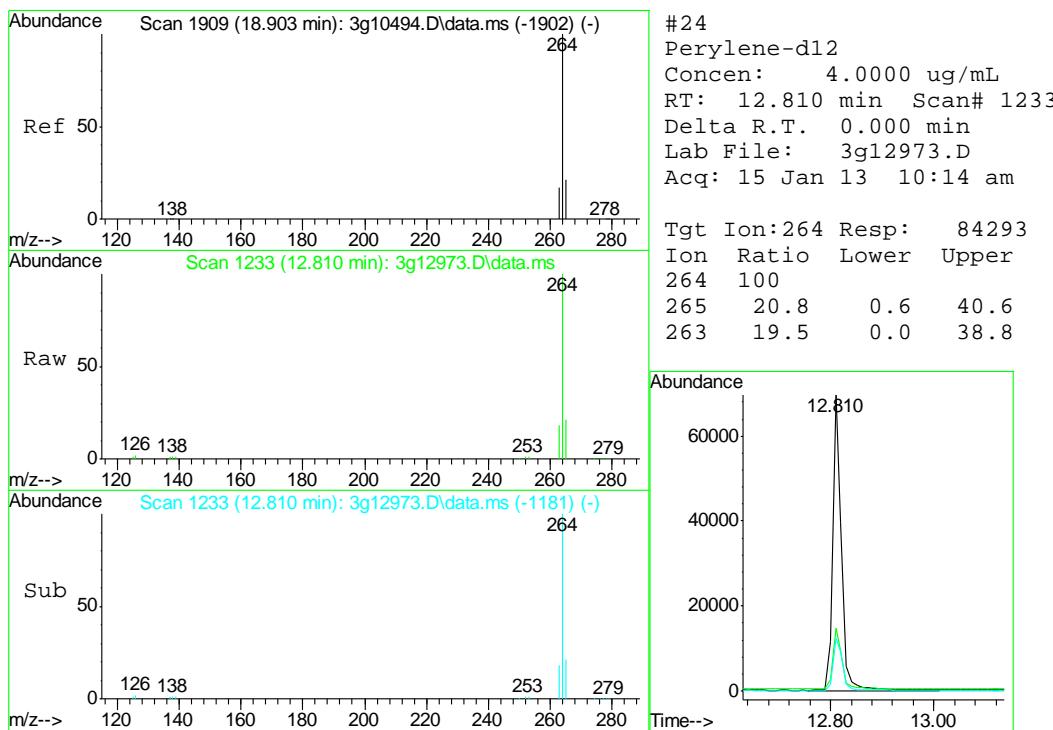
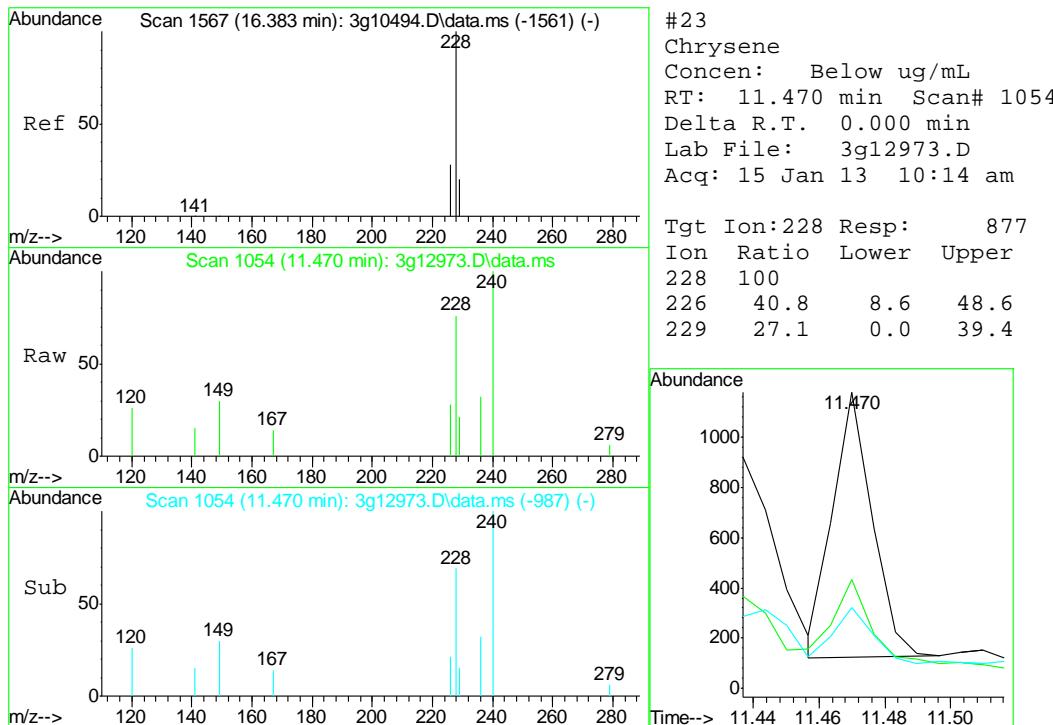
9.2.1

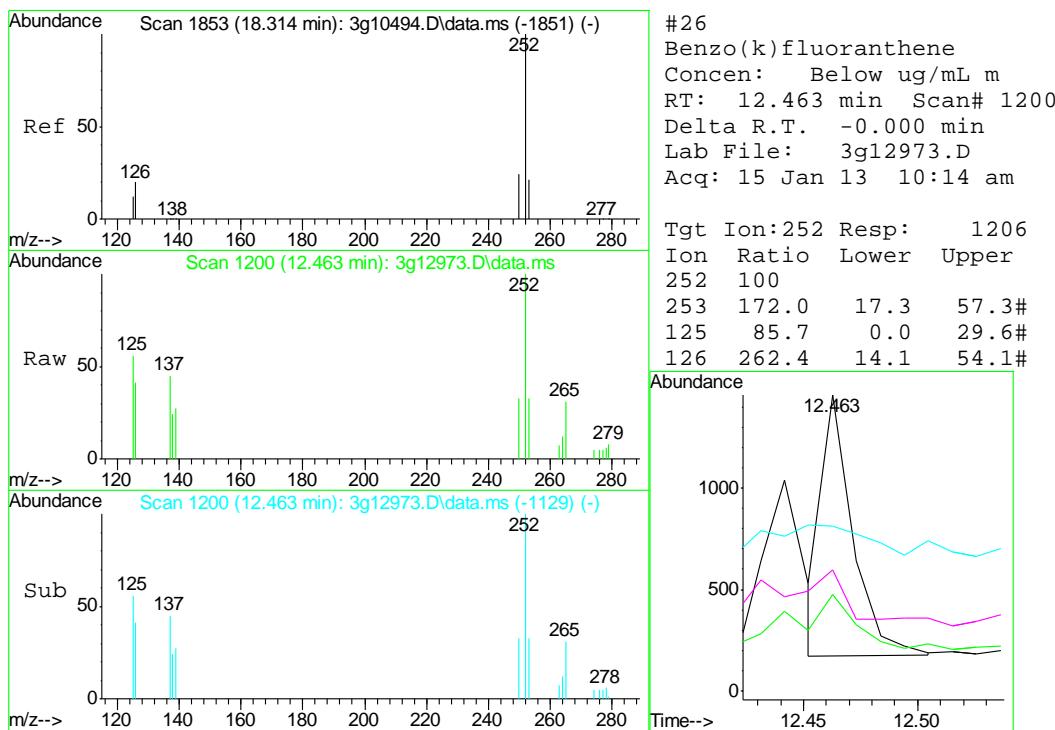
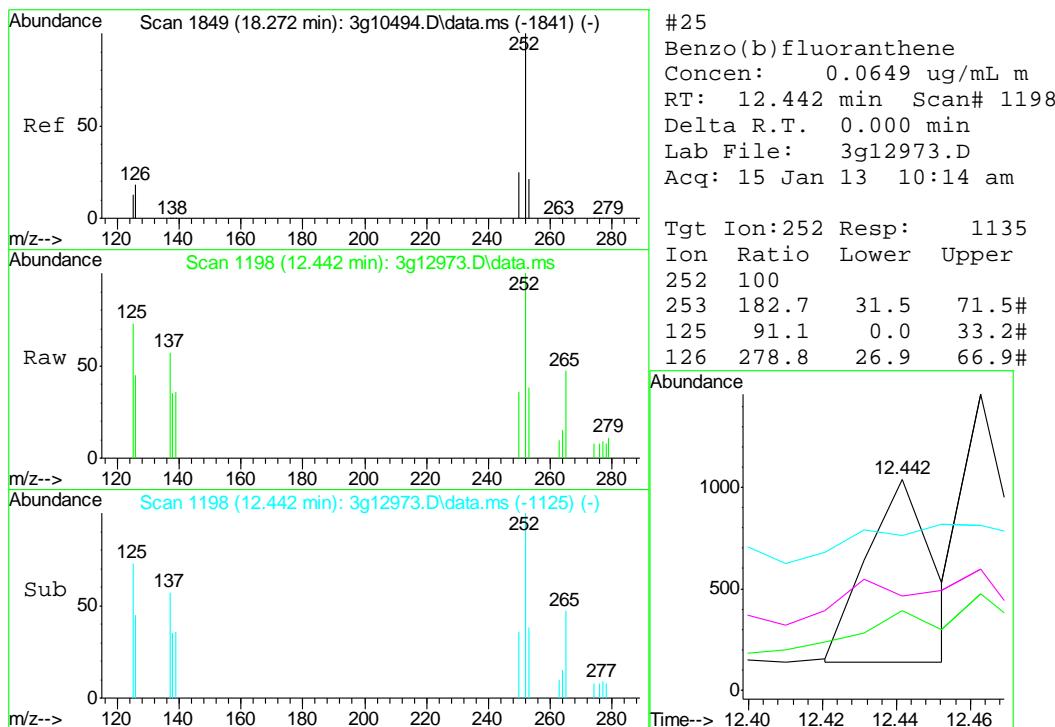
9

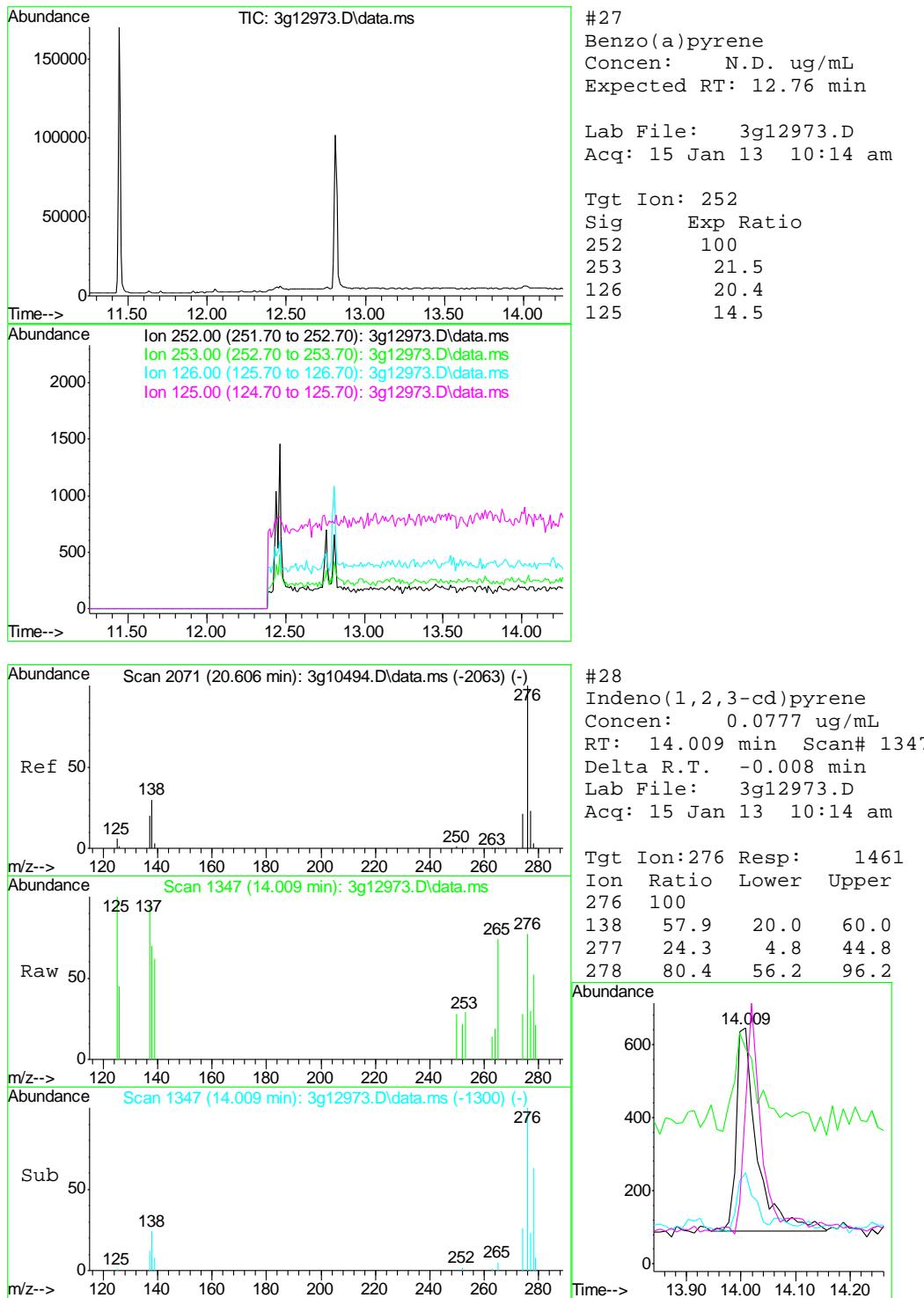


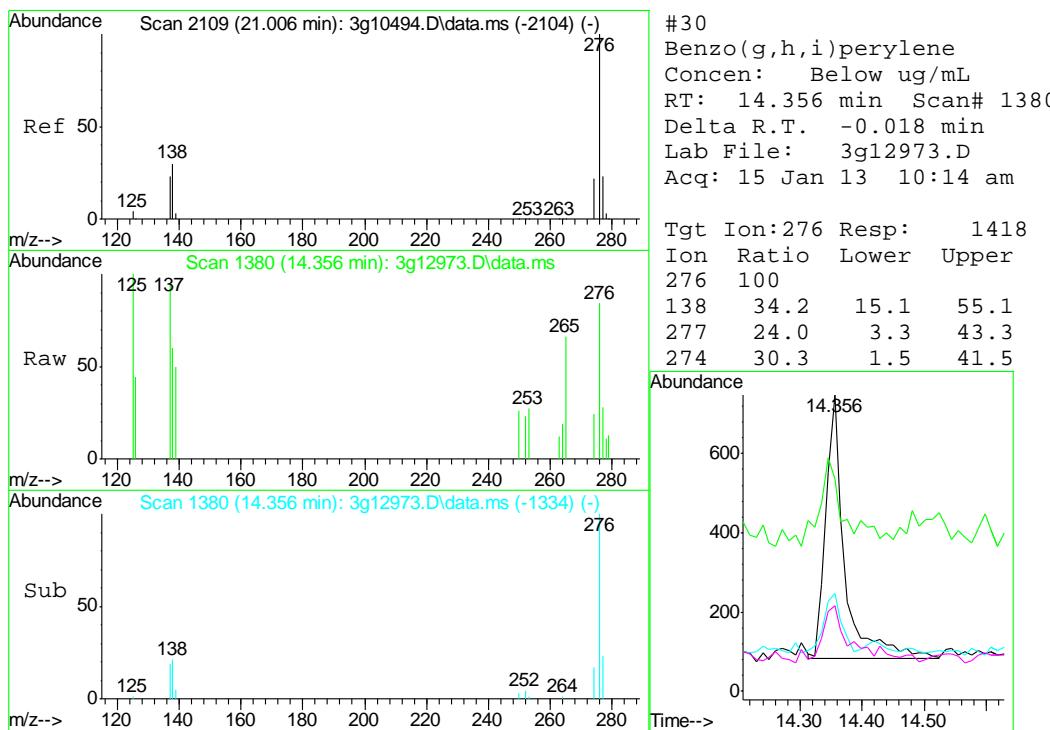
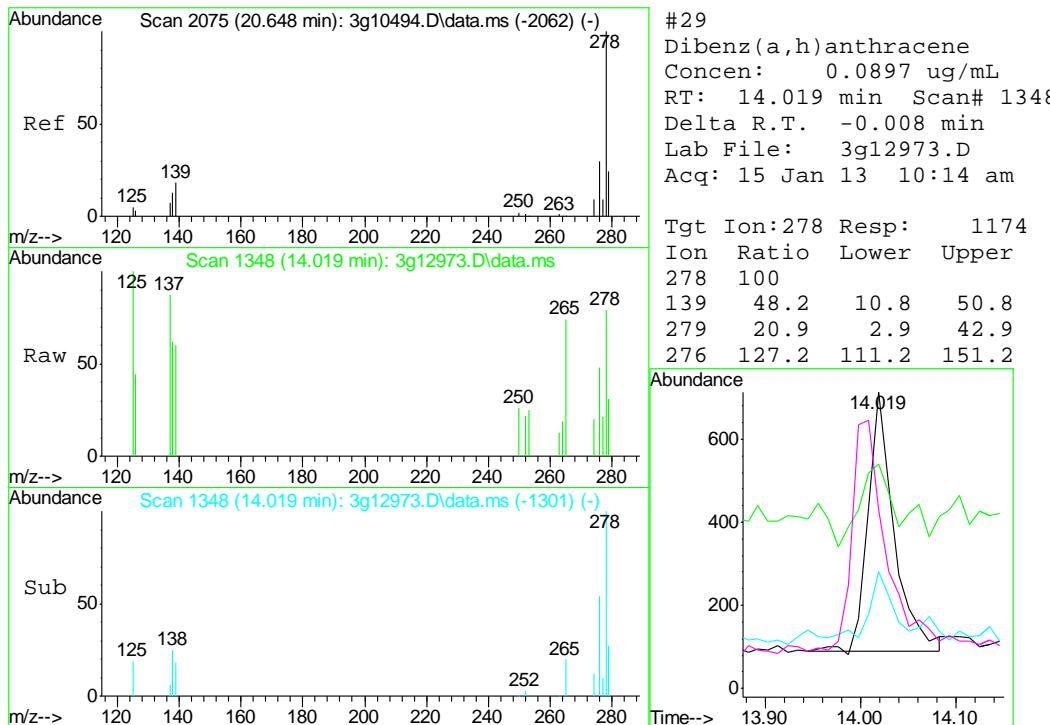














GC Volatiles

QC Data Summaries

Includes the following where applicable:

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

Method Blank Summary

Job Number: D42511
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1044-MB	GB19088.D	1	01/10/13	SK	n/a	n/a	GGB1044

The QC reported here applies to the following samples:

Method: SW846 8015B

D42511-1

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	87% 60-140%

10.1.1

10

Blank Spike Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1044-BS	GB19089.D	1	01/10/13	SK	n/a	n/a	GGB1044

The QC reported here applies to the following samples:

Method: SW846 8015B

D42511-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	123	112	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	104%	60-140%

10.2.1

10

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D42435-1MS	GB19091.D	1	01/10/13	SK	n/a	n/a	GGB1044
D42435-1MSD	GB19092.D	1	01/10/13	SK	n/a	n/a	GGB1044
D42435-1	GB19090.D	1	01/10/13	SK	n/a	n/a	GGB1044

The QC reported here applies to the following samples:

Method: SW846 8015B

D42511-1

CAS No.	Compound	D42435-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		127	140	111	140	111	0	70-130/30
10.3.1										
CAS No.	Surrogate Recoveries	MS	MSD	D42435-1			Limits			
120-82-1	1,2,4-Trichlorobenzene	92%	95%	93%			60-140%			

* = Outside of Control Limits.



GC Volatiles

Raw Data

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson 01/11/13 11:39

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011013\GB19112.D\FID1A.CH Vial: 27
 Signal #2 : Y:\1\DATA\011013\GB19112.D\FID2B.CH
 Acq On : 11 Jan 2013 6:12 am Operator: StephK
 Sample : D42511-1, 50X Inst : GC/MS Ins
 Misc : GC3347,GGB1044,,5.034,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Jan 11 08:57:32 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Jan 10 16:31:50 2013
 Response via : Initial Calibration
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.37	2835281	90.486 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.37	14291134	87.931 %	

Target Compounds

1) H	TVH-Gasoline	7.23	3605558	<MDL	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	7.66	124969	0.315	ug/L
7) T	Ethylbenzene	0.00	0	N.D.	ug/L d
8) T	m,p-Xylene	0.00	0	N.D.	ug/L d
9) T	o-Xylene	0.00	0	N.D.	ug/L d
11) T	Naphthalene	14.55	15343	<MDL	ug/L m

11.11

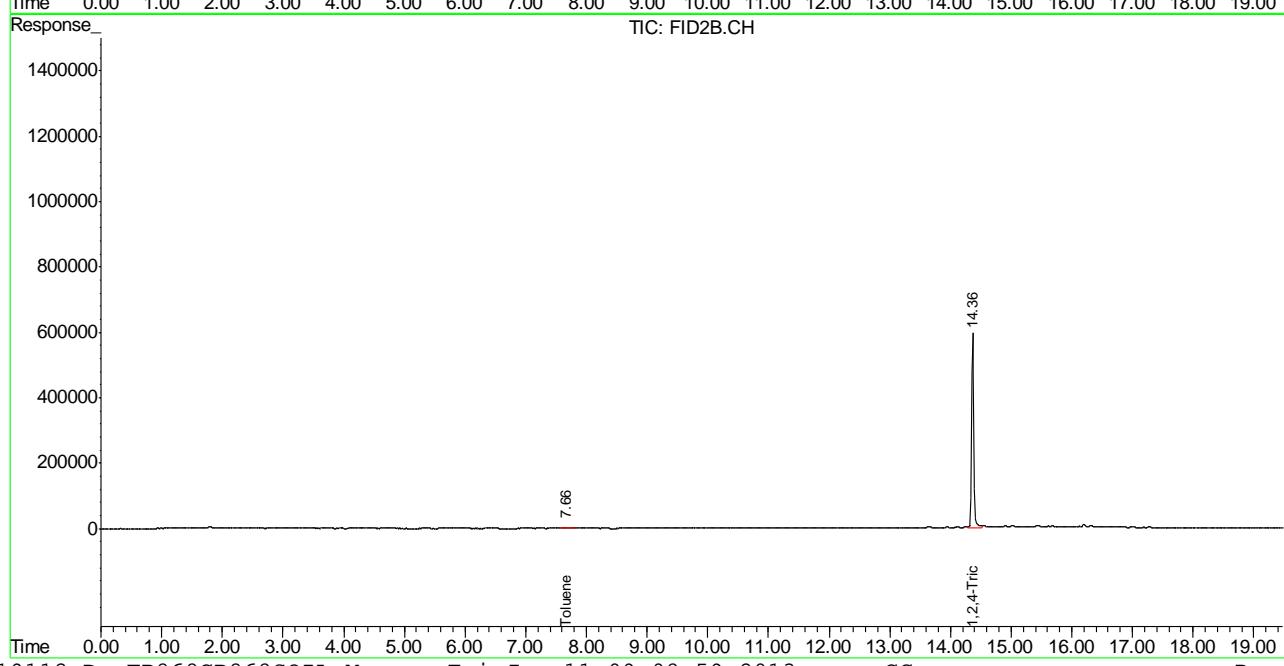
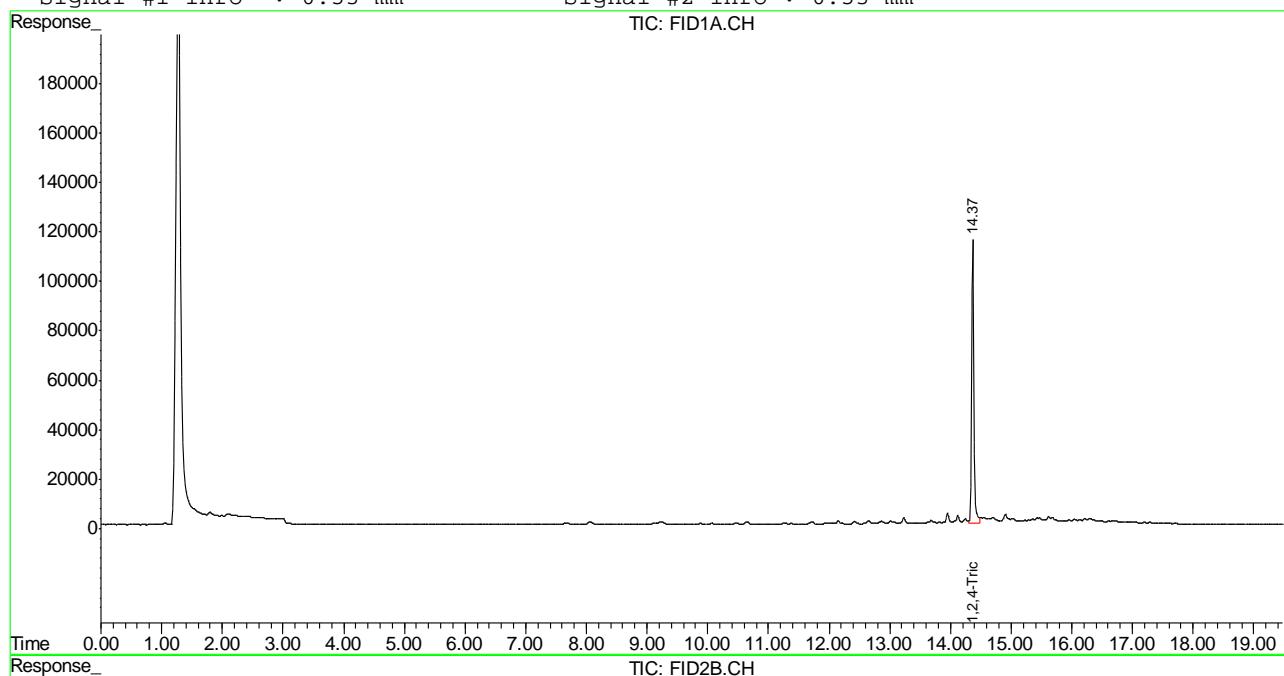
 (f)=RT Delta > 1/2 Window (m)=manual int.
 GB19112.D TB868GB868SOIL.M Fri Jan 11 09:08:59 2013 GC

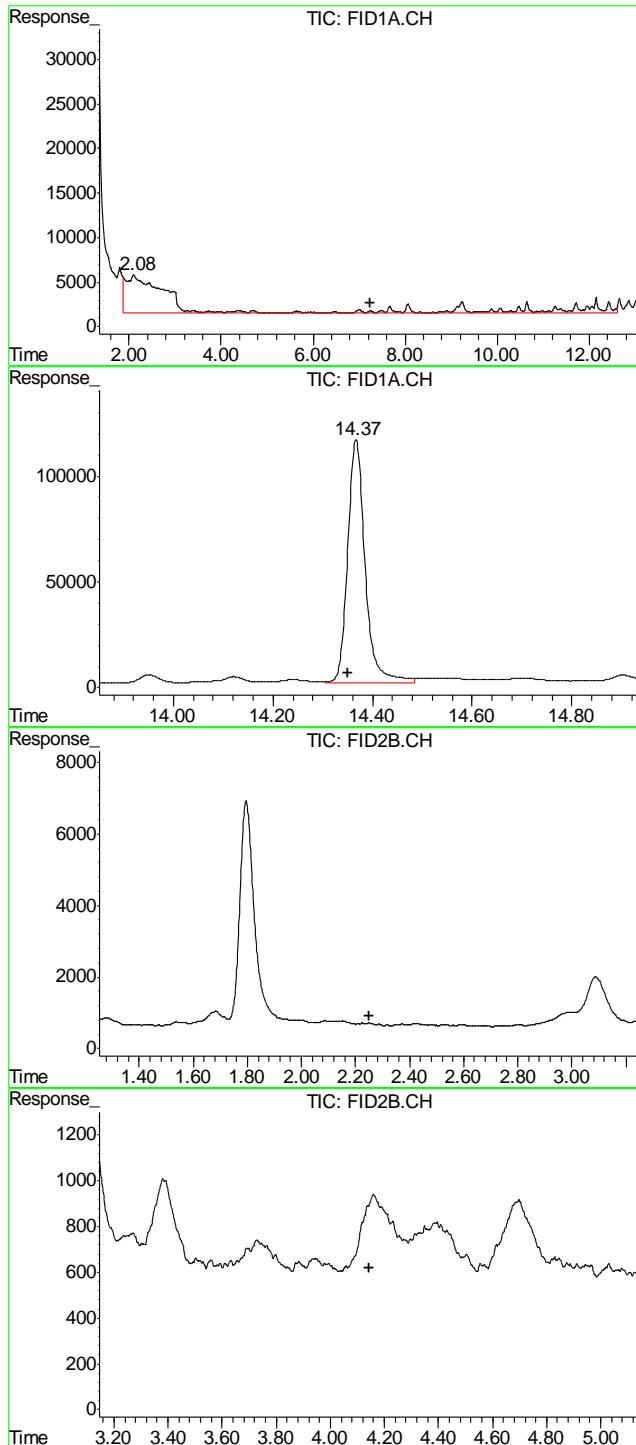
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011013\GB19112.D\FID1A.CH Vial: 27
 Signal #2 : Y:\1\DATA\011013\GB19112.D\FID2B.CH
 Acq On : 11 Jan 2013 6:12 am Operator: StephK
 Sample : D42511-1, 50X Inst : GC/MS Ins
 Misc : GC3347,GGB1044,,5.034,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Jan 11 9:06 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Jan 10 16:31:50 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.: 7.230 min
 Delta R.T.: 0.000 min
 Response: 3605558
 Conc: N.D.

#2 1,2,4-Trichlorobenzene

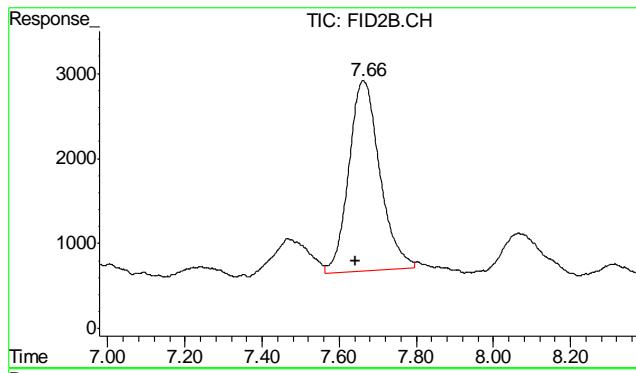
R.T.: 14.366 min
 Delta R.T.: 0.016 min
 Response: 2835281
 Conc: 90.49 % m

#4 Methyl-t-butyl-ether

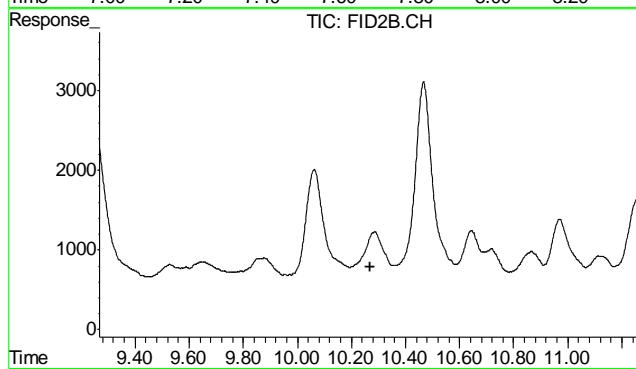
R.T.: 0.000 min
 Exp R.T. : 2.252 min
 Response: 0
 Conc: N.D.

#5 Benzene

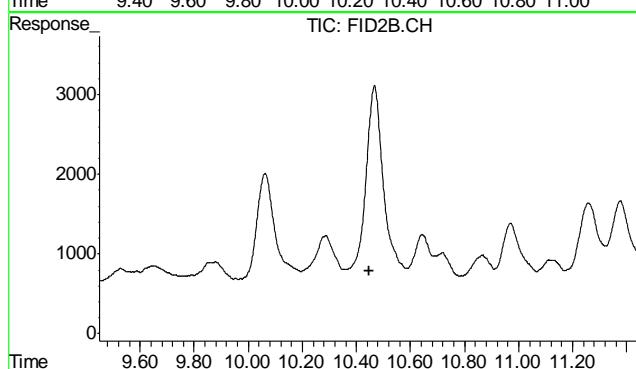
R.T.: 0.000 min
 Exp R.T. : 4.145 min
 Response: 0
 Conc: N.D.



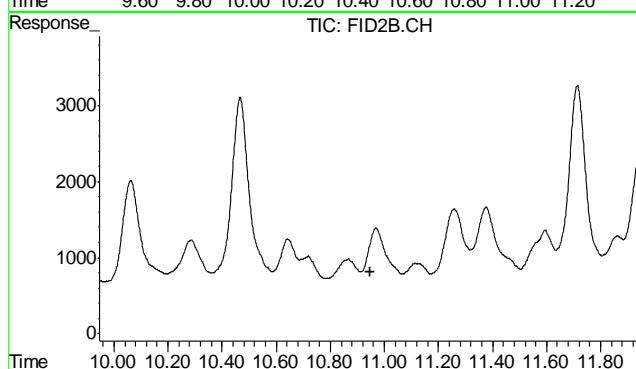
#6 Toluene
R.T.: 7.663 min
Delta R.T.: 0.019 min
Response: 124969
Conc: 0.32 ug/L



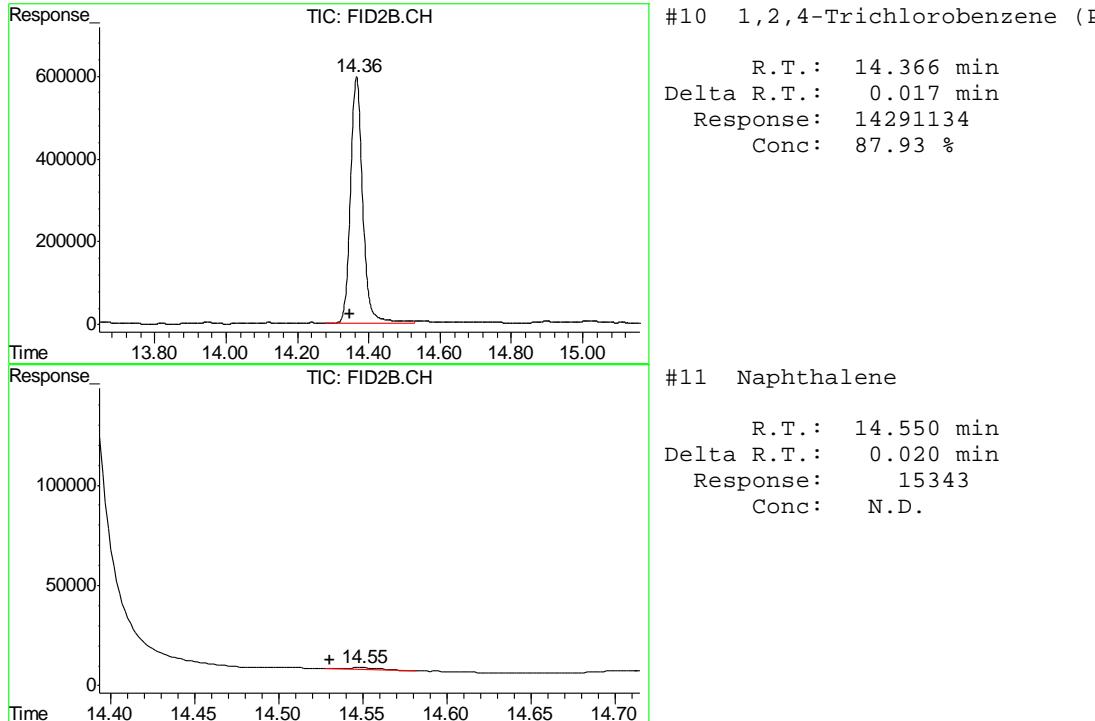
#7 Ethylbenzene
R.T.: 0.000 min
Exp R.T. : 10.268 min
Response: 0
Conc: N.D.



#8 m,p-Xylene
R.T.: 0.000 min
Exp R.T. : 10.449 min
Response: 0
Conc: N.D.



#9 o-Xylene
R.T.: 0.000 min
Exp R.T. : 10.946 min
Response: 0
Conc: N.D.



11.1.1

**Manual Integrations
APPROVED
(compounds with "m" flag)**
**Judy Nelson
01/11/13 11:38**

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011013\GB19088.D\FID1A.CH Vial: 3
 Signal #2 : Y:\1\DATA\011013\GB19088.D\FID2B.CH
 Acq On : 10 Jan 2013 4:00 pm Operator: StephK
 Sample : MB Inst : GC/MS Ins
 Misc : GC3347,GGB1044,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Jan 10 16:32:09 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Jan 10 16:31:50 2013
 Response via : Initial Calibration
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.35	2717831	86.737 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.35	13731194	84.485 %	

Target Compounds

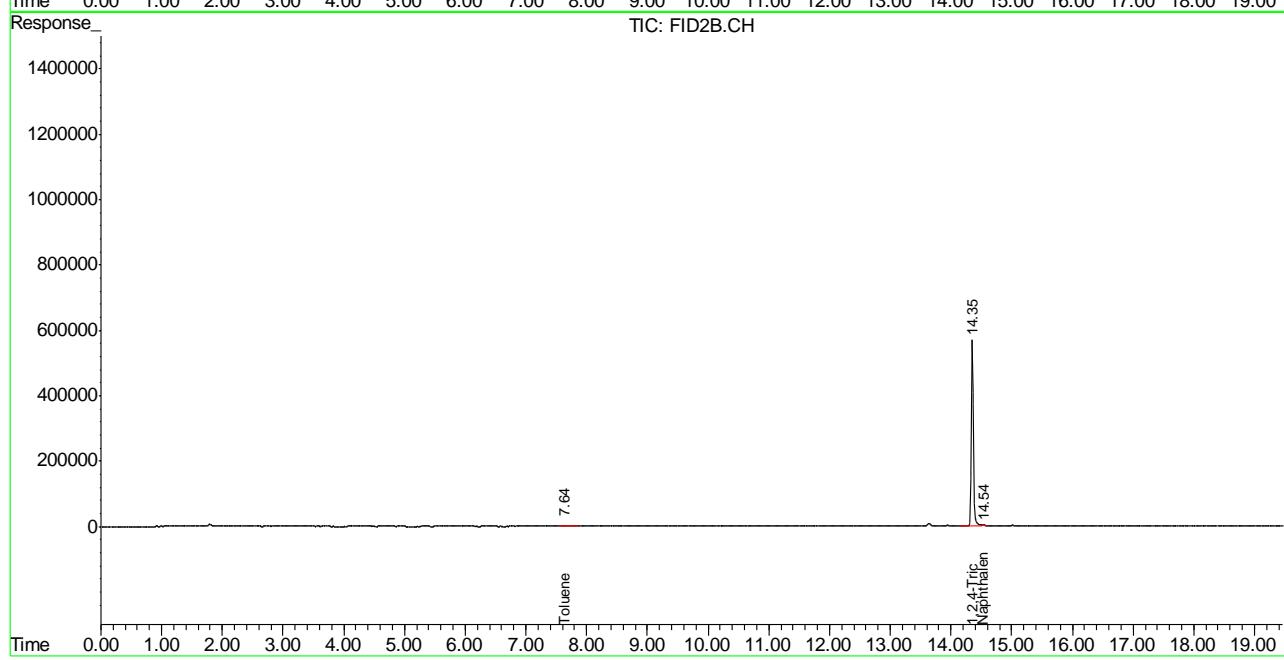
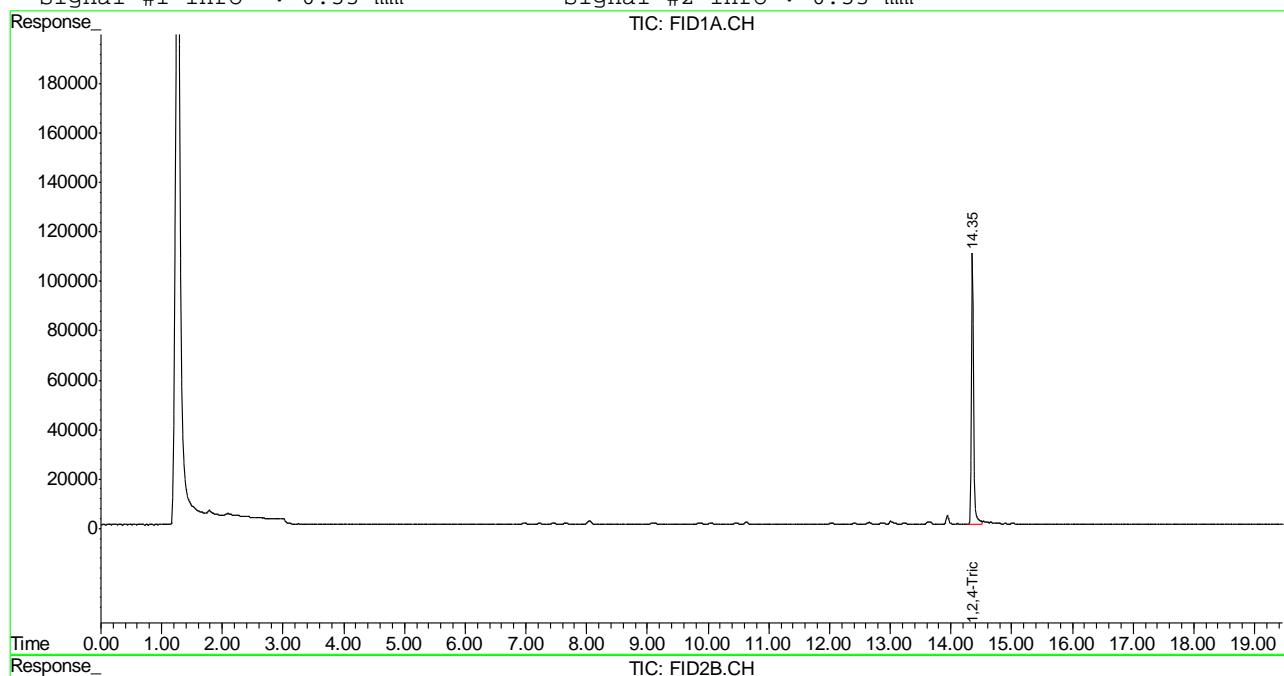
1) H	TVH-Gasoline	7.23	3461894	<MDL	mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D.	ug/L d
5) T	Benzene	0.00	0	N.D.	ug/L d
6) T	Toluene	7.64	126661	0.320	ug/L
7) T	Ethylbenzene	0.00	0	N.D.	ug/L d
8) T	m,p-Xylene	0.00	0	N.D.	ug/L d
9) T	o-Xylene	0.00	0	N.D.	ug/L d
11) T	Naphthalene	14.54	24002	0.122	ug/L m

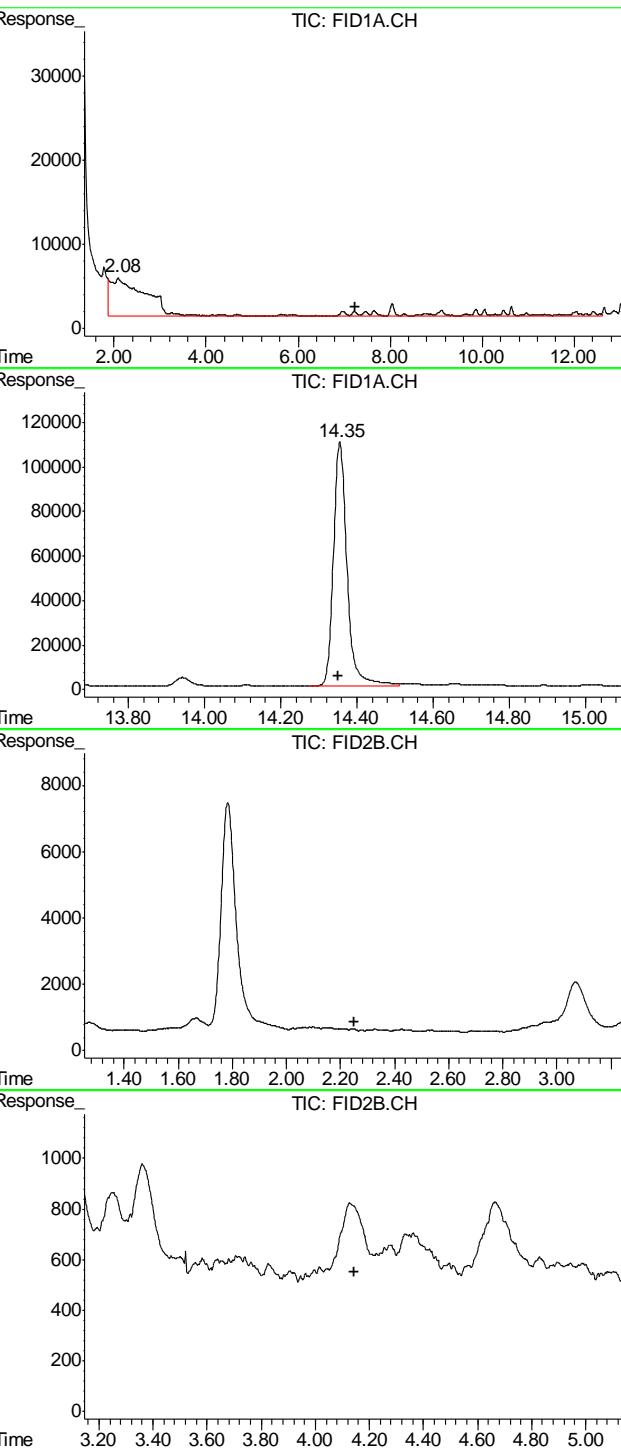
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011013\GB19088.D\FID1A.CH Vial: 3
 Signal #2 : Y:\1\DATA\011013\GB19088.D\FID2B.CH
 Acq On : 10 Jan 2013 4:00 pm Operator: StephK
 Sample : MB Inst : GC/MS Ins
 Misc : GC3347,GGB1044,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Jan 10 16:31 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Thu Jan 10 16:31:50 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.: 7.230 min
 Delta R.T.: 0.000 min
 Response: 3461894
 Conc: N.D.

#2 1,2,4-Trichlorobenzene

R.T.: 14.355 min
 Delta R.T.: 0.004 min
 Response: 2717831
 Conc: 86.74 % m

#4 Methyl-t-butyl-ether

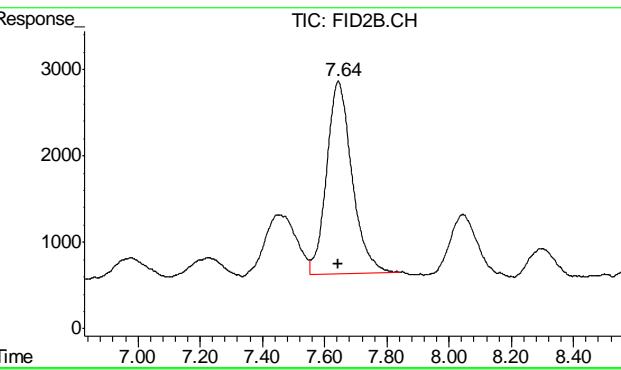
R.T.: 0.000 min
 Exp R.T. : 2.252 min
 Response: 0
 Conc: N.D.

#5 Benzene

R.T.: 0.000 min
 Exp R.T. : 4.145 min
 Response: 0
 Conc: N.D.

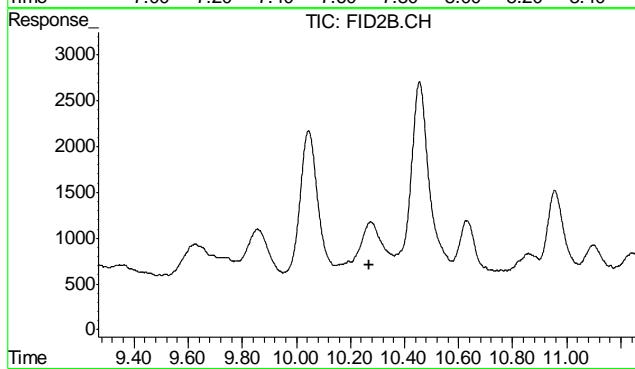
11.2.1

11



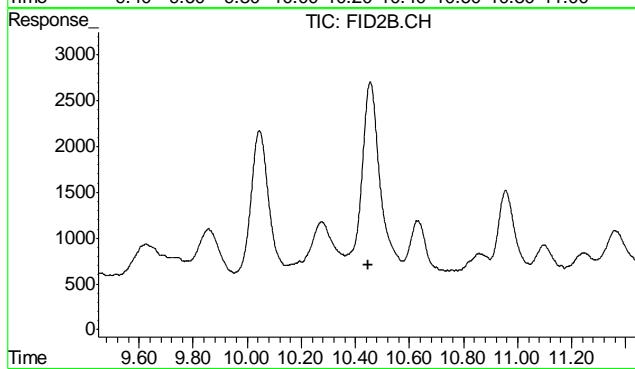
#6 Toluene

R.T.: 7.644 min
Delta R.T.: 0.000 min
Response: 126661
Conc: 0.32 ug/L



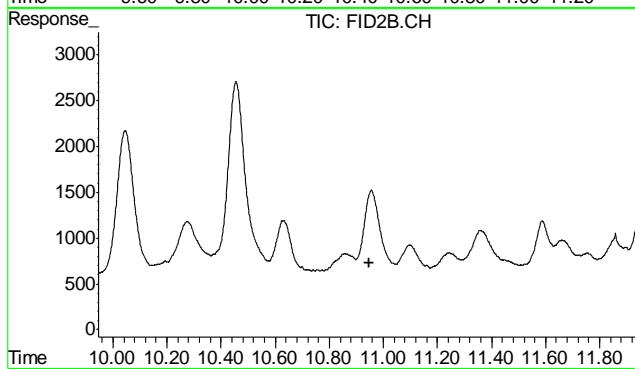
#7 Ethylbenzene

R.T.: 0.000 min
Exp R.T. : 10.268 min
Response: 0
Conc: N.D.



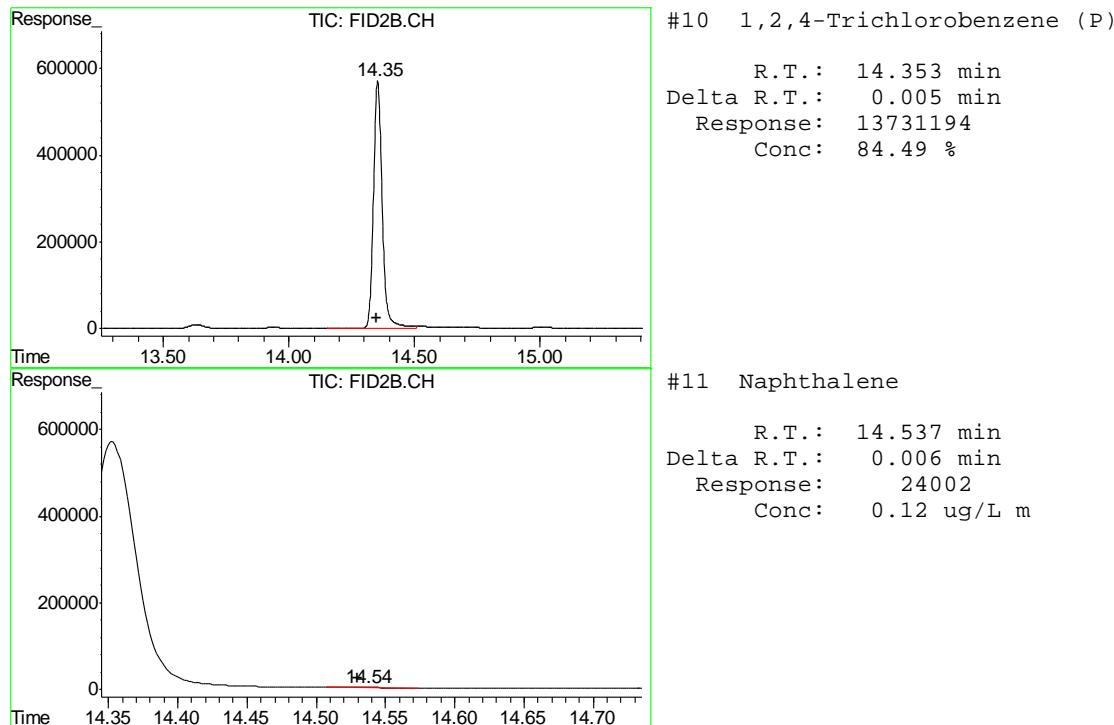
#8 m,p-Xylene

R.T.: 0.000 min
Exp R.T. : 10.449 min
Response: 0
Conc: N.D.



#9 o-Xylene

R.T.: 0.000 min
Exp R.T. : 10.946 min
Response: 0
Conc: N.D.



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

Job Number: D42511
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7222-MB	FH008528.D	1	01/14/13	AV	01/14/13	OP7222	GFH472

The QC reported here applies to the following samples:

Method: SW846-8015B

D42511-1

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	ND	6.7	4.0	mg/kg	

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	91% 35-130%

Blank Spike Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7222-BS	FH008532.D	1	01/14/13	AV	01/14/13	OP7222	GFH472

The QC reported here applies to the following samples:

Method: SW846-8015B

D42511-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	534	80	48-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	88%	35-130%

* = Outside of Control Limits.

12.2.1
12

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42511

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7222-MS	FH008534.D 1		01/14/13	AV	01/14/13	OP7222	GFH472
OP7222-MSD	FH008536.D 1		01/14/13	AV	01/14/13	OP7222	GFH472
D42509-2	FH008572.D 1		01/15/13	AV	01/14/13	OP7222	GFH472

The QC reported here applies to the following samples:

Method: SW846-8015B

D42511-1

CAS No.	Compound	D42509-2		Spike mg/kg	MS mg/kg	MS %	MSD mg/kg	MSD %	RPD	Limits Rec/RPD
		mg/kg	Q							
	TPH-DRO (C10-C28)	46.1		736	647	82	493	61	27	20-168/30
CAS No.	Surrogate Recoveries	MS	MSD	D42509-2		Limits				
84-15-1	o-Terphenyl	79%	62%	74%		35-130%				

* = Outside of Control Limits.

12.3.1
12



GC Semi-volatiles

Raw Data

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH011413.SEC\
 Data File : FH008578.D
 Signal(s) : FID2B.ch
 Acq On : 15 Jan 2013 6:01 am
 Operator : ashleyv
 Sample : D42511-1
 Misc : OP7222,GFH472,30.02,,,1,1
 ALS Vial : 77 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Jan 15 11:54:29 2013
 Quant Method : C:\msdchem\1\METHODS\DRD-GFH464R.M
 Quant Title : DRO-ORO REAR
 QLast Update : Mon Jan 07 08:59:40 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
3) s o-Terphenyl	12.543	1853045135	1347.455	ug/ml
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	10.243	1912608191	1620.074	ug/ml
2) H TPH-DRO (C8-C20)	7.898	936279254	842.351	ug/ml
<hr/>				

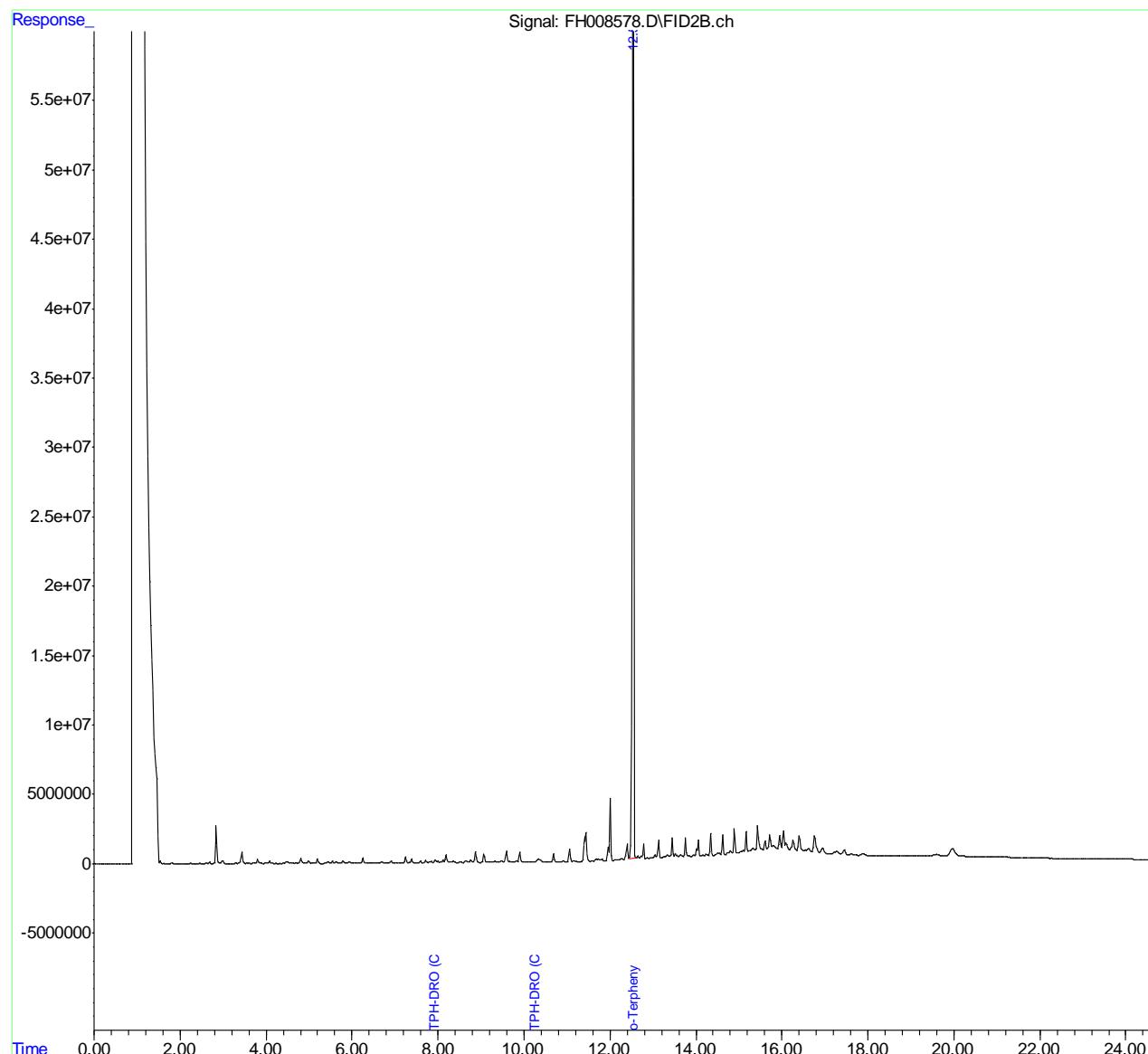
(f)=RT Delta > 1/2 Window (m)=manual int.

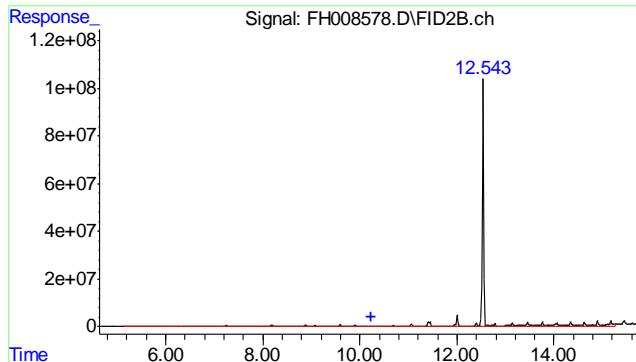
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH011413.SEC\
 Data File : FH008578.D
 Signal(s) : FID2B.ch
 Acq On : 15 Jan 2013 6:01 am
 Operator : ashleyv
 Sample : D42511-1
 Misc : OP7222,GFH472,30.02,,,1,1
 ALS Vial : 77 Sample Multiplier: 1

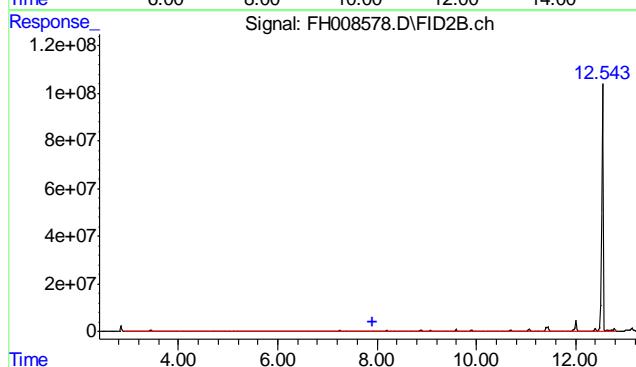
Integration File: events.e
 Quant Time: Jan 15 11:54:29 2013
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH464R.M
 Quant Title : DRO-ORO REAR
 QLast Update : Mon Jan 07 08:59:40 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

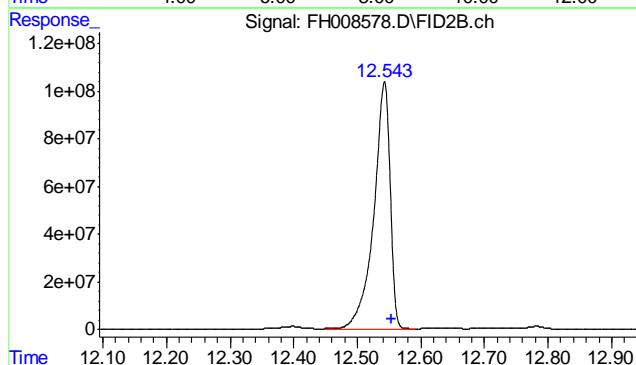




#1 TPH-DRO (C10-C28)
 R.T.: 10.243 min
 Delta R.T.: 0.000 min
 Response: 1912608191
 Conc: 1620.07 ug/ml m



#2 TPH-DRO (C8-C20)
 R.T.: 7.898 min
 Delta R.T.: 0.000 min
 Response: 936279254
 Conc: 842.35 ug/ml m



#3 o-Terphenyl
 R.T.: 12.543 min
 Delta R.T.: -0.010 min
 Response: 1853045135
 Conc: 1347.46 ug/ml

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH011413.SEC\
 Data File : FH008528.D
 Signal(s) : FID2B.ch
 Acq On : 14 Jan 2013 3:50 pm
 Operator : ashleyv
 Sample : OP7222-MB
 Misc : OP7222,GFH472,30.00,,,1,1
 ALS Vial : 53 Sample Multiplier: 1

Integration File: events.e
 Quant Time: Jan 15 08:27:53 2013
 Quant Method : C:\msdchem\1\METHODS\DRD-GFH464R.M
 Quant Title : DRO-ORO REAR
 QLast Update : Mon Jan 07 08:59:40 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
3) s o-Terphenyl	12.557	2504718000	1821.324	ug/ml
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	10.243	70088032	59.368	ug/ml
2) H TPH-DRO (C8-C20)	7.898	20531892	18.472	ug/ml
<hr/>				

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

13.2.1

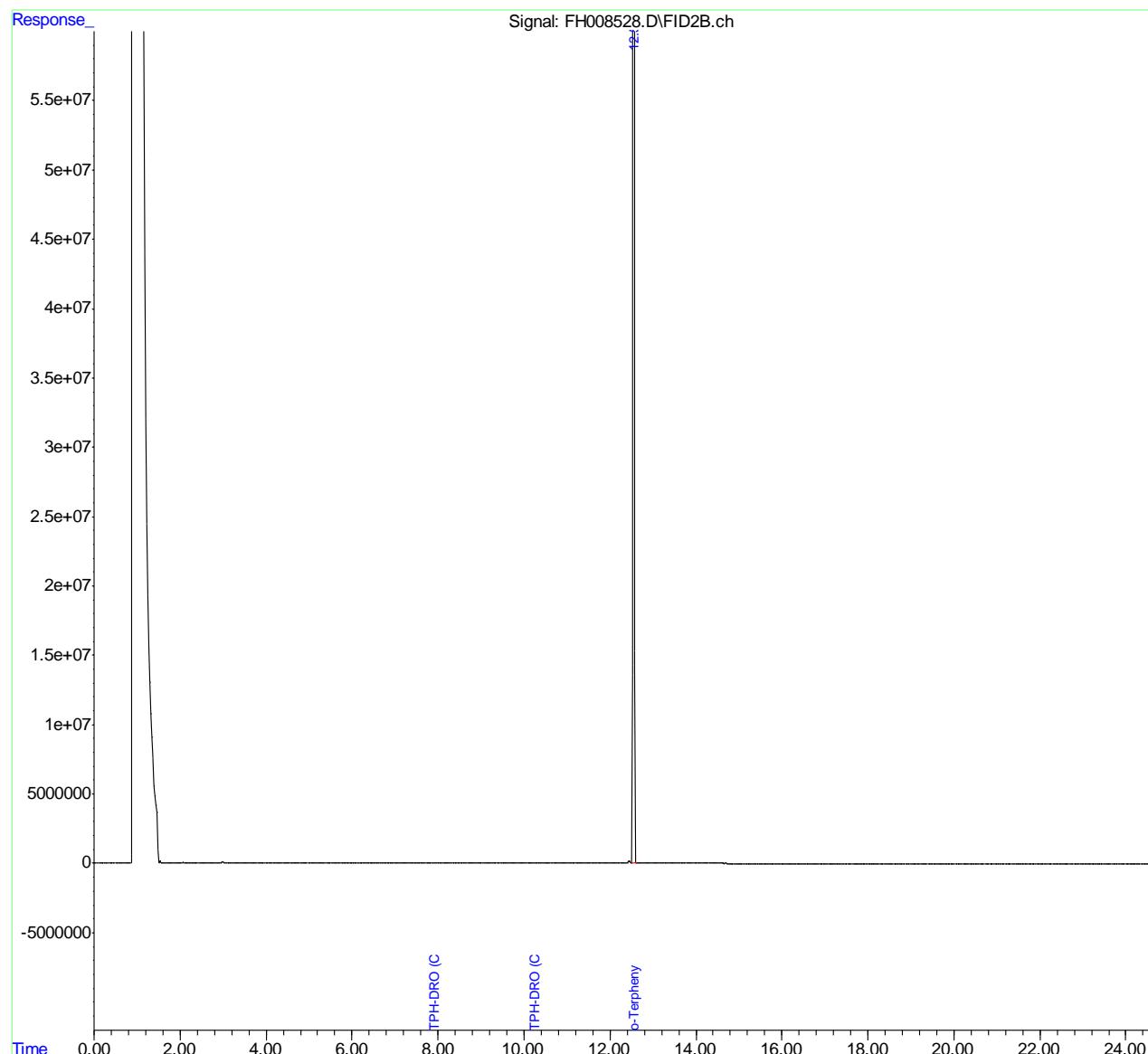
13

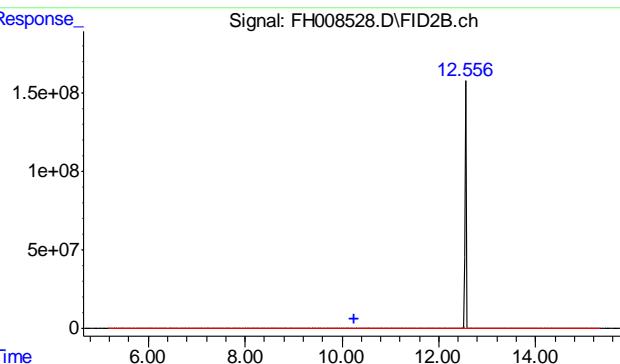
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH011413.SEC\
 Data File : FH008528.D
 Signal(s) : FID2B.ch
 Acq On : 14 Jan 2013 3:50 pm
 Operator : ashleyv
 Sample : OP7222-MB
 Misc : OP7222,GFH472,30.00,,,1,1
 ALS Vial : 53 Sample Multiplier: 1

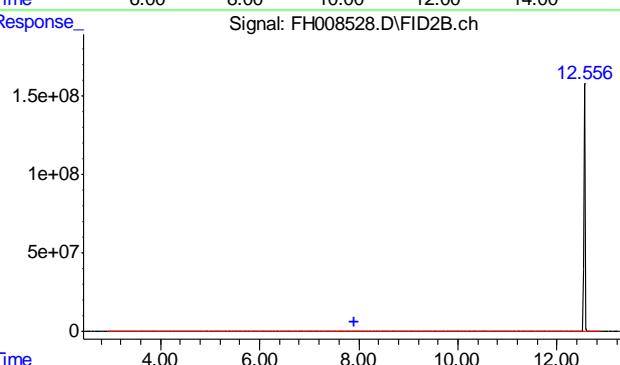
Integration File: events.e
 Quant Time: Jan 15 08:27:53 2013
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH464R.M
 Quant Title : DRO-ORO REAR
 QLast Update : Mon Jan 07 08:59:40 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

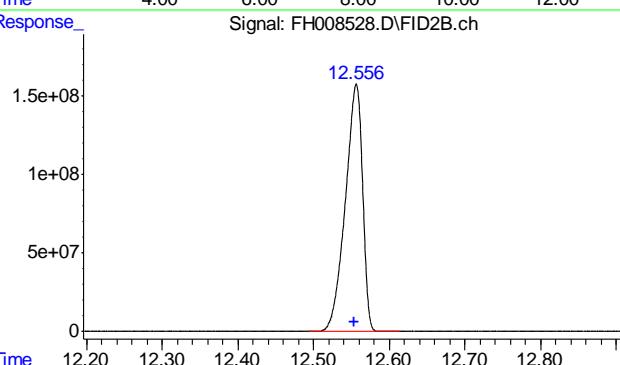




#1 TPH-DRO (C10-C28)
R.T.: 10.243 min
Delta R.T.: 0.000 min
Response: 70088032
Conc: 59.37 ug/ml m



#2 TPH-DRO (C8-C20)
R.T.: 7.898 min
Delta R.T.: 0.000 min
Response: 20531892
Conc: 18.47 ug/ml m



#3 o-Terphenyl
R.T.: 12.557 min
Delta R.T.: 0.004 min
Response: 2504718000
Conc: 1821.32 ug/ml



Metals Analysis

QC Data Summaries

Includes the following where applicable:

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9237
Matrix Type: AQUEOUS

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

Prep Date:

01/11/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	48	130		
Antimony	150	8.5	18		
Arsenic	130	22	42		
Barium	50	.5	9		
Beryllium	50	6.5	16		
Boron	250	5	22		
Cadmium	50	3	3		
Calcium	2000	27	80	-27	<2000
Chromium	50	1.5	2.8		
Cobalt	25	2	2.1		
Copper	50	6	15		
Iron	350	6	100		
Lead	250	9.5	15		
Lithium	10	2.5			
Magnesium	1000	33	110	-38	<1000
Manganese	25	6	6		
Molybdenum	50	11	11		
Nickel	150	2.5	2.9		
Phosphorus	500	70	300		
Potassium	5000	310	750		
Selenium	250	24	55		
Silicon	250	15			
Silver	150	2	4.9		
Sodium	2000	30	490	-180	<2000
Strontium	25	.2	7.5		
Thallium	50	15	43		
Tin	250	60			
Titanium	50	.5			
Uranium	250	11	23		
Vanadium	50	1	2.4		
Zinc	150	2.5	12		

Associated samples MP9237: D42511-1A

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9237
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: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9237
 Matrix Type: AQUEOUS

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

Prep Date:

01/11/13

Metal	D42427-2 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	241000	164000	125000	-61.6N(a) 75-125
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	1310	130000	125000	103.0 75-125
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	266000	295000	125000	23.2N(a) 75-125
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9237: D42511-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9237
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
(a) Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9237
 Matrix Type: AQUEOUS

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

Prep Date: 01/11/13

Metal	D42427-2 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic						
Barium						
Beryllium						
Boron						
Cadmium						
Calcium	241000	166000	125000	-60.0N(a	1.2	20
Chromium						
Cobalt						
Copper						
Iron						
Lead						
Lithium						
Magnesium	1310	129000	125000	102.2	0.8	20
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium						
Silicon						
Silver						
Sodium	266000	292000	125000	20.8N(a)	1.0	20
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP9237: D42511-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9237
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
(a) Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42511
 Account: XTOKWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9237
 Matrix Type: AQUEOUS

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

Prep Date: 01/11/13

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	126000	125000	100.8	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9237: D42511-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9237
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: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9237
 Matrix Type: AQUEOUS

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

Prep Date: 01/11/13

Metal	D42427-2	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	48100	6440	86.6*(a)	0-10	
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	262	1240	375.2(b)	0-10	
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	53100	34700	34.7*(a)	0-10	
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP9237: D42511-1A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9237
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60
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: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9242
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

01/14/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	.96	.57		
Antimony	3.0	.17	.12		
Arsenic	2.5	.44	.56		
Barium	1.0	.01	.11	0.25	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	0.020	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.040	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.13	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	0.020	<5.0
Lithium	0.20	.05	.054		
Magnesium	20	.65	.98		
Manganese	0.50	.12	.022		
Molybdenum	1.0	.21	.08		
Nickel	3.0	.05	.026	-0.020	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	-0.46	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	0.030	<3.0
Sodium	40	.59	1.9		
Strontium	5.0	.004	.017		
Thallium	1.0	.29	.53		
Tin	5.0	1.2	2		
Titanium	1.0	.01	.038		
Uranium	5.0	.22	.26		
Vanadium	1.0	.02	.036		
Zinc	3.0	.05	.37	0.090	<3.0

Associated samples MP9242: D42511-1

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9242
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9242
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 01/14/13

Metal	D42510-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	4420	4470	235	21.3 (a) 75-125
Beryllium				
Boron				
Cadmium	0.0	49.1	58.7	83.7 75-125
Calcium				
Chromium	28.6	73.0	58.7	83.2 75-125
Cobalt				
Copper	15.4	67.2	58.7	88.3 75-125
Iron				
Lead	12.0	107	117	81.0 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	19.2	77.2	58.7	98.9 75-125
Phosphorus				
Potassium				
Selenium	0.0	105	117	89.5 75-125
Silicon				
Silver	0.0	21.0	23.5	89.5 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	37.6	82.7	58.7	76.9 75-125

Associated samples MP9242: D42511-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9242
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9242
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

01/14/13

Metal	D42510-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic						
Barium	4420	4930	232	219.5(a)	9.8	20
Beryllium						
Boron						
Cadmium	0.0	48.6	58.1	83.7	1.0	20
Calcium						
Chromium	28.6	74.5	58.1	86.6	2.0	20
Cobalt						
Copper	15.4	67.2	58.1	89.2	0.0	20
Iron						
Lead	12.0	105	116	80.1	1.9	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	19.2	77.0	58.1	99.5	0.3	20
Phosphorus						
Potassium						
Selenium	0.0	104	116	89.5	1.0	20
Silicon						
Silver	0.0	20.8	23.2	89.5	1.0	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	37.6	81.9	58.1	76.3	1.0	20

Associated samples MP9242: D42511-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9242
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9242
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 01/14/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	198	200	99.0	80-120
Beryllium				
Boron				
Cadmium	46.5	50	93.0	80-120
Calcium				
Chromium	49.9	50	99.8	80-120
Cobalt				
Copper	47.4	50	94.8	80-120
Iron				
Lead	96.0	100	96.0	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	47.5	50	95.0	80-120
Phosphorus				
Potassium				
Selenium	97.5	100	97.5	80-120
Silicon				
Silver	19.8	20	99.0	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	49.2	50	98.4	80-120

Associated samples MP9242: D42511-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9242
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9242
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: ug/l

Prep Date: 01/14/13

Metal	D42510-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	32800	36700	2.5	0-10
Beryllium				
Boron				
Cadmium	0.00	0.00	NC	0-10
Calcium				
Chromium	206	243	17.8*(a)	0-10
Cobalt				
Copper	121	142	8.3	0-10
Iron				
Lead	119	122	19.0*(a)	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	135	155	5.2	0-10
Phosphorus				
Potassium				
Selenium	0.00	0.00	NC	0-10
Silicon				
Silver	0.00	0.00	NC	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	320	412	28.5*(a)	0-10

Associated samples MP9242: D42511-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9242
Matrix Type: SOLID

Methods: SW846 6010C
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested
(a) Serial dilution indicates possible matrix interference.

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D42511
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9243
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date:

01/14/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	0.013	<0.10
Barium	1.0	.0065	.037		
Beryllium	0.10	.016	.09		
Boron	20	1.2	1.2		
Cadmium	0.050	.014	.021		
Calcium	200	7.9	8		
Chromium	1.0	.033	.19		
Cobalt	0.10	.0012	.015		
Copper	1.0	.017	.065		
Iron	20	.8	5		
Lead	0.25	.0011	.024		
Magnesium	50	.44	.85		
Manganese	0.50	.0043	.02		
Molybdenum	0.50	.018	.018		
Nickel	1.0	.0049	.011		
Phosphorus	30	1.4	3.6		
Potassium	100	9.8	10		
Selenium	0.20	.029	.14		
Silver	0.050	.0009	.0065		
Sodium	250	1.5	2.3		
Strontium	10	.036	.036		
Thallium	0.10	.00095	.0095		
Tin	5.0	.023	.34		
Titanium	1.0	.044	.1		
Uranium	0.25	.00085	.001		
Vanadium	2.0	.12	.21		
Zinc	5.0	.033	.35		

Associated samples MP9243: D42511-1

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: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9243
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 01/14/13

Metal	D42510-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	11.1	116	117	89.4 75-125
Barium	anr			
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium	anr			
Silver	anr			
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9243: D42511-1

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: D42511
 Account: XTOKWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9243
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 01/14/13

Metal	D42510-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	11.1	123	116	96.3	5.9	20
Barium						
Beryllium						
Boron						
Cadmium		anr				
Calcium						
Chromium		anr				
Cobalt						
Copper						
Iron						
Lead		anr				
Magnesium						
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium		anr				
Silver		anr				
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP9243: D42511-1

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: D42511
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9243
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 01/14/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	101	100	101.0	80-120
Barium	anr			
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium	anr			
Silver	anr			
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9243: D42511-1

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

14.3.3
14

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42511
 Account: XTOKWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9243
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: ug/l

Prep Date: 01/14/13

Metal	D42510-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	94.9	94.9	0.0	0-10
Barium	anr			
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium	anr			
Silver	anr			
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9243: D42511-1

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

14.3.4
14

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9244
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 01/15/13

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.0009	0.0018	<0.10

Associated samples MP9244: D42511-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42511
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9244
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 01/15/13

Metal	D42445-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.054	0.45	0.393	100.7 75-125

Associated samples MP9244: D42511-1

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: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9244
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

01/15/13

Metal	D42445-1 Original	MSD HGWSR1	Spikelot % Rec	MSD RPD	QC Limit
Mercury	0.054	0.45	0.393	100.7	0.0 20

Associated samples MP9244: D42511-1

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: D42511
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9244
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 01/15/13

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	Limits
Mercury	0.40	0.4	100.0	80-120

Associated samples MP9244: D42511-1

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

14.4.3
14



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: D42511
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP9086/GN18425	1.0	0.0	mg/kg	92.9	91.5	98.5	80-120%
Specific Conductivity	GP9098/GN18435	1.0	<1.0	umhos/cm	9992	10500	105.2	90-110%
pH	GN18407			su	8.00	7.99	99.9	99.3-100.7%

Associated Samples:

Batch GP9086: D42511-1

Batch GP9098: D42511-1

Batch GN18407: D42511-1

(*) Outside of QC limits

DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D42511
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP9086/GN18425 GN18410	D42556-1 D42511-1	mg/kg mv	0.0 137	0.0 137	0.0 0.0	0-20% 0-20%

Associated Samples:
Batch GP9086: D42511-1
Batch GN18410: D42511-1
(*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D42511
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP9086/GN18425	D42556-1	mg/kg	0.0	40.0	35.0	87.5	75-125%

Associated Samples:

Batch GP9086: D42511-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D42511
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP9086/GN18425	D42556-1	mg/kg	0.0	40.0	33.9	3.1	20%

Associated Samples:

Batch GP9086: D42511-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits