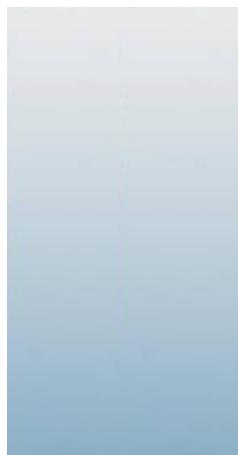




09/14/12



Technical Report for

XTO Energy

PCU 197-36A

1203-02

Accutest Job Number: D38480

Sampling Date: 09/05/12

Report to:

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Total number of pages in report: 222



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)

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

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Sample Summary

XTO Energy

Job No: D38480

PCU 197-36A

Project No: 1203-02

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D38480-1	09/05/12	10:10 DS	09/07/12	SO	Soil	CUT 1 POST SOLIDIFICATION
D38480-1A	09/05/12	10:10 DS	09/07/12	SO	Soil	CUT 1 POST SOLIDIFICATION
D38480-2	09/05/12	10:00 DS	09/07/12	SO	Soil	CUT 2 POST SOLIDIFICATION
D38480-2A	09/05/12	10:00 DS	09/07/12	SO	Soil	CUT 2 POST SOLIDIFICATION
D38480-3	09/05/12	09:50 DS	09/07/12	SO	Soil	CUT 3 POST SOLIDIFICATION
D38480-3A	09/05/12	09:50 DS	09/07/12	SO	Soil	CUT 3 POST SOLIDIFICATION

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



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: XTO Energy

Job No D38480

Site: PCU 197-36A

Report Date 9/14/2012 4:49:18 PM

On 09/07/2012, 3 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 4.5 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D38480 was assigned to the project. The lab sample IDs, client sample IDs, 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: V3V1182
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D38454-1MS, D38454-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

Extractables by GCMS By Method SW846 8270C BY SIM

Matrix SO	Batch ID: OP6602
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D38513-1MS, D38513-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

Volatiles by GC By Method SW846 8015B

Matrix SO	Batch ID: GGB958
------------------	-------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D38454-1MS, D38454-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

Extractables by GC By Method SW846-8015B

Matrix SO	Batch ID: OP6603
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D38483-1MS, D38483-1MSD were used as the QC samples indicated.

Metals By Method SW846 6010C

Matrix AQ

Batch ID: MP8372

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D38518-1AMS, D38518-1AMSD, D38518-1ASDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Sodium 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 Magnesium are outside control limits for sample MP8372-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).

Matrix SO

Batch ID: MP8358

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D38480-1MS, D38480-1MSD, D38480-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Copper are outside control limits. Spike recovery indicates possible matrix interference.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The RPD(s) for the MS and MSD recoveries of Barium are outside control limits for sample MP8358-S2. High RPD due to possible sample matrix or nonhomogeneity.
- The serial dilution RPD(s) for Cadmium, Selenium, Silver, Copper are outside control limits for sample MP8358-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP8358-SD1 for Copper: Serial dilution indicates possible matrix interference.

Metals By Method SW846 6020A

Matrix SO

Batch ID: MP8359

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D38480-1MSD, D38480-1SDL, D38480-1MS were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Arsenic are outside control limits. Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.

Metals By Method SW846 7471B

Matrix SO

Batch ID: MP8357

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D38479-1MS, D38479-1MSD were used as the QC samples for the metals analysis.

Wet Chemistry By Method ASTM D1498-76M

Matrix SO

Batch ID: GN16683

- Sample(s) D38518-2DUP were used as the QC samples for the Redox Potential Vs H₂ analysis.

Wet Chemistry By Method SM19 2540B M

Matrix SO

Batch ID: GN16673

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

Wet Chemistry By Method SM2510B-1997 MOD

Matrix SO

Batch ID: GP8183

- 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 SW846 3060/7196A M

Matrix SO

Batch ID: R14352

- The data for SW846 3060/7196A M meets quality control requirements.
- D38480-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Matrix SO

Batch ID: R14353

- The data for SW846 3060/7196A M meets quality control requirements.
- D38480-2 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Matrix SO

Batch ID: R14354

- The data for SW846 3060/7196A M meets quality control requirements.
- D38480-3 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Wet Chemistry By Method SW846 3060A/7196A

Matrix SO

Batch ID: GP8138

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D38513-1MS, D38513-1MSD, D38513-1DUP were used as the QC samples for the Chromium, Hexavalent analysis.
- The duplicate RPD(s) for Chromium, Hexavalent are outside control limits for sample GP8138-D1. RPD acceptable due to low duplicate and sample concentrations.

Wet Chemistry By Method SW846 9045D

Matrix SO

Batch ID: GN16667

- The following samples were run outside of holding time for method SW846 9045D: D38480-1, D38480-2, D38480-3

Wet Chemistry By Method USDA HANDBOOK 60

Matrix SO

Batch ID: MP8372

- All samples 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 3

Job Number: D38480
Account: XTO Energy
Project: PCU 197-36A
Collected: 09/05/12

3

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

D38480-1 CUT 1 POST SOLIDIFICATION

Toluene	0.490	0.15	0.075	mg/kg	SW846 8260B
Ethylbenzene	0.131 J	0.15	0.029	mg/kg	SW846 8260B
Xylene (total)	0.567	0.30	0.15	mg/kg	SW846 8260B
Naphthalene	0.170	0.015	0.013	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	68.0	15	7.5	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	842	17	11	mg/kg	SW846-8015B
Arsenic	16.8	0.13		mg/kg	SW846 6020A
Barium	1270	1.3		mg/kg	SW846 6010C
Chromium	28.9	1.3		mg/kg	SW846 6010C
Copper	28.8	1.3		mg/kg	SW846 6010C
Lead	15.0	6.3		mg/kg	SW846 6010C
Nickel	22.4	3.8		mg/kg	SW846 6010C
Zinc	51.8	3.8		mg/kg	SW846 6010C
Specific Conductivity	11400	1.0		umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent ^a	28.9	2.3		mg/kg	SW846 3060/7196A M
Redox Potential Vs H2	146			mv	ASTM D1498-76M
pH	12.43			su	SW846 9045D

D38480-1A CUT 1 POST SOLIDIFICATION

Calcium	15.9	2.0	mg/l	SW846 6010C
Sodium	1360	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	93.8		ratio	USDA HANDBOOK 60

D38480-2 CUT 2 POST SOLIDIFICATION

Benzene	0.109	0.082	0.041	mg/kg	SW846 8260B
Toluene	0.240	0.16	0.082	mg/kg	SW846 8260B
Ethylbenzene	0.0507 J	0.16	0.031	mg/kg	SW846 8260B
Xylene (total)	0.270 J	0.33	0.16	mg/kg	SW846 8260B
Benzo(a)anthracene	0.0258	0.011	0.0057	mg/kg	SW846 8270C BY SIM
Benzo(a)pyrene	0.0275	0.011	0.0057	mg/kg	SW846 8270C BY SIM
Chrysene	0.0850	0.011	0.0057	mg/kg	SW846 8270C BY SIM
Fluoranthene	0.0377	0.011	0.0057	mg/kg	SW846 8270C BY SIM
Fluorene	0.131	0.011	0.0057	mg/kg	SW846 8270C BY SIM
Indeno(1,2,3-cd)pyrene	0.0110	0.011	0.0057	mg/kg	SW846 8270C BY SIM
Naphthalene	0.838	0.015	0.014	mg/kg	SW846 8270C BY SIM
Pyrene	0.0729	0.011	0.0057	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	27.3	16	8.2	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	253	18	11	mg/kg	SW846-8015B
Arsenic	11.0	0.13		mg/kg	SW846 6020A
Barium	2590	1.3		mg/kg	SW846 6010C
Chromium	17.3	1.3		mg/kg	SW846 6010C

Summary of Hits

Page 2 of 3

Job Number: D38480
Account: XTO Energy
Project: PCU 197-36A
Collected: 09/05/12

3

Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
Copper		163	1.3		mg/kg	SW846 6010C
Lead		33.3	6.4		mg/kg	SW846 6010C
Nickel		29.3	3.9		mg/kg	SW846 6010C
Zinc		100	3.9		mg/kg	SW846 6010C
Specific Conductivity		11900	1.0		umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent ^a		17.3	2.3		mg/kg	SW846 3060/7196A M
Redox Potential Vs H2		160			mv	ASTM D1498-76M
pH		11.74			su	SW846 9045D

D38480-2A CUT 2 POST SOLIDIFICATION

Calcium	86.6	2.0		mg/l	SW846 6010C
Sodium	2270	2.0		mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	67.0			ratio	USDA HANDBOOK 60

D38480-3 CUT 3 POST SOLIDIFICATION

Benzene	0.0957	0.072	0.036	mg/kg	SW846 8260B
Toluene	0.267	0.14	0.072	mg/kg	SW846 8260B
Ethylbenzene	0.0273 J	0.14	0.027	mg/kg	SW846 8260B
Xylene (total)	0.327	0.29	0.14	mg/kg	SW846 8260B
Benzo(a)anthracene	0.0310	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Benzo(a)pyrene	0.0548	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Benzo(b)fluoranthene	0.0440	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Benzo(k)fluoranthene	0.0098 J	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Chrysene	0.108	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Dibenzo(a,h)anthracene	0.0141	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Fluoranthene	0.0342	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Fluorene	0.139	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Indeno(1,2,3-cd)pyrene	0.0139	0.010	0.0053	mg/kg	SW846 8270C BY SIM
Naphthalene	0.482	0.014	0.013	mg/kg	SW846 8270C BY SIM
Pyrene	0.0689	0.010	0.0053	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	44.4	14	7.2	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	434	16	11	mg/kg	SW846-8015B
Arsenic	9.1	0.12		mg/kg	SW846 6020A
Barium	5630	5.9		mg/kg	SW846 6010C
Chromium	16.6	1.2		mg/kg	SW846 6010C
Copper	28.0	1.2		mg/kg	SW846 6010C
Lead	28.8	5.9		mg/kg	SW846 6010C
Nickel	20.7	18		mg/kg	SW846 6010C
Zinc	56.2	18		mg/kg	SW846 6010C
Specific Conductivity	11400	1.0		umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent ^a	16.6	2.2		mg/kg	SW846 3060/7196A M
Redox Potential Vs H2	145			mv	ASTM D1498-76M
pH	12.45			su	SW846 9045D

Summary of Hits

Page 3 of 3

Job Number: D38480
Account: XTO Energy
Project: PCU 197-36A
Collected: 09/05/12

3

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

D38480-3A CUT 3 POST SOLIDIFICATION

Calcium	2.95	2.0	mg/l	SW846 6010C
Sodium	2100	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	336		ratio	USDA HANDBOOK 60

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

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



4

Sample Results

Report of Analysis

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 1 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-1	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	79.4
Method:	SW846 8260B		
Project:	PCU 197-36A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3V20315.D	1	09/08/12	BD	n/a	n/a	V3V1182
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.06 g	5.0 ml	100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	0.075	0.038	mg/kg	
108-88-3	Toluene	0.490	0.15	0.075	mg/kg	
100-41-4	Ethylbenzene	0.131	0.15	0.029	mg/kg	J
1330-20-7	Xylene (total)	0.567	0.30	0.15	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	95%		64-130%
460-00-4	4-Bromofluorobenzene	103%		62-131%
17060-07-0	1,2-Dichloroethane-D4	115%		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

Page 1 of 1

Client Sample ID:	CUT 1 POST SOLIDIFICATION			Date Sampled:	09/05/12
Lab Sample ID:	D38480-1			Date Received:	09/07/12
Matrix:	SO - Soil			Percent Solids:	79.4
Method:	SW846 8270C BY SIM SW846 3546				
Project:	PCU 197-36A				
File ID	DF	Analyzed	By	Prep Date	Prep Batch
Run #1 3G11215.D	1	09/13/12	DC	09/11/12	OP6602
Run #2					E3G522
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	0.0055	mg/kg	
120-12-7	Anthracene	ND	0.010	0.0055	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.010	0.0055	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.010	0.0055	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.010	0.0055	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.010	0.0055	mg/kg	
218-01-9	Chrysene	ND	0.010	0.0055	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.010	0.0055	mg/kg	
206-44-0	Fluoranthene	ND	0.010	0.0055	mg/kg	
86-73-7	Fluorene	ND	0.010	0.0055	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.010	0.0055	mg/kg	
91-20-3	Naphthalene	0.170	0.015	0.013	mg/kg	
129-00-0	Pyrene	ND	0.010	0.0055	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	64%		10-145%
321-60-8	2-Fluorobiphenyl	76%		10-130%
1718-51-0	Terphenyl-d14	65%		22-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 1 POST SOLIDIFICATION**Lab Sample ID:** D38480-1**Date Sampled:** 09/05/12**Matrix:** SO - Soil**Date Received:** 09/07/12**Method:** SW846 8015B**Percent Solids:** 79.4**Project:** PCU 197-36A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB17471.D	1	09/10/12	SK	n/a	n/a	GGB958
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	68.0	15	7.5	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1		Run# 2	Limits	
120-82-1	1,2,4-Trichlorobenzene	88%			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

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 1 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-1	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	79.4
Method:	SW846-8015B SW846 3546		
Project:	PCU 197-36A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FD17346.D	1	09/12/12	AW	09/11/12	OP6603	GFD891
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	2.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	842	17	11	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	74%		43-136%		

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

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

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 1 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-1	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	79.4
Project:	PCU 197-36A		

Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	16.8	0.13	mg/kg	5	09/11/12	09/14/12 JB	SW846 6020A ³	SW846 3050B ⁶
Barium	1270	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Cadmium	< 1.3	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Chromium	28.9	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Copper	28.8	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Lead	15.0	6.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Mercury	< 0.12	0.12	mg/kg	1	09/11/12	09/11/12 JM	SW846 7471B ¹	SW846 7471B ⁴
Nickel	22.4	3.8	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Selenium	< 6.3	6.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Silver	< 3.8	3.8	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Zinc	51.8	3.8	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵

- (1) Instrument QC Batch: MA2789
- (2) Instrument QC Batch: MA2795
- (3) Instrument QC Batch: MA2802
- (4) Prep QC Batch: MP8357
- (5) Prep QC Batch: MP8358
- (6) Prep QC Batch: MP8359

RL = Reporting Limit

Report of Analysis

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Client Sample ID: CUT 1 POST SOLIDIFICATION**Lab Sample ID:** D38480-1**Matrix:** SO - Soil**Date Sampled:** 09/05/12**Date Received:** 09/07/12**Percent Solids:** 79.4**Project:** PCU 197-36A**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	11400	1.0	umhos/cm	1	09/14/12	JK	SM2510B-1997 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	09/11/12	CJ	SW846 3060A/7196A
Chromium, Trivalent ^a	28.9	2.3	mg/kg	1	09/12/12 18:04	JB	SW846 3060/7196A M
Redox Potential Vs H2	146		mv	1	09/10/12	CT	ASTM D1498-76M
Solids, Percent	79.4		%	1	09/10/12	SWT	SM19 2540B M
pH	12.43		su	1	09/07/12 14:30	JD	SW846 9045D

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

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 1 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-1A	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	79.4
Project:	PCU 197-36A		

SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	15.9	2.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	< 1.0	1.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Sodium	1360	2.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA2796

(2) Prep QC Batch: MP8372

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 1 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-1A	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	79.4
Project:	PCU 197-36A		

General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	93.8		ratio	1	09/12/12 17:47	JB	USDA HANDBOOK 60

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

RL = Reporting Limit

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 2 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-2	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	75.5
Method:	SW846 8260B		
Project:	PCU 197-36A		
Run #1	File ID 3V20316.D	DF 1	Analyzed 09/08/12
Run #2			By BD n/a
			Prep Date n/a
			Prep Batch n/a
			Analytical Batch V3V1182
Run #1	Initial Weight 5.06 g	Final Volume 5.0 ml	Methanol Aliquot 100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	0.109	0.082	0.041	mg/kg	
108-88-3	Toluene	0.240	0.16	0.082	mg/kg	
100-41-4	Ethylbenzene	0.0507	0.16	0.031	mg/kg	J
1330-20-7	Xylene (total)	0.270	0.33	0.16	mg/kg	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	96%		64-130%
460-00-4	4-Bromofluorobenzene	104%		62-131%
17060-07-0	1,2-Dichloroethane-D4	112%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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4.3
4

Client Sample ID:	CUT 2 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-2	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	75.5
Method:	SW846 8270C BY SIM	SW846 3546	
Project:	PCU 197-36A		
File ID	DF	Analyzed	By
Run #1	3G11216.D	1	09/13/12 DC
Run #2			
Initial Weight	Final Volume		
Run #1	30.1 g	1.0 ml	
Run #2			

COGCC Table 910-1 PAH List

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	0.011	0.0057	mg/kg	
120-12-7	Anthracene	ND	0.011	0.0057	mg/kg	
56-55-3	Benzo(a)anthracene	0.0258	0.011	0.0057	mg/kg	
50-32-8	Benzo(a)pyrene	0.0275	0.011	0.0057	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.011	0.0057	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.011	0.0057	mg/kg	
218-01-9	Chrysene	0.0850	0.011	0.0057	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.011	0.0057	mg/kg	
206-44-0	Fluoranthene	0.0377	0.011	0.0057	mg/kg	
86-73-7	Fluorene	0.131	0.011	0.0057	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	0.0110	0.011	0.0057	mg/kg	
91-20-3	Naphthalene	0.838	0.015	0.014	mg/kg	
129-00-0	Pyrene	0.0729	0.011	0.0057	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	70%		10-145%
321-60-8	2-Fluorobiphenyl	72%		10-130%
1718-51-0	Terphenyl-d14	77%		22-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 2 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-2	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	75.5
Method:	SW846 8015B		
Project:	PCU 197-36A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB17472.D	1	09/10/12	SK	n/a	n/a	GGB958
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	27.3	16	8.2	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	93%		60-140%		

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

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

Accutest Laboratories

Report of Analysis

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4.3
4**Client Sample ID:** CUT 2 POST SOLIDIFICATION**Lab Sample ID:** D38480-2**Date Sampled:** 09/05/12**Matrix:** SO - Soil**Date Received:** 09/07/12**Method:** SW846-8015B SW846 3546**Percent Solids:** 75.5**Project:** PCU 197-36A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FD17348.D	1	09/12/12	AW	09/11/12	OP6603	GFD891
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	2.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	253	18	11	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	77%		43-136%		

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

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

Report of Analysis

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Client Sample ID:	CUT 2 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-2	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	75.5
Project:	PCU 197-36A		

Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	11.0	0.13	mg/kg	5	09/11/12	09/14/12 JB	SW846 6020A ³	SW846 3050B ⁶
Barium	2590	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Cadmium	< 1.3	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Chromium	17.3	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Copper	163	1.3	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Lead	33.3	6.4	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Mercury	< 0.12	0.12	mg/kg	1	09/11/12	09/11/12 JM	SW846 7471B ¹	SW846 7471B ⁴
Nickel	29.3	3.9	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Selenium	< 6.4	6.4	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Silver	< 3.9	3.9	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵
Zinc	100	3.9	mg/kg	1	09/11/12	09/12/12 JB	SW846 6010C ²	SW846 3050B ⁵

- (1) Instrument QC Batch: MA2789
- (2) Instrument QC Batch: MA2795
- (3) Instrument QC Batch: MA2802
- (4) Prep QC Batch: MP8357
- (5) Prep QC Batch: MP8358
- (6) Prep QC Batch: MP8359

RL = Reporting Limit

Report of Analysis

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Client Sample ID: CUT 2 POST SOLIDIFICATION**Lab Sample ID:** D38480-2**Matrix:** SO - Soil**Project:** PCU 197-36A**Date Sampled:** 09/05/12**Date Received:** 09/07/12**Percent Solids:** 75.5**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	11900	1.0	umhos/cm	1	09/14/12	JK	SM2510B-1997 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	09/11/12	CJ	SW846 3060A/7196A
Chromium, Trivalent ^a	17.3	2.3	mg/kg	1	09/12/12 18:57	JB	SW846 3060/7196A M
Redox Potential Vs H2	160		mv	1	09/10/12	CT	ASTM D1498-76M
Solids, Percent	75.5		%	1	09/10/12	SWT	SM19 2540B M
pH	11.74		su	1	09/07/12 14:30	JD	SW846 9045D

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

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 2 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-2A	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	75.5
Project:	PCU 197-36A		

SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	86.6	2.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	< 1.0	1.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Sodium	2270	2.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA2796

(2) Prep QC Batch: MP8372

RL = Reporting Limit

Report of Analysis

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Client Sample ID:	CUT 2 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-2A	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	75.5
Project:	PCU 197-36A		

General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	67.0		ratio	1	09/12/12 17:57	JB	USDA HANDBOOK 60

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

RL = Reporting Limit

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 3 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-3	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	81.6
Method:	SW846 8260B		
Project:	PCU 197-36A		
File ID	DF	Analyzed	By
Run #1	3V20317.D	1	09/08/12
Run #2			BD
		n/a	
		n/a	
		V3V1182	
Initial Weight	Final Volume	Methanol Aliquot	
Run #1	5.08 g	5.0 ml	100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	0.0957	0.072	0.036	mg/kg	
108-88-3	Toluene	0.267	0.14	0.072	mg/kg	
100-41-4	Ethylbenzene	0.0273	0.14	0.027	mg/kg	J
1330-20-7	Xylene (total)	0.327	0.29	0.14	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	95%		64-130%
460-00-4	4-Bromofluorobenzene	104%		62-131%
17060-07-0	1,2-Dichloroethane-D4	109%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 3 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-3	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	81.6
Method:	SW846 8270C BY SIM	SW846 3546	
Project:	PCU 197-36A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G11217.D	1	09/13/12	DC	09/11/12	OP6602	E3G522
Run #2							

	Initial Weight	Final Volume
Run #1	30.2 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	0.0310	0.010	0.0053	mg/kg	
50-32-8	Benzo(a)pyrene	0.0548	0.010	0.0053	mg/kg	
205-99-2	Benzo(b)fluoranthene	0.0440	0.010	0.0053	mg/kg	
207-08-9	Benzo(k)fluoranthene	0.0098	0.010	0.0053	mg/kg	J
218-01-9	Chrysene	0.108	0.010	0.0053	mg/kg	
53-70-3	Dibenz(a,h)anthracene	0.0141	0.010	0.0053	mg/kg	
206-44-0	Fluoranthene	0.0342	0.010	0.0053	mg/kg	
86-73-7	Fluorene	0.139	0.010	0.0053	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	0.0139	0.010	0.0053	mg/kg	
91-20-3	Naphthalene	0.482	0.014	0.013	mg/kg	
129-00-0	Pyrene	0.0689	0.010	0.0053	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	83%		10-145%
321-60-8	2-Fluorobiphenyl	66%		10-130%
1718-51-0	Terphenyl-d14	77%		22-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 3 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-3	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	81.6
Method:	SW846 8015B		
Project:	PCU 197-36A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB17474.D	1	09/10/12	SK	n/a	n/a	GGB958
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	44.4	14	7.2	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	88%		60-140%		

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

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

Accutest Laboratories

Report of Analysis

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Client Sample ID:	CUT 3 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-3	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	81.6
Method:	SW846-8015B SW846 3546		
Project:	PCU 197-36A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FD17350.D	1	09/12/12	AW	09/11/12	OP6603	GFD891
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	2.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	434	16	11	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	91%		43-136%		

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

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

Report of Analysis

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Client Sample ID:	CUT 3 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-3	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	81.6
Project:	PCU 197-36A		

Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	9.1	0.12	mg/kg	5	09/11/12	09/14/12	JB	SW846 6020A ⁴
Barium	5630	5.9	mg/kg	5	09/11/12	09/13/12	JM	SW846 6010C ³
Cadmium	< 1.2	1.2	mg/kg	1	09/11/12	09/12/12	JB	SW846 6010C ²
Chromium	16.6	1.2	mg/kg	1	09/11/12	09/12/12	JB	SW846 6010C ²
Copper	28.0	1.2	mg/kg	1	09/11/12	09/12/12	JB	SW846 6010C ²
Lead	28.8	5.9	mg/kg	1	09/11/12	09/12/12	JB	SW846 6010C ²
Mercury	< 0.12	0.12	mg/kg	1	09/11/12	09/11/12	JM	SW846 7471B ¹
Nickel	20.7	18	mg/kg	5	09/11/12	09/13/12	JM	SW846 6010C ³
Selenium	< 5.9	5.9	mg/kg	1	09/11/12	09/12/12	JB	SW846 6010C ²
Silver	< 3.6	3.6	mg/kg	1	09/11/12	09/12/12	JB	SW846 6010C ²
Zinc	56.2	18	mg/kg	5	09/11/12	09/13/12	JM	SW846 6010C ³

- (1) Instrument QC Batch: MA2789
- (2) Instrument QC Batch: MA2795
- (3) Instrument QC Batch: MA2799
- (4) Instrument QC Batch: MA2802
- (5) Prep QC Batch: MP8357
- (6) Prep QC Batch: MP8358
- (7) Prep QC Batch: MP8359

RL = Reporting Limit

Report of Analysis

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Client Sample ID: CUT 3 POST SOLIDIFICATION**Lab Sample ID:** D38480-3**Matrix:** SO - Soil**Project:** PCU 197-36A**Date Sampled:** 09/05/12**Date Received:** 09/07/12**Percent Solids:** 81.6**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	11400	1.0	umhos/cm	1	09/14/12	JK	SM2510B-1997 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	09/11/12	CJ	SW846 3060A/7196A
Chromium, Trivalent ^a	16.6	2.2	mg/kg	1	09/12/12 19:26	JB	SW846 3060/7196A M
Redox Potential Vs H2	145		mv	1	09/10/12	CT	ASTM D1498-76M
Solids, Percent	81.6		%	1	09/10/12	SWT	SM19 2540B M
pH	12.45		su	1	09/07/12 14:30	JD	SW846 9045D

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

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 3 POST SOLIDIFICATION	Date Sampled:	09/05/12
Lab Sample ID:	D38480-3A	Date Received:	09/07/12
Matrix:	SO - Soil	Percent Solids:	81.6
Project:	PCU 197-36A		

SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	2.95	2.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	< 1.0	1.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Sodium	2100	2.0	mg/l	1	09/11/12	09/12/12 JB	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA2796

(2) Prep QC Batch: MP8372

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 3 POST SOLIDIFICATION**Lab Sample ID:** D38480-3A**Matrix:** SO - Soil**Project:** PCU 197-36A**Date Sampled:** 09/05/12**Date Received:** 09/07/12**Percent Solids:** 81.6**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	336		ratio	1	09/12/12 18:08	JB	USDA HANDBOOK 60

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

RL = Reporting Limit



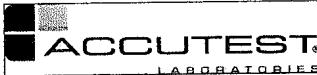
Misc. Forms

5

Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody



CHAIN OF CUSTODY

PAGE 1 OF 1

D38480: Chain of Custody

Page 1 of 2



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D38480

Client: KRW CONSULTING

Immediate Client Services Action Required: No

Date / Time Received: 9/7/2012 12:30:00 PM

No. Coolers:

1

Client Service Action Required at Login: No

Project: PCU 197-36A XTO

Airbill #'s: HDCO

Cooler Security Y or N

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

Cooler Temperature Y or N

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

Quality Control Preservation Y or N N/A

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

Sample Integrity - Documentation

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

Sample Integrity - Condition

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

Sample Integrity - Instructions

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

Comments

Accutest Laboratories
V:(303) 425-6021

4036 Youngfield Street
F: (303) 425-6854

Wheat Ridge, CO
www.accutest.com

5.1

5

D38480: 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: D38480
Account: XTOKRWR XTO Energy
Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1182-MB	3V20302.D	1	09/07/12	BD	n/a	n/a	V3V1182

The QC reported here applies to the following samples:

Method: SW846 8260B

D38480-1, D38480-2, D38480-3

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	102%
460-00-4	4-Bromofluorobenzene	97%
17060-07-0	1,2-Dichloroethane-D4	115%

Blank Spike Summary

Job Number: D38480
Account: XTOKWR XTO Energy
Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1182-BS	3V20303.D	1	09/07/12	BD	n/a	n/a	V3V1182

The QC reported here applies to the following samples:

Method: SW846 8260B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	48.5	97	70-130
100-41-4	Ethylbenzene	50	48.6	97	70-130
108-88-3	Toluene	50	45.6	91	70-130
1330-20-7	Xylene (total)	150	148	99	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	100%	64-130%
460-00-4	4-Bromofluorobenzene	102%	62-131%
17060-07-0	1,2-Dichloroethane-D4	111%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D38454-1MS	3V20305.D	1	09/07/12	BD	n/a	n/a	V3V1182
D38454-1MSD	3V20306.D	1	09/07/12	BD	n/a	n/a	V3V1182
D38454-1	3V20304.D	1	09/07/12	BD	n/a	n/a	V3V1182

The QC reported here applies to the following samples:

Method: SW846 8260B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	D38454-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	ND		3820	2930	77	3610	95	21	64-139/30
100-41-4	Ethylbenzene	ND		3820	3060	80	3740	98	20	68-136/30
108-88-3	Toluene	108	J	3820	2760	69	3350	85	19	60-130/30
1330-20-7	Xylene (total)	ND		11500	9580	84	11600	101	19	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D38454-1	Limits
2037-26-5	Toluene-D8	94%	93%	95%	64-130%
460-00-4	4-Bromofluorobenzene	109%	107%	101%	62-131%
17060-07-0	1,2-Dichloroethane-D4	109%	109%	111%	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\V3090712.S\
 Data File : 3V20315.D
 Acq On : 8 Sep 2012 12:15 am
 Operator : BRETD
 Sample : D38480-1
 Misc : MS4630,V3V1182,5.064,,100,5,1
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Sep 08 11:23:40 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1160TVH1160SOIL.M
 Quant Title : 8260
 QLast Update : Fri Aug 24 10:57:50 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.860	168	192918	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.655	114	320306	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.293	117	345598	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.282	152	201499	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.248	102	25013	57.59	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	115.18%	
61) Toluene-d8	14.051	98	426392	47.26	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	94.52%	
69) 4-Bromofluorobenzene	16.243	95	181735	51.47	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	102.94%	

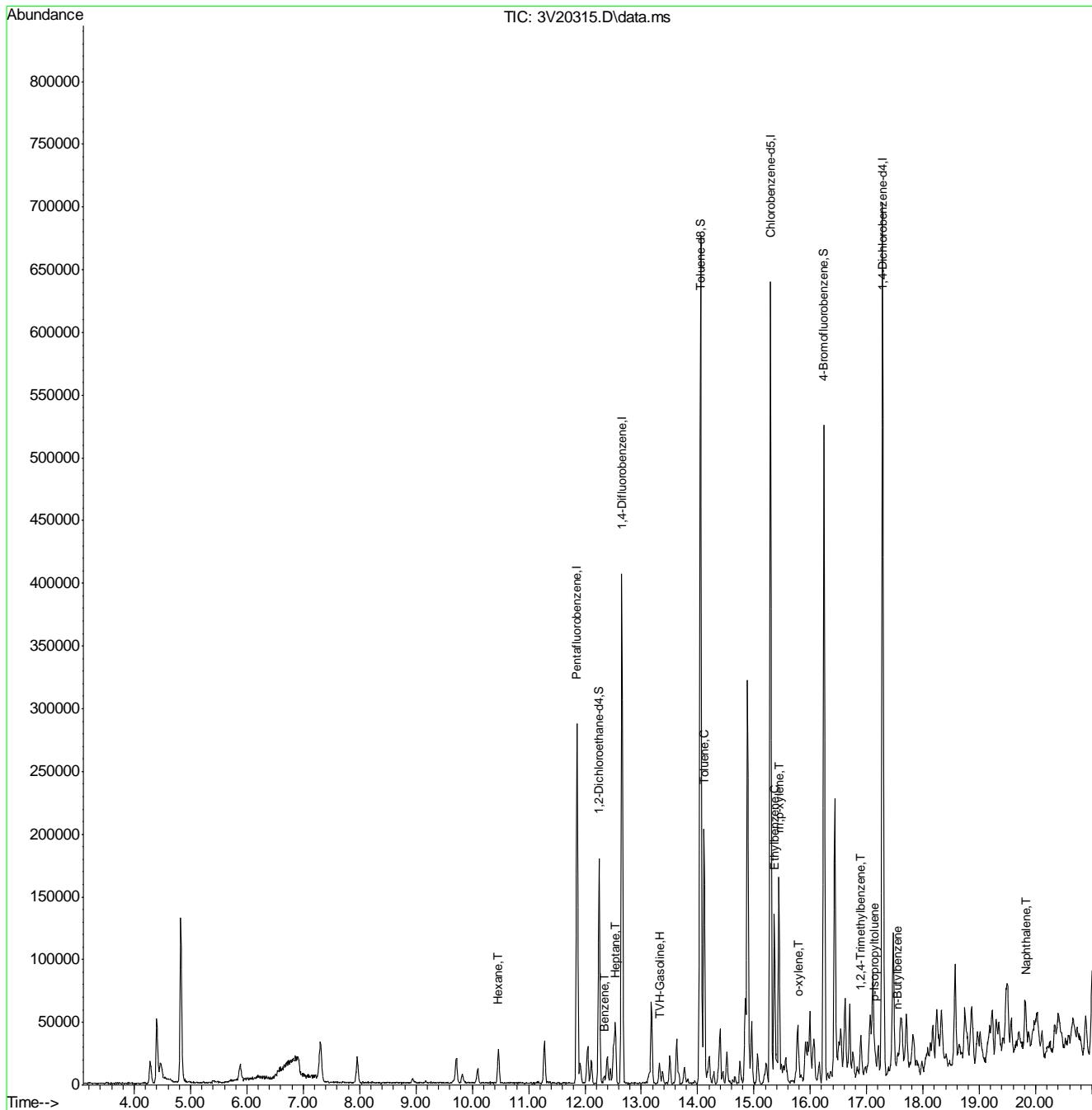
Target Compounds						Qvalue
1) TVH-Gasoline	13.329	TIC	4159379m	146.11	ug/l	
41) Hexane	10.467	57	14000	2.78	ug/l	100
43) Heptane	12.540	43	20614	3.21	ug/l	94
50) Benzene	12.347	78	3885	0.34	ug/l	100
62) Toluene	14.109	92	57037	6.52	ug/l	98
66) Ethylbenzene	15.360	91	28347	1.74	ug/l	94
72) m,p-xylene	15.444	106	44624	6.87	ug/l	92
73) o-xylene	15.794	106	3392	0.68	ug/l	80
82) 1,2,4-Trimethylbenzene	16.894	105	9623	0.71	ug/l	89
86) p-Isopropyltoluene	17.148	119	7683	0.55	ug/l #	88
88) n-Butylbenzene	17.539	91	4721	0.35	ug/l #	87
91) Naphthalene	19.840	128	4708	0.46	ug/l	100

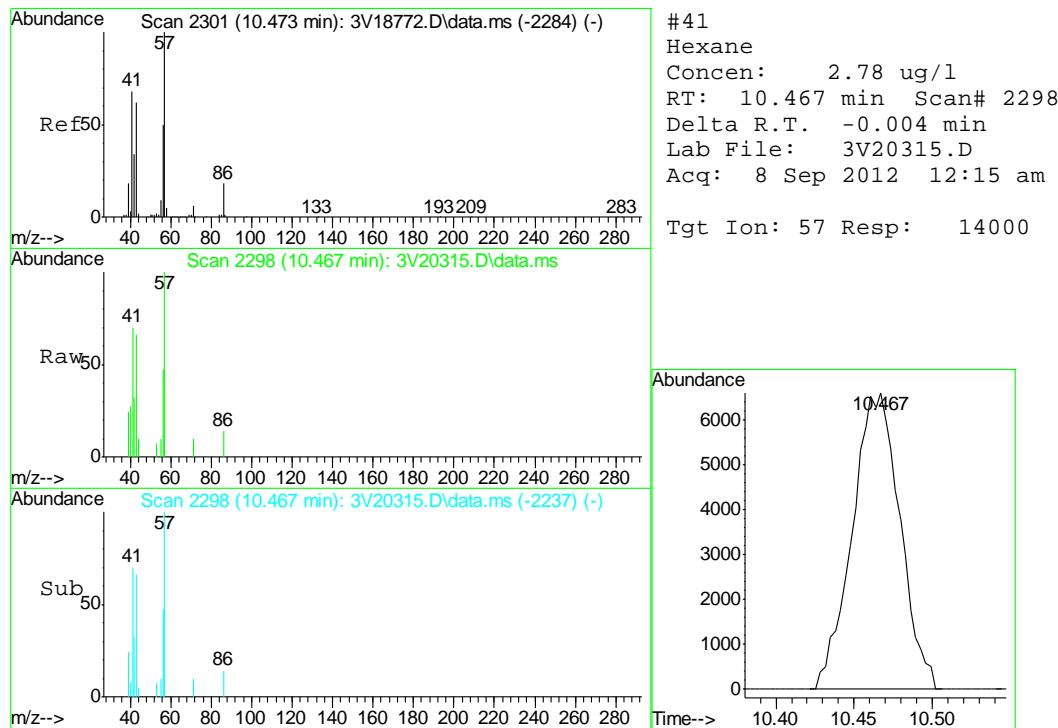
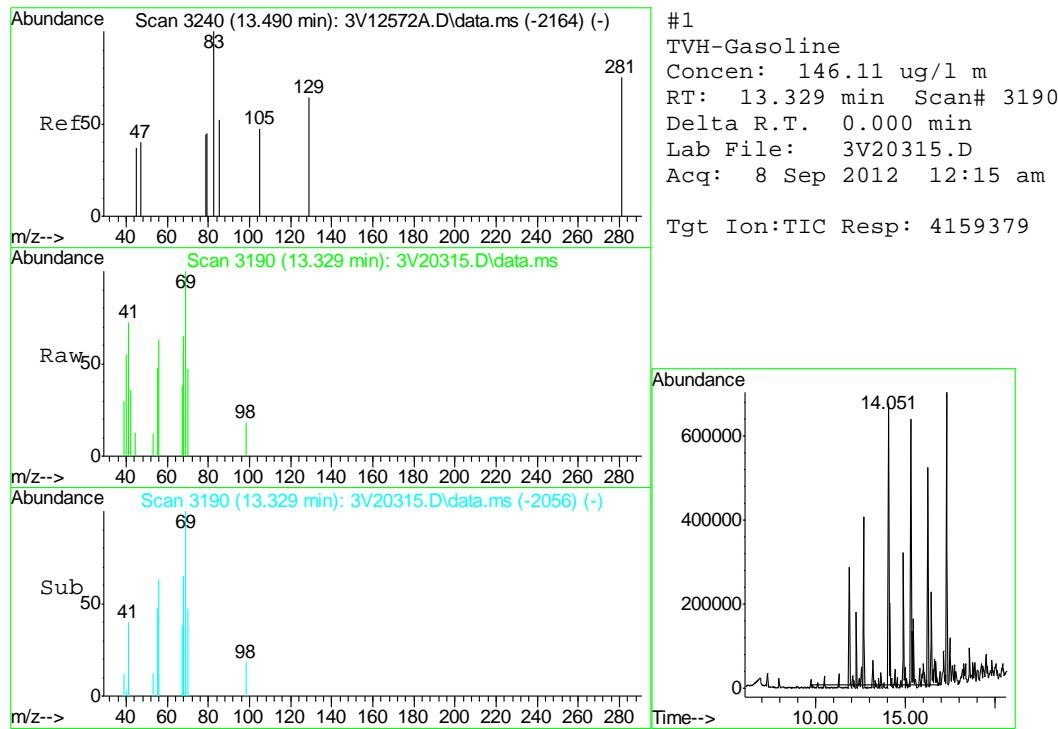
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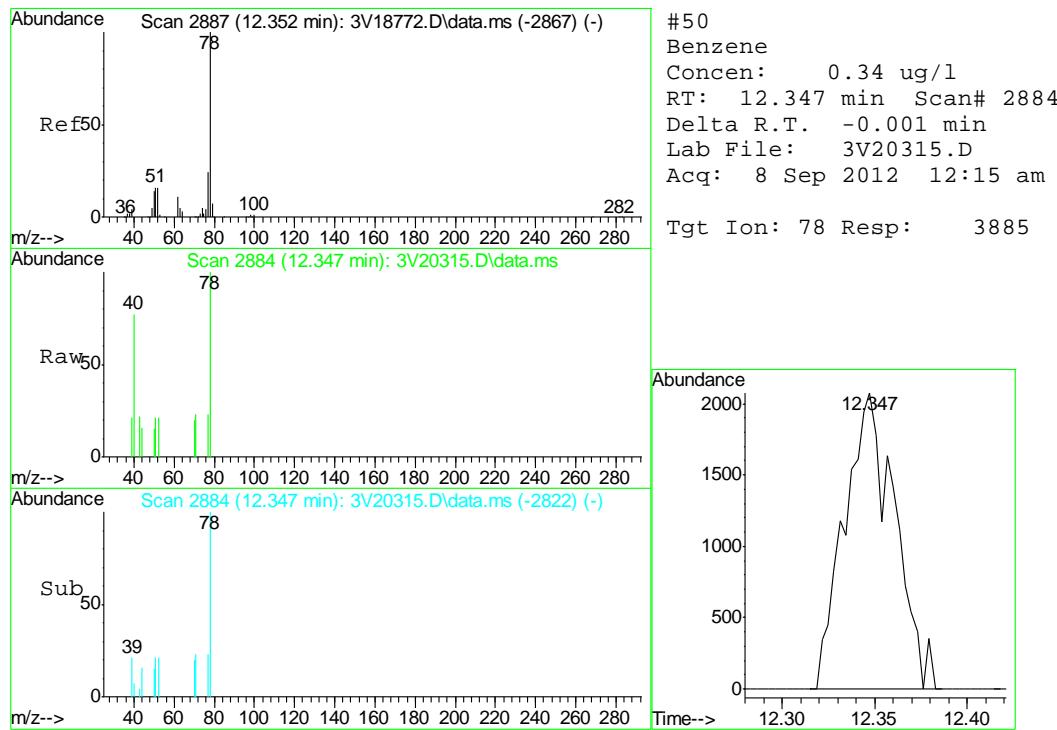
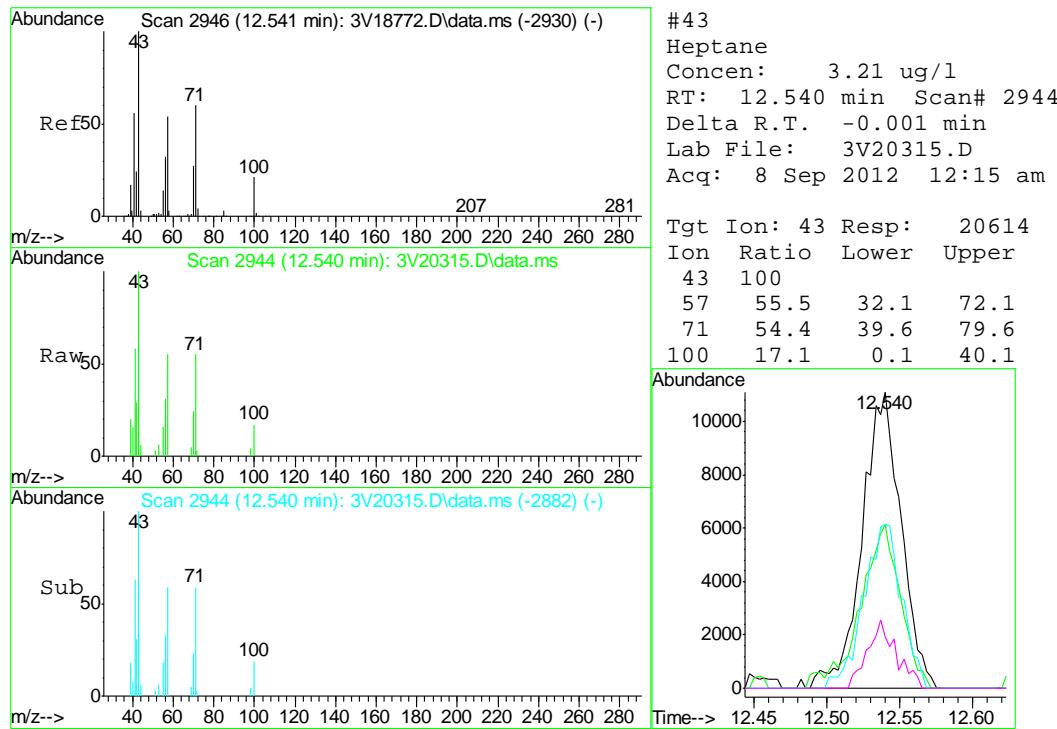
Quantitation Report (QT Reviewed)

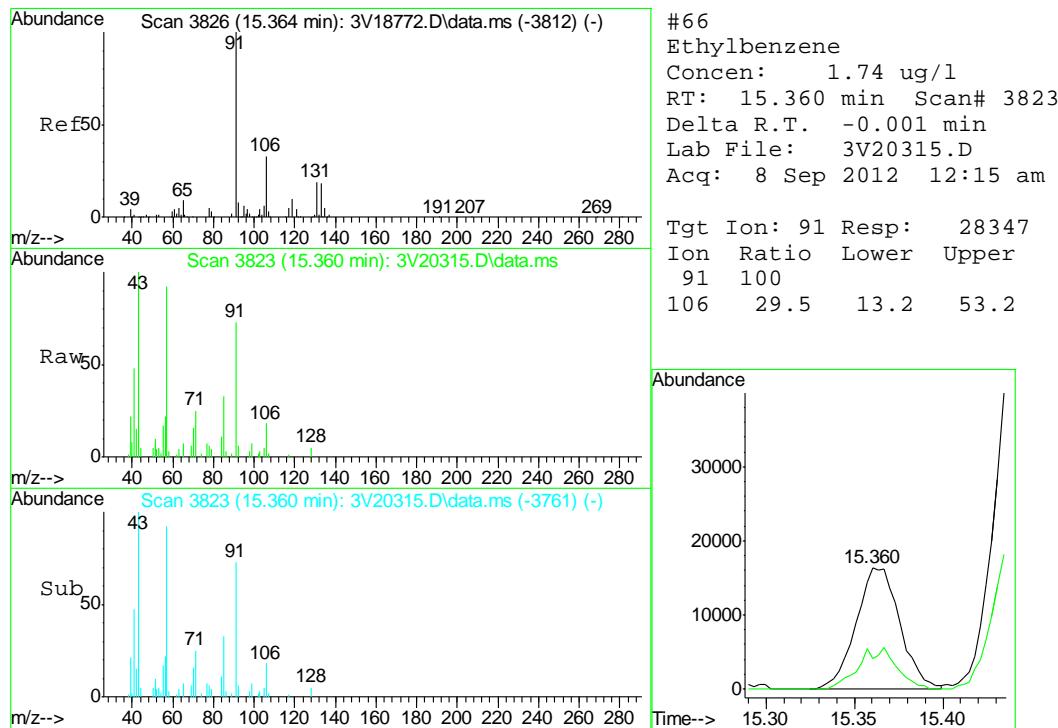
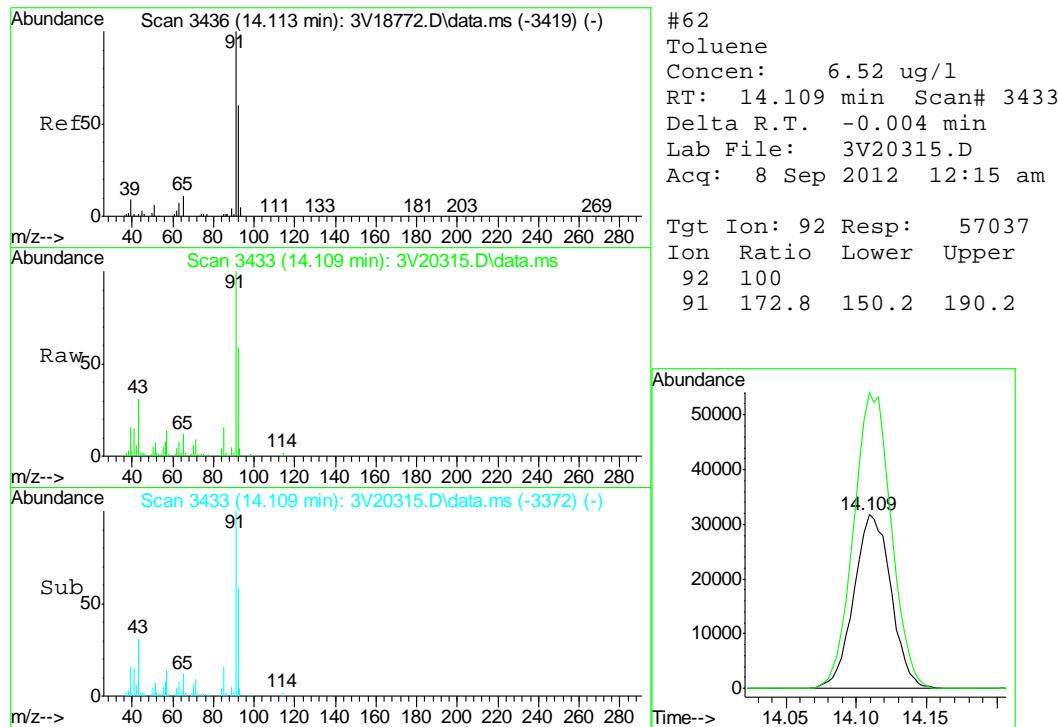
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 Operator : BRETD
 Sample : D38480-1
 Misc : MS4630,V3V1182,5.064,,100,5,1
 ALS Vial : 17 Sample Multiplier: 1

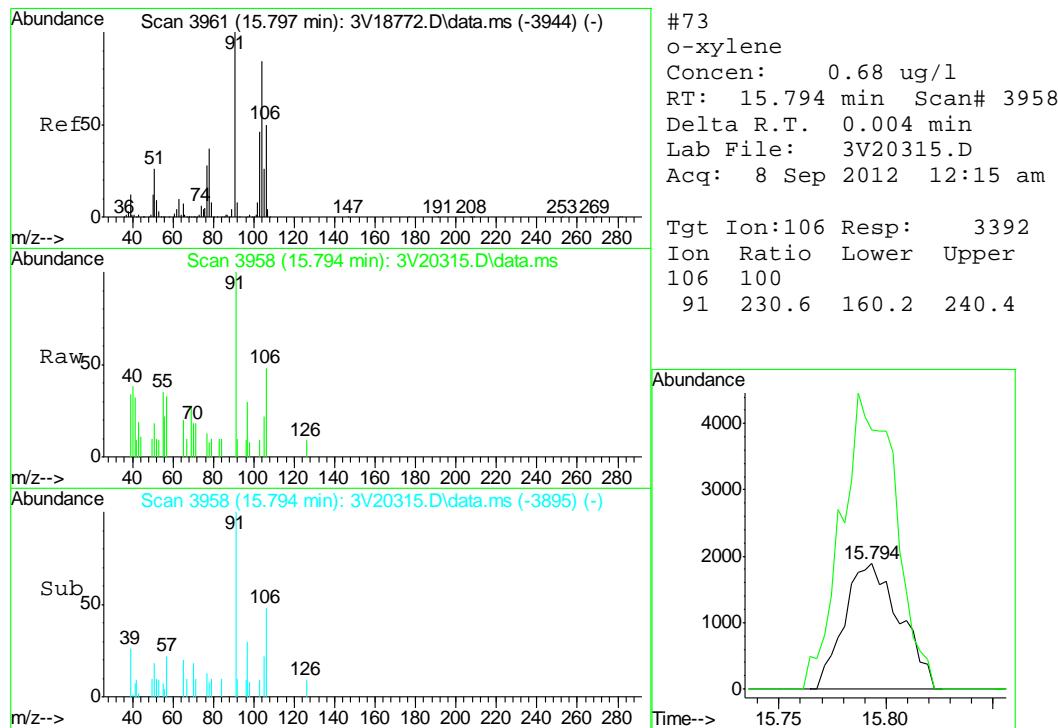
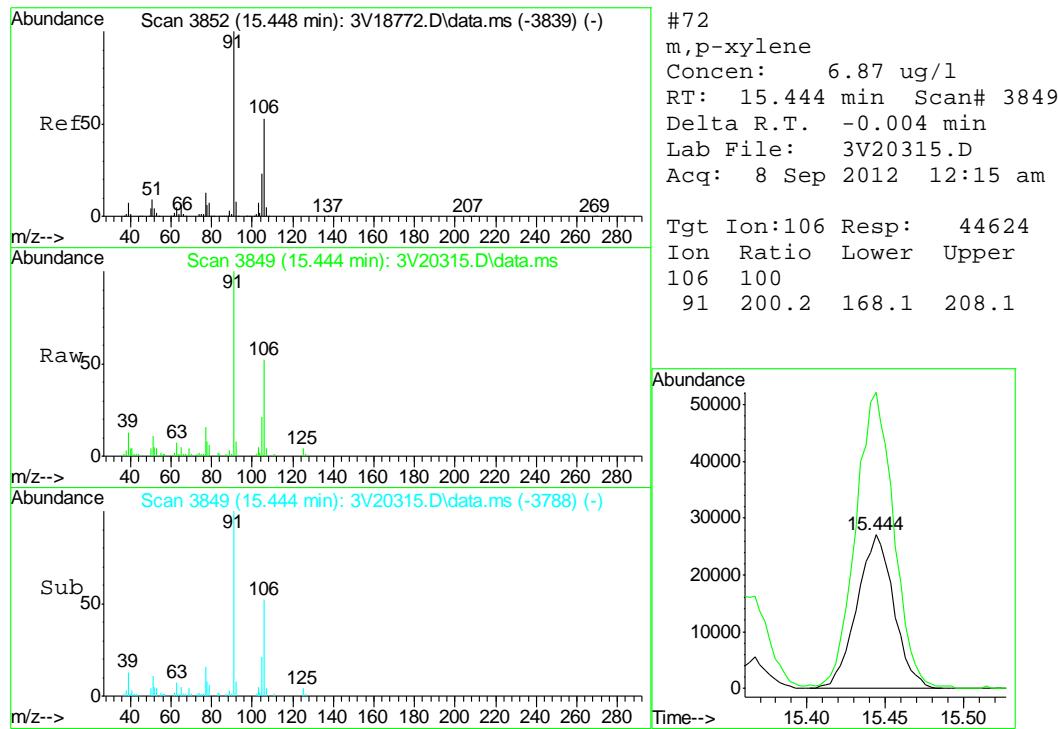
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 Quant Title : 8260
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 Response via : Initial Calibration

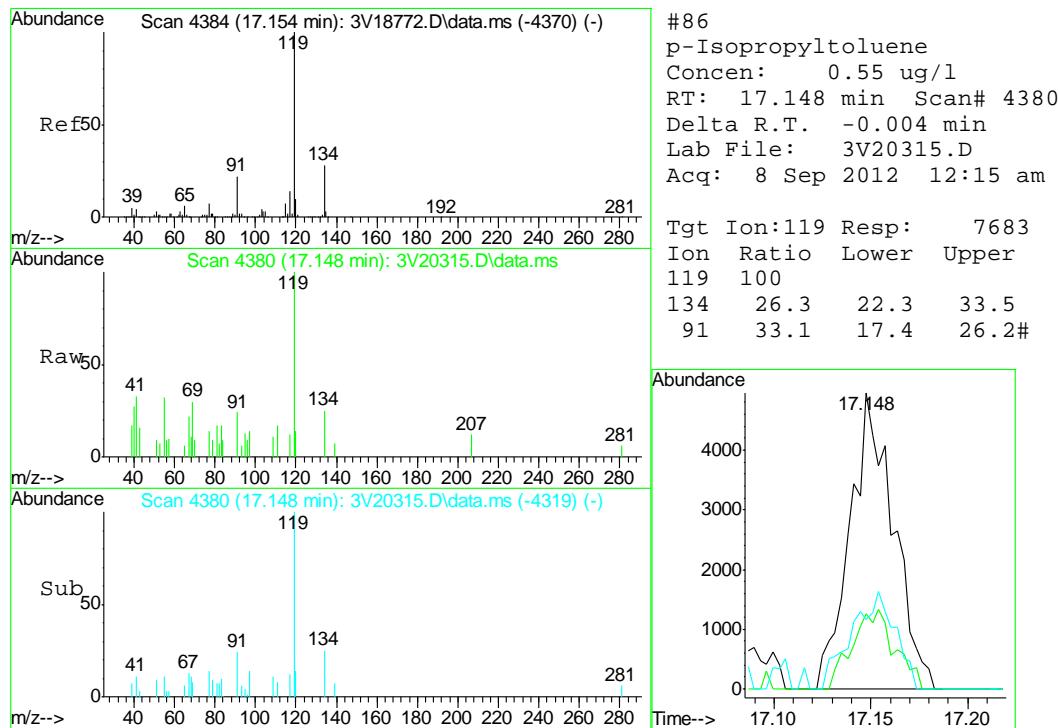
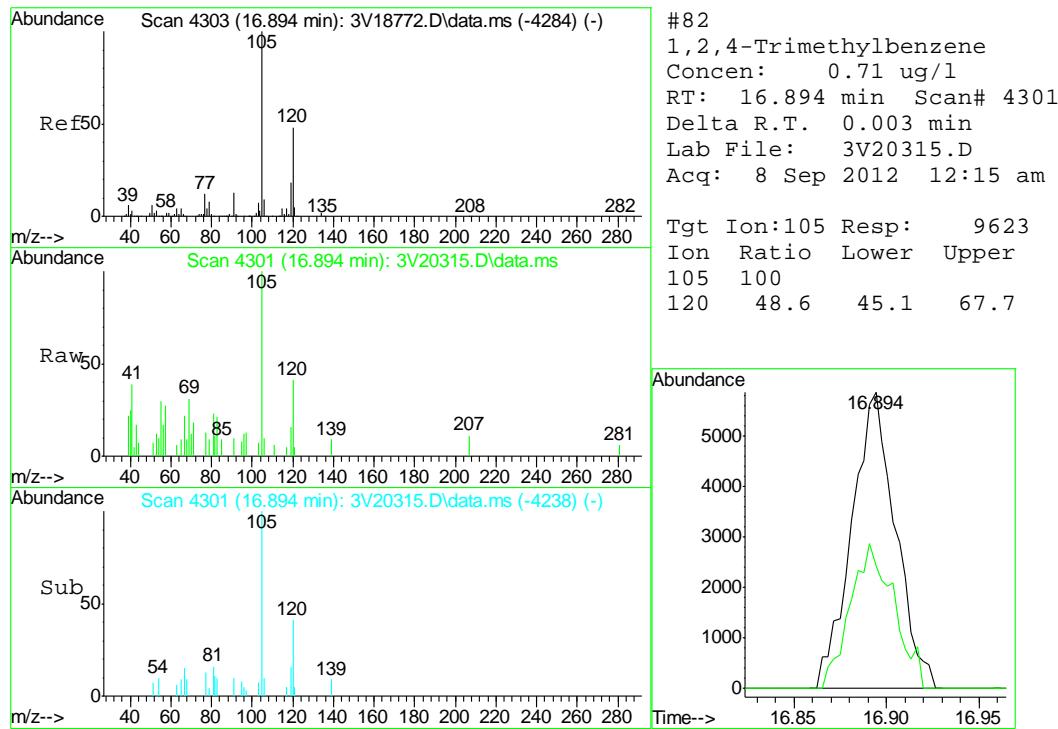


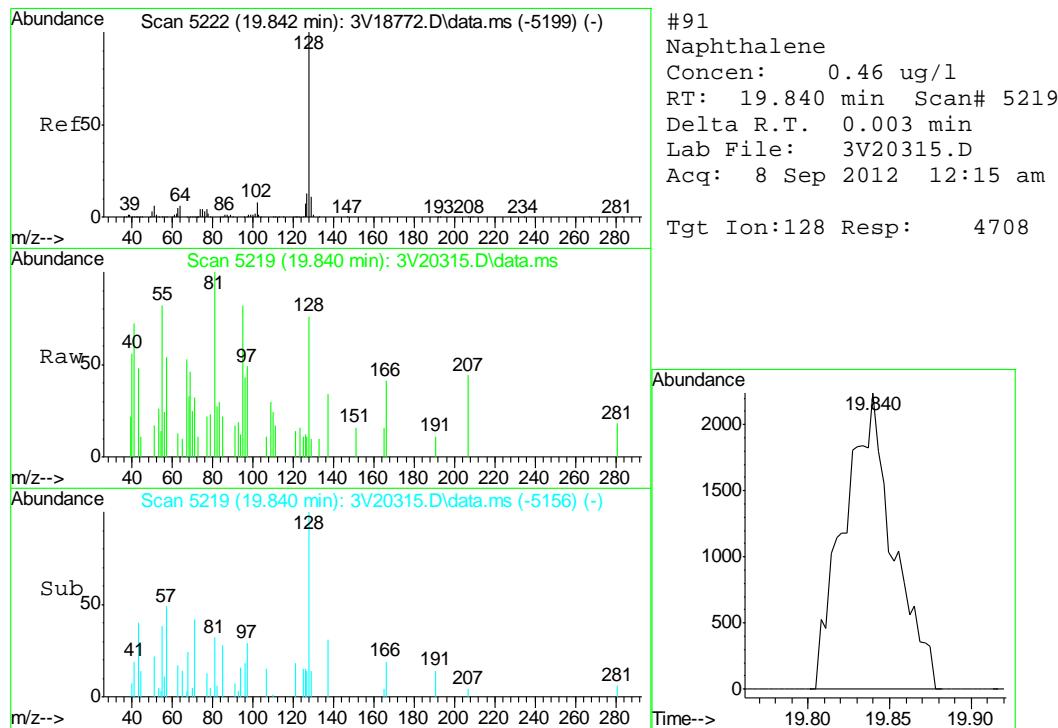
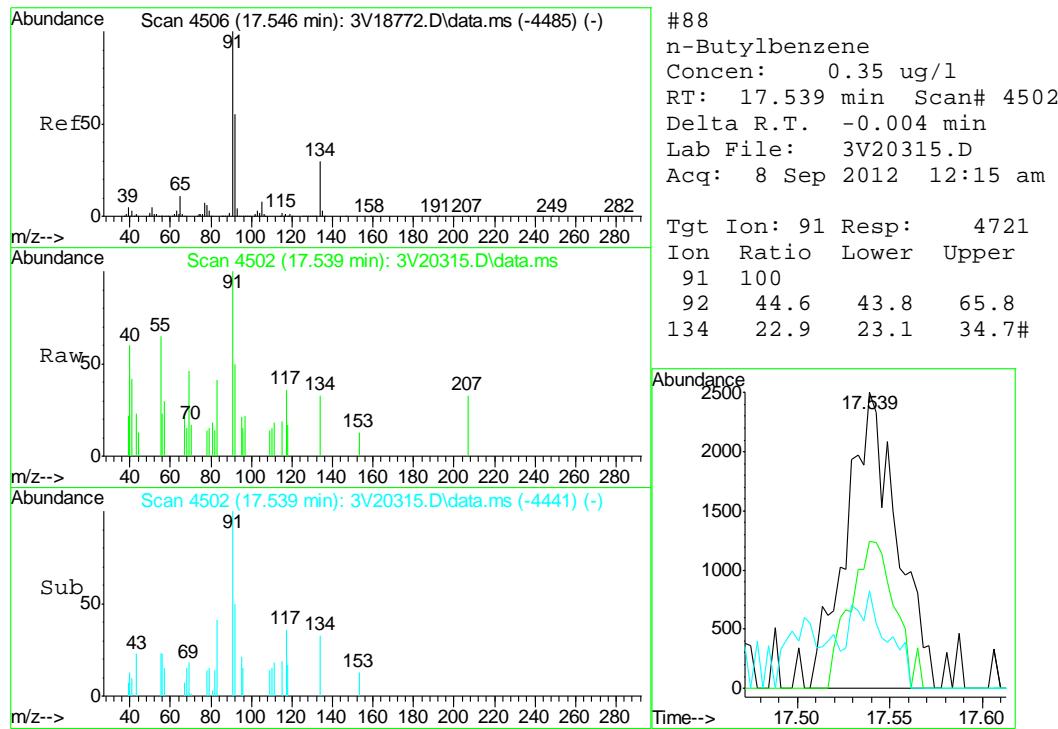












Judy Nelson
 09/10/12 16:37

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3090712.S\
 Data File : 3V20316.D
 Acq On : 8 Sep 2012 12:46 am
 Operator : BRETD
 Sample : D38480-2
 Misc : MS4630,V3V1182,5.062,,100,5,1
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Sep 08 11:25:50 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1160TVH1160SOIL.M
 Quant Title : 8260
 QLast Update : Fri Aug 24 10:57:50 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.860	168	200277	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.656	114	329230	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.294	117	350955	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.283	152	204961	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.252	102	25231	55.95	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	111.90%	
61) Toluene-d8	14.048	98	438053	47.81	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	95.62%	
69) 4-Bromofluorobenzene	16.243	95	185826	51.83	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	103.66%	

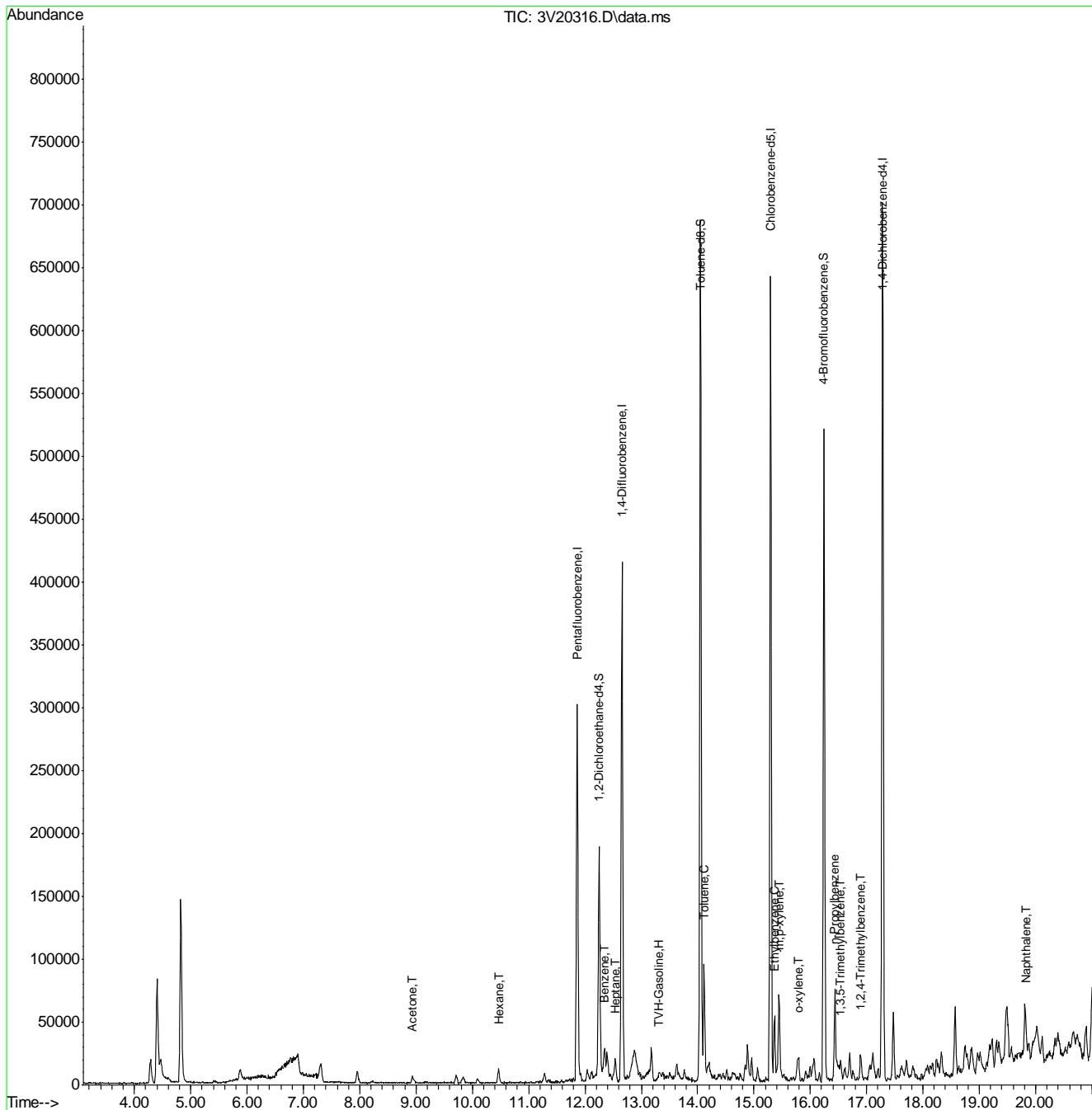
Target Compounds						Qvalue
1) TVH-Gasoline	13.329	TIC	1484752m	52.16	ug/l	
15) Acetone	8.937	58	2228	1.21	ug/l	# 51
41) Hexane	10.461	57	5795	1.12	ug/l	100
43) Heptane	12.540	43	7788	1.18	ug/l	93
50) Benzene	12.348	78	15941	1.34	ug/l	100
62) Toluene	14.109	92	26096	2.94	ug/l	96
66) Ethylbenzene	15.364	91	10250	0.62	ug/l	95
72) m,p-xylene	15.444	106	16544	2.51	ug/l	# 78
73) o-xylene	15.791	106	4177	0.81	ug/l	94
77) n-Propylbenzene	16.426	91	4725	0.24	ug/l	# 84
80) 1,3,5-Trimethylbenzene	16.532	105	5059m	0.37	ug/l	
82) 1,2,4-Trimethylbenzene	16.891	105	9456	0.68	ug/l	88
91) Naphthalene	19.840	128	8208	0.78	ug/l	100

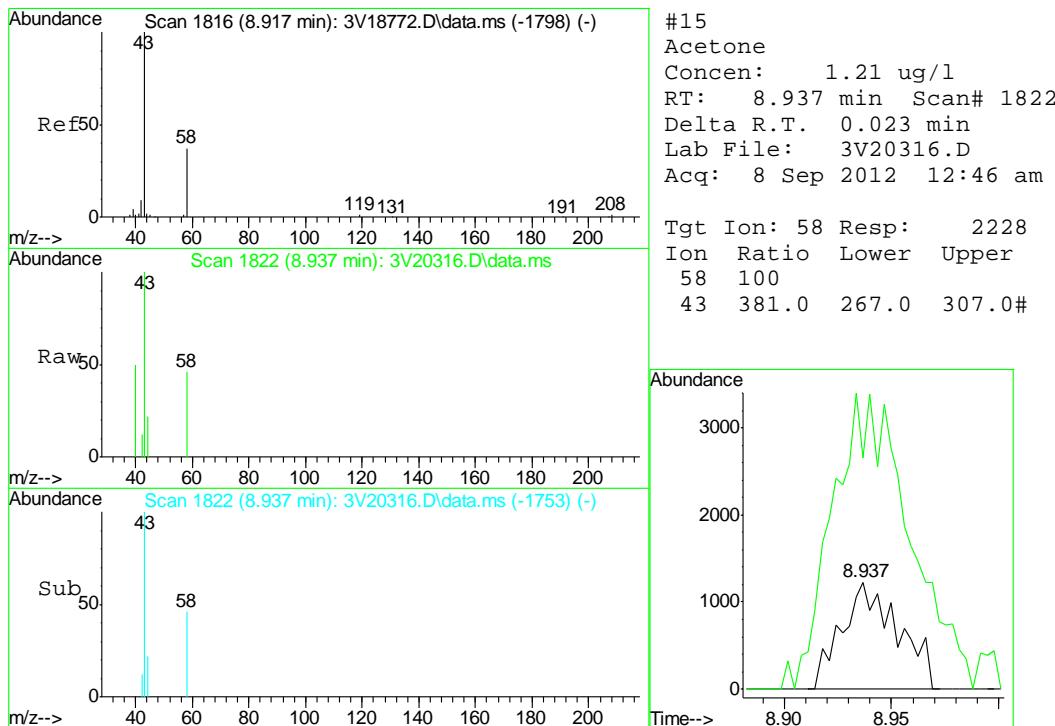
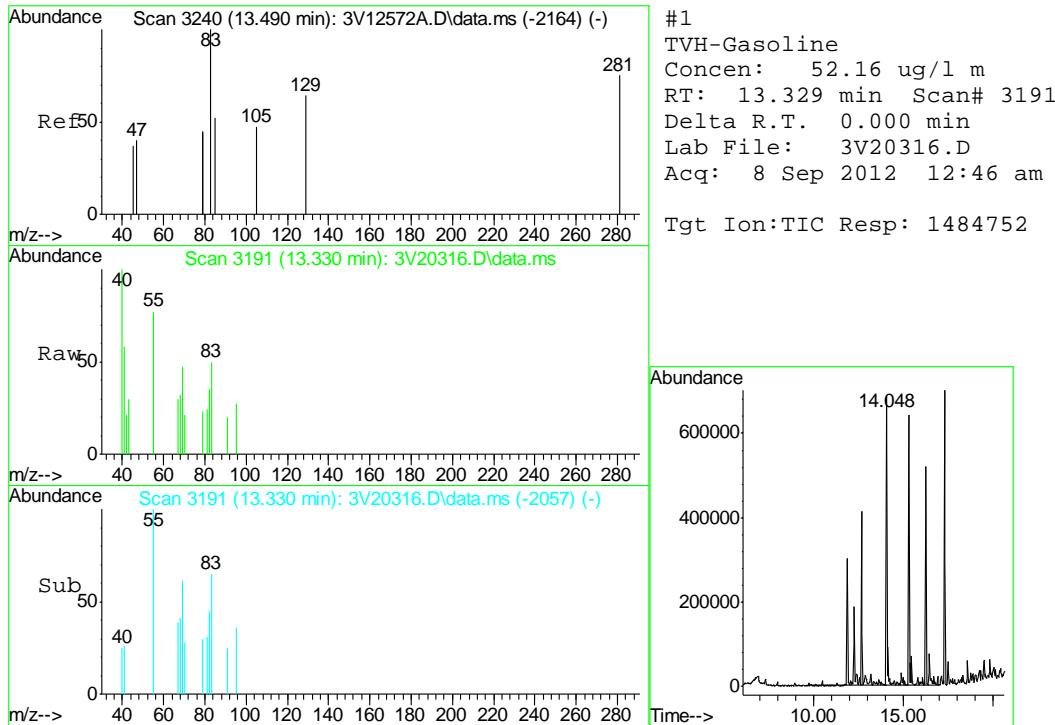
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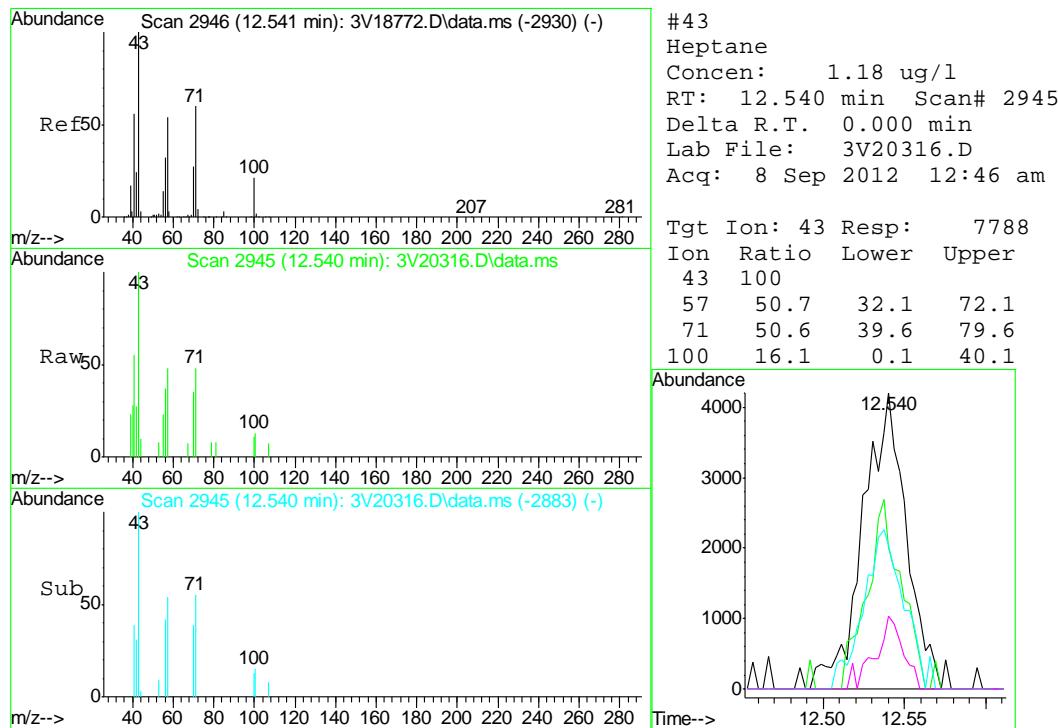
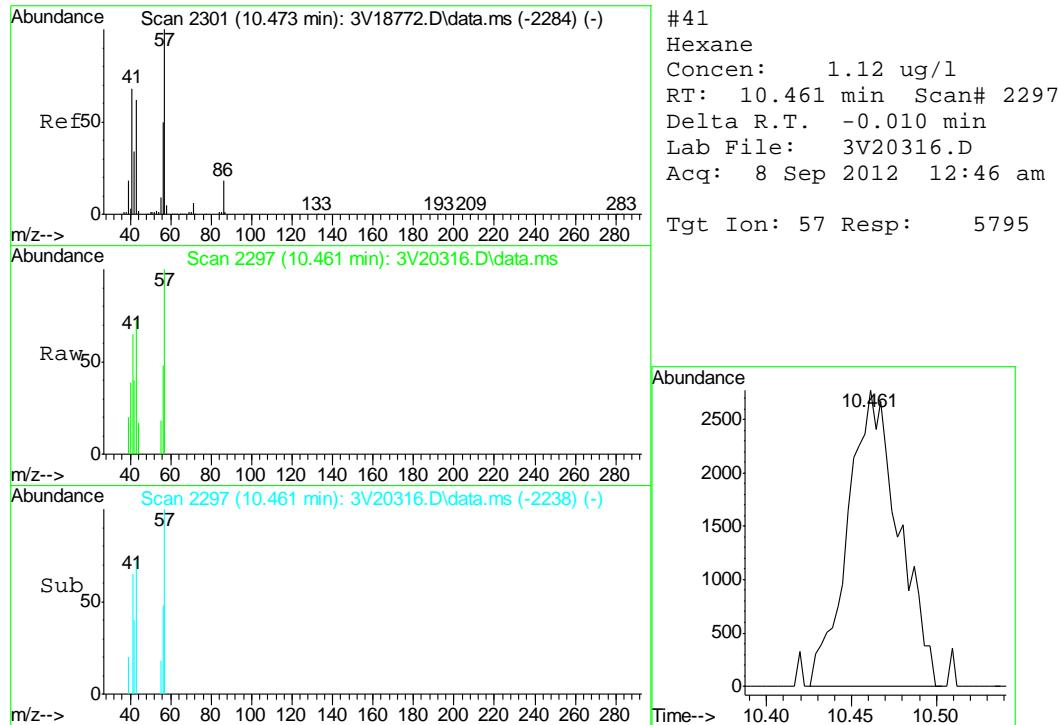
Quantitation Report (QT Reviewed)

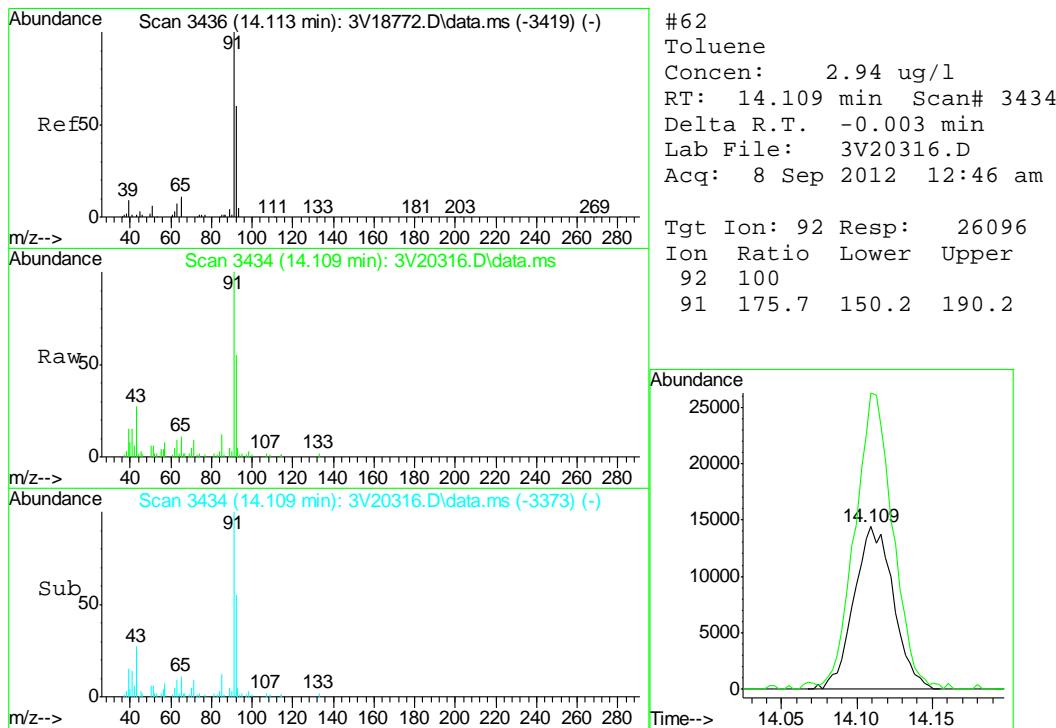
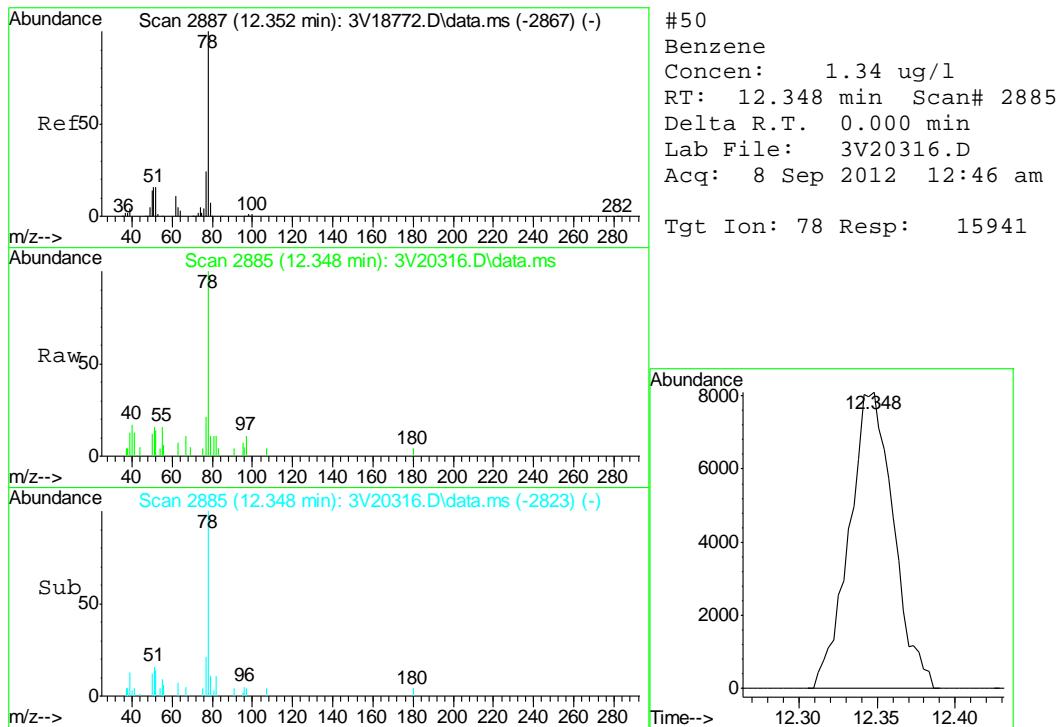
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 ALS Vial : 18 Sample Multiplier: 1

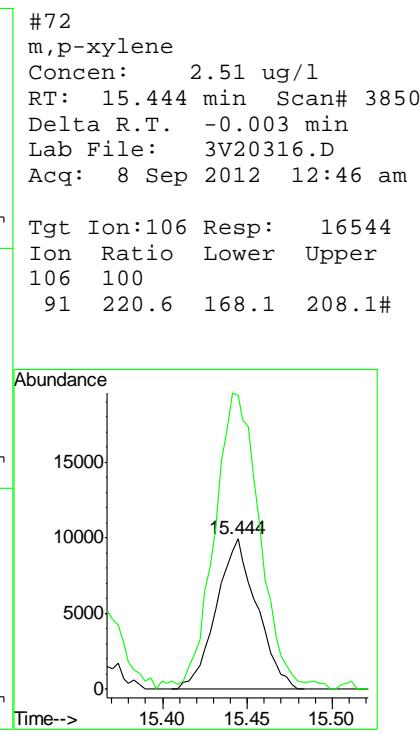
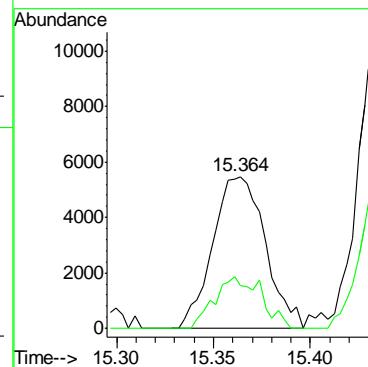
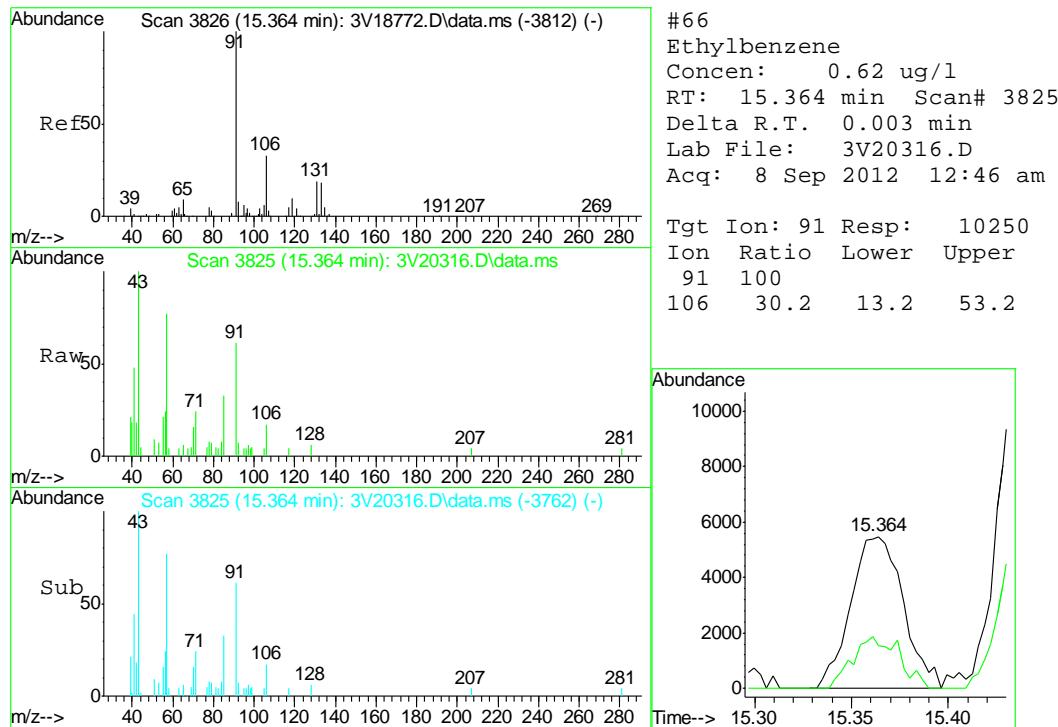
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 Response via : Initial Calibration

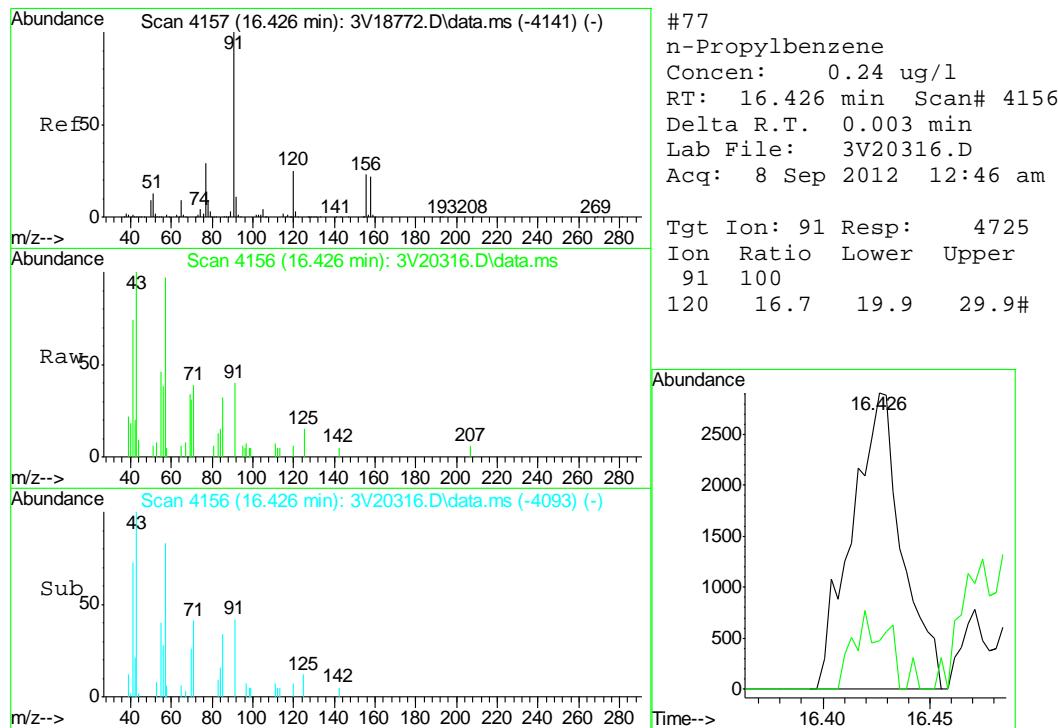
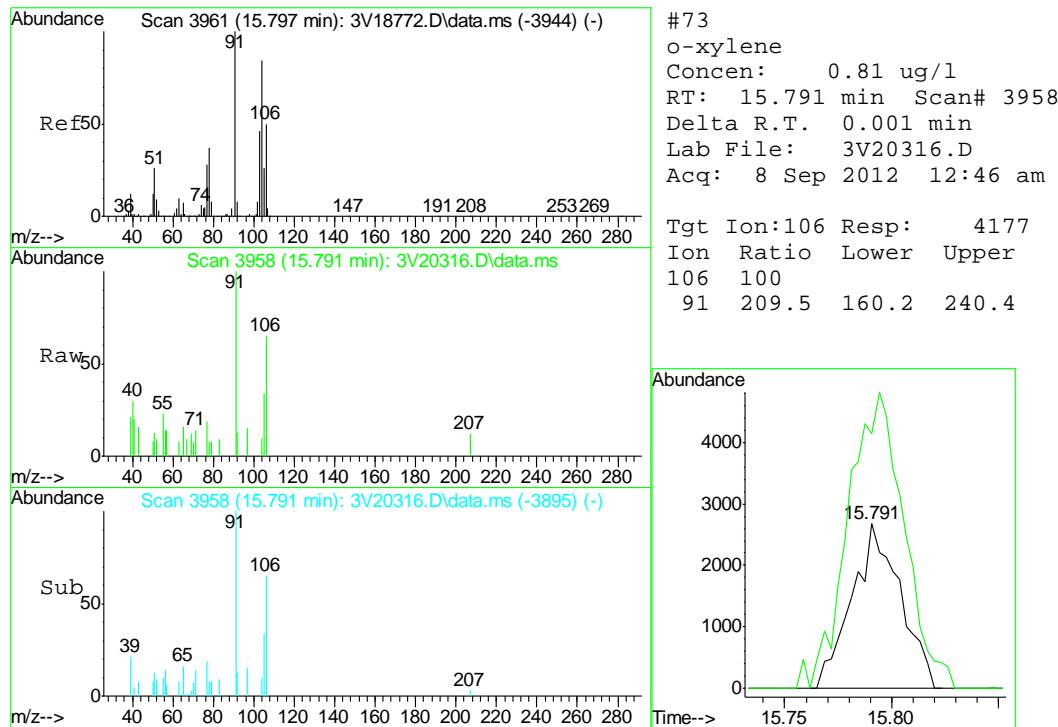


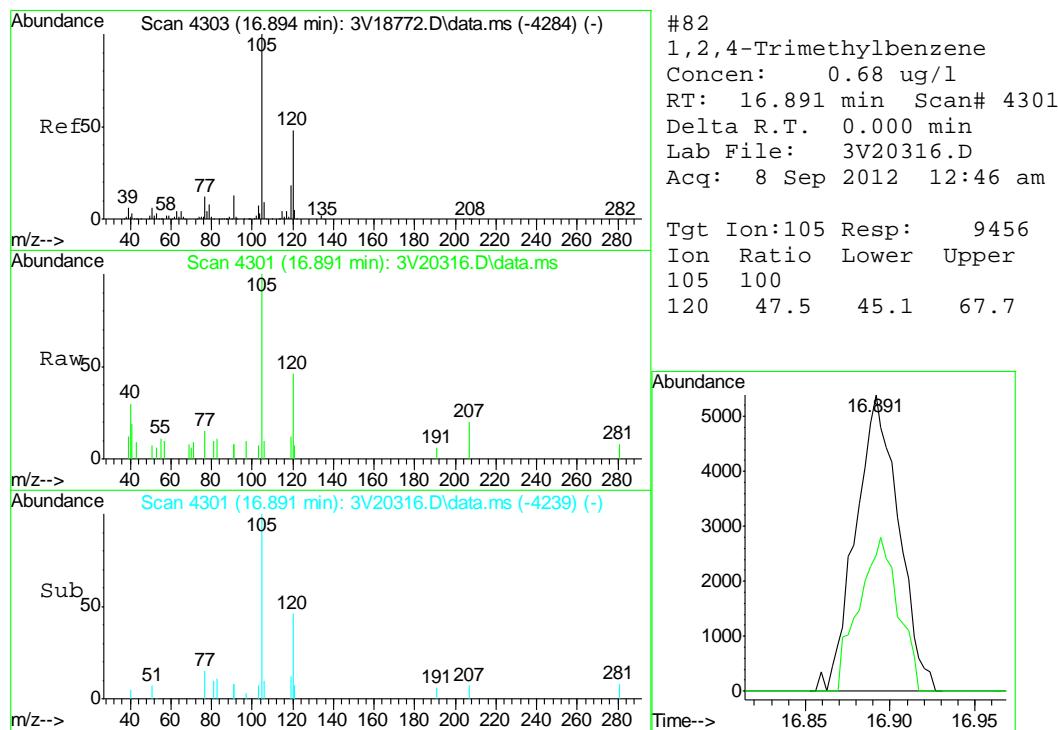
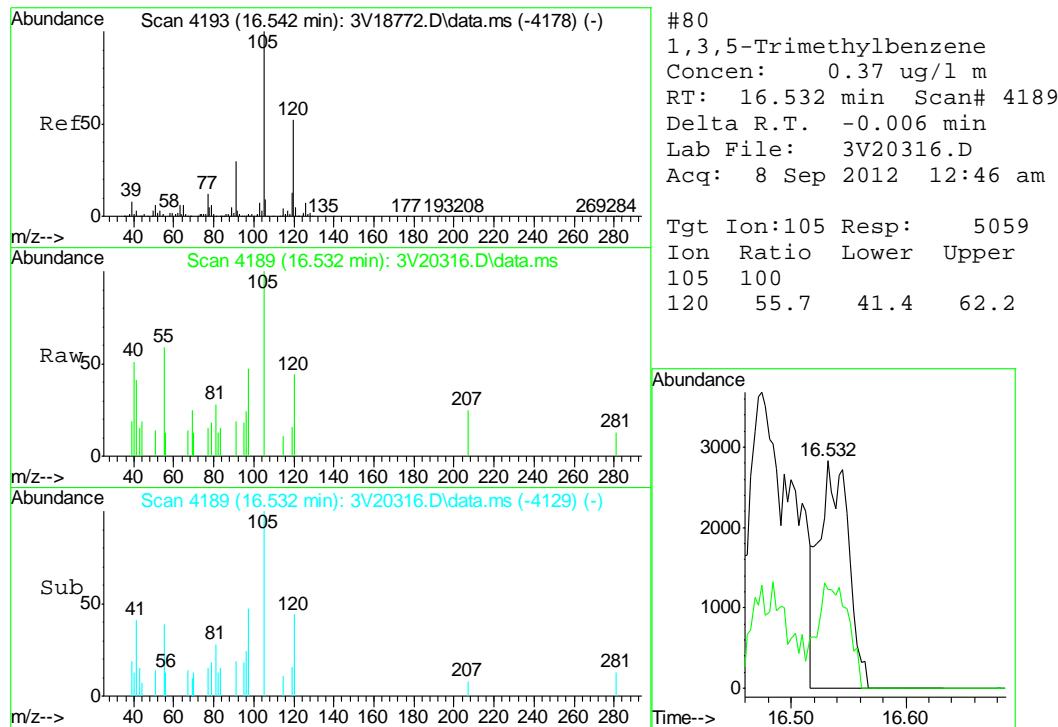


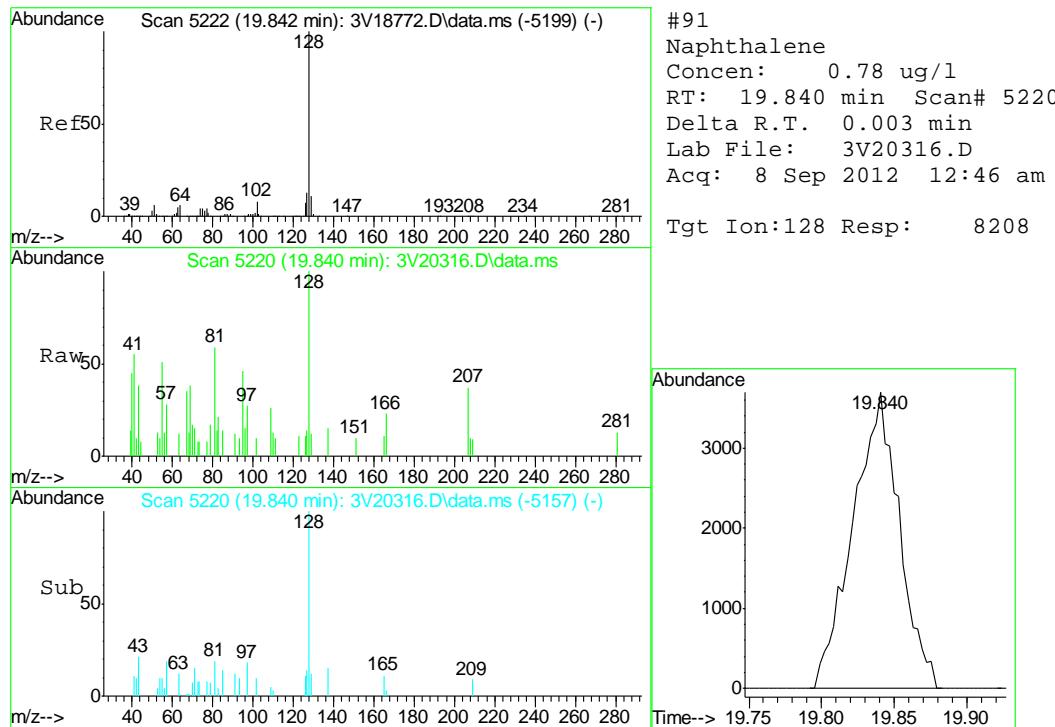












Quantitation Report (QT Reviewed)

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 Data File : 3V20317.D
 Acq On : 8 Sep 2012 1:17 am
 Operator : BRETD
 Sample : D38480-3
 Misc : MS4630,V3V1182,5.078,,100,5,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Sep 08 11:27:31 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1160TVH1160SOIL.M
 Quant Title : 8260
 QLast Update : Fri Aug 24 10:57:50 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.863	168	204526	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.659	114	333739	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.293	117	357536	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.282	152	207505	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.248	102	25137	54.59	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	109.18%	
61) Toluene-d8	14.051	98	444203	47.59	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	95.18%	
69) 4-Bromofluorobenzene	16.243	95	190372	52.12	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	104.24%	

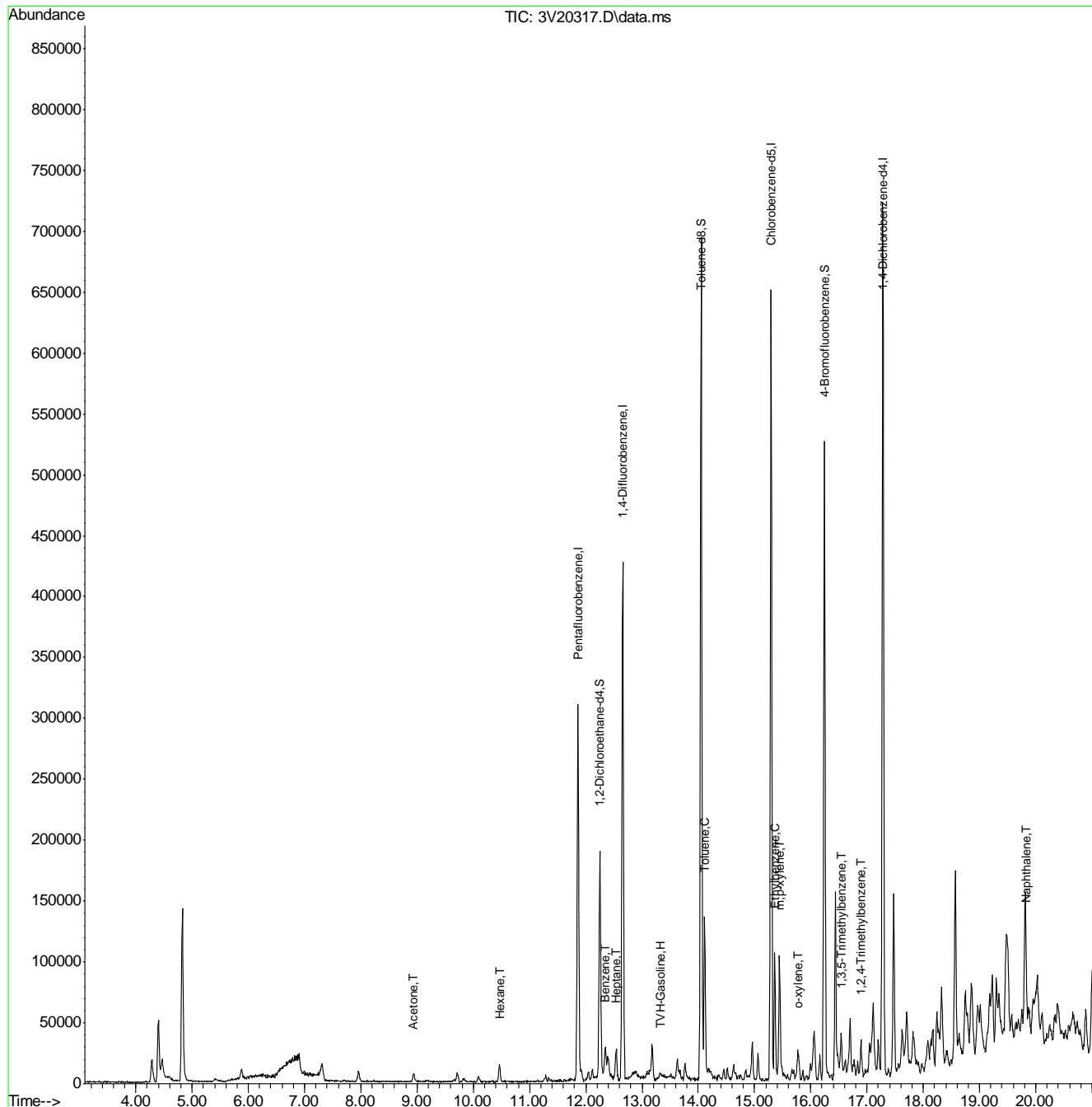
Target Compounds					Qvalue
1) TVH-Gasoline	13.329	TIC	2100016m	73.77	ug/l
15) Acetone	8.936	58	3167	4.45	ug/l # 71
41) Hexane	10.464	57	6518	1.24	ug/l 100
43) Heptane	12.533	43	12109	1.81	ug/l 95
50) Benzene	12.347	78	16105	1.34	ug/l 100
62) Toluene	14.112	92	33715	3.72	ug/l 98
66) Ethylbenzene	15.364	91	6397	0.38	ug/l 88
72) m,p-xylene	15.447	106	26731	3.98	ug/l 90
73) o-xylene	15.787	106	3013	0.59	ug/l # 71
80) 1,3,5-Trimethylbenzene	16.538	105	9176	0.66	ug/l 88
82) 1,2,4-Trimethylbenzene	16.894	105	13094	0.93	ug/l 87
91) Naphthalene	19.837	128	5929	0.56	ug/l 100

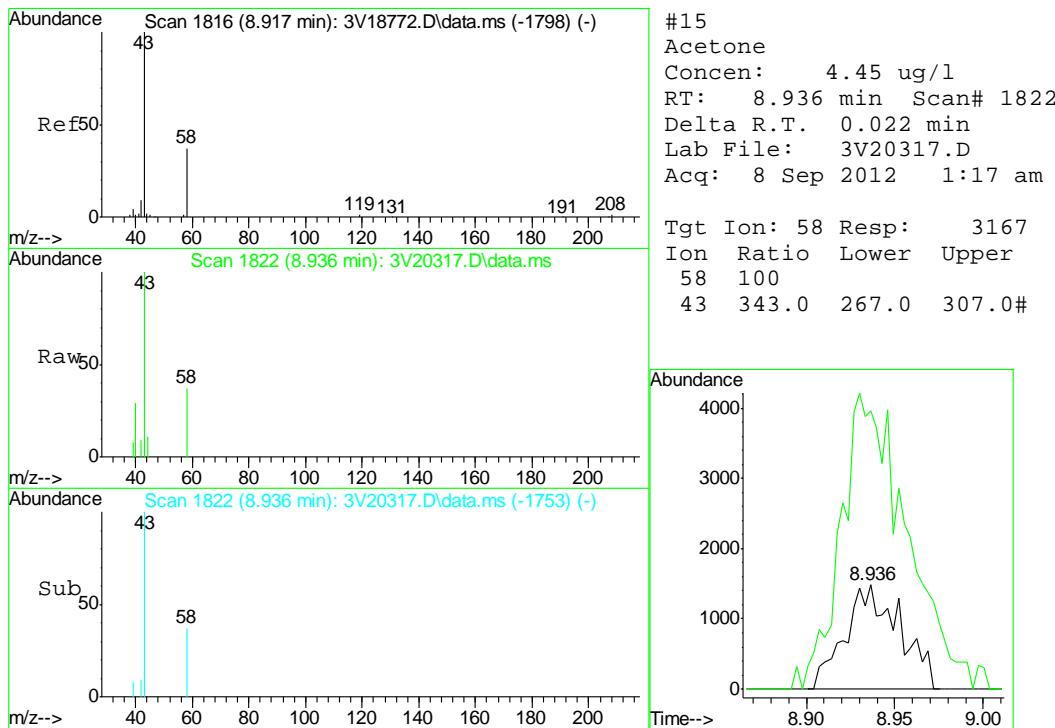
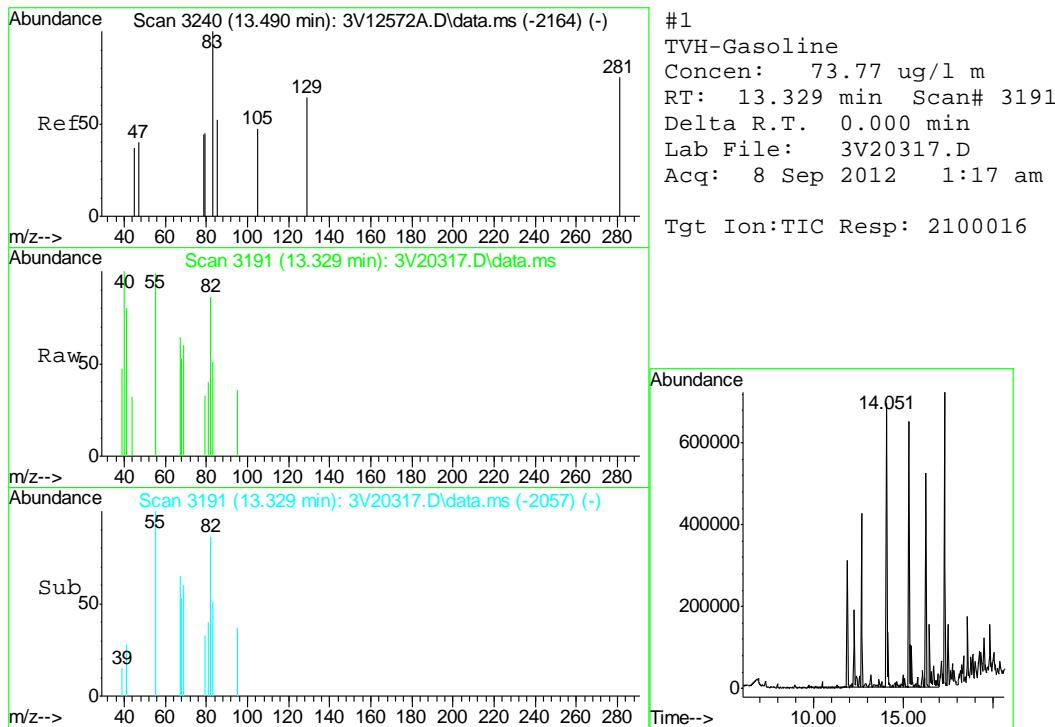
(#) = qualifier out of range (m) = manual integration (+) = signals summed

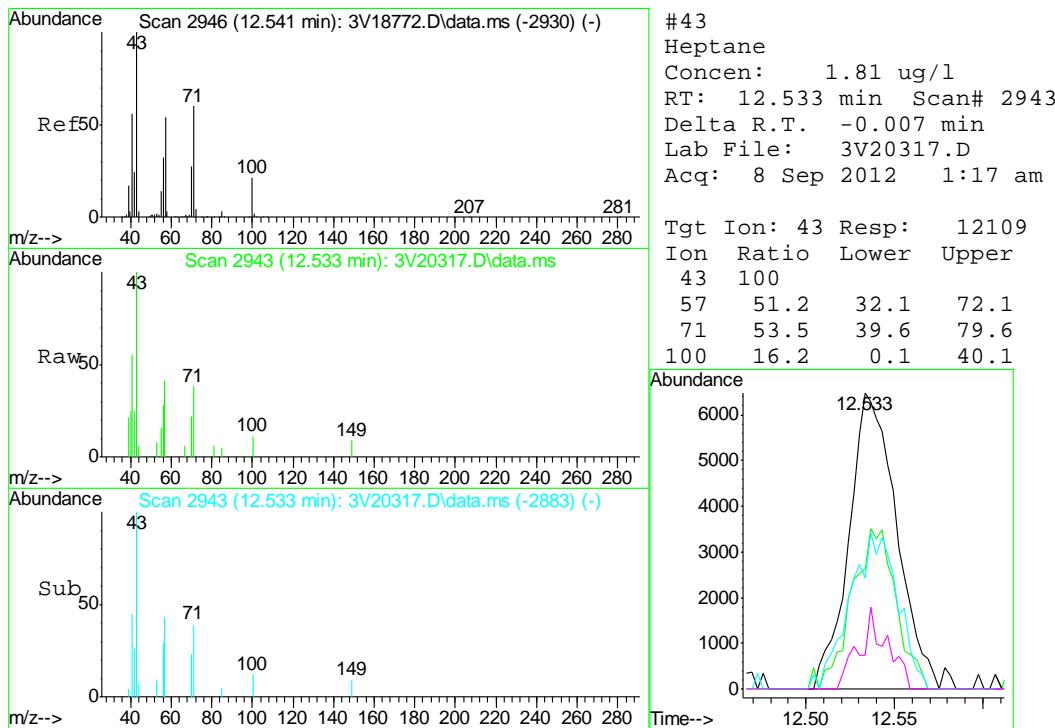
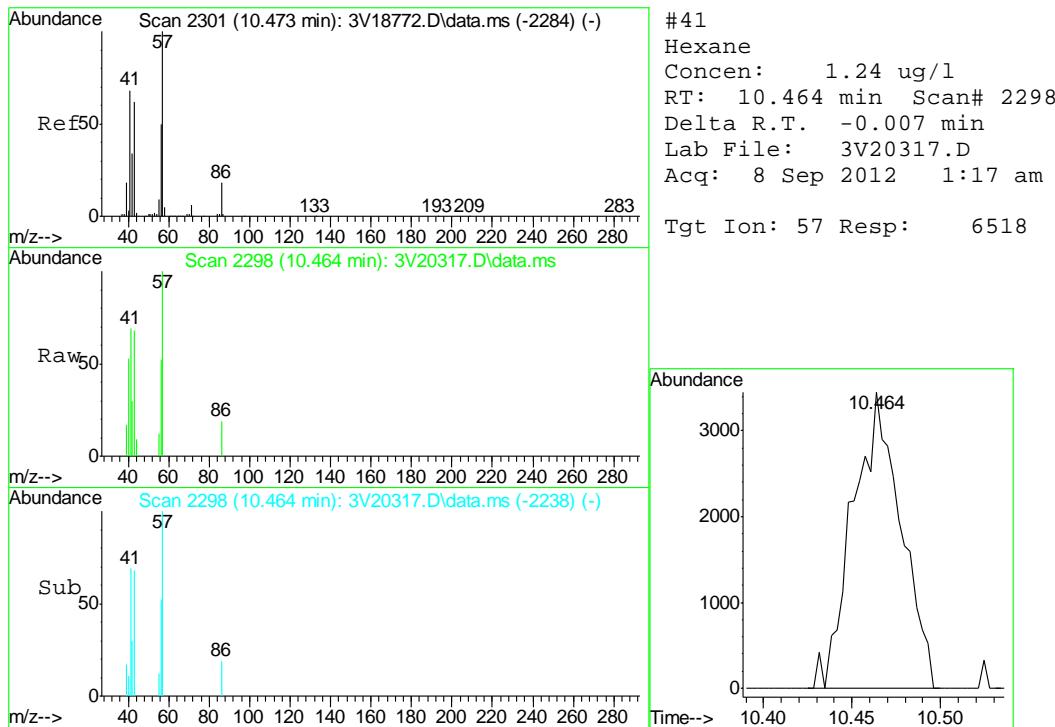
Quantitation Report (QT Reviewed)

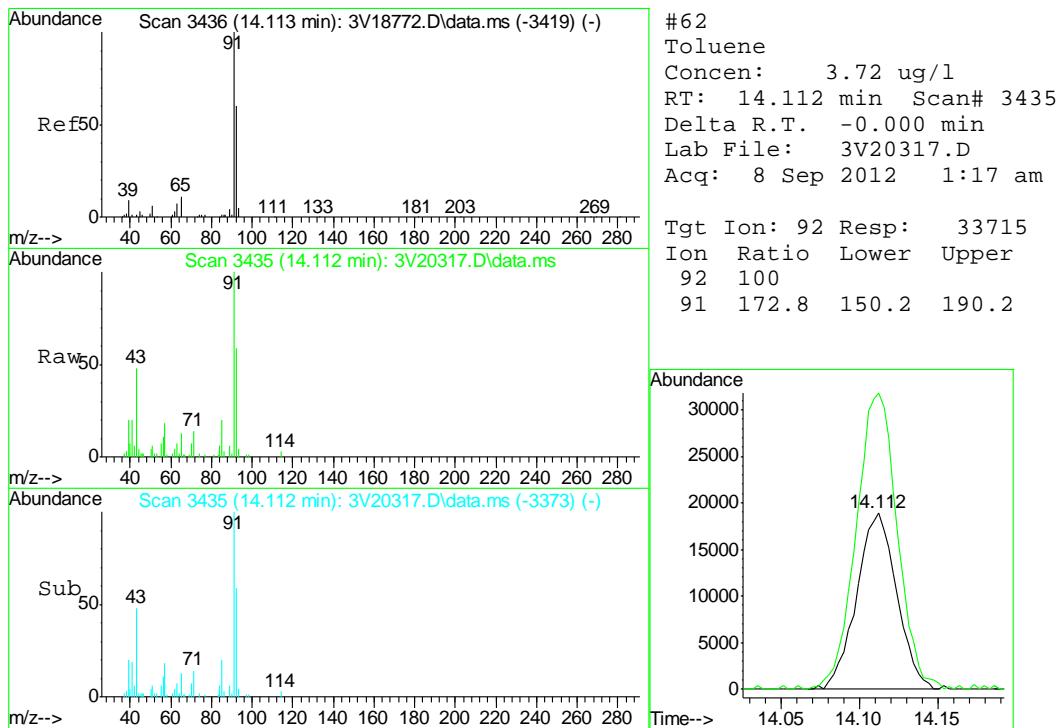
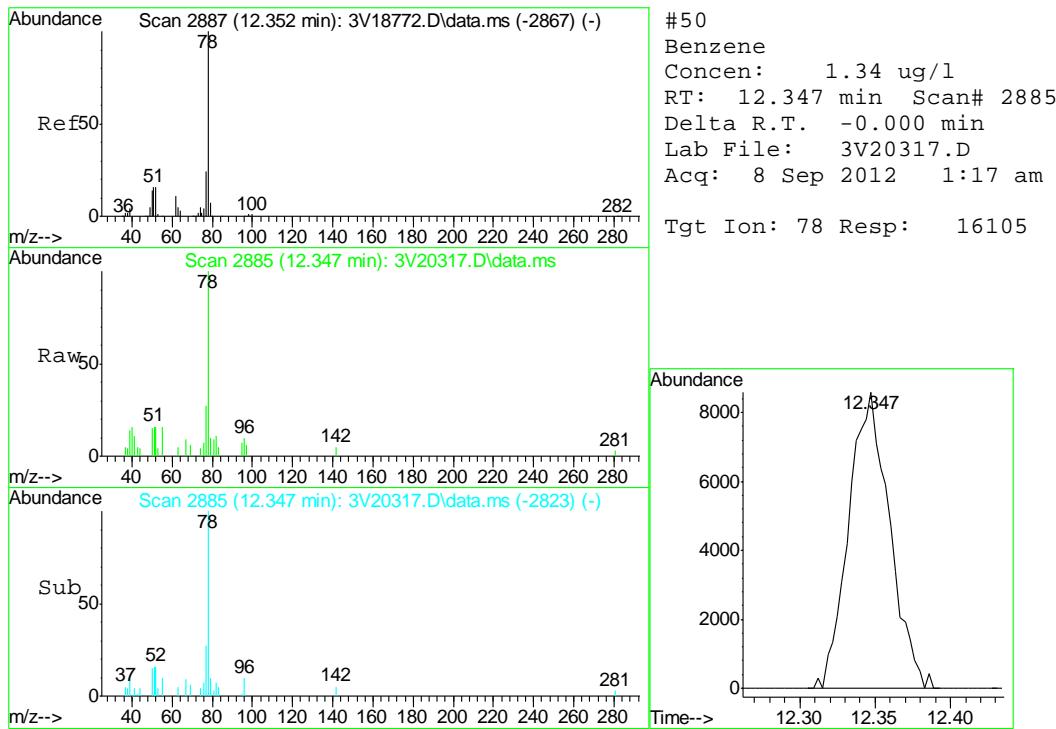
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 Data File : 3V20317.D
 Acq On : 8 Sep 2012 1:17 am
 Operator : BRETD
 Sample : D38480-3
 Misc : MS4630,V3V1182,5.078,,100,5,1
 ALS Vial : 19 Sample Multiplier: 1

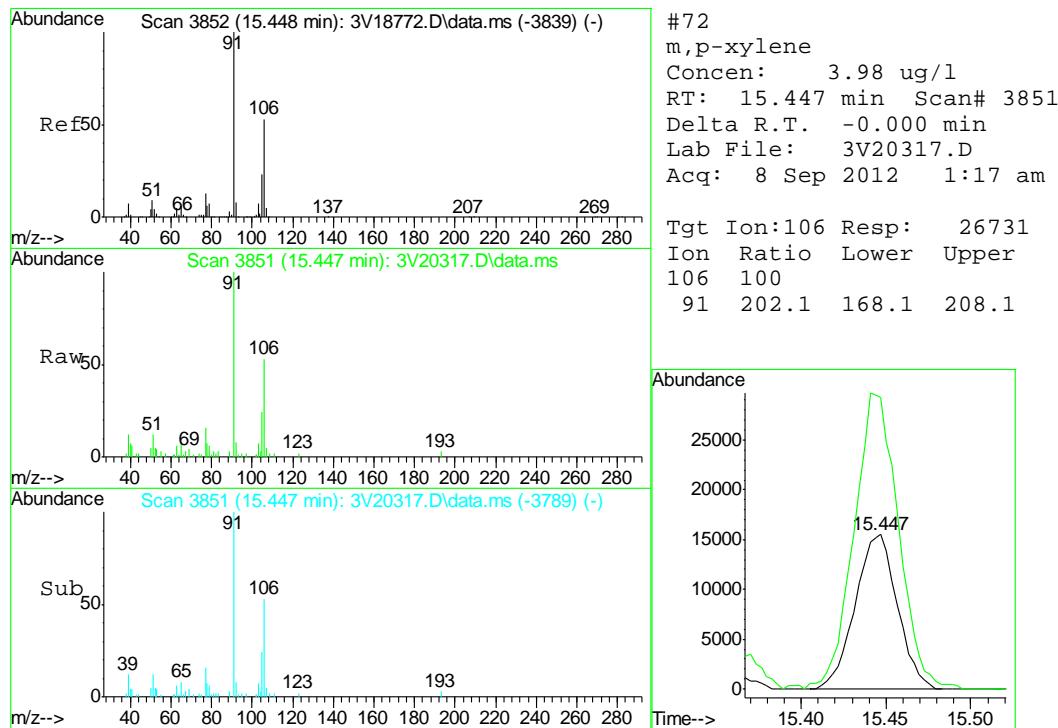
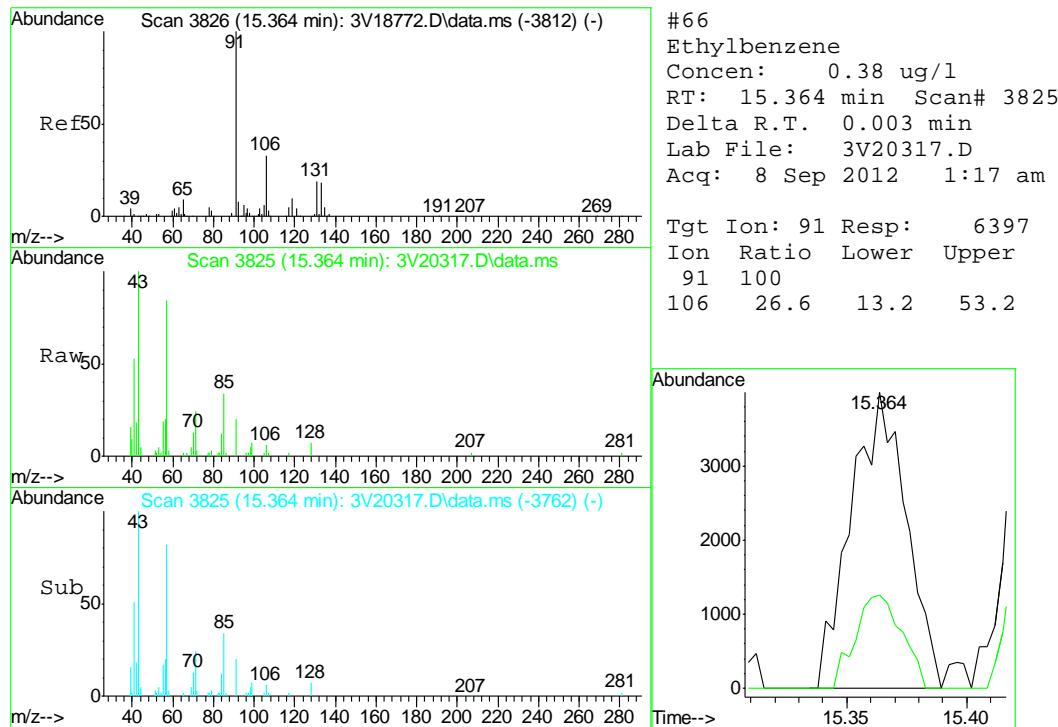
Quant Time: Sep 08 11:27:31 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1160TVH1160SOIL.M
 Quant Title : 8260
 QLast Update : Fri Aug 24 10:57:50 2012
 Response via : Initial Calibration

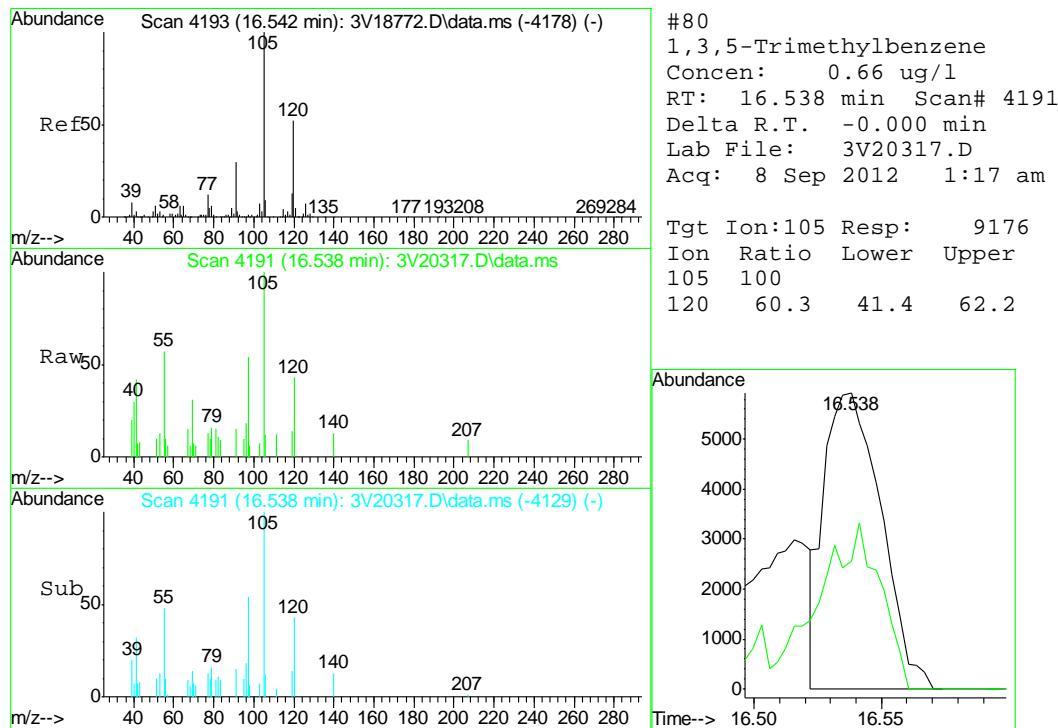
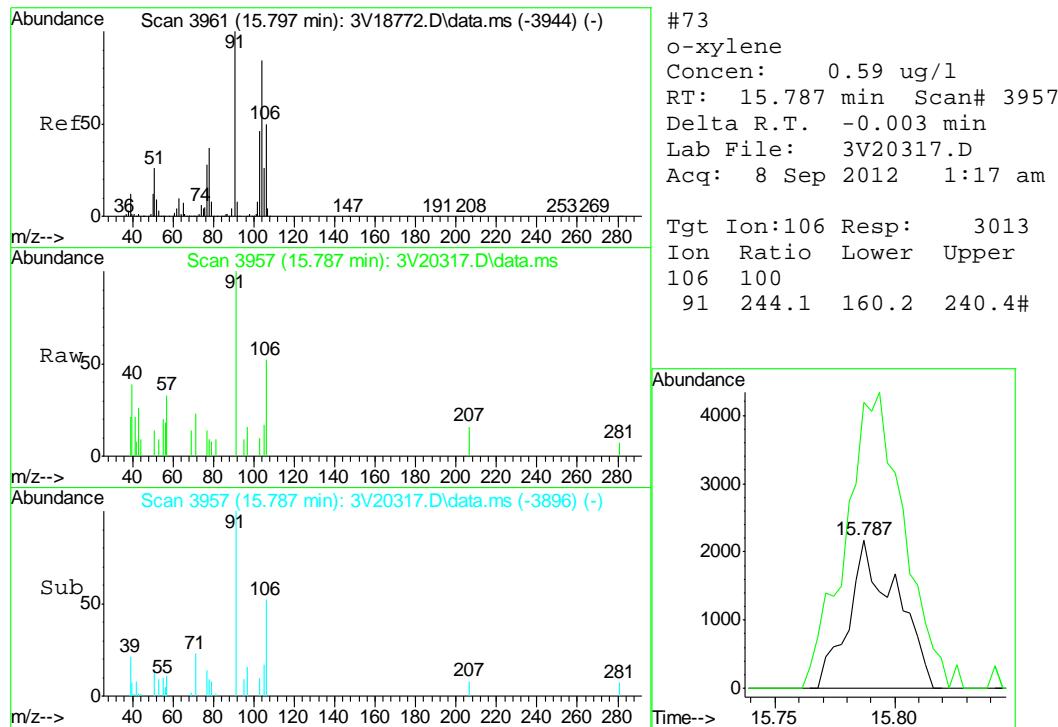


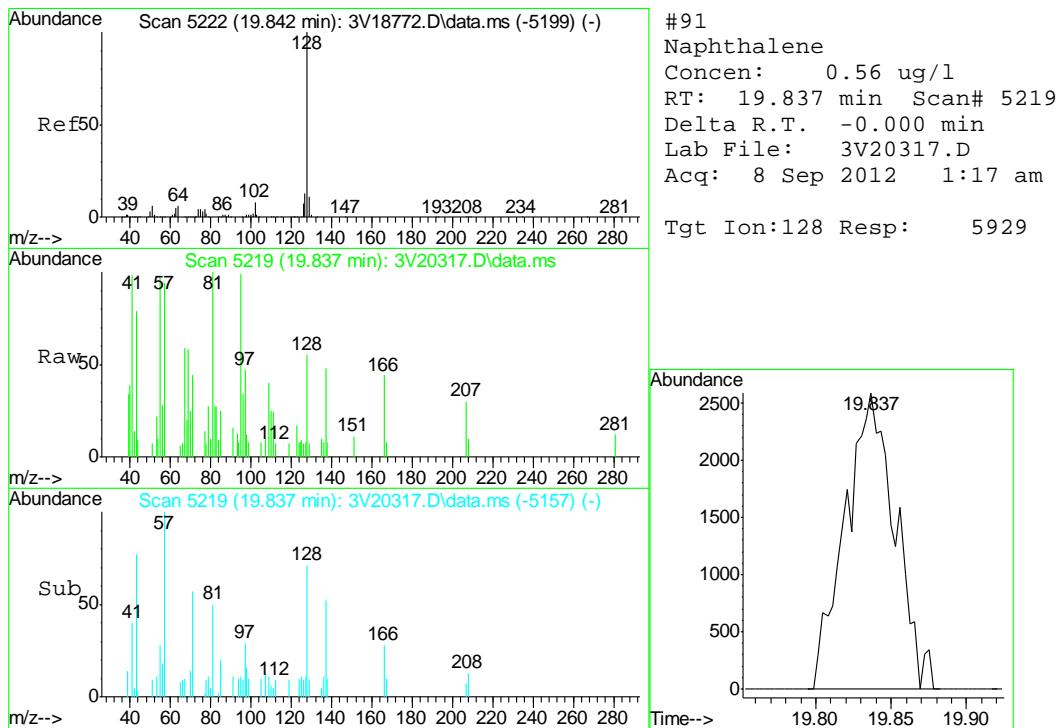
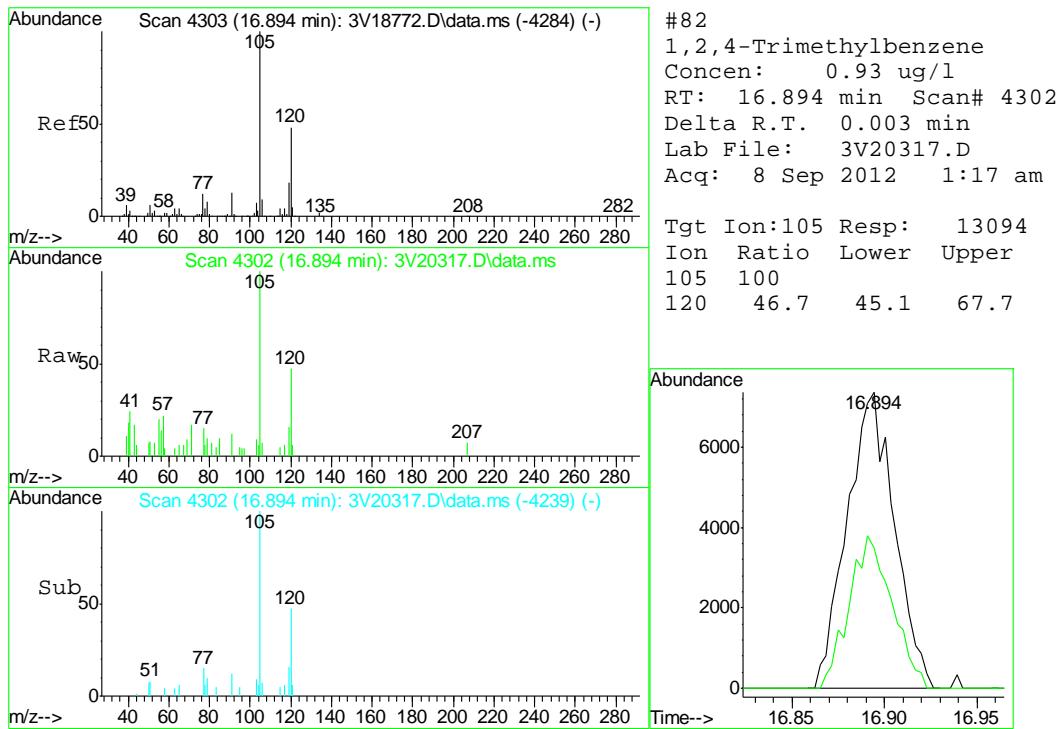












Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3090712.S\
 Data File : 3V20302.D
 Acq On : 7 Sep 2012 5:31 pm
 Operator : BRETD
 Sample : MB
 Misc : MS4630,V3V1182,5.00,,100,5,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Sep 08 10:51:42 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1160TVH1160SOIL.M
 Quant Title : 8260
 QLast Update : Fri Aug 24 10:57:50 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.863	168	214790	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.659	114	349912	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.294	117	344887	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.283	152	191769	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.255	102	27742	57.36	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	114.72%
61) Toluene-d8	14.052	98	460707	51.17	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	102.34%
69) 4-Bromofluorobenzene	16.243	95	170109	48.28	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	96.56%

Target Compounds					Qvalue
1) TVH-Gasoline	13.329	TIC	1798871m	63.19	ug/l
62) Toluene	14.110	92	2192	0.25	ug/l
91) Naphthalene	19.847	128	8104	0.82	ug/l
					100

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

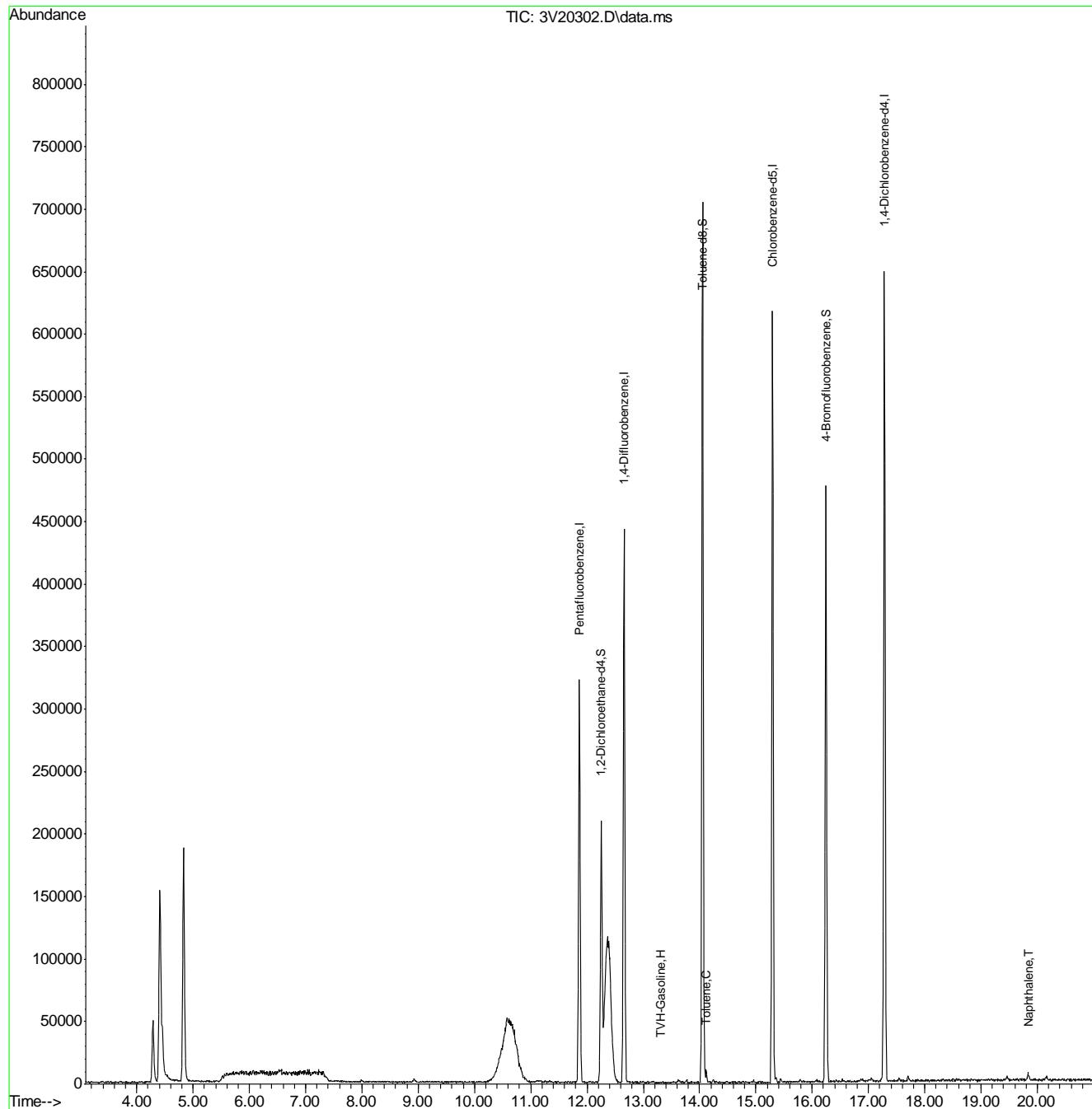
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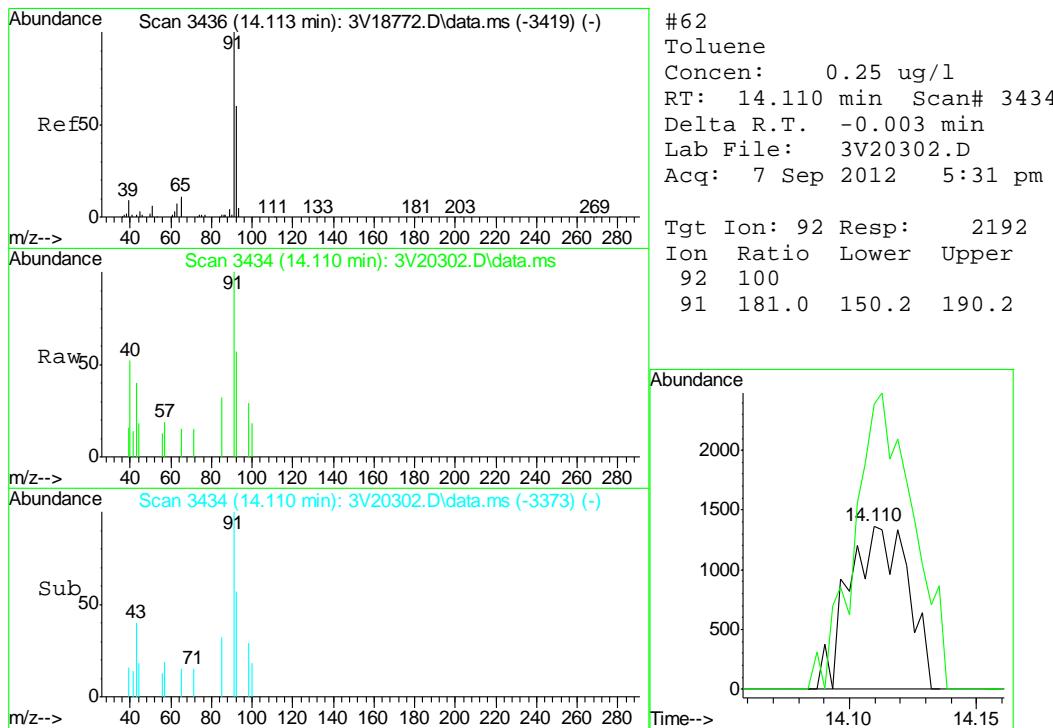
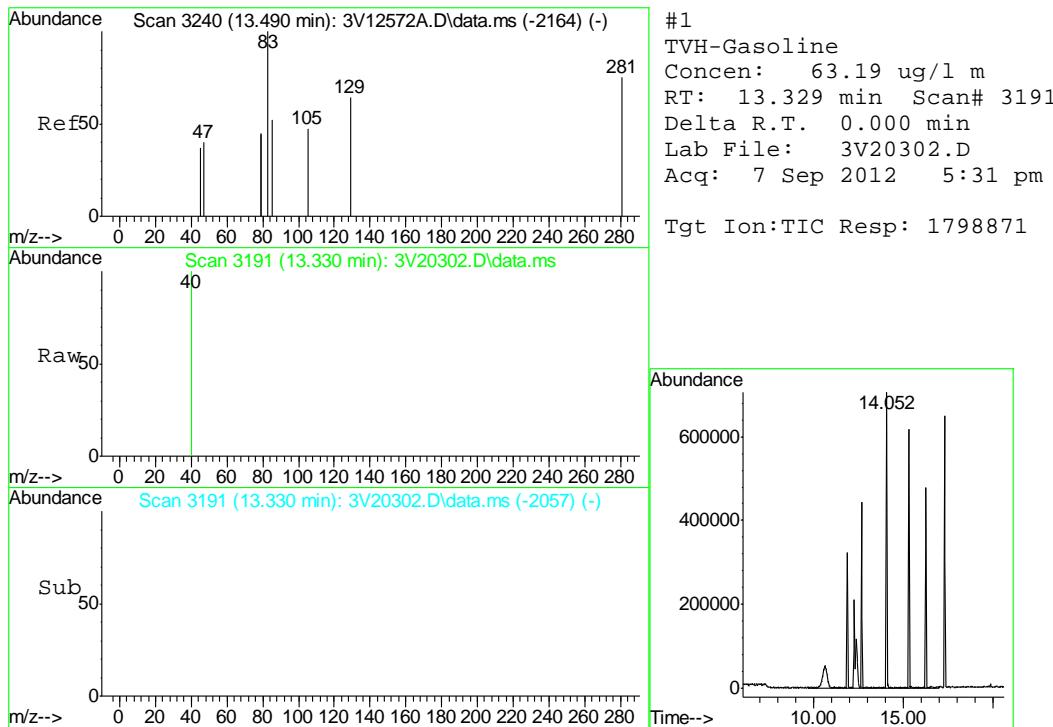
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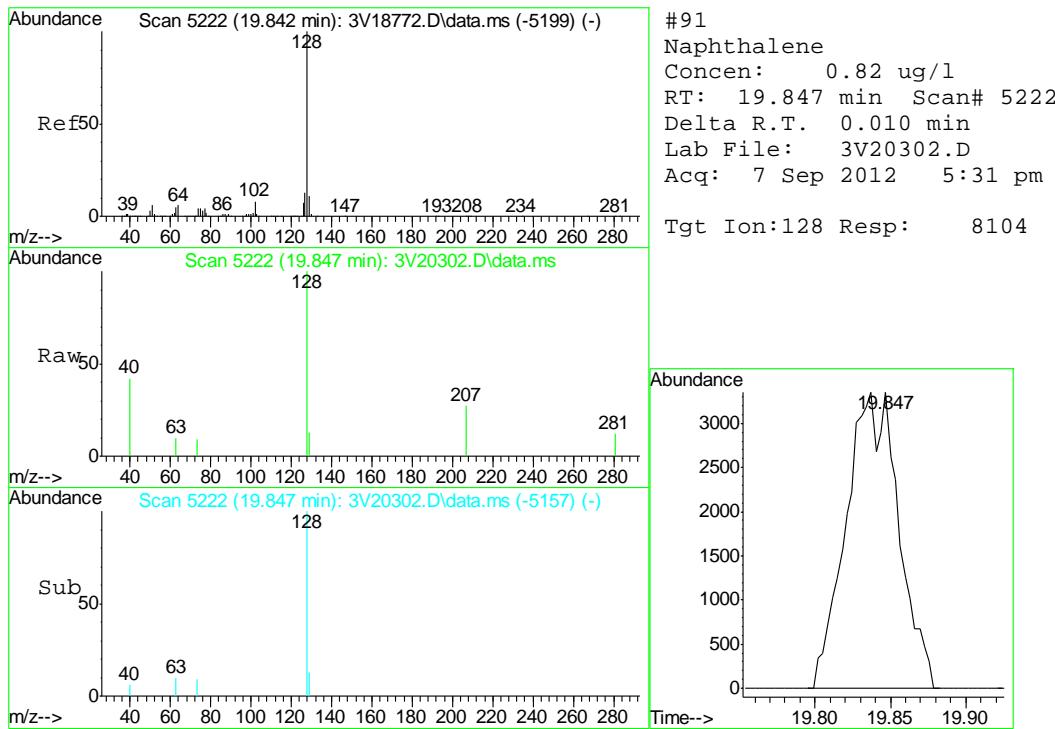
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3090712.S\
 Data File : 3V20302.D
 Acq On : 7 Sep 2012 5:31 pm
 Operator : BRETD
 Sample : MB
 Misc : MS4630,V3V1182,5.00,,100,5,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Sep 08 10:51:42 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1160TVH1160SOIL.M
 Quant Title : 8260
 QLast Update : Fri Aug 24 10:57:50 2012
 Response via : Initial Calibration









GC/MS Semi-volatiles

QC Data Summaries

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Includes the following where applicable:

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

Method Blank Summary

Page 1 of 1

Job Number: D38480
Account: XTOKRWR XTO Energy
Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6602-MB	3G11149.D	1	09/11/12	DC	09/11/12	OP6602	E3G518

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D38480-1, D38480-2, D38480-3

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	
50-32-8	Benzo(a)pyrene	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	
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	96%
321-60-8	2-Fluorobiphenyl	94%
1718-51-0	Terphenyl-d14	95%

Blank Spike Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6602-BS	3G11150.D	1	09/11/12	DC	09/11/12	OP6602	E3G518

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D38480-1, D38480-2, D38480-3

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	77.8	93	34-130
120-12-7	Anthracene	83.3	81.6	98	35-130
56-55-3	Benzo(a)anthracene	83.3	79.3	95	36-130
50-32-8	Benzo(a)pyrene	83.3	71.7	86	36-130
205-99-2	Benzo(b)fluoranthene	83.3	89.7	108	35-130
207-08-9	Benzo(k)fluoranthene	83.3	68.2	82	37-130
218-01-9	Chrysene	83.3	73.0	88	40-130
53-70-3	Dibenzo(a,h)anthracene	83.3	63.5	76	32-130
206-44-0	Fluoranthene	83.3	75.7	91	38-130
86-73-7	Fluorene	83.3	76.3	92	35-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	62.7	75	28-130
91-20-3	Naphthalene	83.3	77.0	92	35-130
129-00-0	Pyrene	83.3	81.8	98	37-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	89%	10-145%
321-60-8	2-Fluorobiphenyl	87%	10-130%
1718-51-0	Terphenyl-d14	87%	22-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6602-MS	3G11152.D	1	09/11/12	DC	09/11/12	OP6602	E3G518
OP6602-MSD	3G11153.D	1	09/11/12	DC	09/11/12	OP6602	E3G518
D38513-1	3G11151.D	1	09/11/12	DC	09/11/12	OP6602	E3G518

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D38480-1, D38480-2, D38480-3

CAS No.	Compound	D38513-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND		92.5	83.8	91	79.6	86	5	10-155/30
120-12-7	Anthracene	ND		92.5	91.1	98	88.8	96	3	10-155/30
56-55-3	Benzo(a)anthracene	ND		92.5	90.0	97	89.7	97	0	10-175/30
50-32-8	Benzo(a)pyrene	ND		92.5	77.4	84	78.0	84	1	10-164/30
205-99-2	Benzo(b)fluoranthene	ND		92.5	97.3	105	98.1	106	1	10-165/30
207-08-9	Benzo(k)fluoranthene	ND		92.5	71.3	77	67.7	73	5	10-178/30
218-01-9	Chrysene	ND		92.5	81.5	88	80.1	87	2	10-147/30
53-70-3	Dibenzo(a,h)anthracene	ND		92.5	70.0	76	69.8	76	0	10-144/30
206-44-0	Fluoranthene	ND		92.5	88.0	95	86.8	94	1	10-207/30
86-73-7	Fluorene	ND		92.5	86.5	94	82.3	89	5	10-163/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND		92.5	68.6	74	68.8	74	0	10-180/30
91-20-3	Naphthalene	ND		92.5	81.2	88	78.5	85	3	10-198/30
129-00-0	Pyrene	ND		92.5	94.3	102	93.1	101	1	10-189/30

CAS No.	Surrogate Recoveries	MS	MSD	D38513-1	Limits
4165-60-0	Nitrobenzene-d5	77%	77%	51%	10-145%
321-60-8	2-Fluorobiphenyl	75%	74%	49%	10-130%
1718-51-0	Terphenyl-d14	75%	78%	74%	22-130%

* = Outside of Control Limits.

8.3.1
8



GC/MS Semi-volatiles

Raw Data

Judy Nelson
 09/13/12 14:22

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\091212\
 Data File : 3g11215.D
 Acq On : 13 Sep 2012 3:58 am
 Operator : DONC
 Sample : D38480-1
 Misc : OP6602,E3G522,30.04,,,1,1
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Sep 13 13:09:16 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Sep 06 09:42:23 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.921	136	189738	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.639	164	98450	4.0000	ug/mL	0.00
15) Phenanthrene-d10	9.121	188	180850	4.0000	ug/mL	0.00
19) Chrysene-d12	11.759	240	108763	4.0000	ug/mL	0.00
24) Perylene-d12	13.199	264	47457	4.0000	ug/mL	0.02

System Monitoring Compounds

2) Nitrobenzene-d5	5.223	82	601257	32.2082	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	64.42%
7) 2-Fluorobiphenyl	6.978	172	1558338	38.0521	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	76.10%
21) Terphenyl-d14	10.712	244	536591	32.7432	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	65.48%

Target Compounds

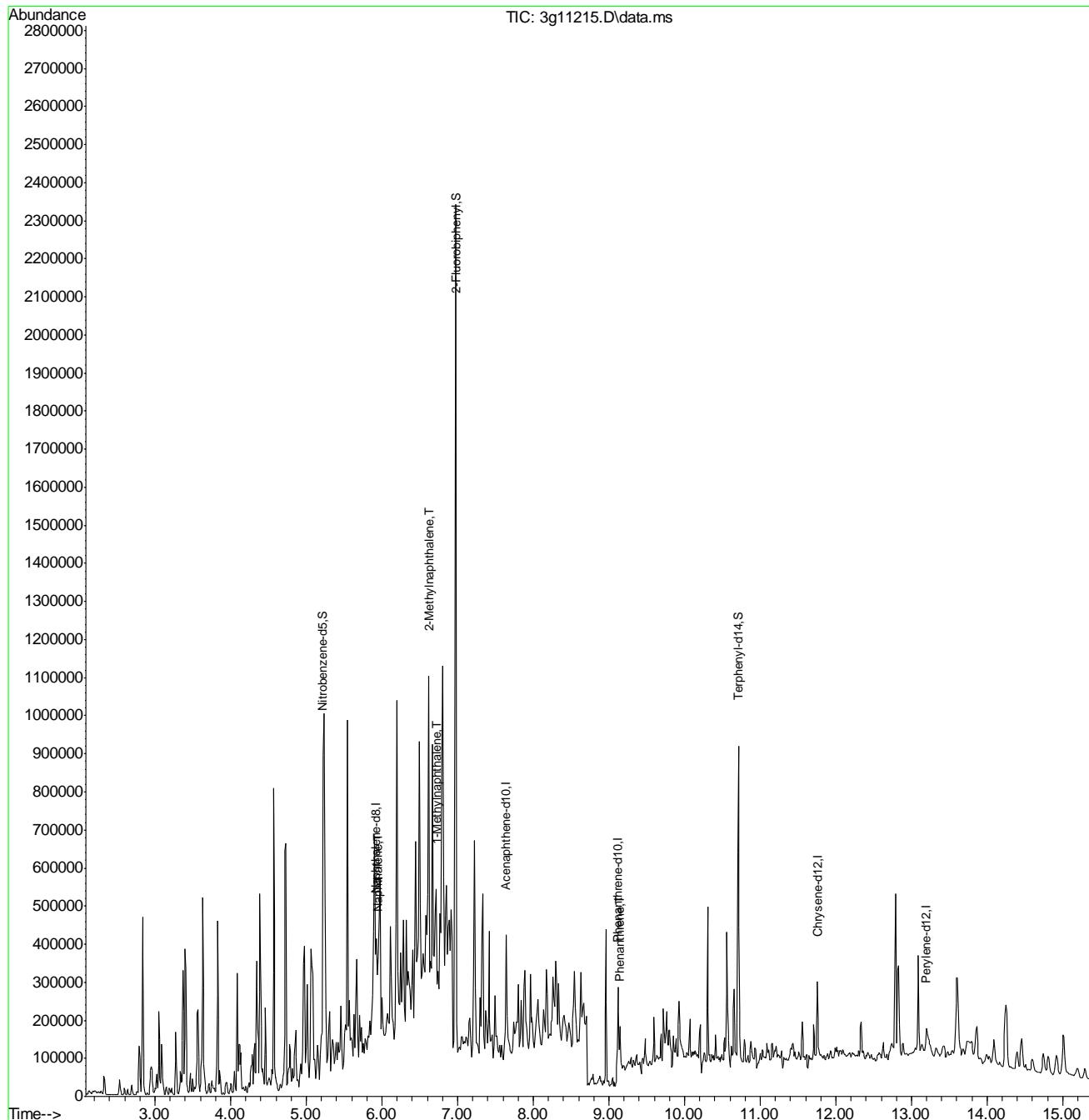
				Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D. d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.946	128	213583	4.0450 ug/mL# 65
8) 2-Methylnaphthalene	6.620	142	383150	13.1640 ug/mL 94
9) 1-Methylnaphthalene	6.719	142	139593m	4.6345 ug/mL
10) Acenaphthylene	0.000	152	0	N.D. d
11) Acenaphthene	0.000	154	0	N.D. d
12) Dibenzofuran	0.000	168	0	N.D. d
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	9.144	178	96156	1.5156 ug/mL# 12
17) Anthracene	0.000	178	0	N.D. d
18) Fluoranthene	0.000	202	0	N.D. d
20) Pyrene	0.000	202	0	N.D. d
22) Benzo(a)anthracene	0.000	228	0	N.D. d
23) Chrysene	0.000	228	0	N.D. d
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d
26) Benzo(k)fluoranthene	0.000	252	0	N.D. d
27) Benzo(a)pyrene	0.000	252	0	N.D. d
28) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D. d
29) Dibenz(a,h)anthracene	0.000	278	0	N.D. d
30) Benzo(g,h,i)perylene	0.000	276	0	N.D. d

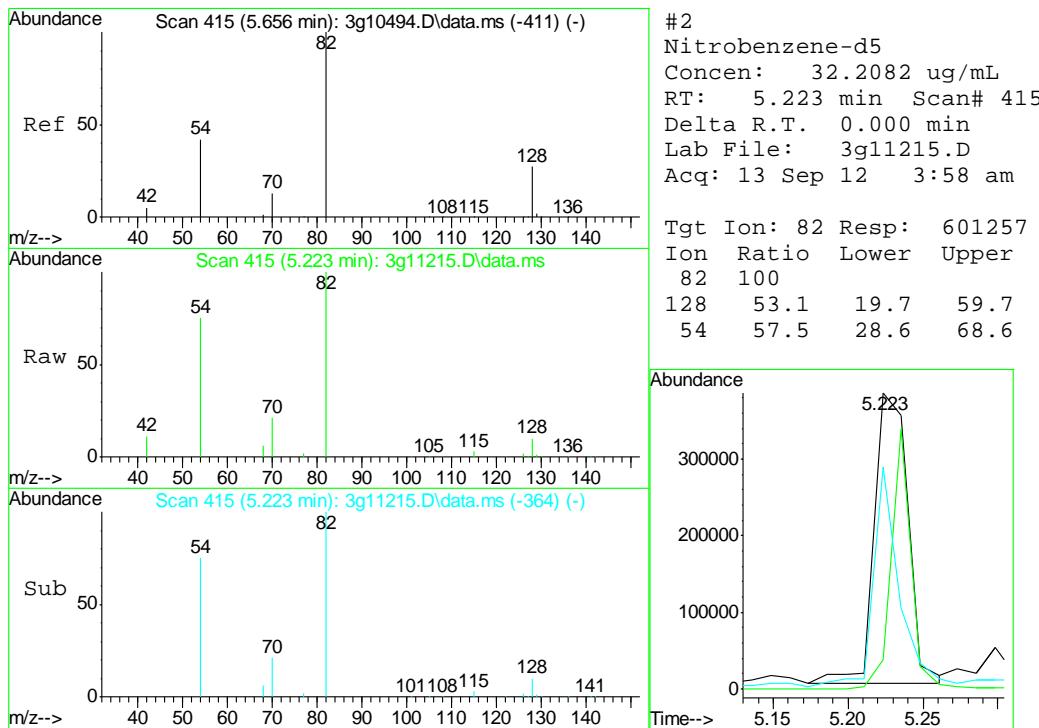
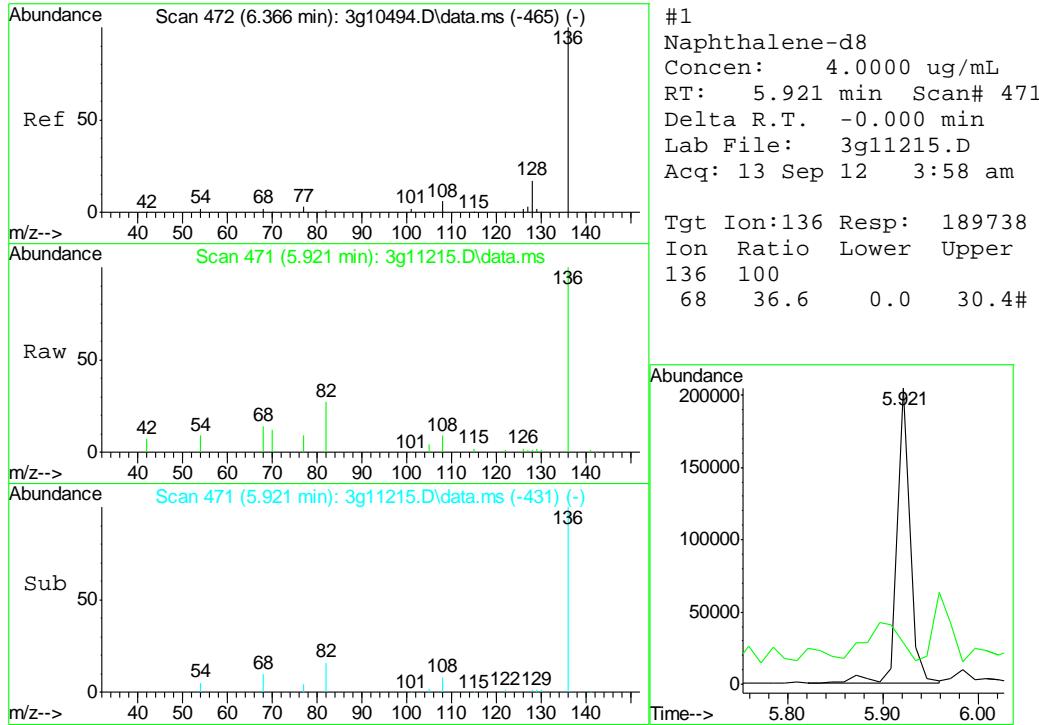
(#) = qualifier out of range (m) = manual integration (+) = signals summed

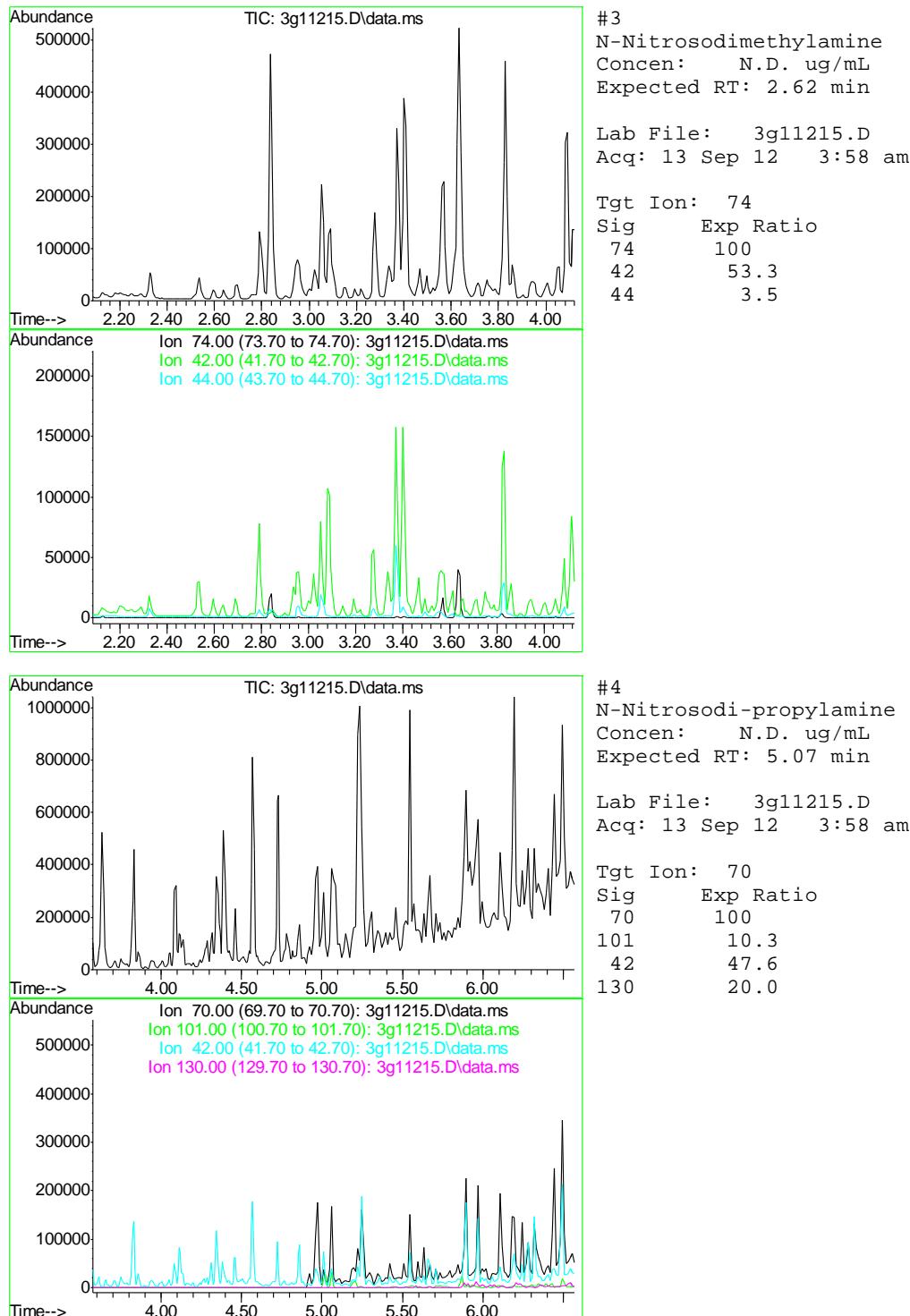
Quantitation Report (QT Reviewed)

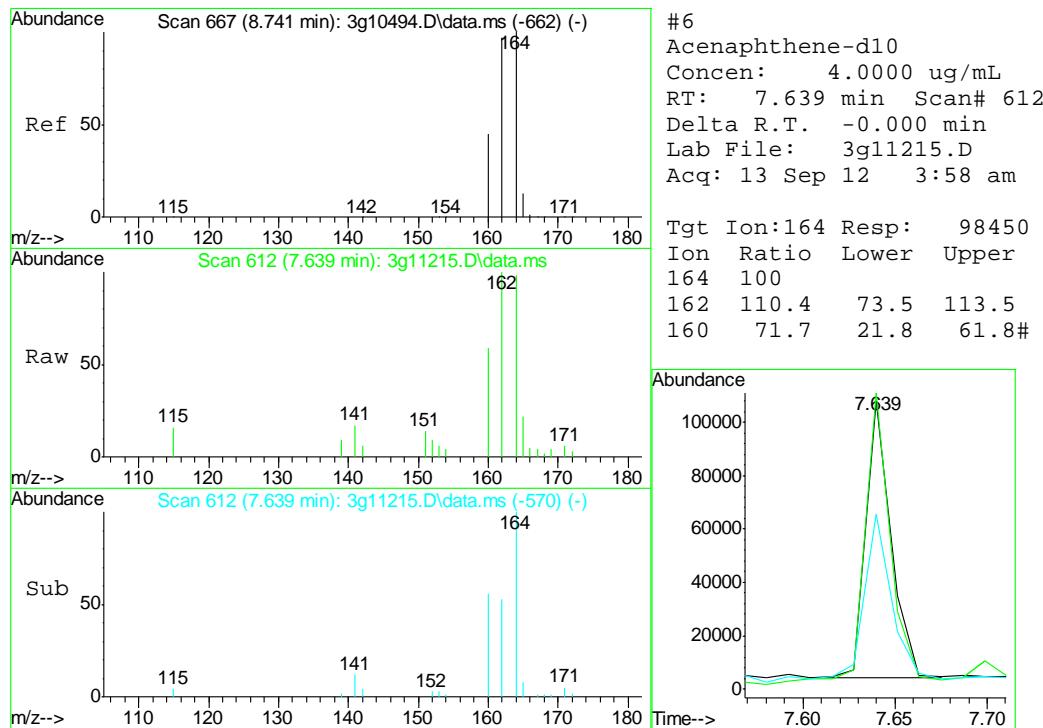
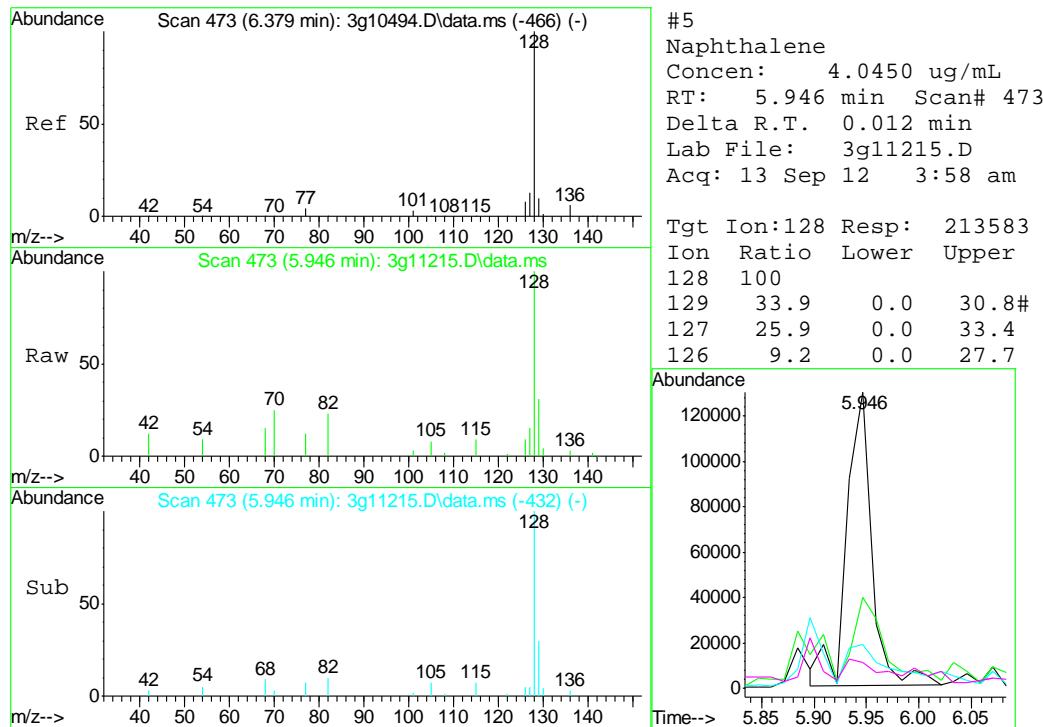
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 Operator : DONC
 Sample : D38480-1
 Misc : OP6602,E3G522,30.04,,,1,1
 ALS Vial : 26 Sample Multiplier: 1

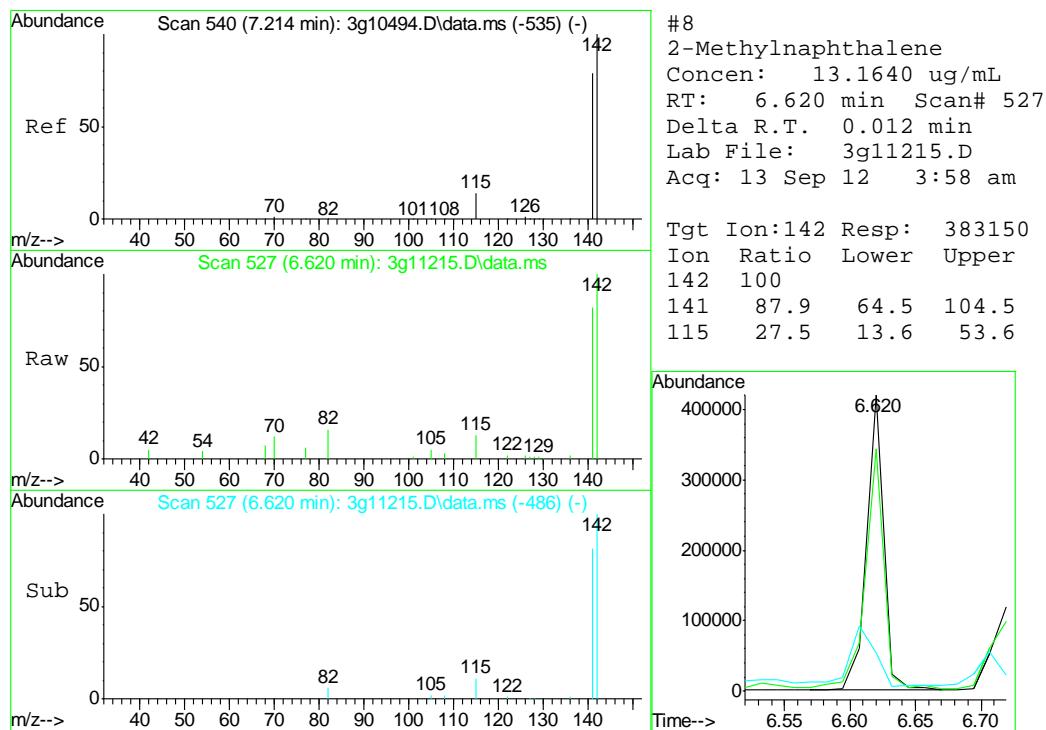
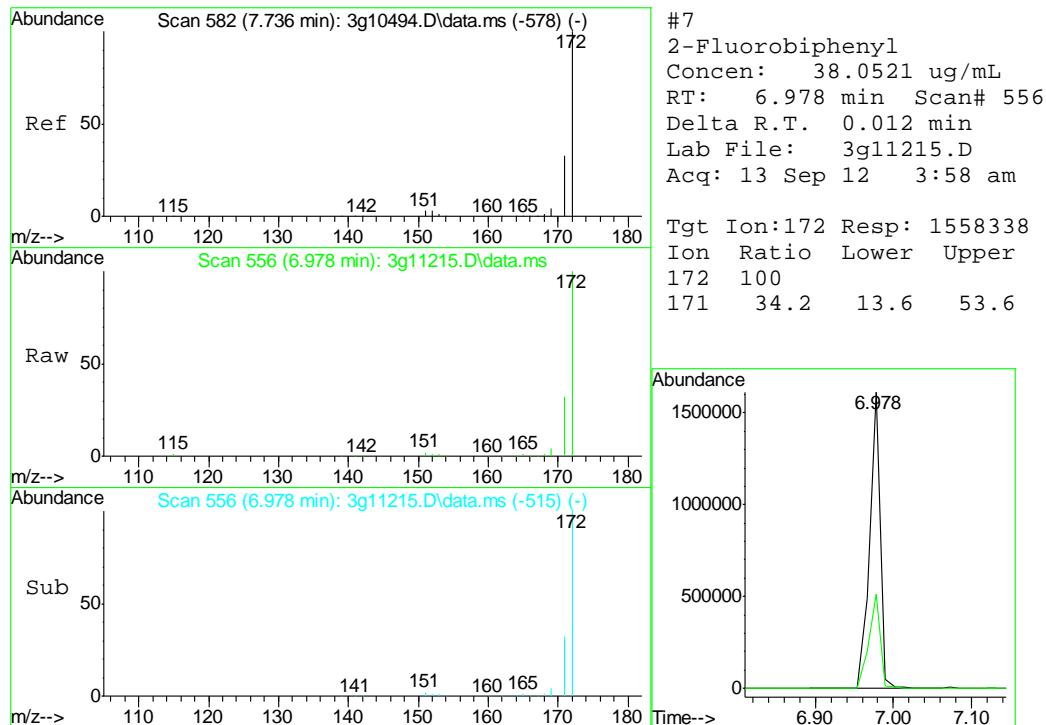
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 Quant Title : PAHSIM BASE
 QLast Update : Thu Sep 06 09:42:23 2012
 Response via : Initial Calibration

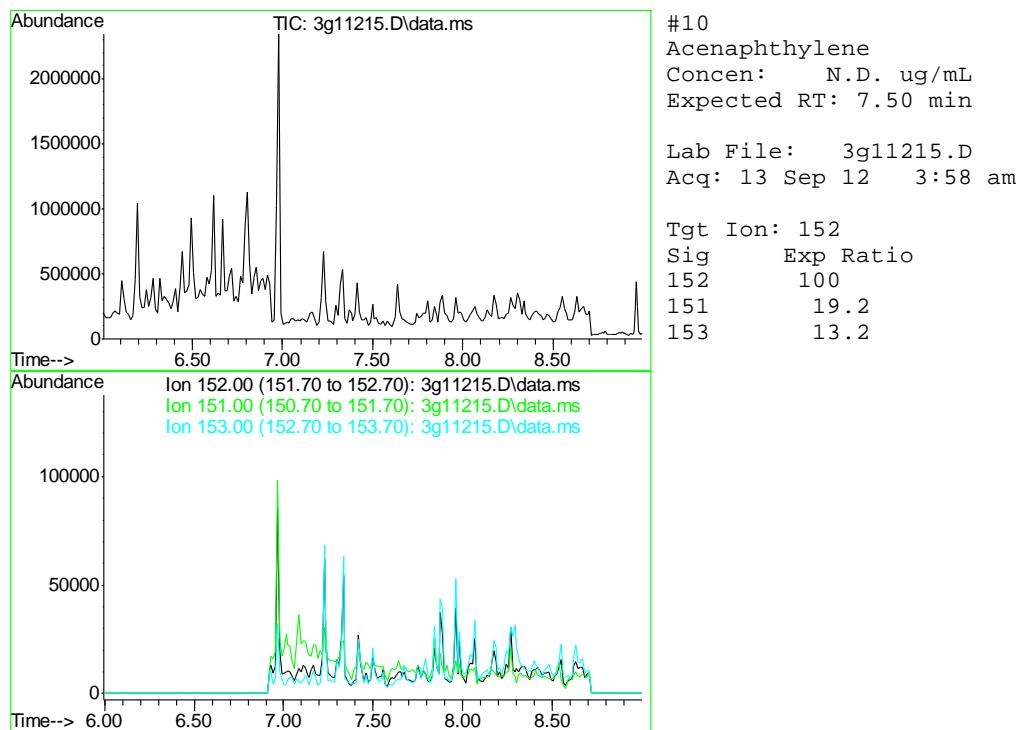
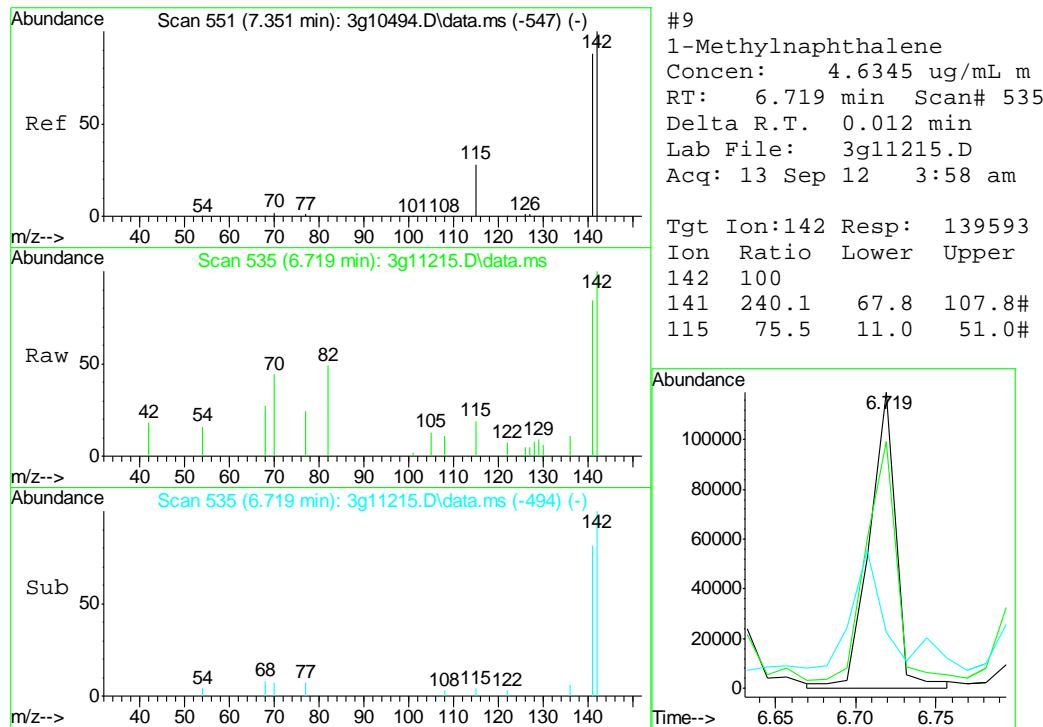


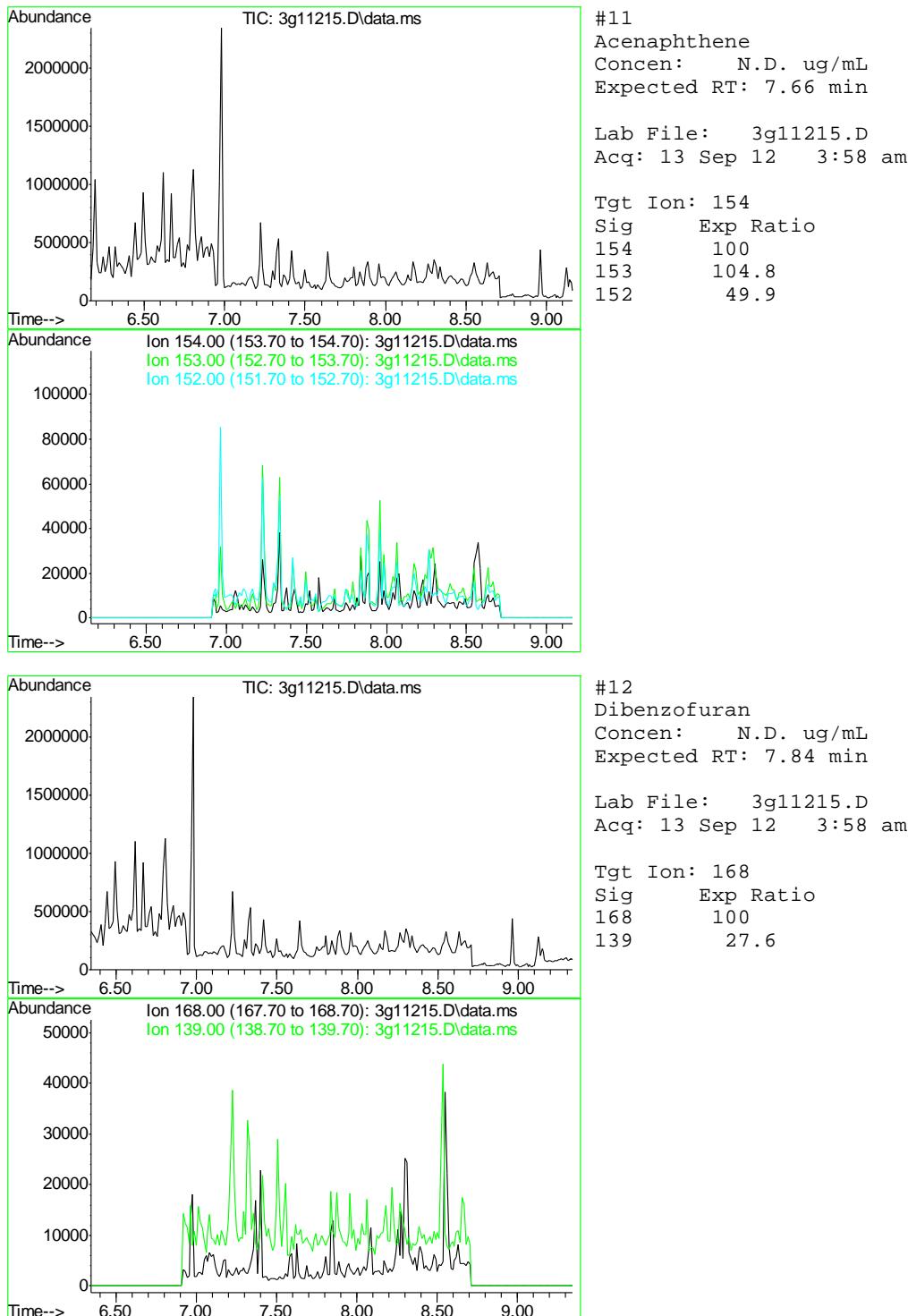


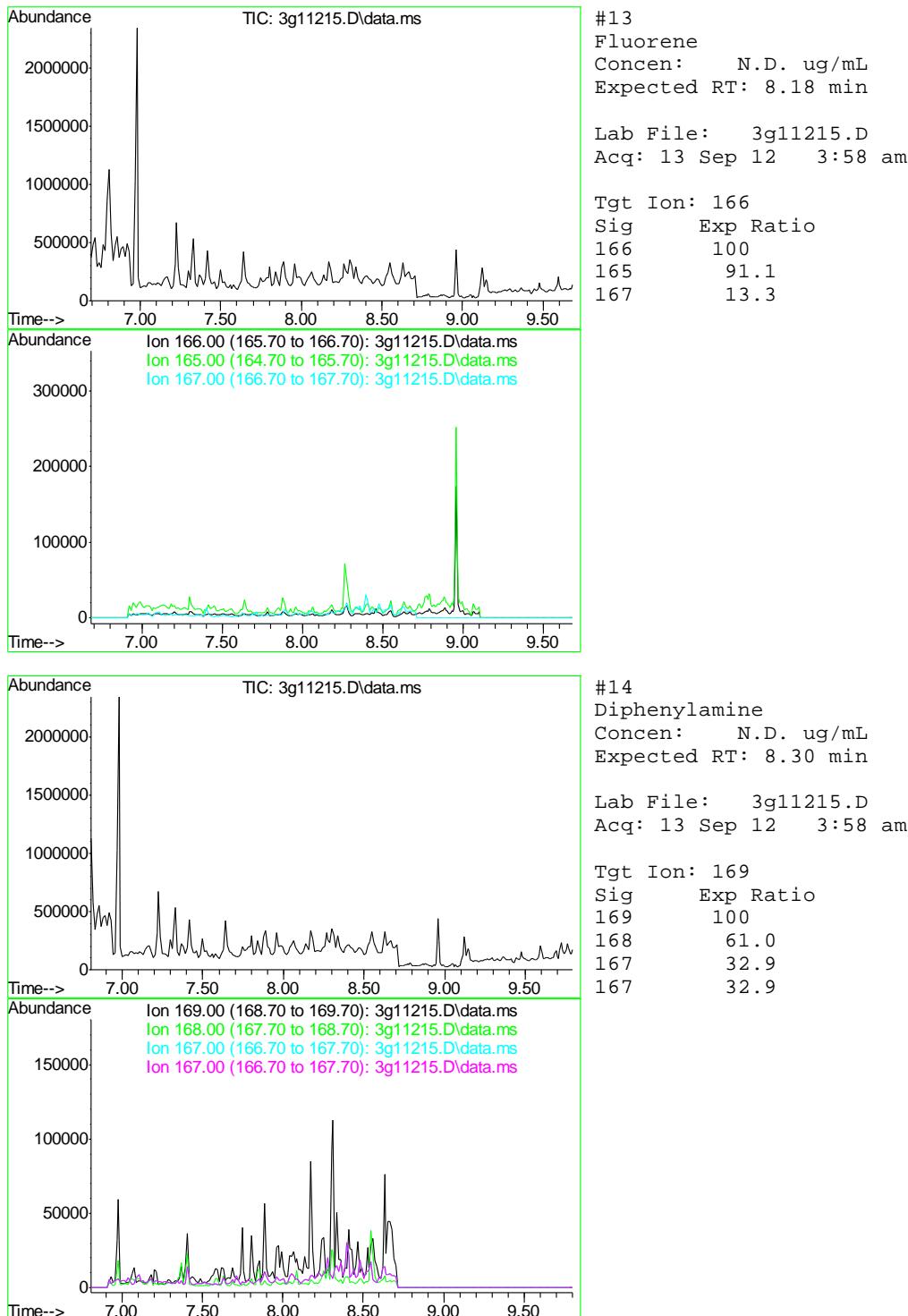


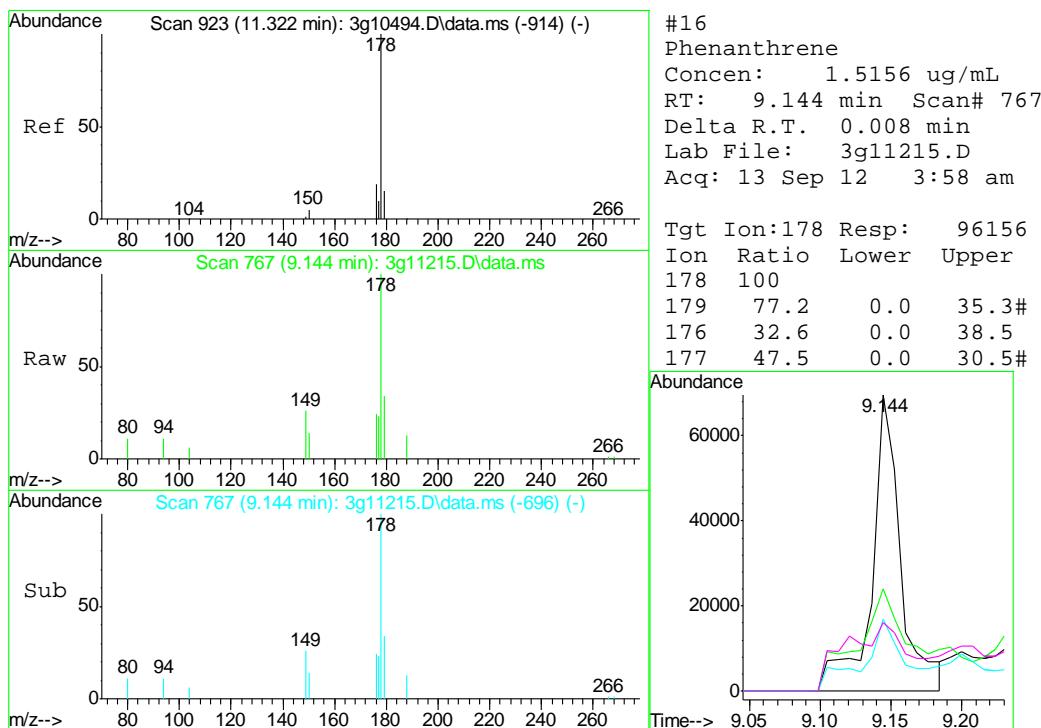
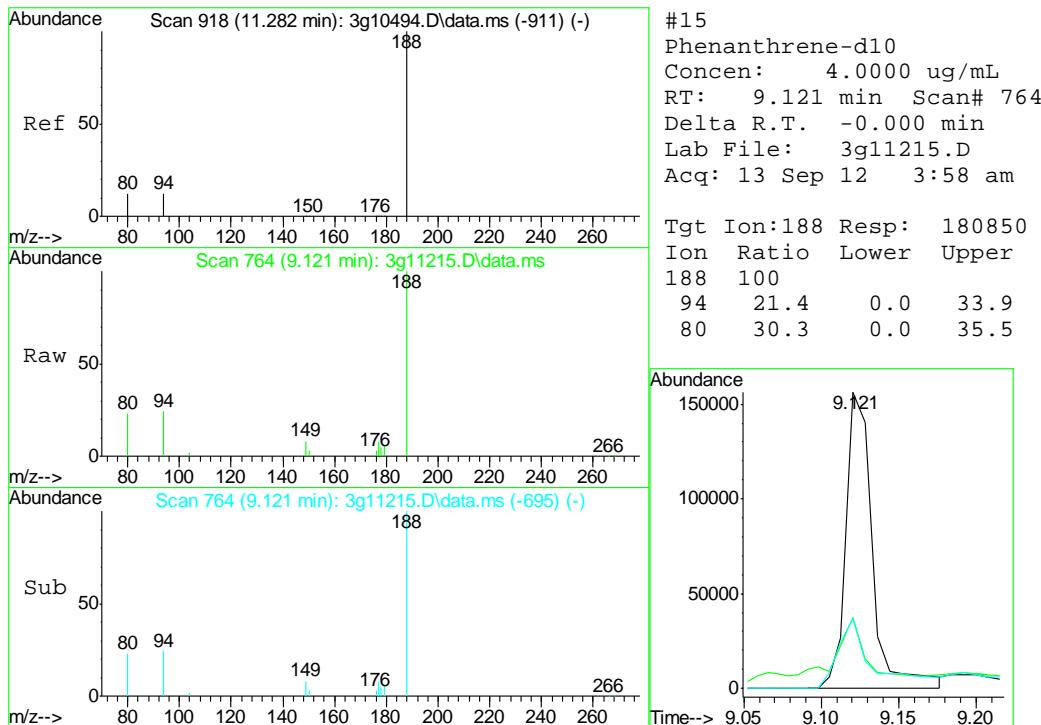


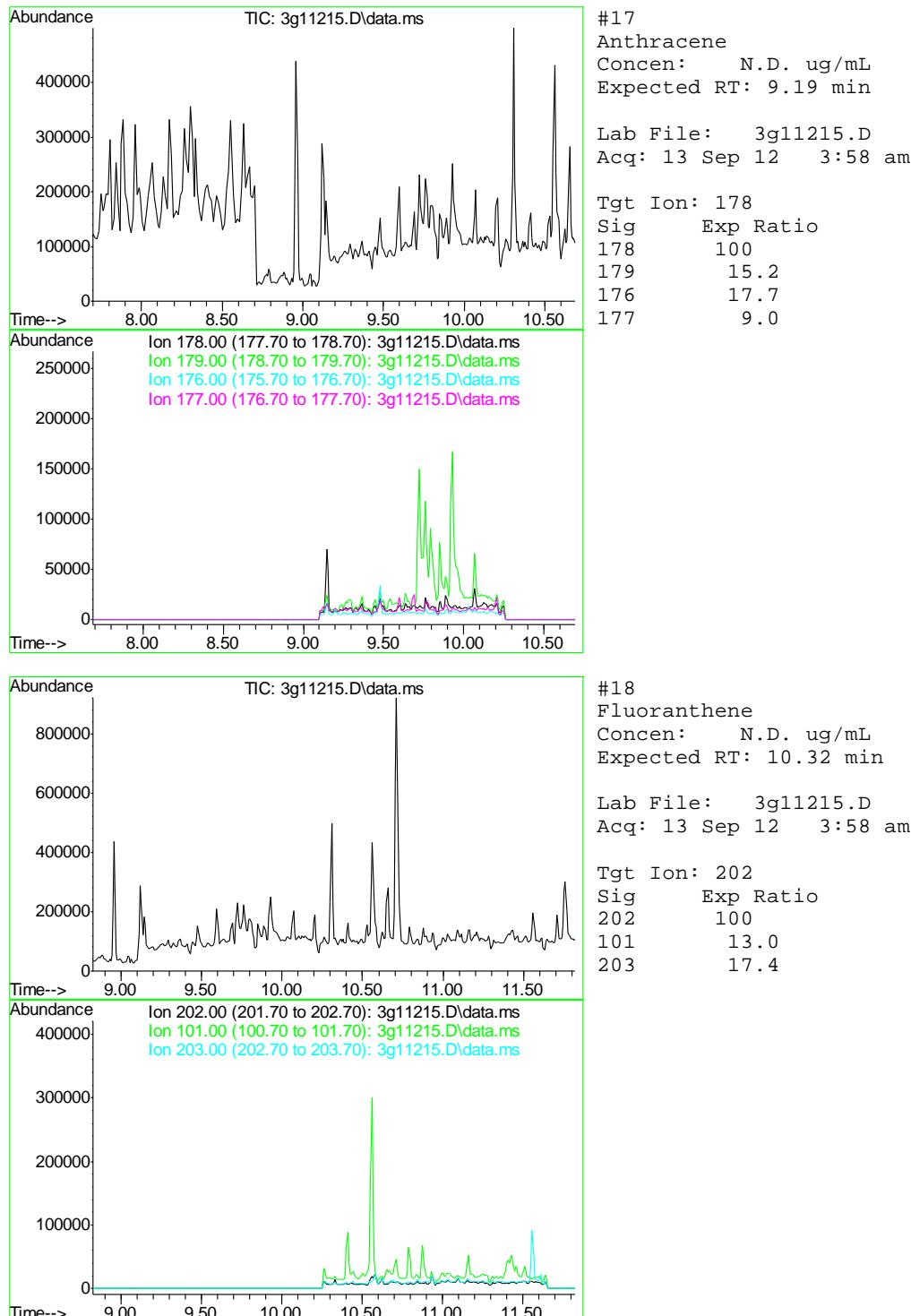


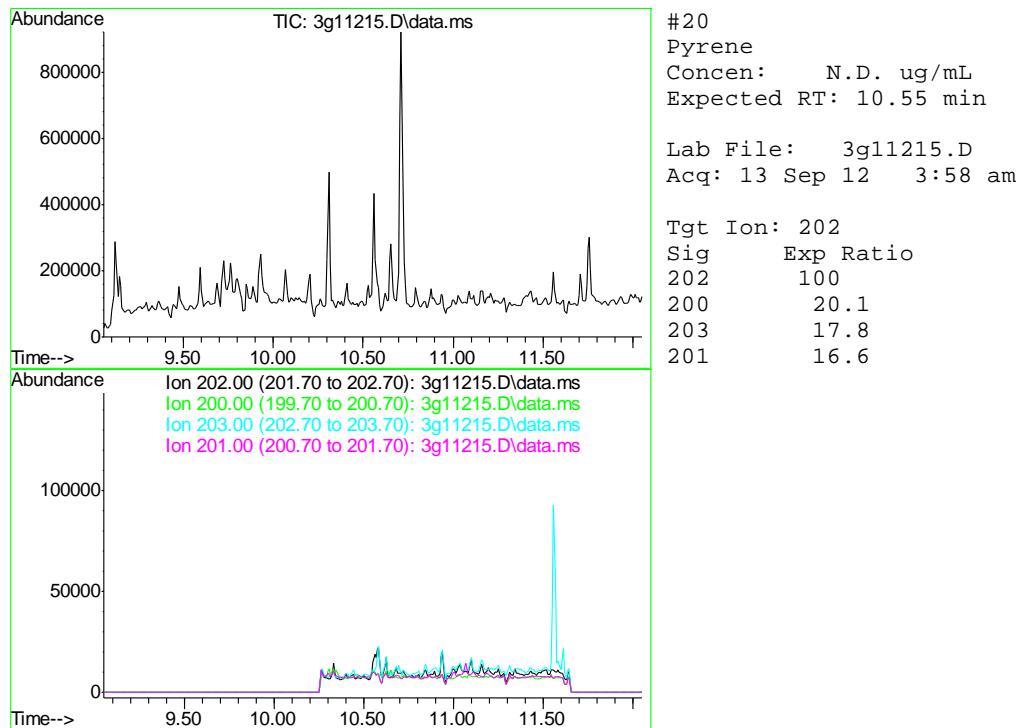
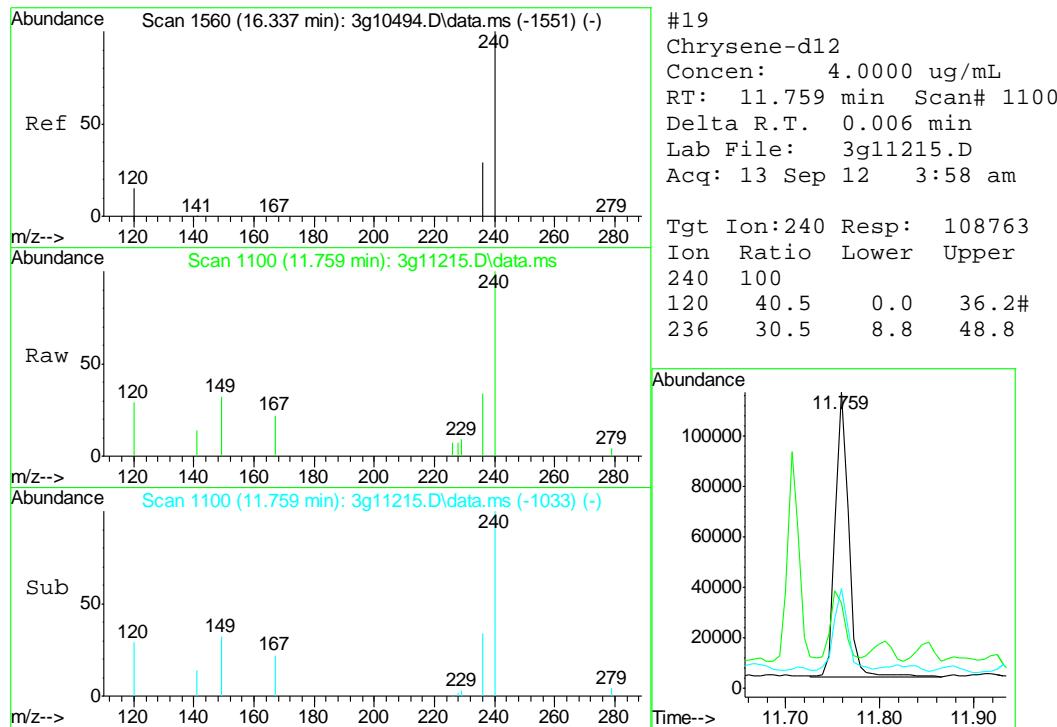


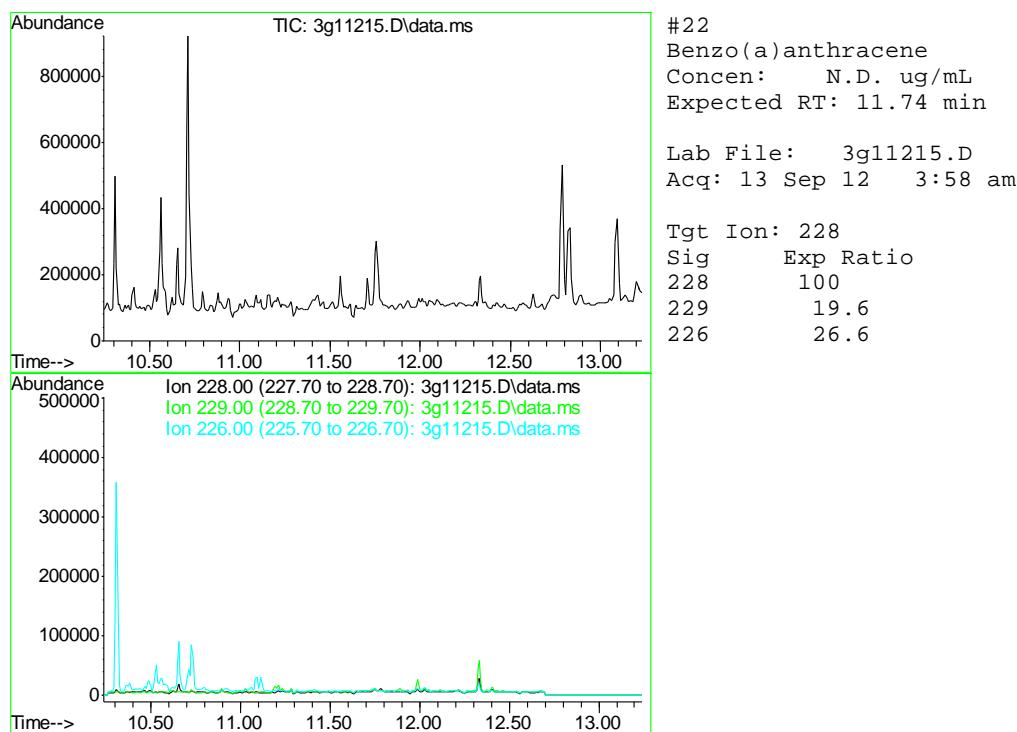
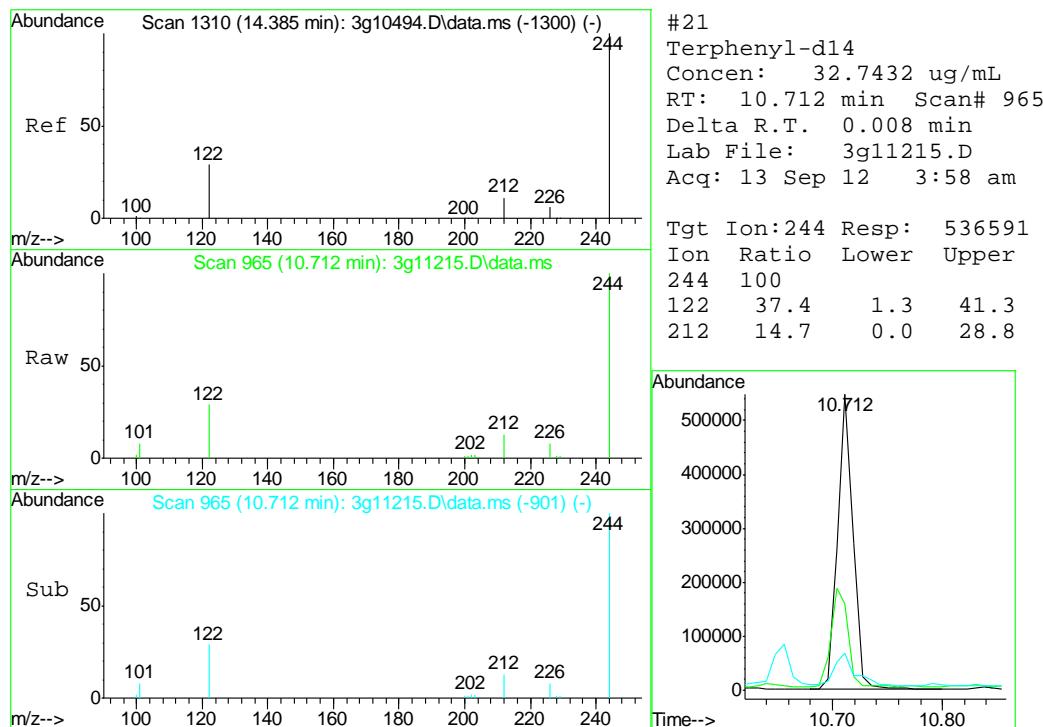


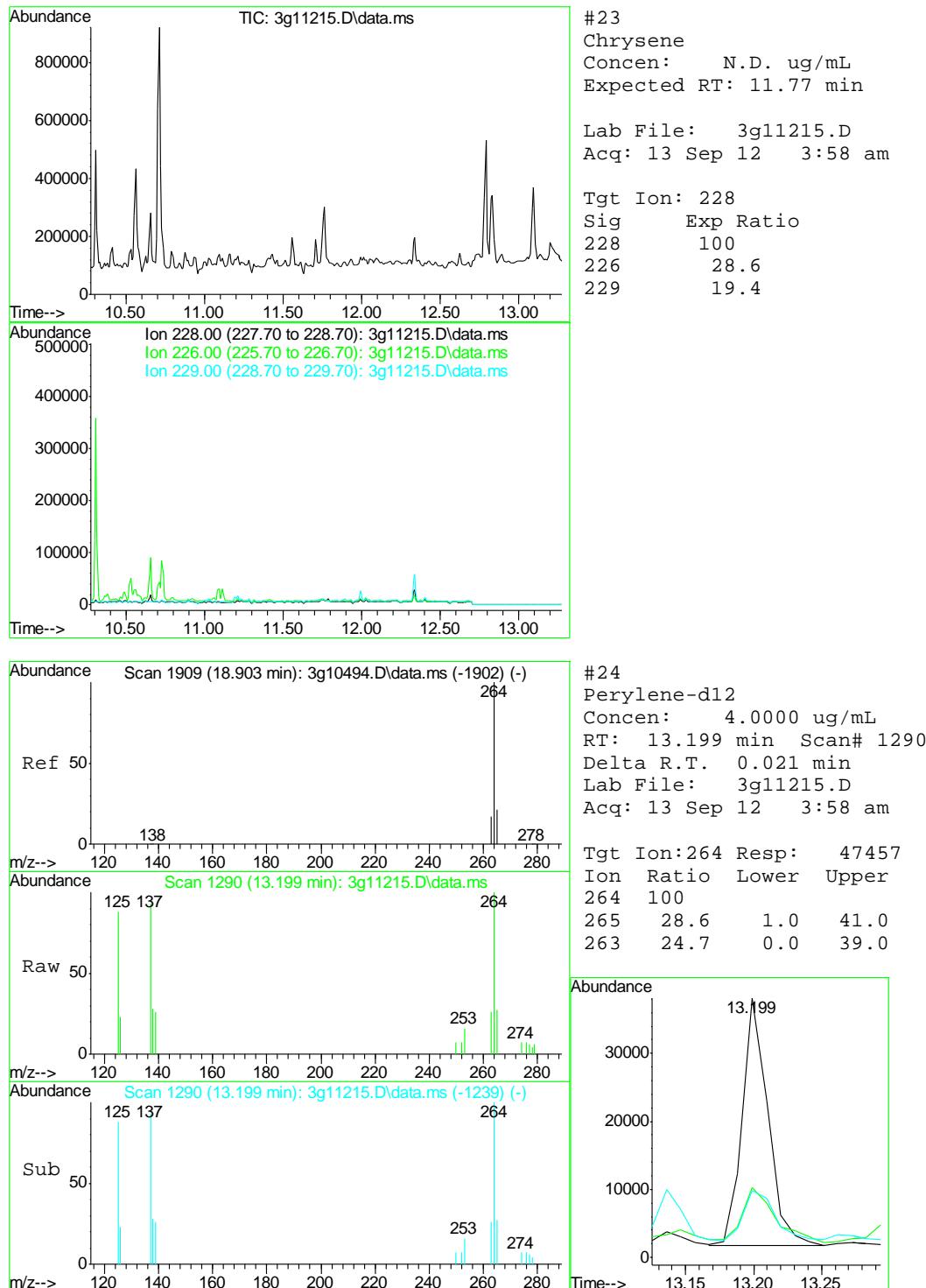


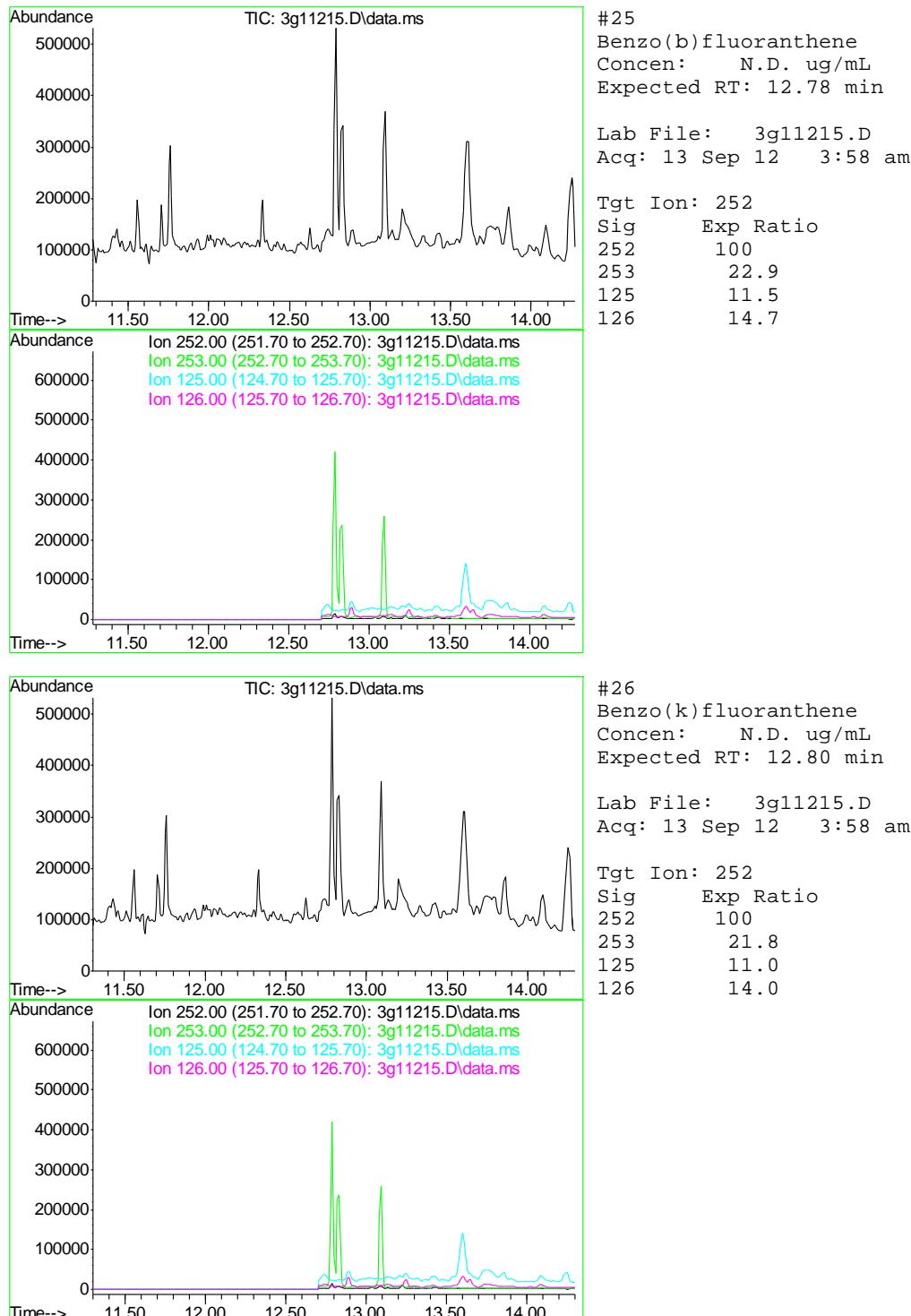


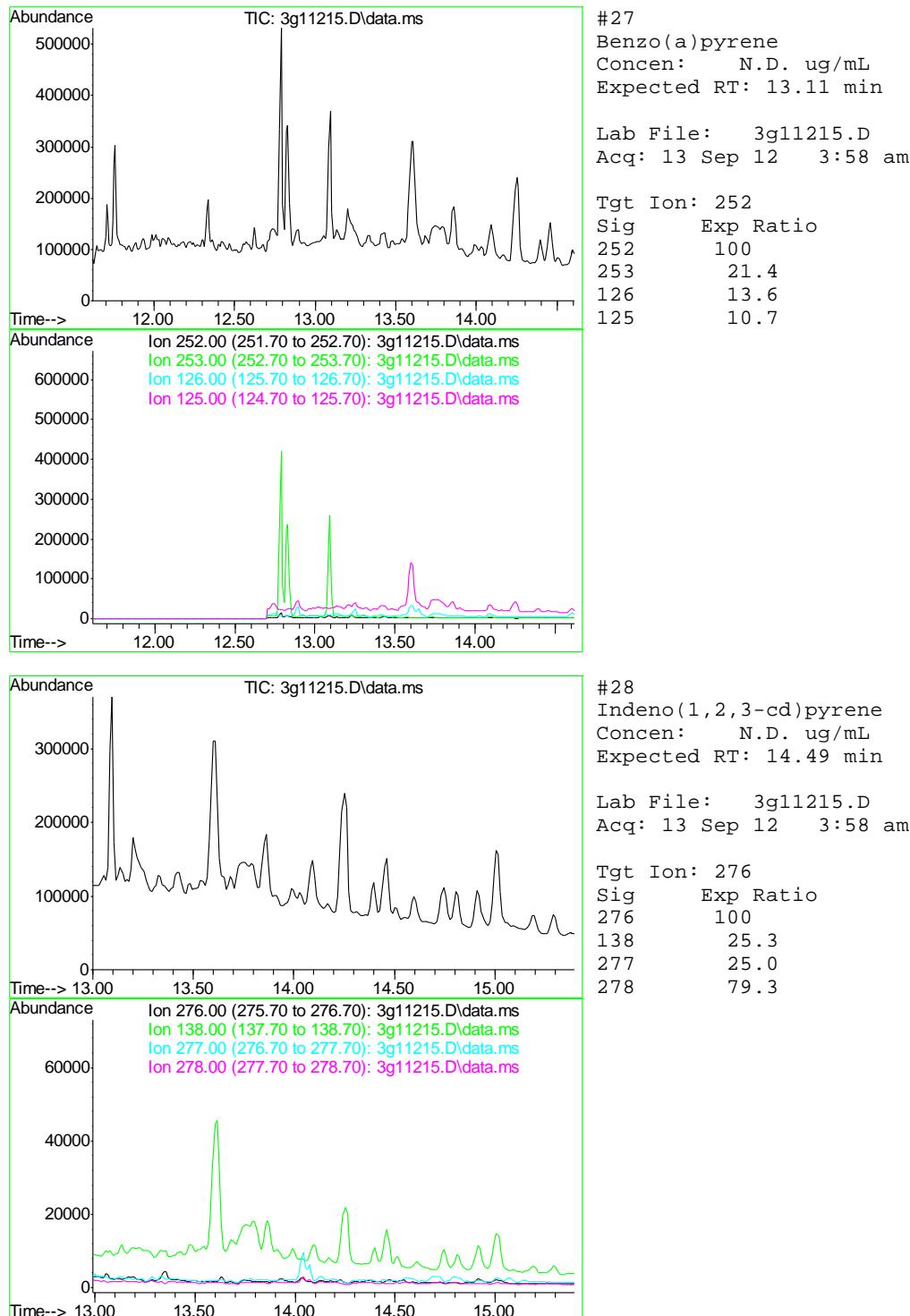


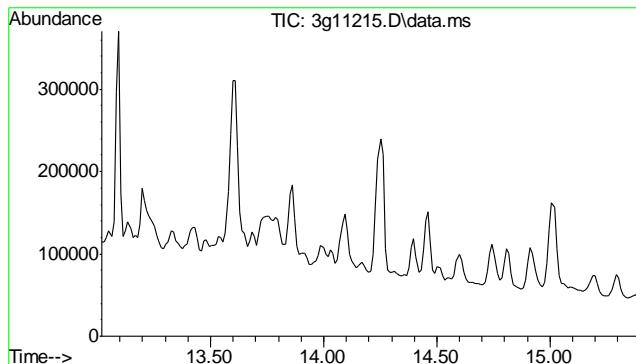








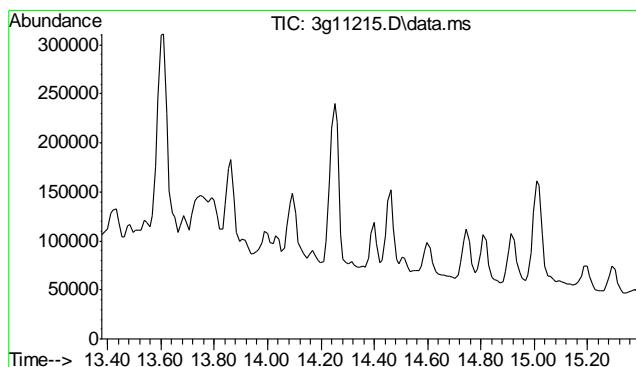
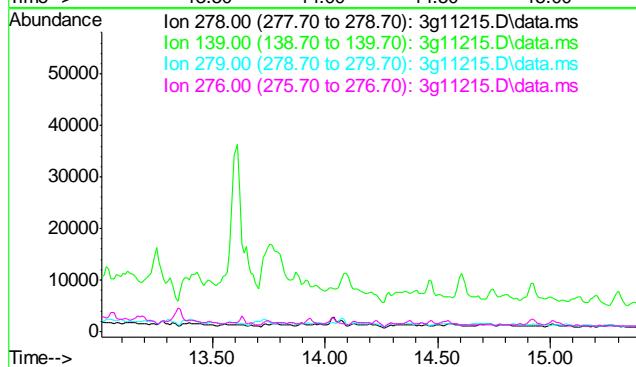




#29
Dibenz(a,h)anthracene
Concen: N.D. ug/mL
Expected RT: 14.51 min

Lab File: 3g11215.D
Acq: 13 Sep 12 3:58 am

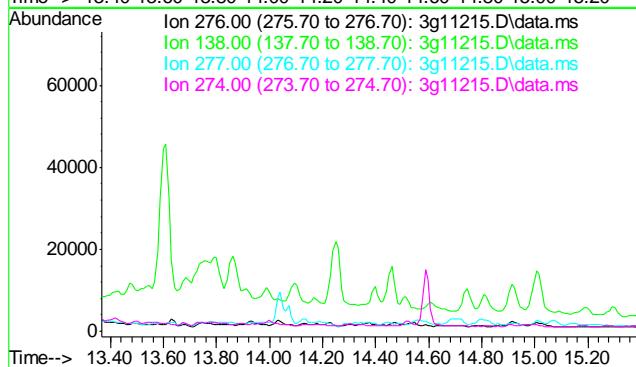
Tgt Ion: 278
Sig Exp Ratio
278 100
139 18.4
279 23.1
276 126.1



#30
Benzo(g,h,i)perylene
Concen: N.D. ug/mL
Expected RT: 14.87 min

Lab File: 3g11215.D
Acq: 13 Sep 12 3:58 am

Tgt Ion: 276
Sig Exp Ratio
276 100
138 21.3
277 23.4
274 21.3



**Manual Integrations
APPROVED
(compounds with "m" flag)**
**Judy Nelson
09/13/12 14:22**

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\091212\
Data File : 3g11216.D
Acq On : 13 Sep 2012 4:22 am
Operator : DONC
Sample : D38480-2
Misc : OP6602,E3G522,30.05,,,1,1
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Sep 13 13:11:55 2012
Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
Quant Title : PAHSIM BASE
QLast Update : Thu Sep 06 09:42:23 2012
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.922	136	198213	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.640	164	109737	4.0000	ug/mL	0.00
15) Phenanthrene-d10	9.121	188	166923	4.0000	ug/mL	0.00
19) Chrysene-d12	11.759	240	103113	4.0000	ug/mL	0.00
24) Perylene-d12	13.188	264	57247	4.0000	ug/mL	0.01

System Monitoring Compounds

2) Nitrobenzene-d5	5.236	82	683832	35.0653	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	= 70.14%	
7) 2-Fluorobiphenyl	6.978	172	1641448	35.9589	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	= 71.92%	
21) Terphenyl-d14	10.712	244	597509	38.4583	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 76.92%	

Target Compounds

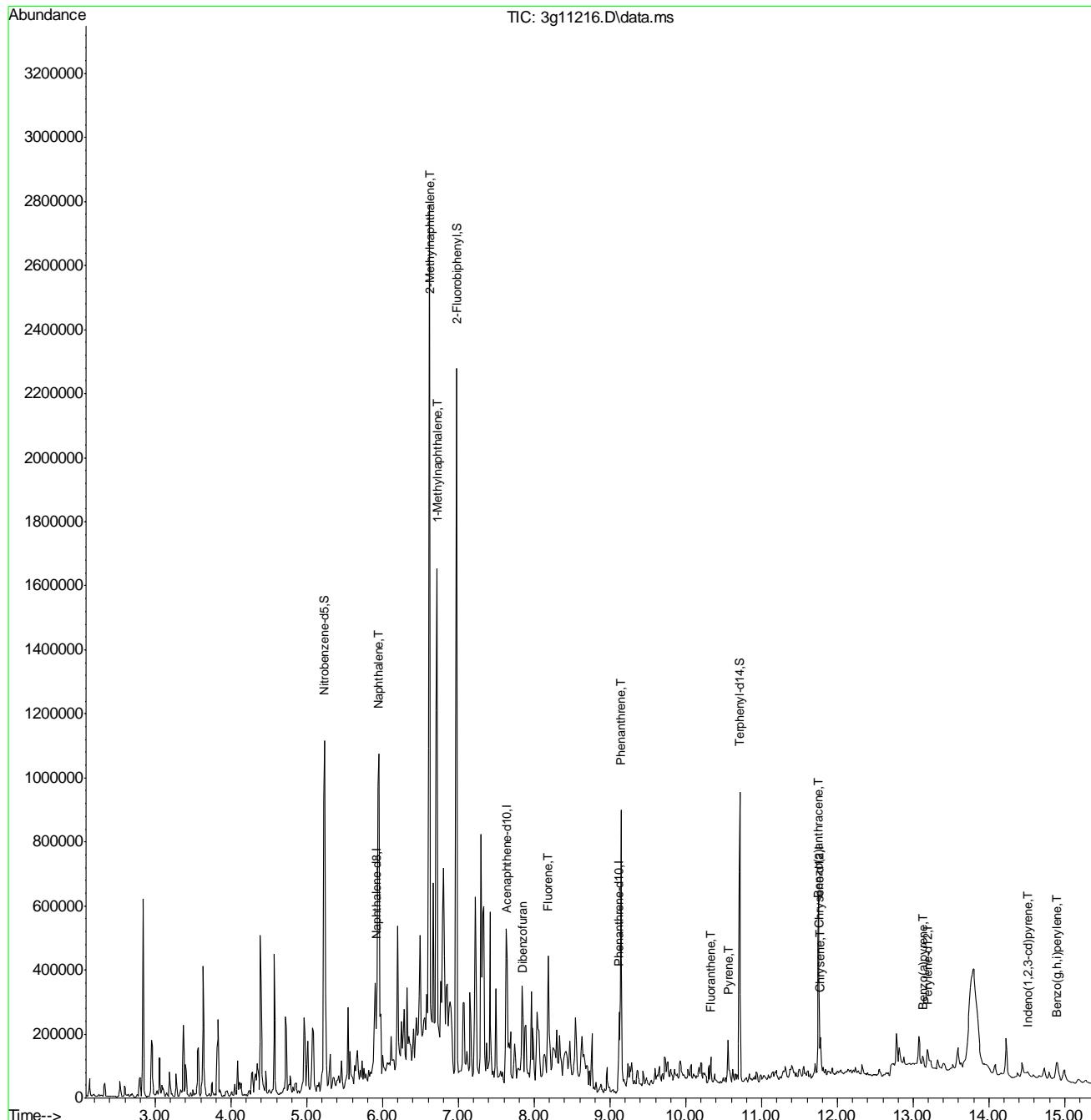
				Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D. d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.946	128	1048115	19.0015 ug/mL 93
8) 2-Methylnaphthalene	6.620	142	1250975	38.5594 ug/mL 96
9) 1-Methylnaphthalene	6.719	142	832403	24.7935 ug/mL 99
10) Acenaphthylene	0.000	152	0	N.D. d
11) Acenaphthene	0.000	154	0	N.D. d
12) Dibenzofuran	7.852	168	140367	2.5062 ug/mL 85
13) Fluorene	8.183	166	132917	2.9657 ug/mL# 70
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	9.145	178	528163	9.0193 ug/mL 89
17) Anthracene	0.000	178	0	N.D. d
18) Fluoranthene	10.332	202	58840	0.8547 ug/mL 56
20) Pyrene	10.561	202	81603m	1.6532 ug/mL
22) Benzo(a)anthracene	11.746	228	25631	0.5853 ug/mL# 66
23) Chrysene	11.779	228	89394	1.9284 ug/mL# 72
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d
26) Benzo(k)fluoranthene	0.000	252	0	N.D. d
27) Benzo(a)pyrene	13.125	252	24979	0.6250 ug/mL# 68
28) Indeno(1,2,3-cd)pyrene	14.513	276	10764	0.2497 ug/mL# 72
29) Dibenz(a,h)anthracene	0.000	278	0	N.D. d
30) Benzo(g,h,i)perylene	14.903	276	34406	0.9346 ug/mL# 45

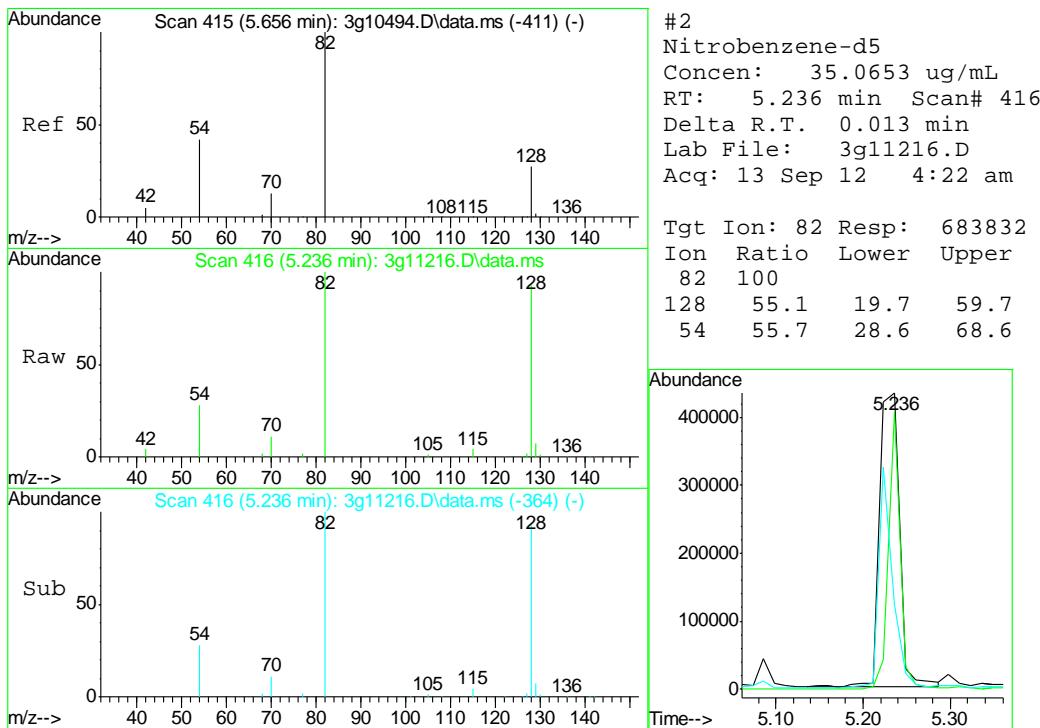
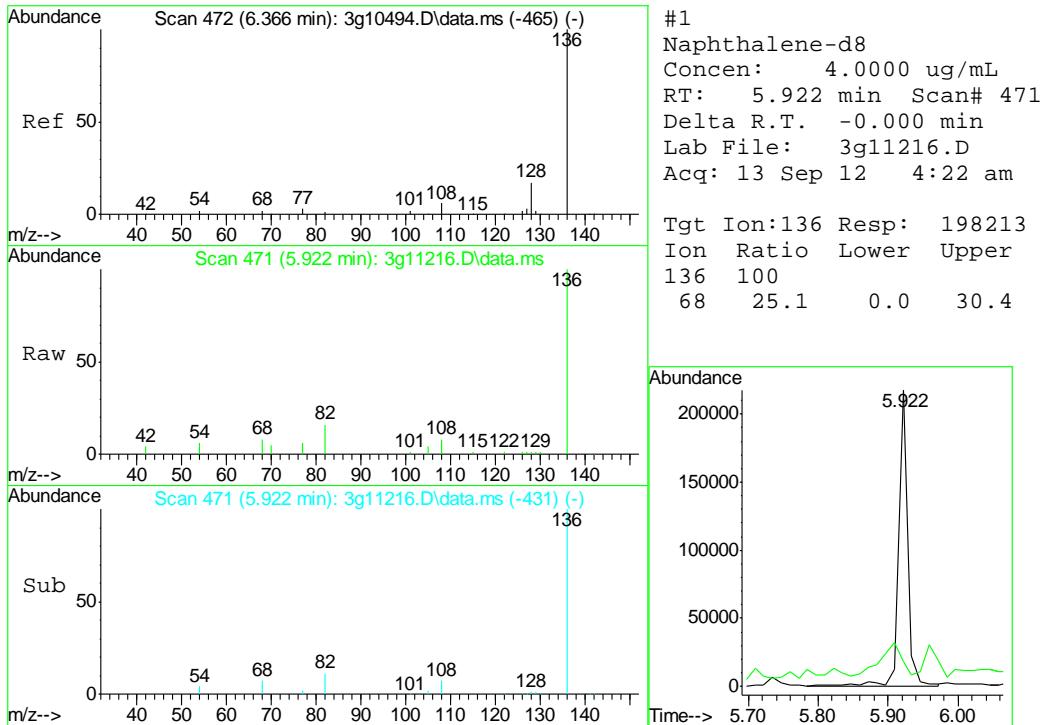
(#) = qualifier out of range (m) = manual integration (+) = signals summed

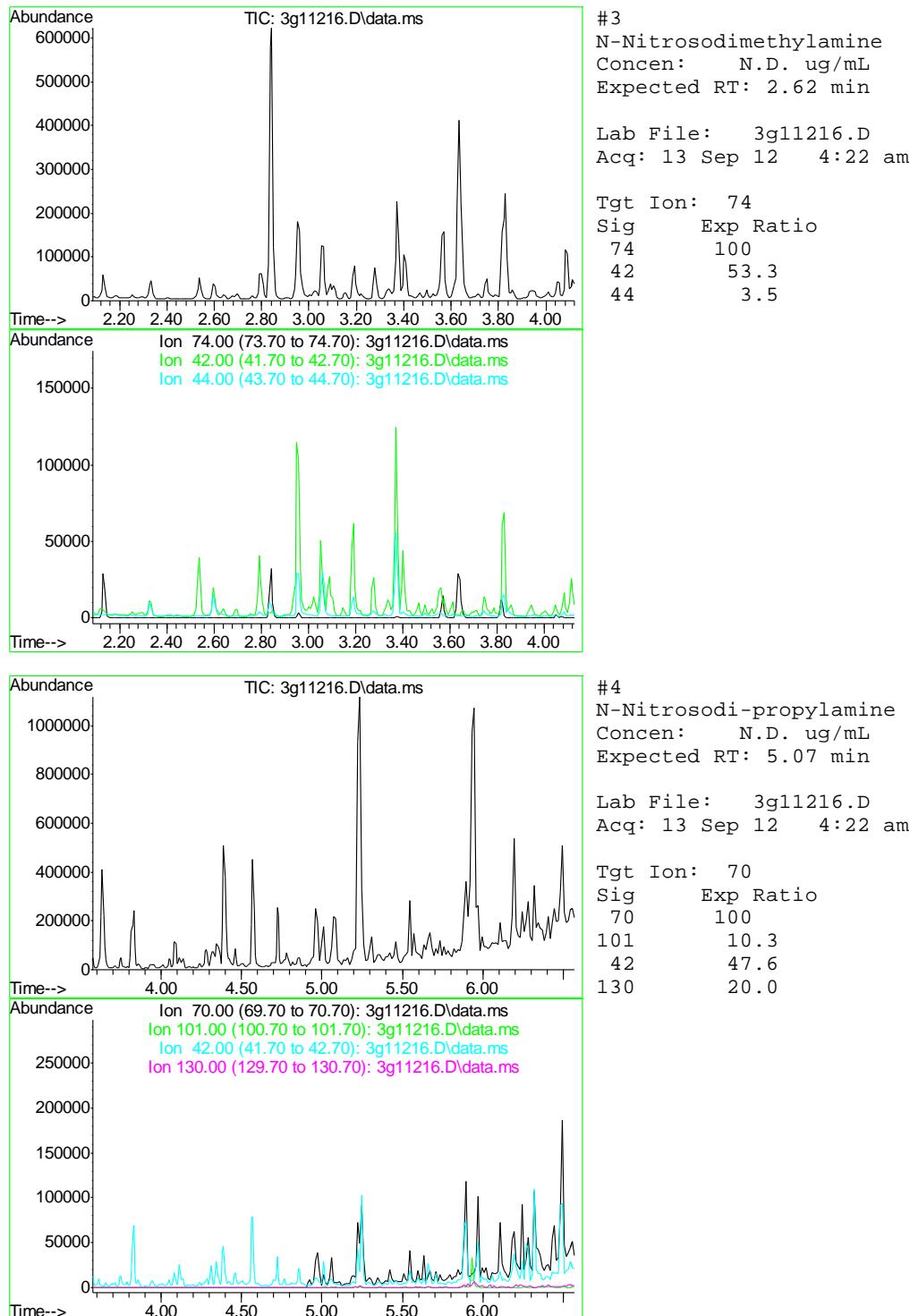
Quantitation Report (QT Reviewed)

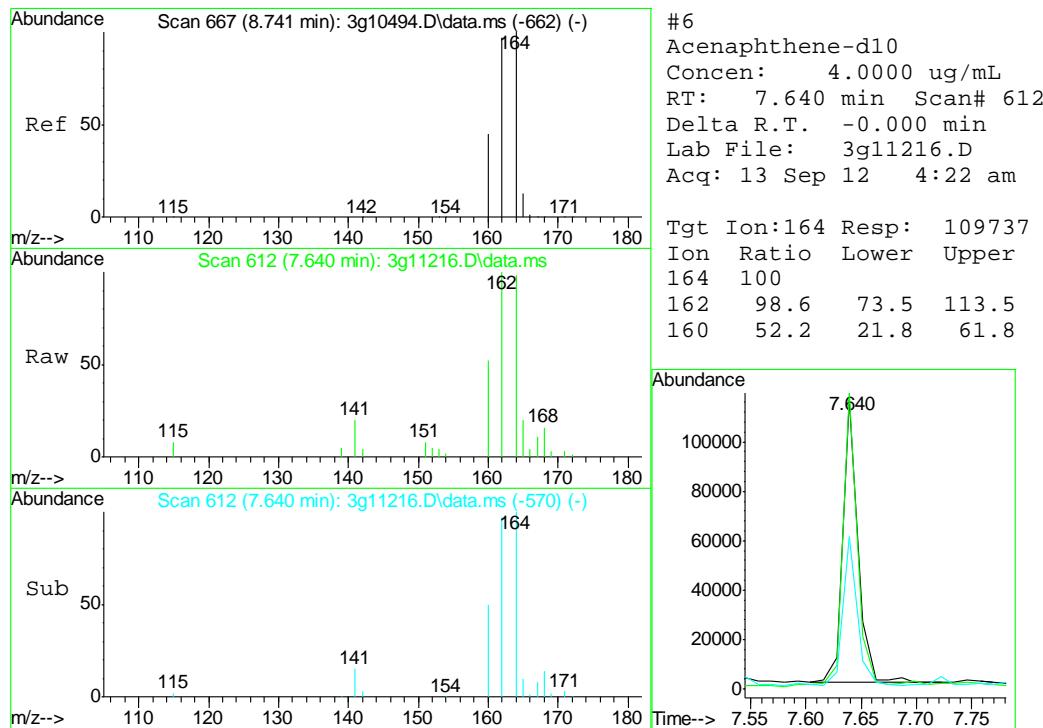
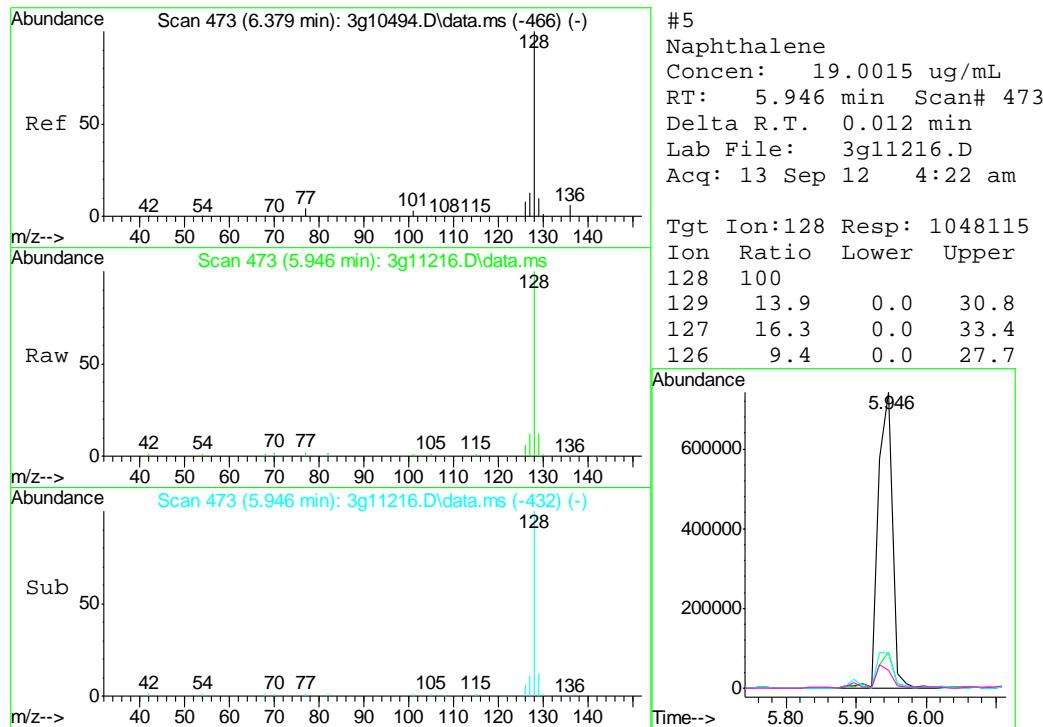
Data Path : C:\msdchem\1\DATA\091212\
 Data File : 3g11216.D
 Acq On : 13 Sep 2012 4:22 am
 Operator : DONC
 Sample : D38480-2
 Misc : OP6602,E3G522,30.05,,,1,1
 ALS Vial : 27 Sample Multiplier: 1

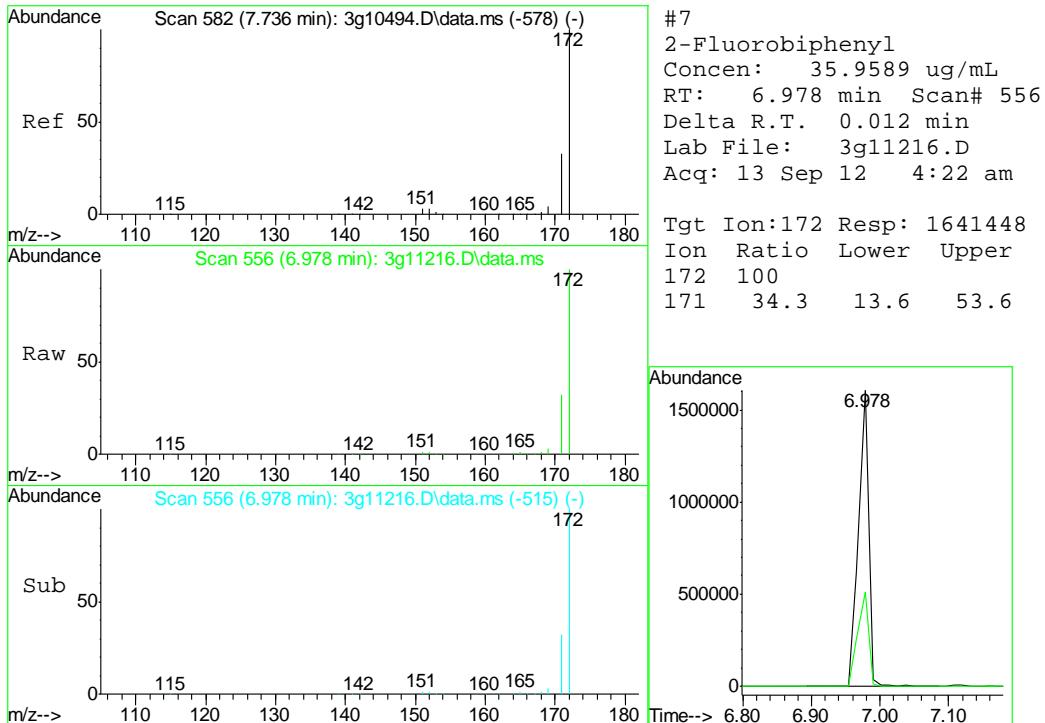
Quant Time: Sep 13 13:11:55 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Sep 06 09:42:23 2012
 Response via : Initial Calibration





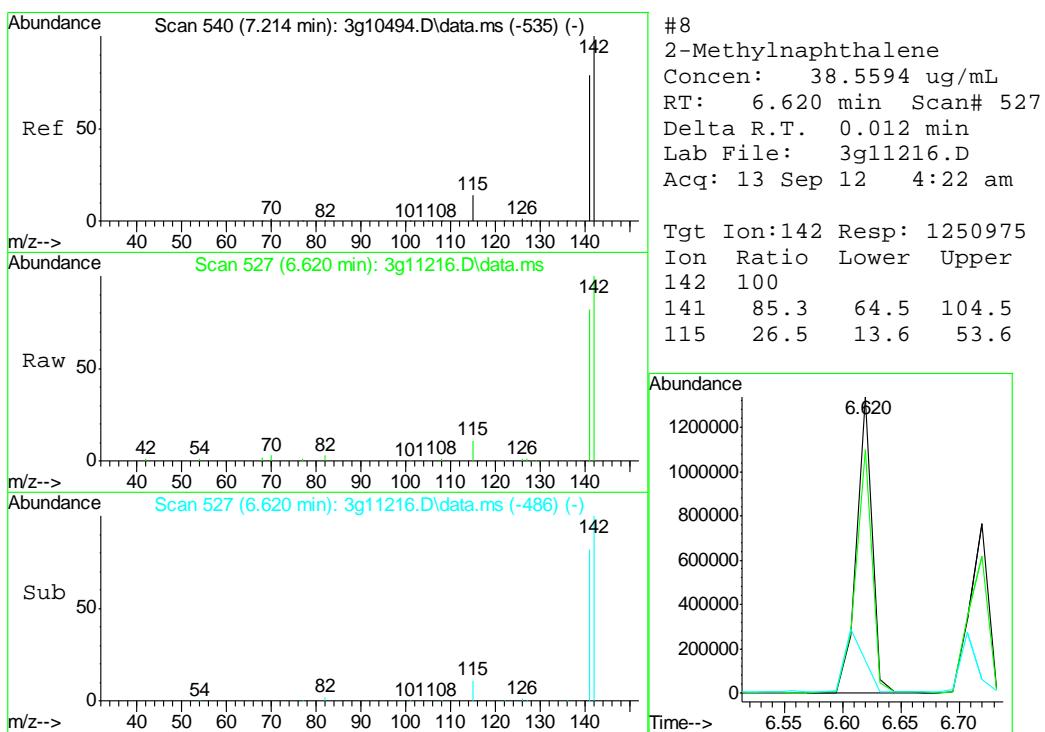


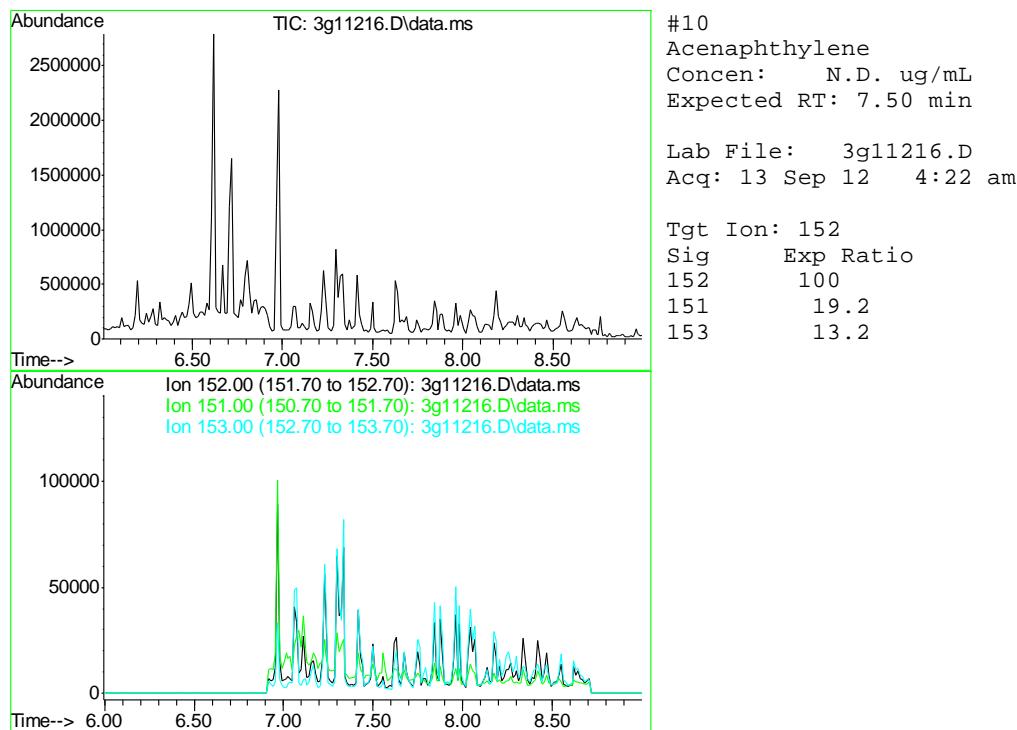
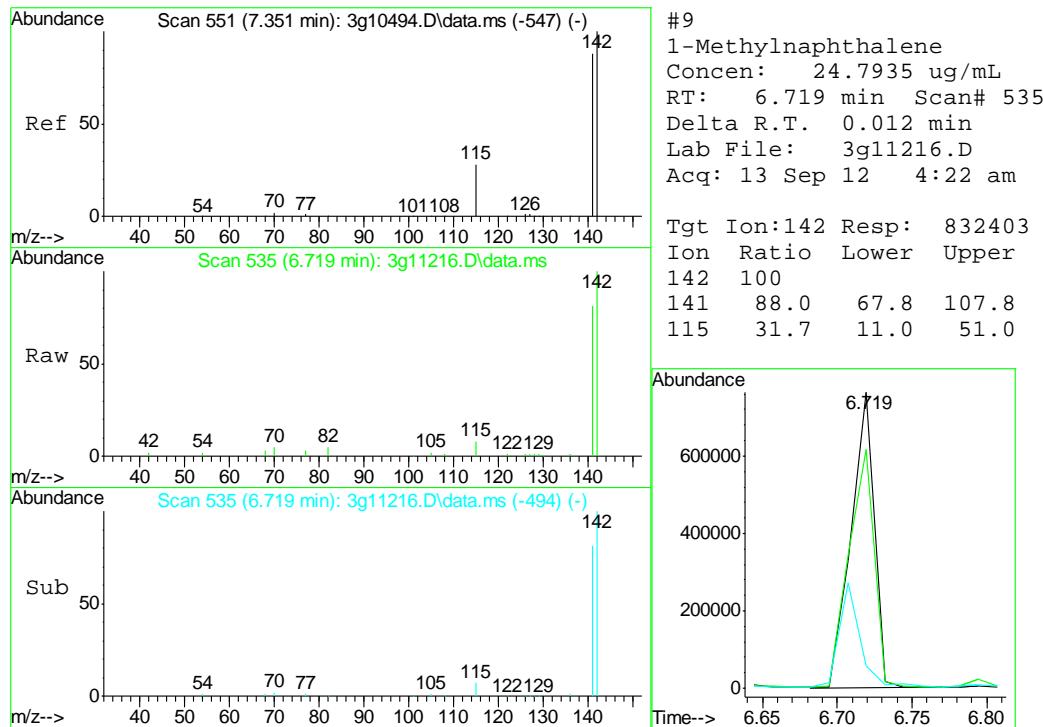


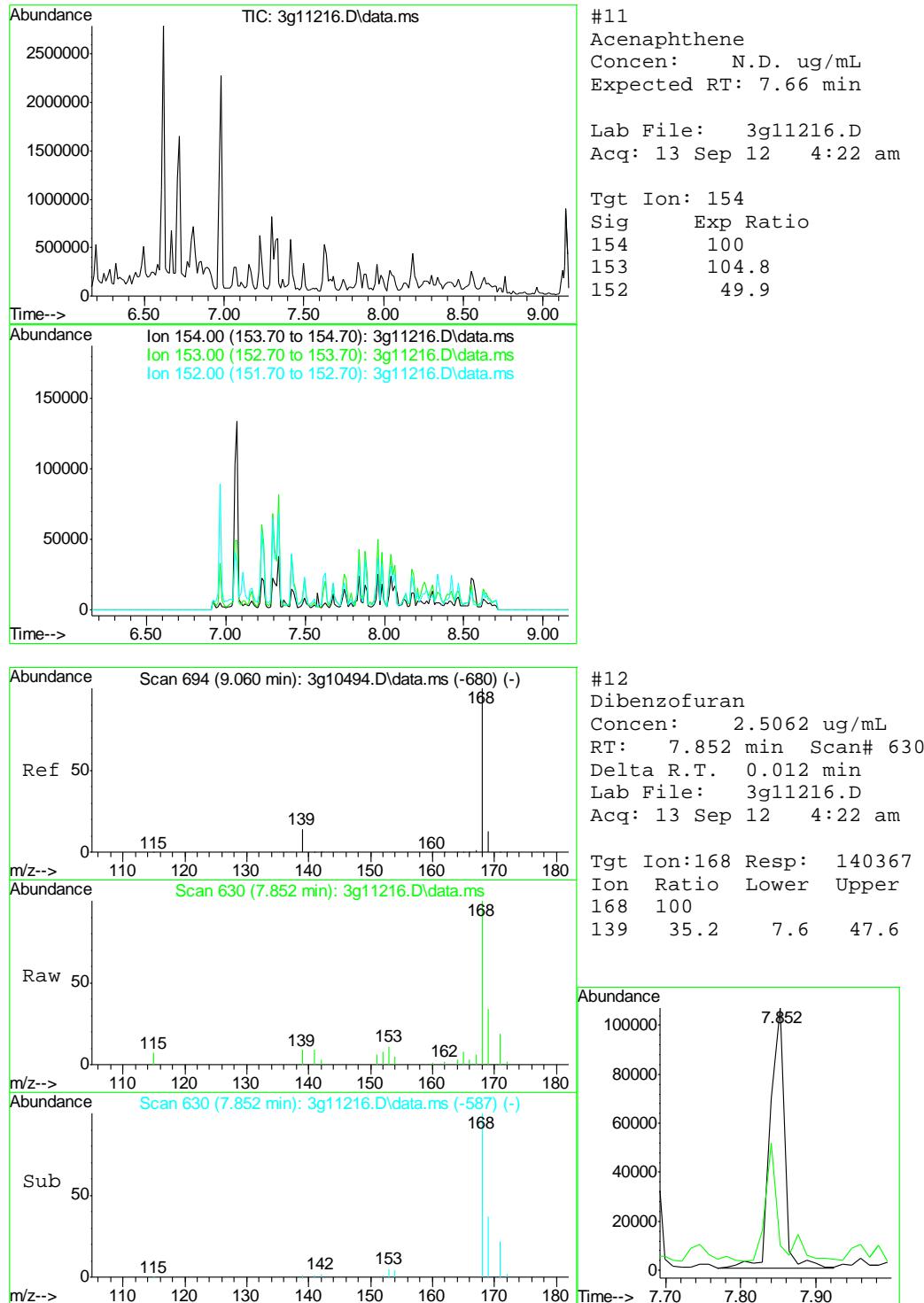


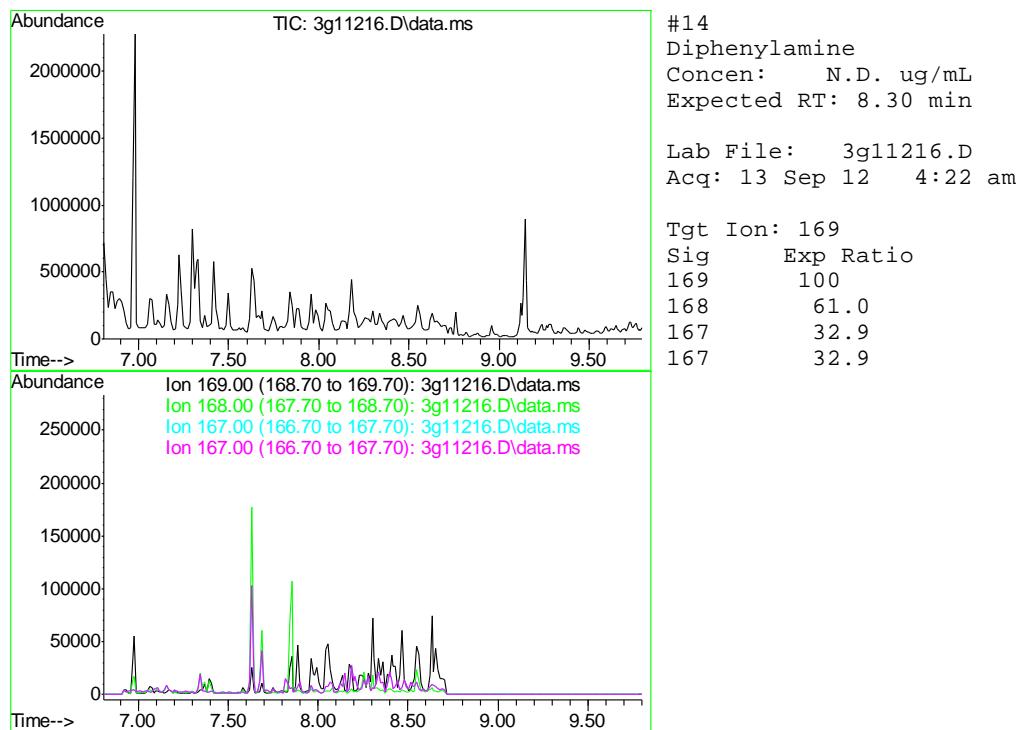
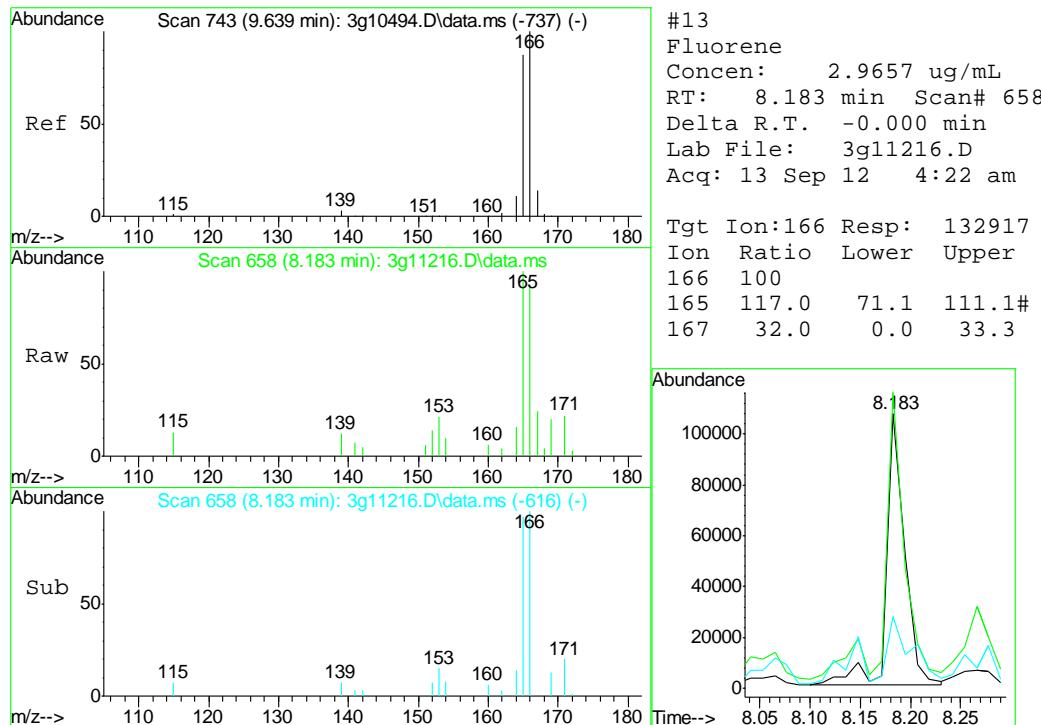
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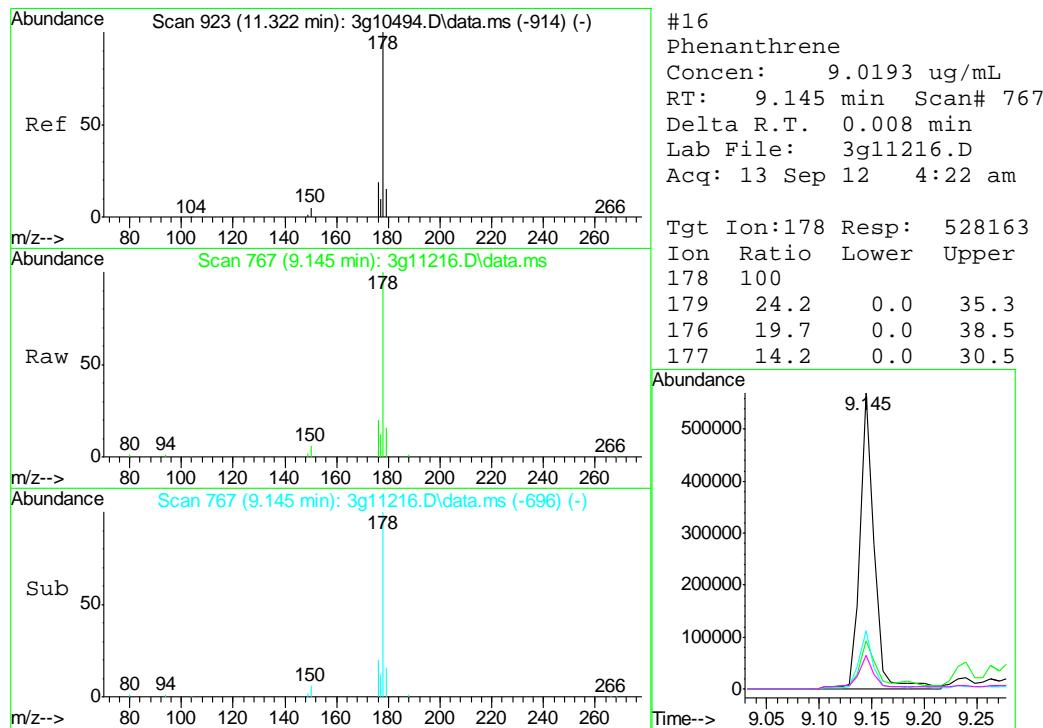
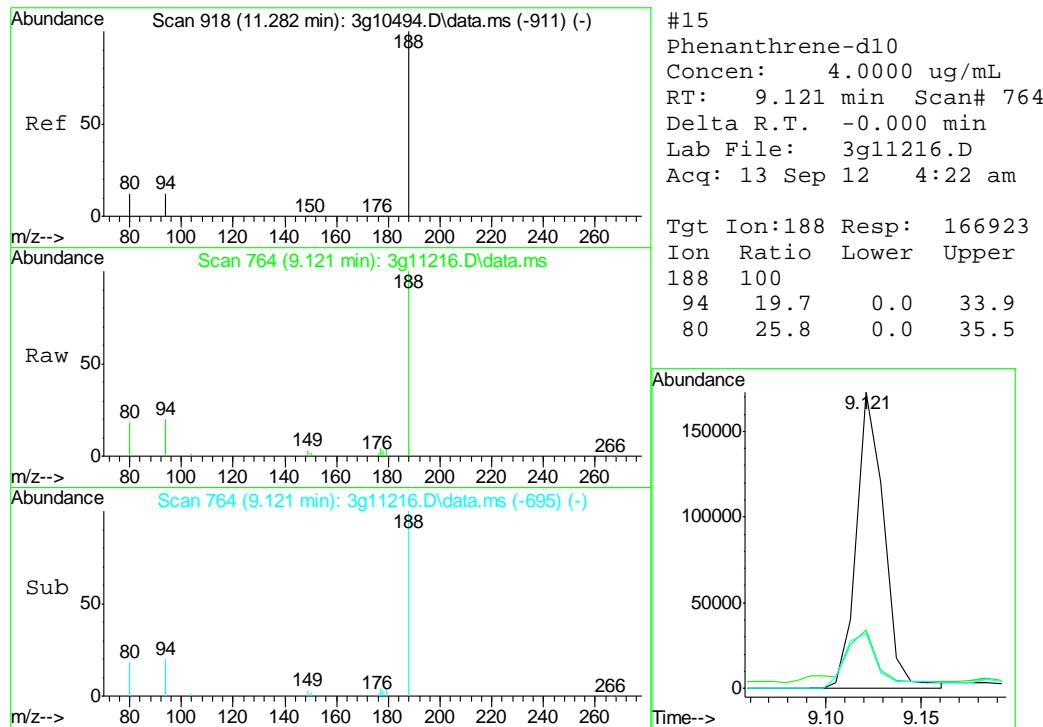
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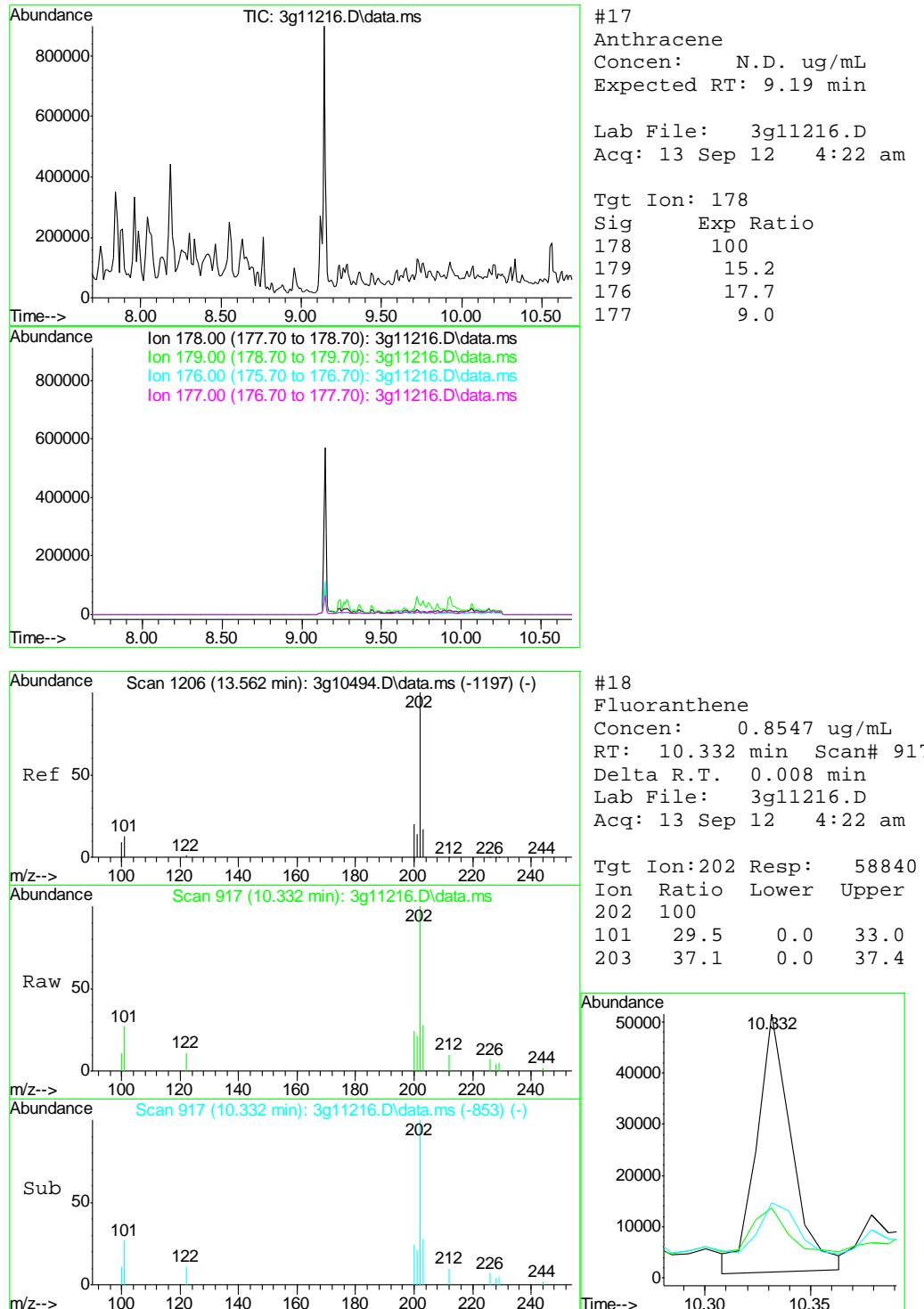


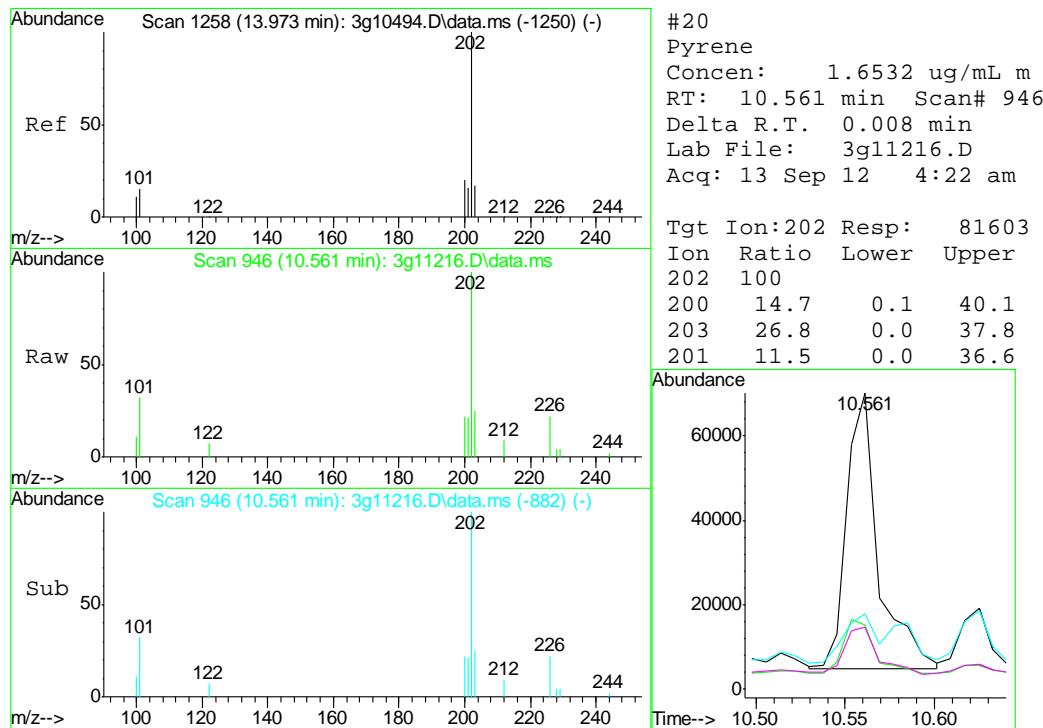
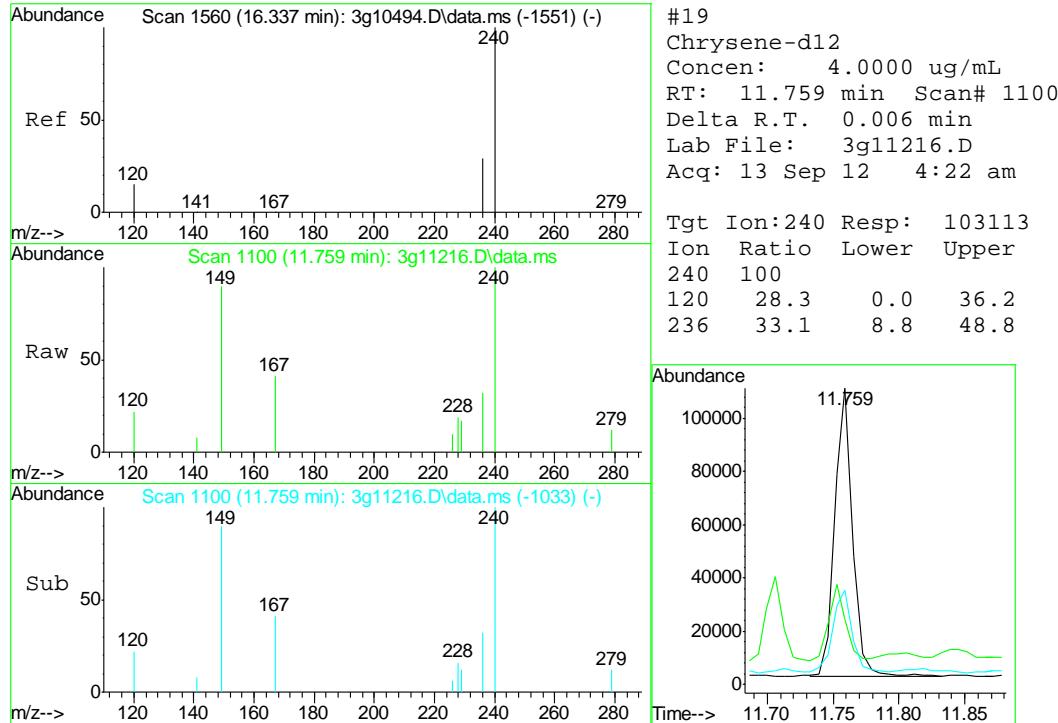


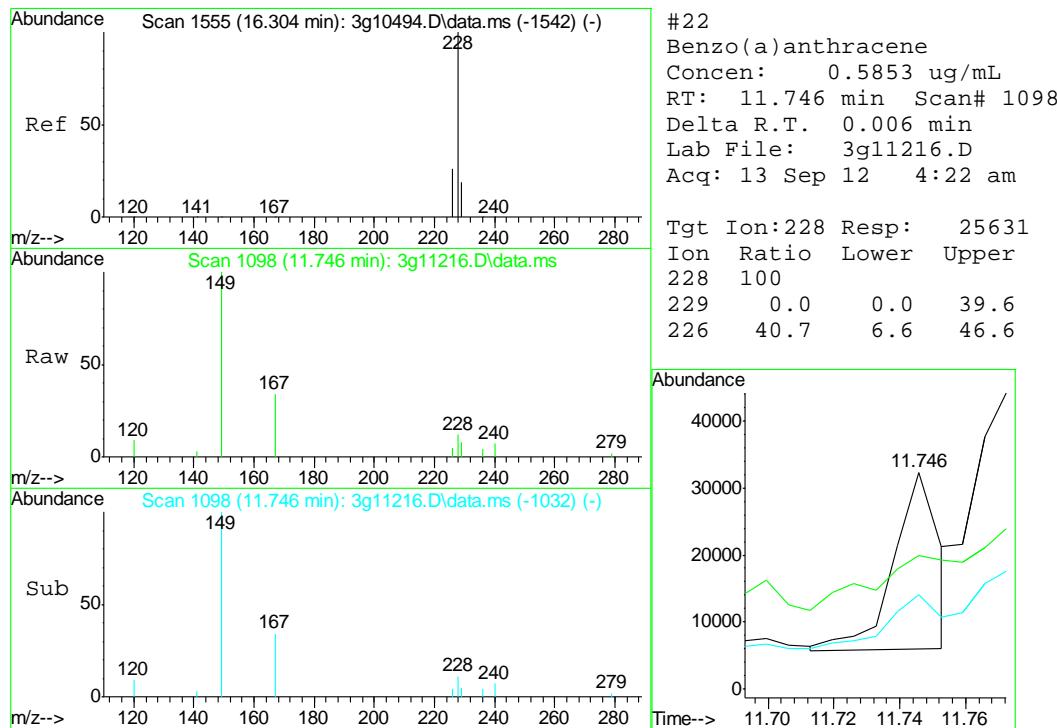
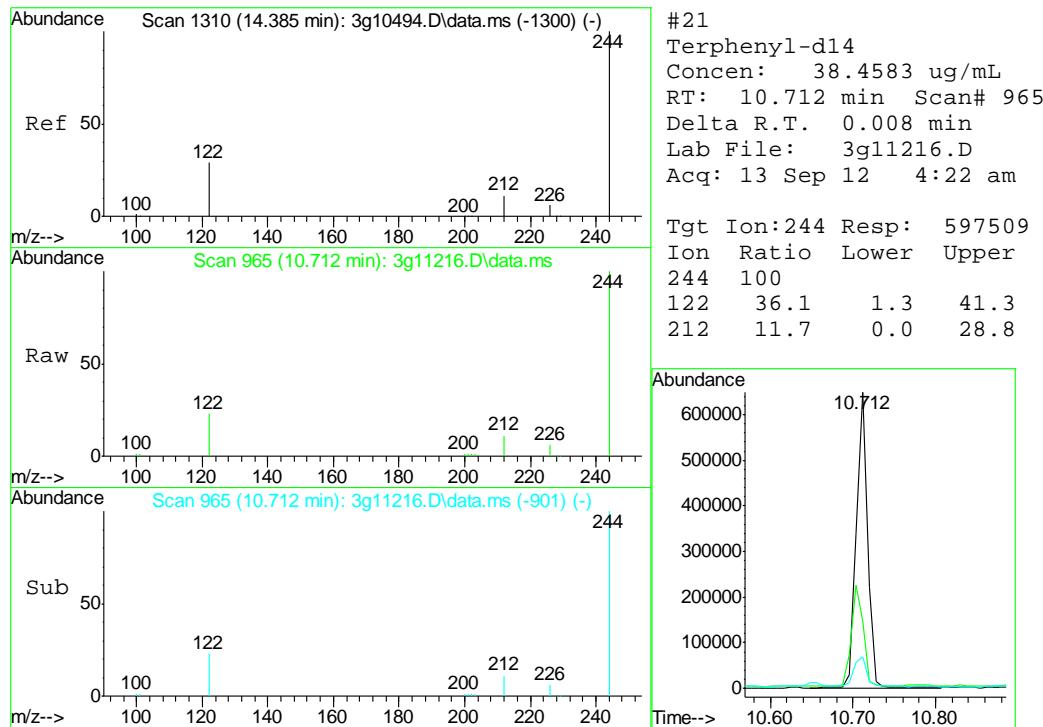


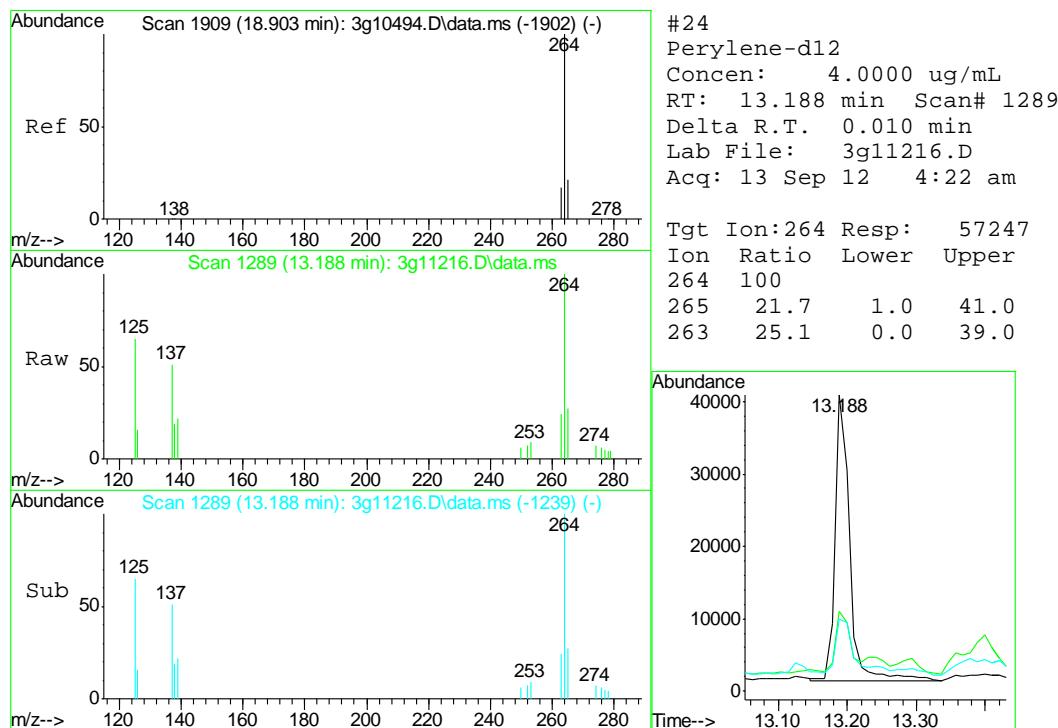
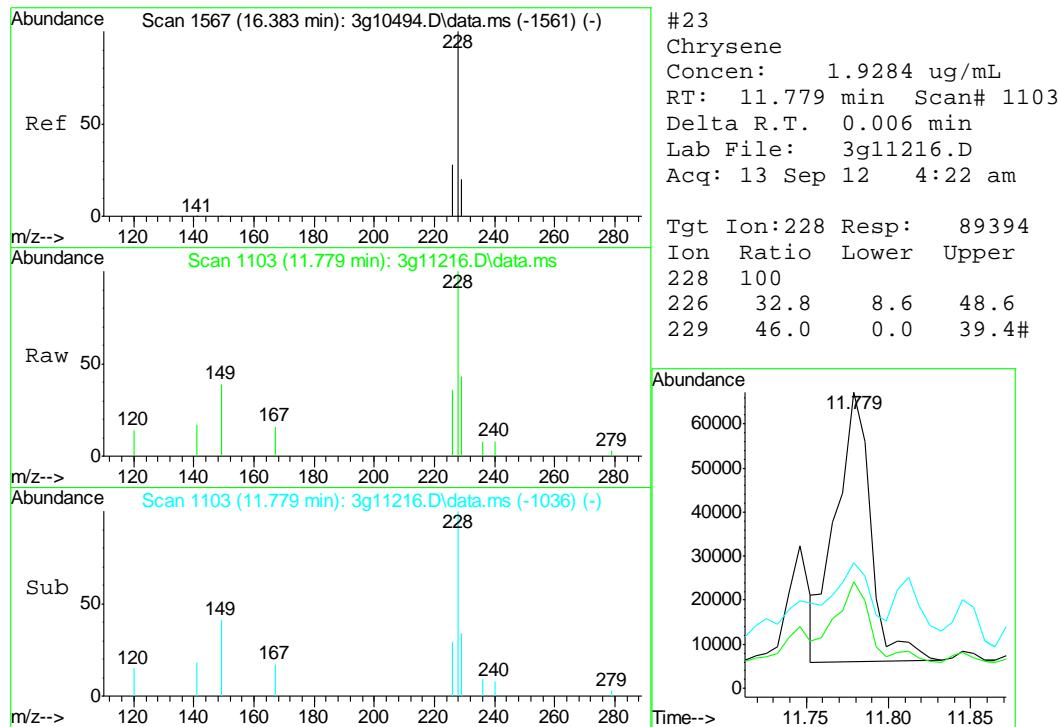


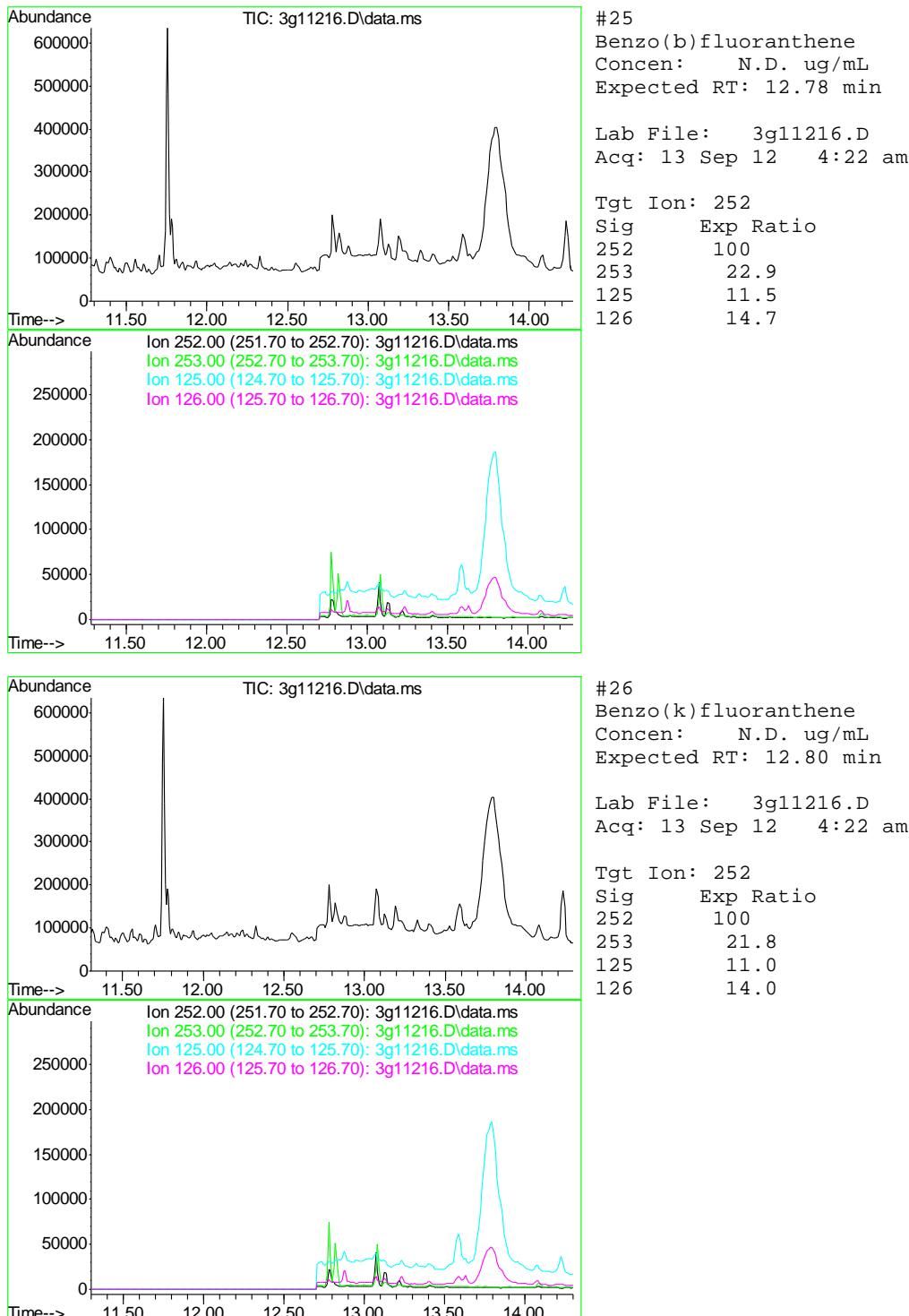


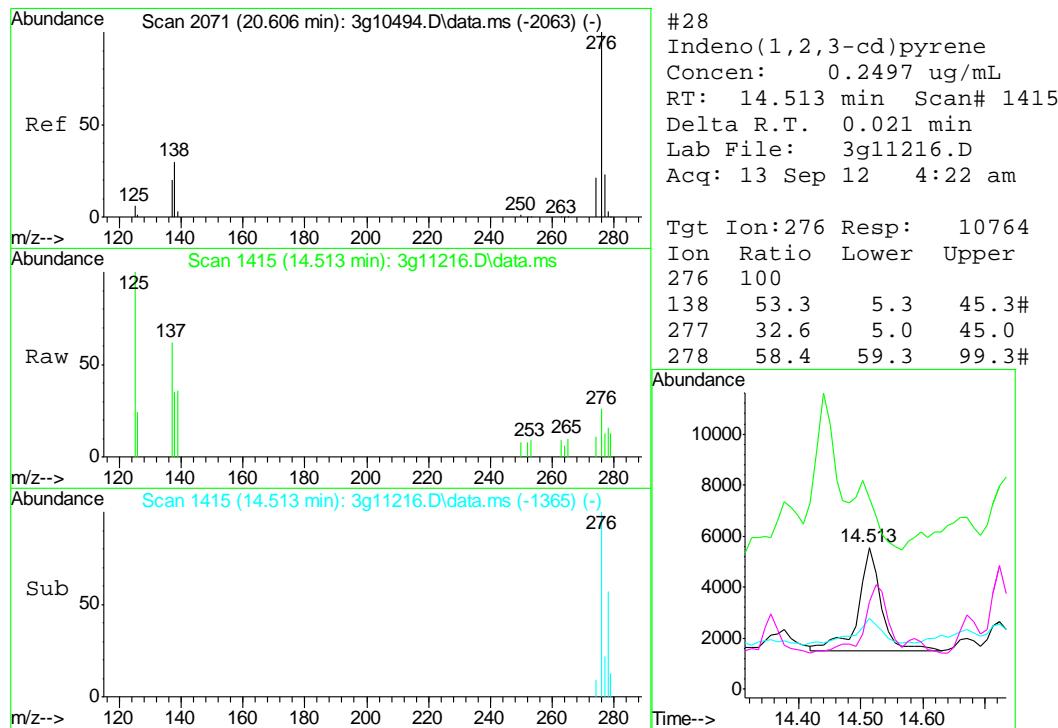
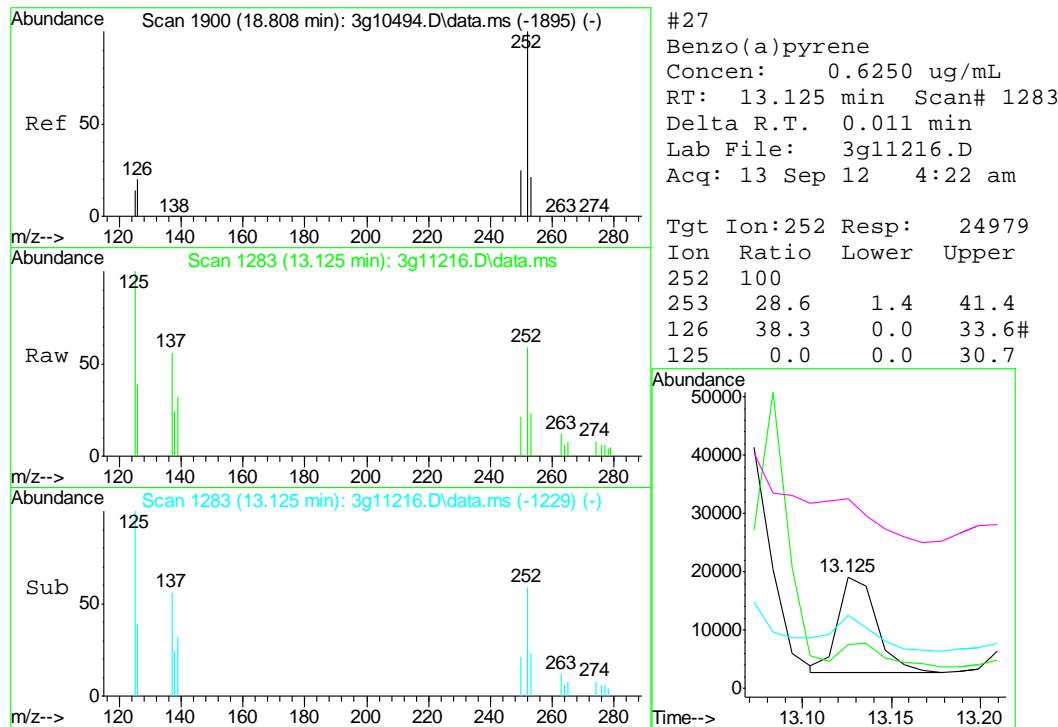


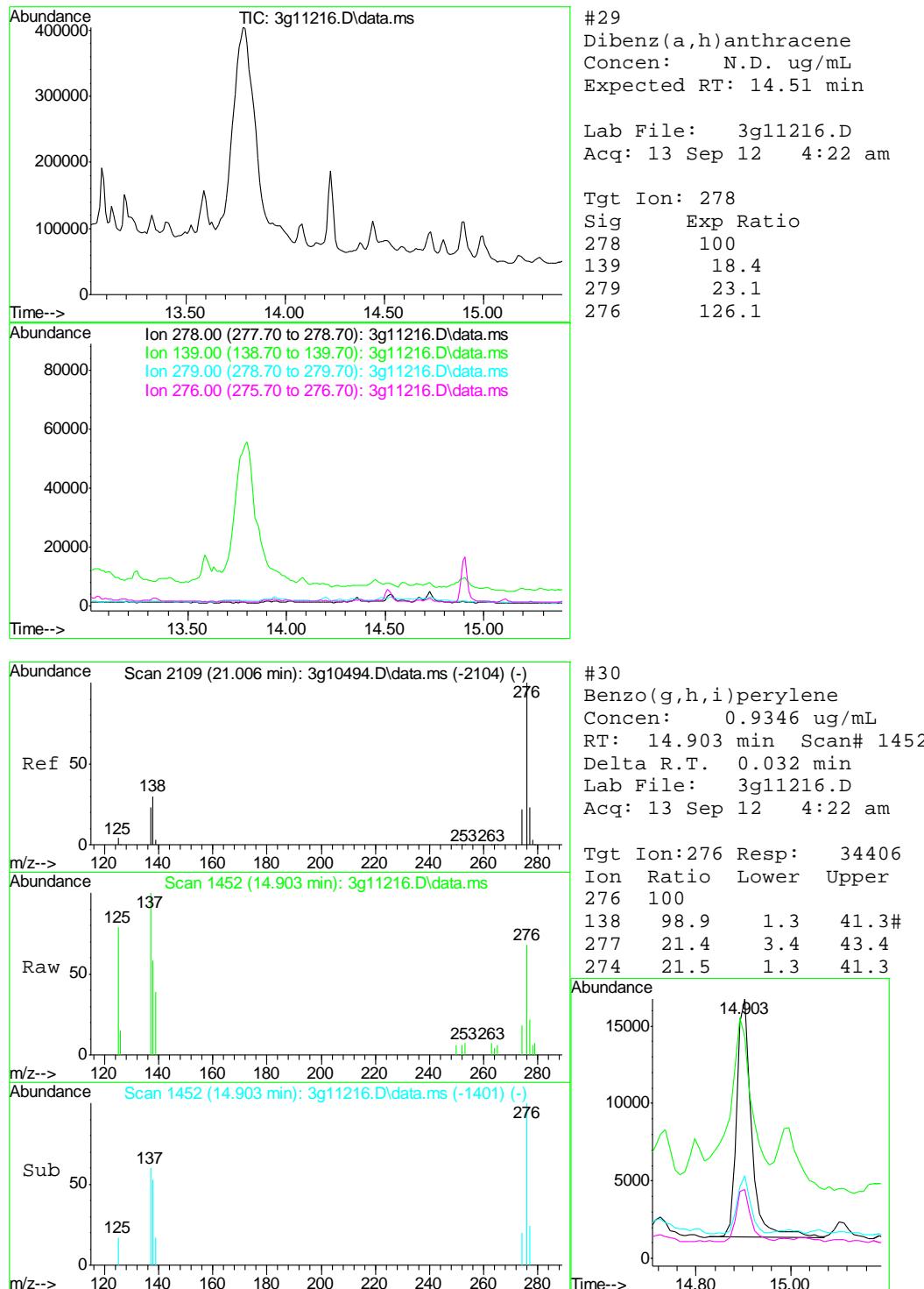












Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson
09/13/12 14:22

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\091212\
 Data File : 3g11217.D
 Acq On : 13 Sep 2012 4:45 am
 Operator : DONC
 Sample : D38480-3
 Misc : OP6602,E3G522,30.17,,,1,1
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Sep 13 13:14:54 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Sep 06 09:42:23 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.921	136	175953	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.639	164	109721	4.0000	ug/mL	0.00
15) Phenanthrene-d10	9.121	188	164625	4.0000	ug/mL	0.00
19) Chrysene-d12	11.759	240	101610	4.0000	ug/mL	0.00
24) Perylene-d12	13.188	264	58034	4.0000	ug/mL	0.01

System Monitoring Compounds

2) Nitrobenzene-d5	5.236	82	720817	41.6379	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	83.28%
7) 2-Fluorobiphenyl	6.978	172	1508827	33.0585	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	66.12%
21) Terphenyl-d14	10.712	244	588930	38.4668	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	76.94%

Target Compounds

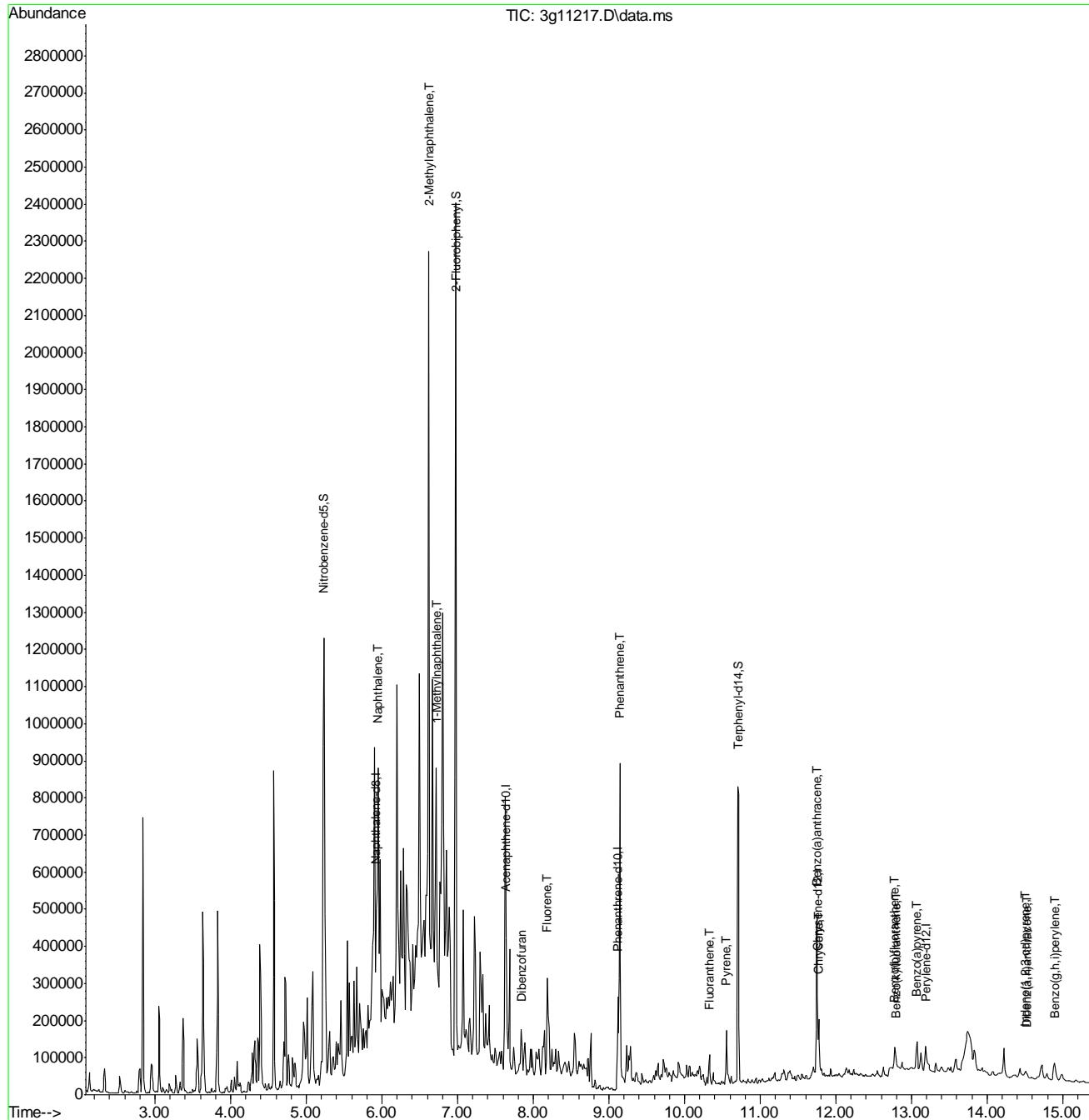
				Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D. d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.946	128	581572	11.8773 ug/mL 97
8) 2-Methylnaphthalene	6.620	142	828664	25.5460 ug/mL 94
9) 1-Methylnaphthalene	6.719	142	253019	7.5374 ug/mL 96
10) Acenaphthylene	0.000	152	0	N.D. d
11) Acenaphthene	0.000	154	0	N.D. d
12) Dibenzofuran	7.852	168	35806	0.6394 ug/mL 88
13) Fluorene	8.183	166	153129	3.4171 ug/mL# 88
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	9.144	178	483183	8.3663 ug/mL 93
17) Anthracene	0.000	178	0	N.D. d
18) Fluoranthene	10.332	202	57097	0.8410 ug/mL 68
20) Pyrene	10.553	202	82508	1.6963 ug/mL 91
22) Benzo(a)anthracene	11.746	228	32939	0.7633 ug/mL 87
23) Chrysene	11.779	228	121299m	2.6553 ug/mL
25) Benzo(b)fluoranthene	12.778	252	41820m	1.0834 ug/mL
26) Benzo(k)fluoranthene	12.799	252	10297m	0.2413 ug/mL
27) Benzo(a)pyrene	13.073	252	54669	1.3494 ug/mL 64
28) Indeno(1,2,3-cd)pyrene	14.503	276	14953	0.3421 ug/mL 92
29) Dibenz(a,h)anthracene	14.524	278	11882	0.3461 ug/mL# 82
30) Benzo(g,h,i)perylene	14.892	276	41127m	1.1020 ug/mL

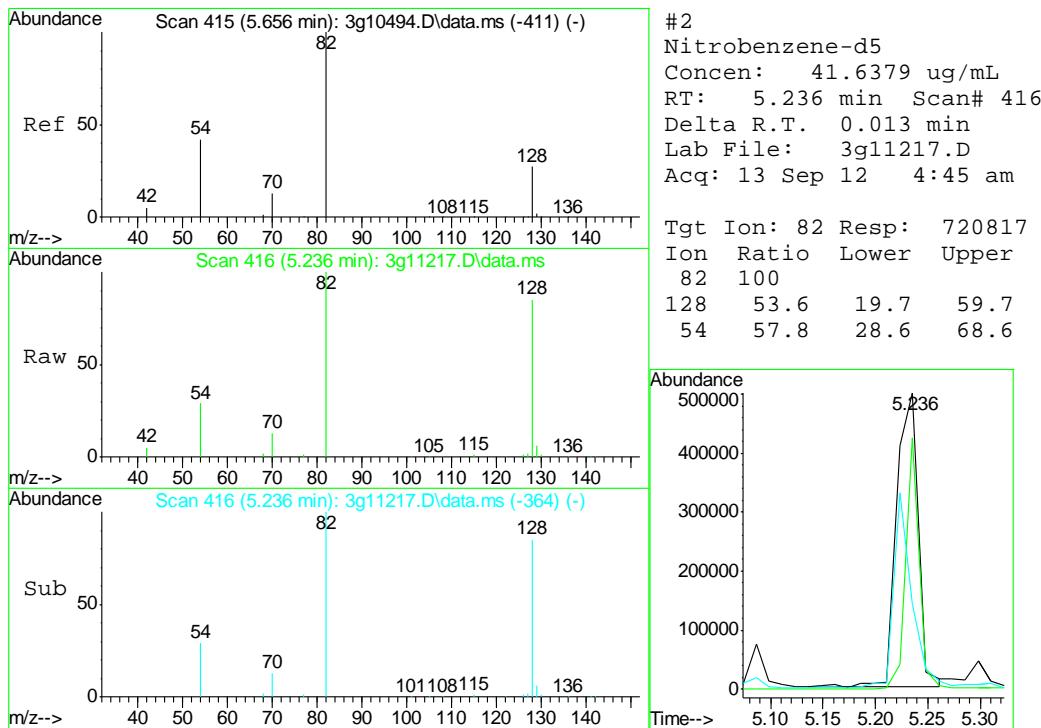
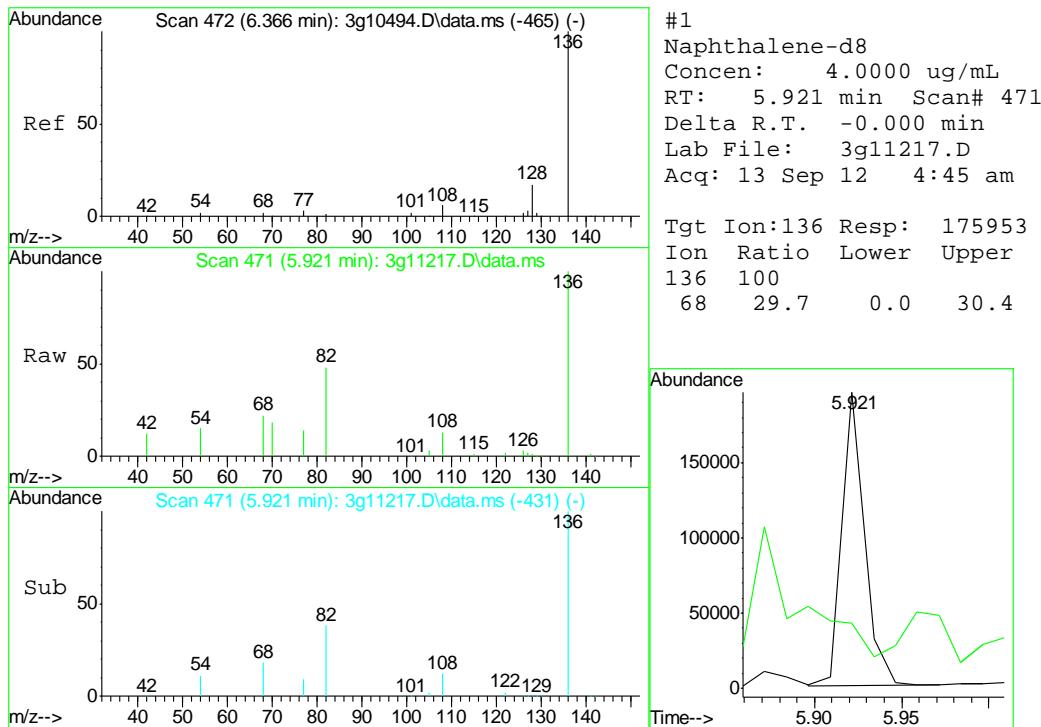
(#) = qualifier out of range (m) = manual integration (+) = signals summed

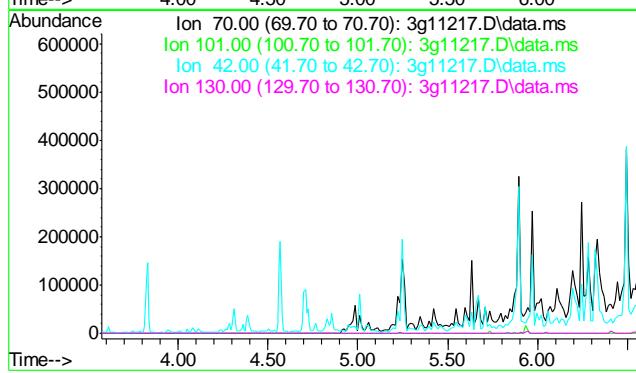
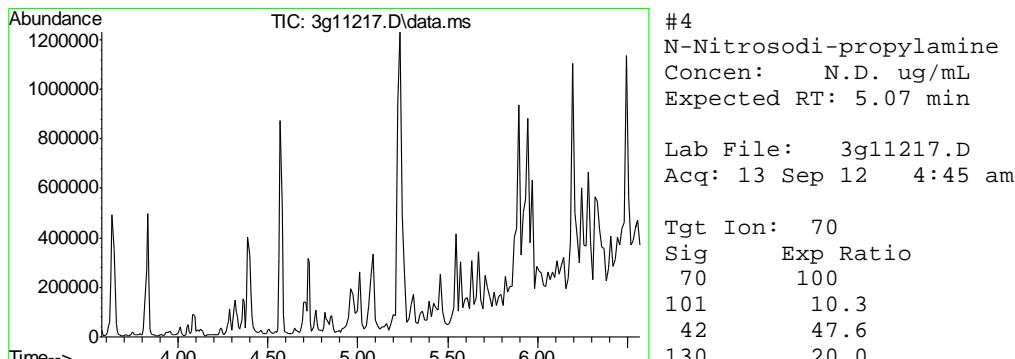
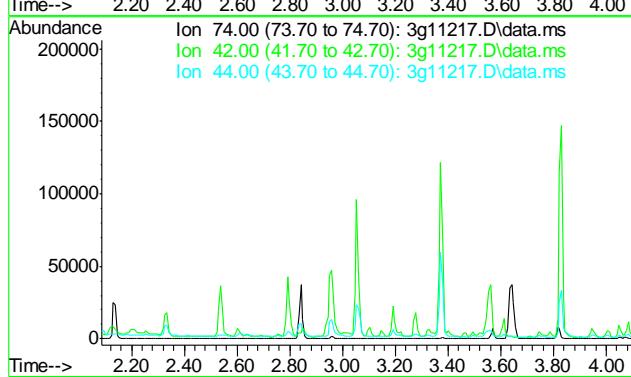
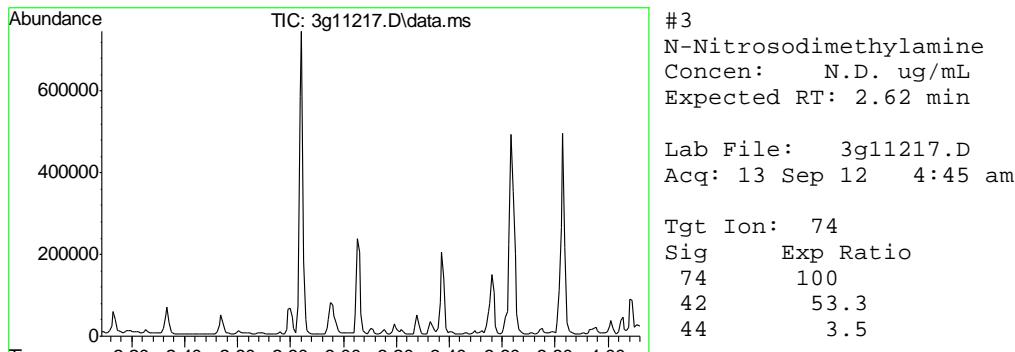
Quantitation Report (QT Reviewed)

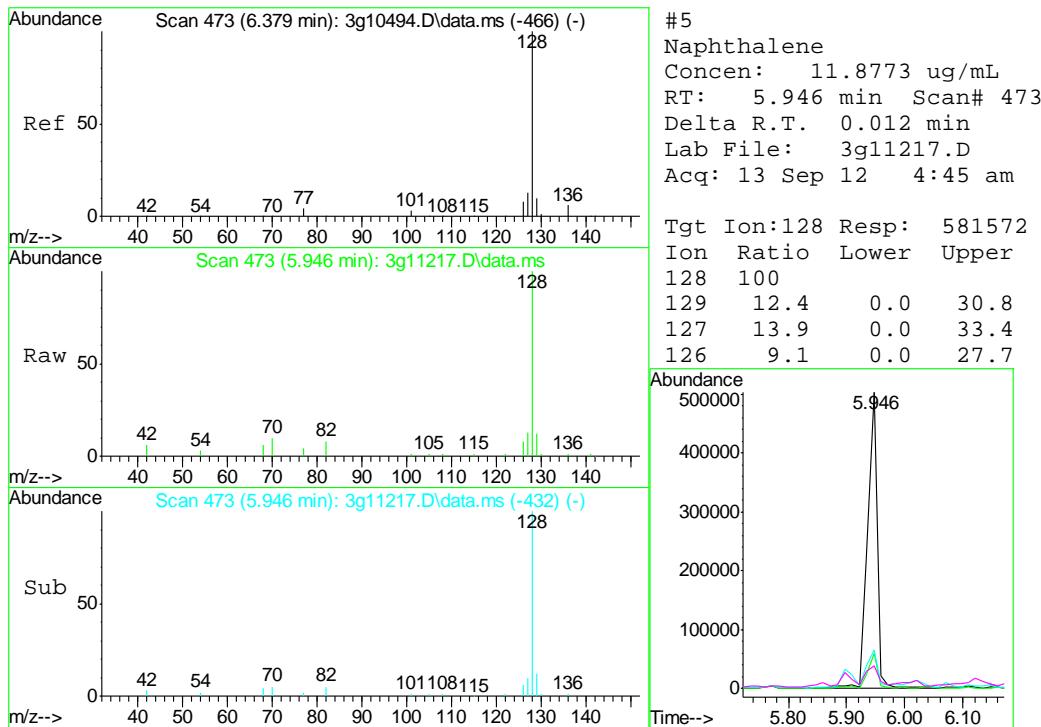
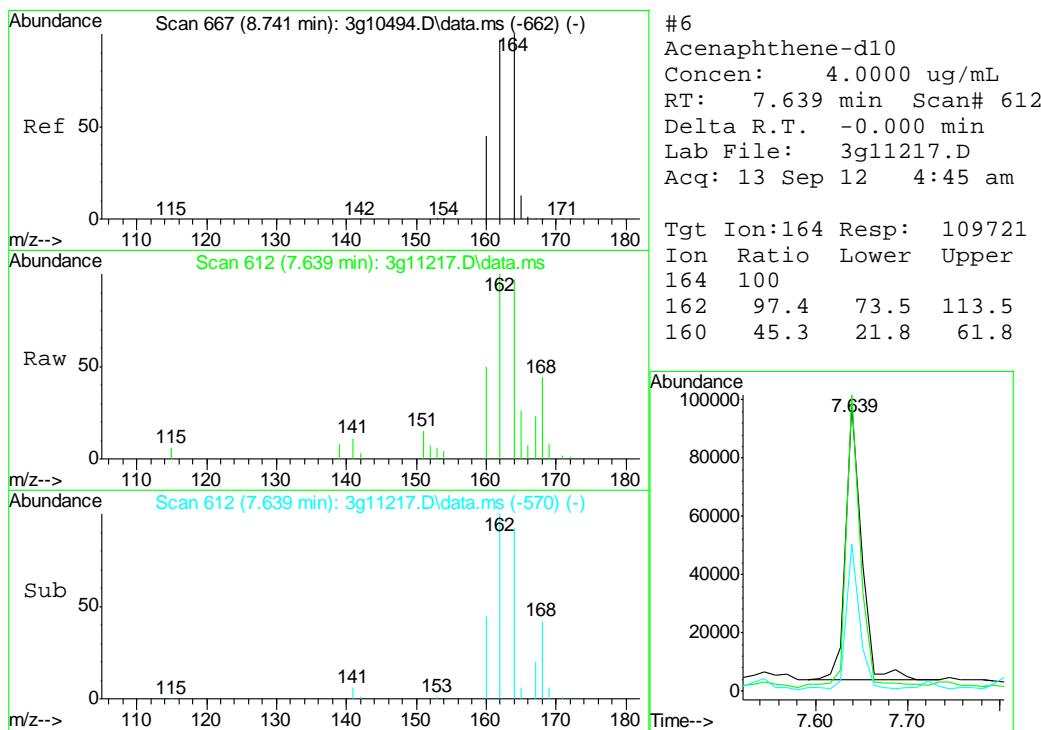
Data Path : C:\msdchem\1\DATA\091212\
 Data File : 3g11217.D
 Acq On : 13 Sep 2012 4:45 am
 Operator : DONC
 Sample : D38480-3
 Misc : OP6602,E3G522,30.17,,,1,1
 ALS Vial : 28 Sample Multiplier: 1

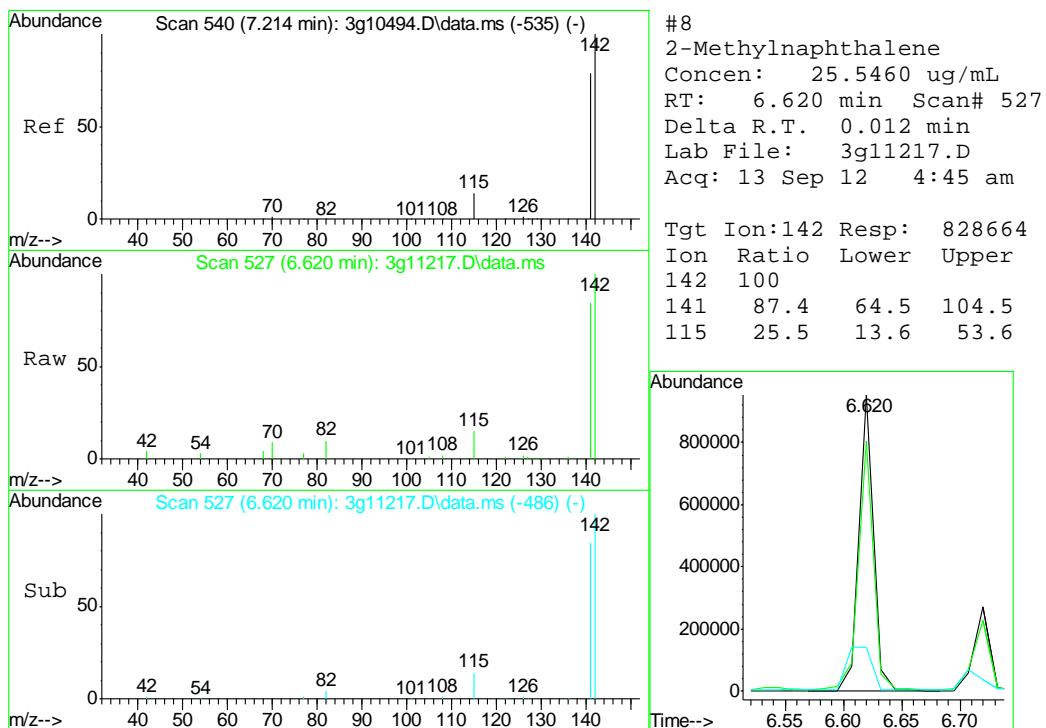
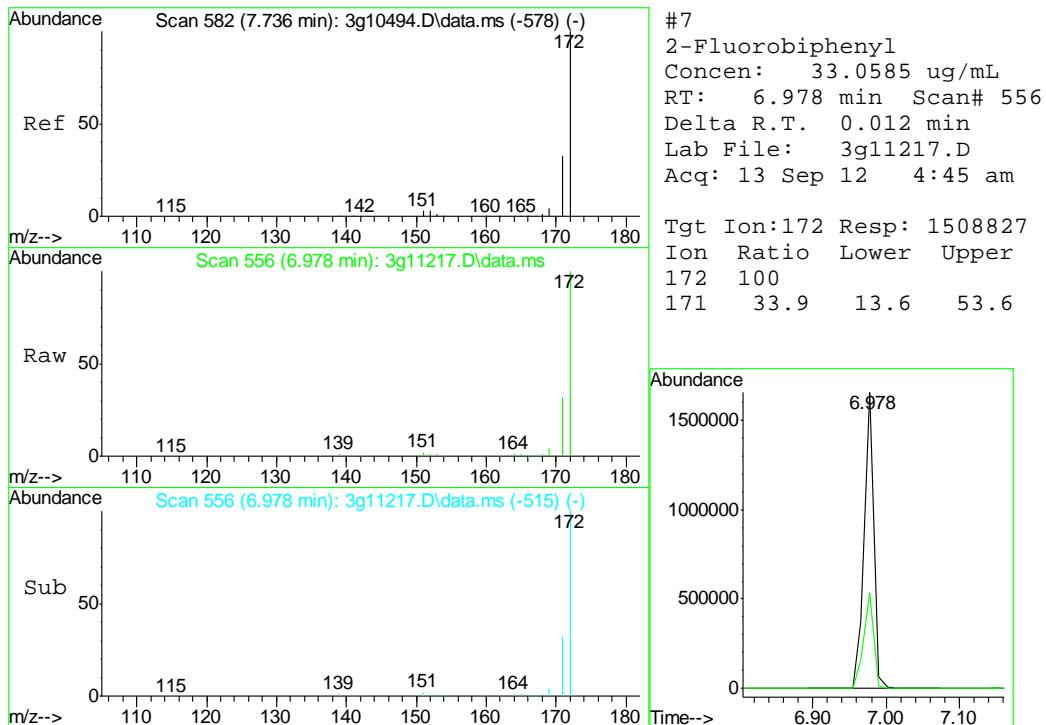
Quant Time: Sep 13 13:14:54 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Sep 06 09:42:23 2012
 Response via : Initial Calibration

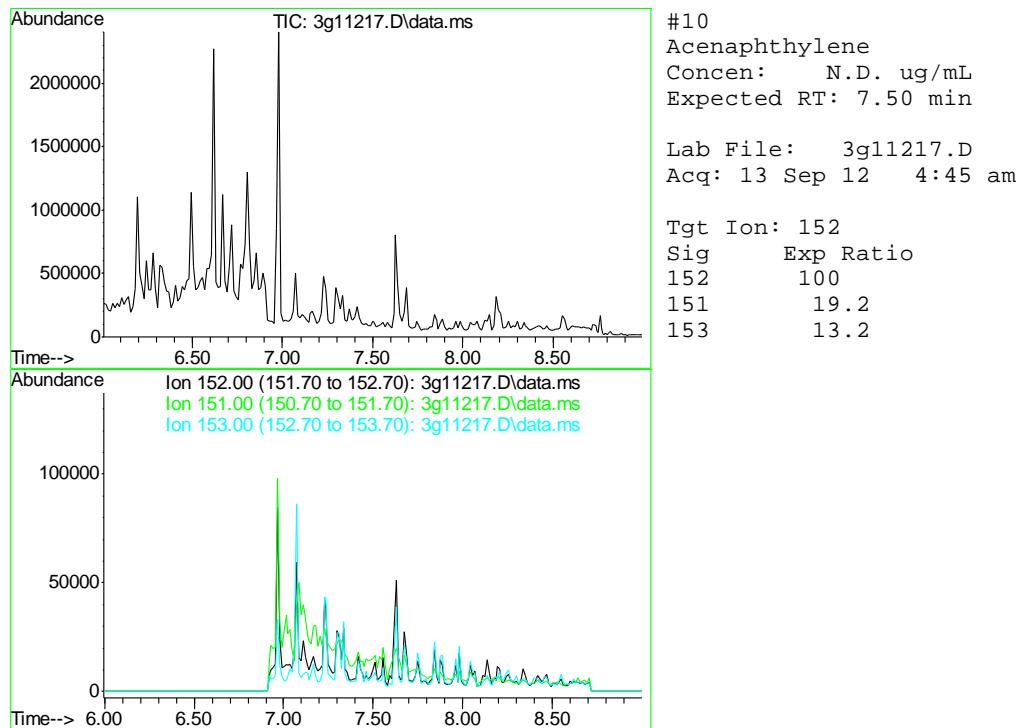
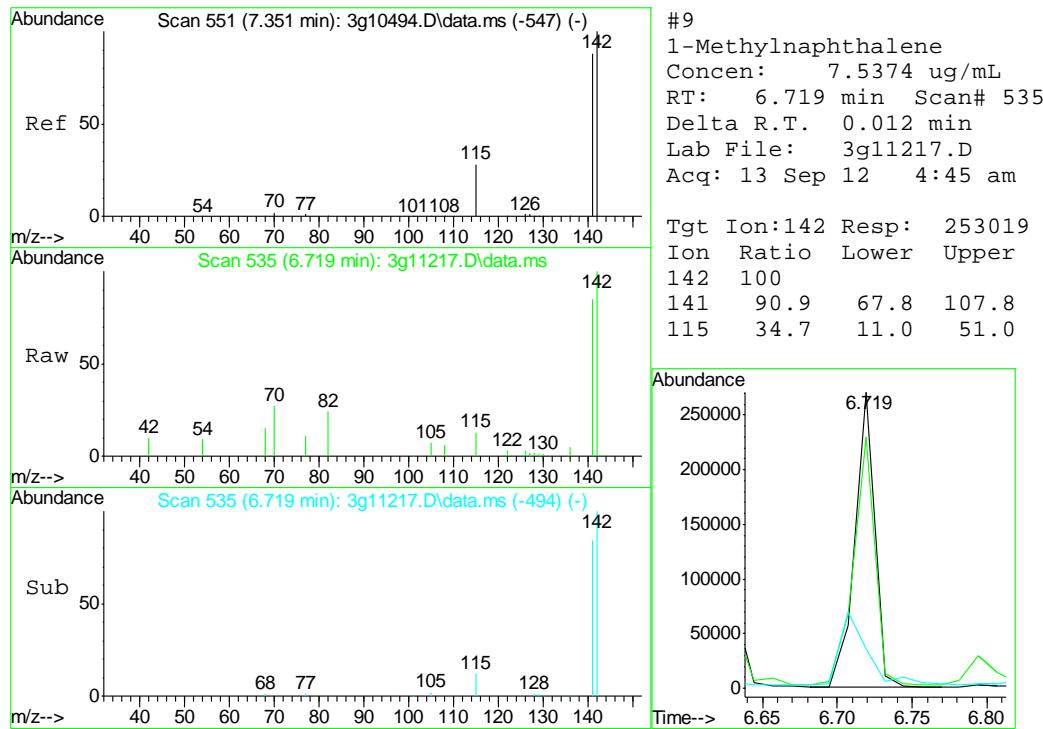


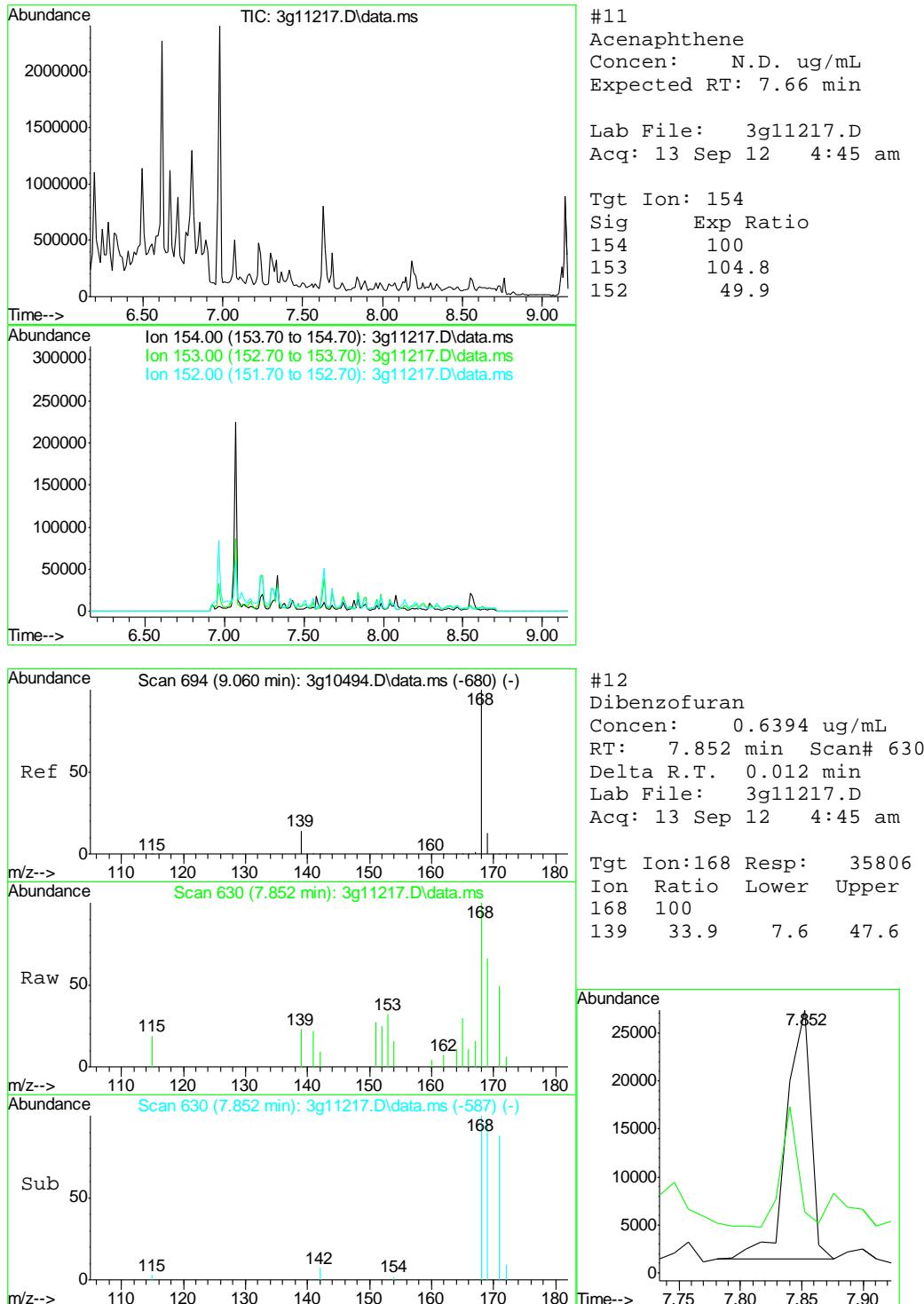


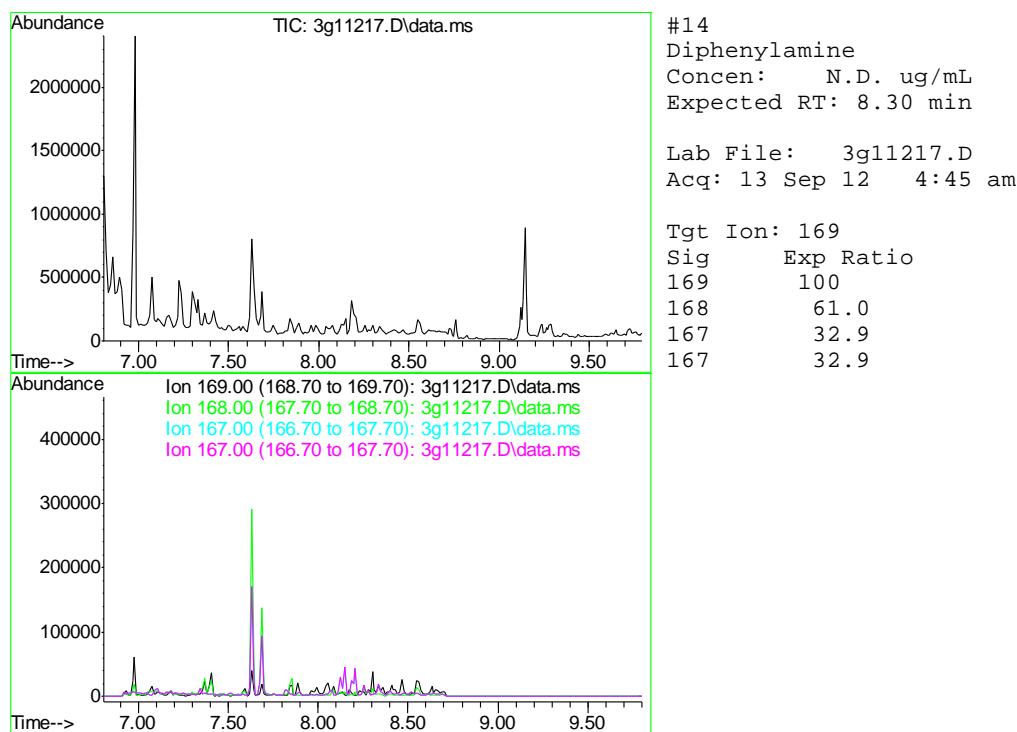
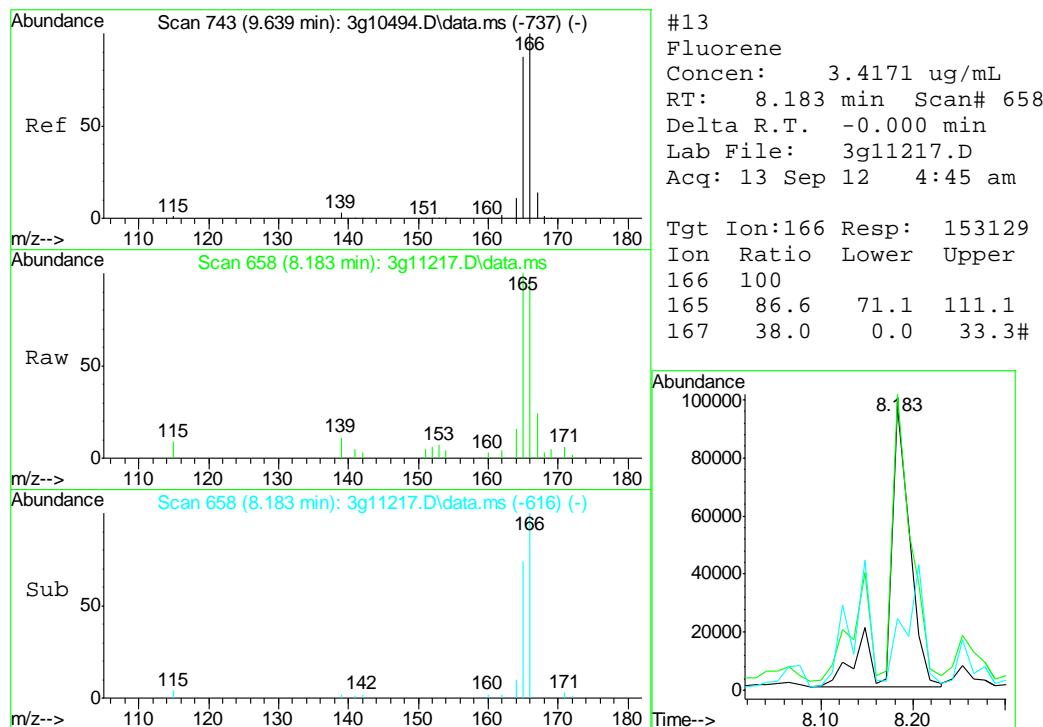


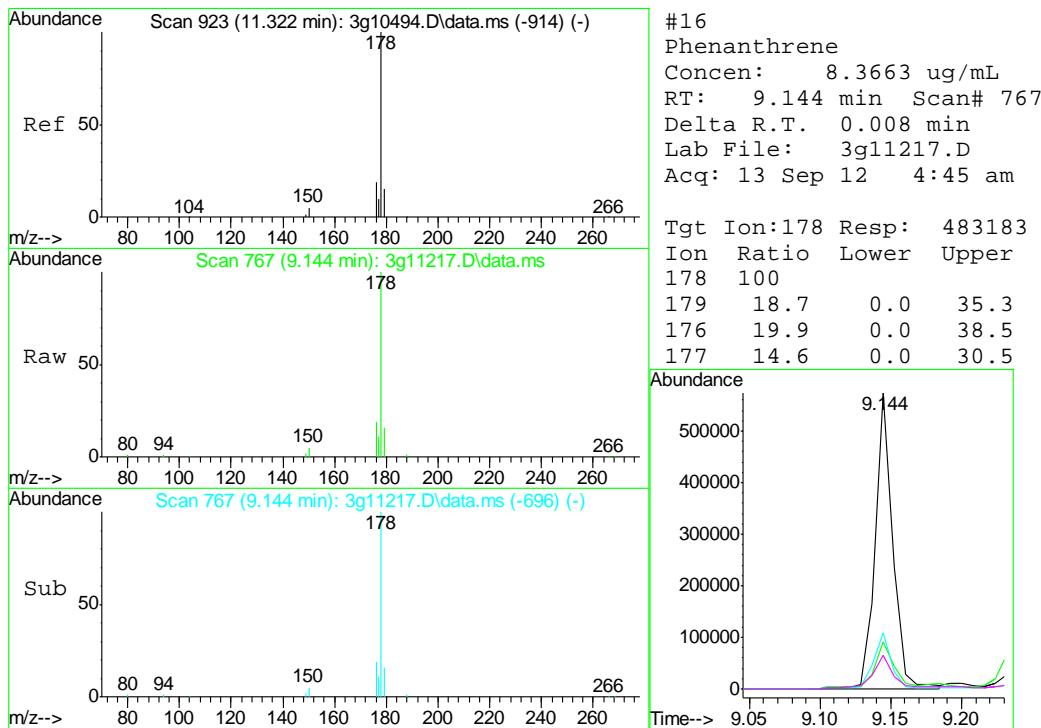
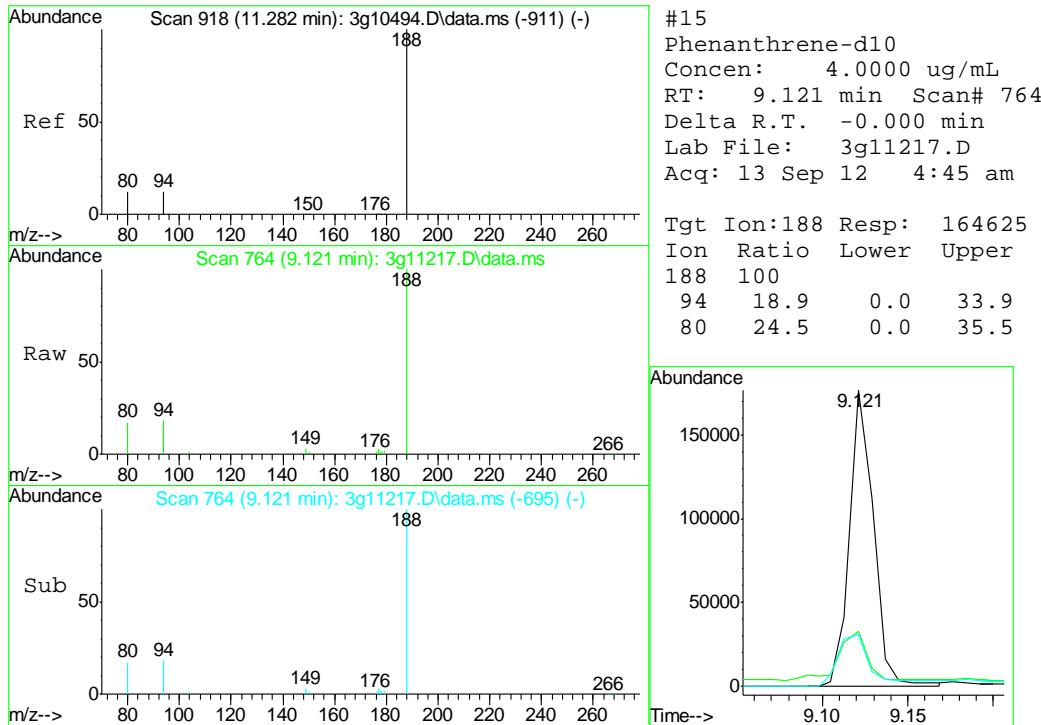
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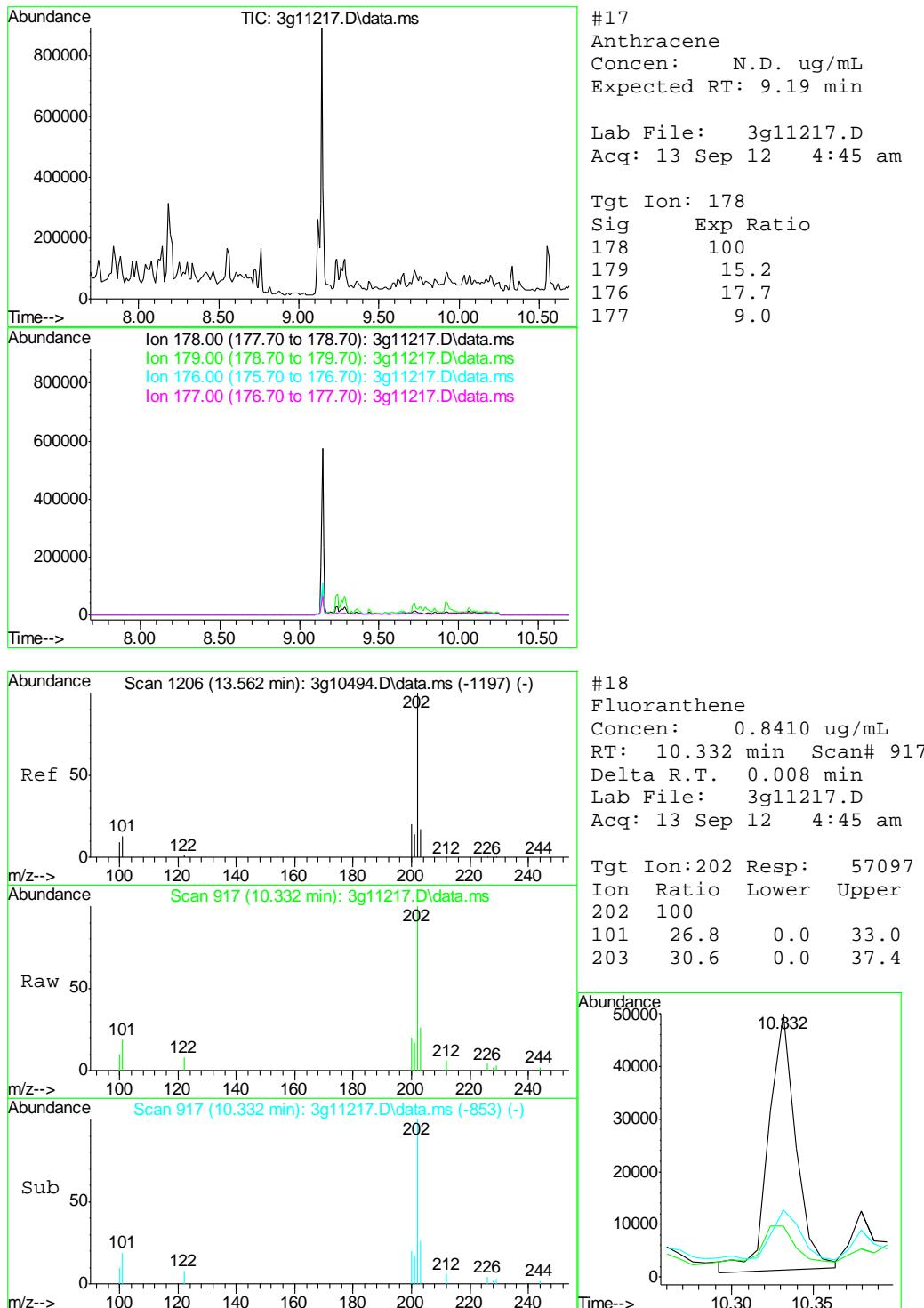


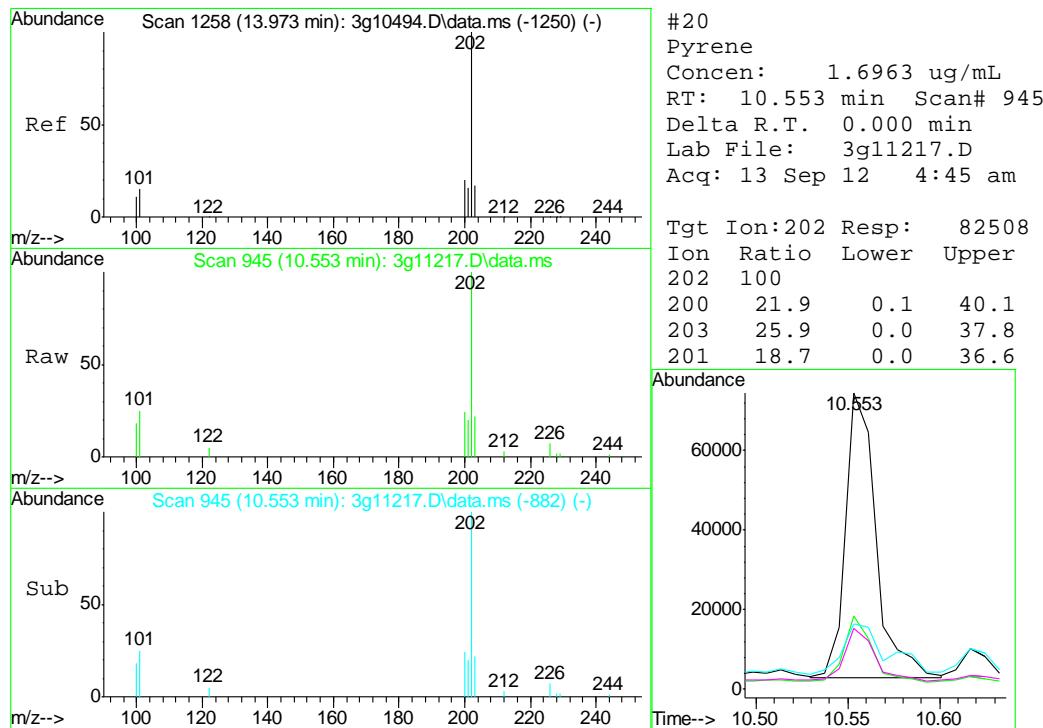
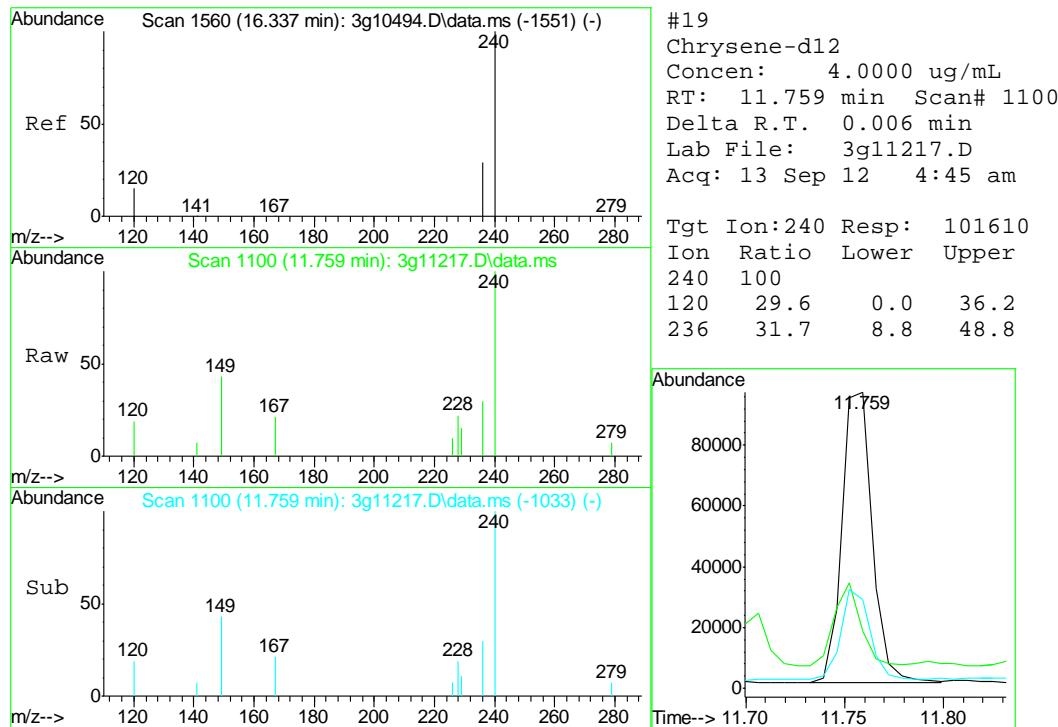


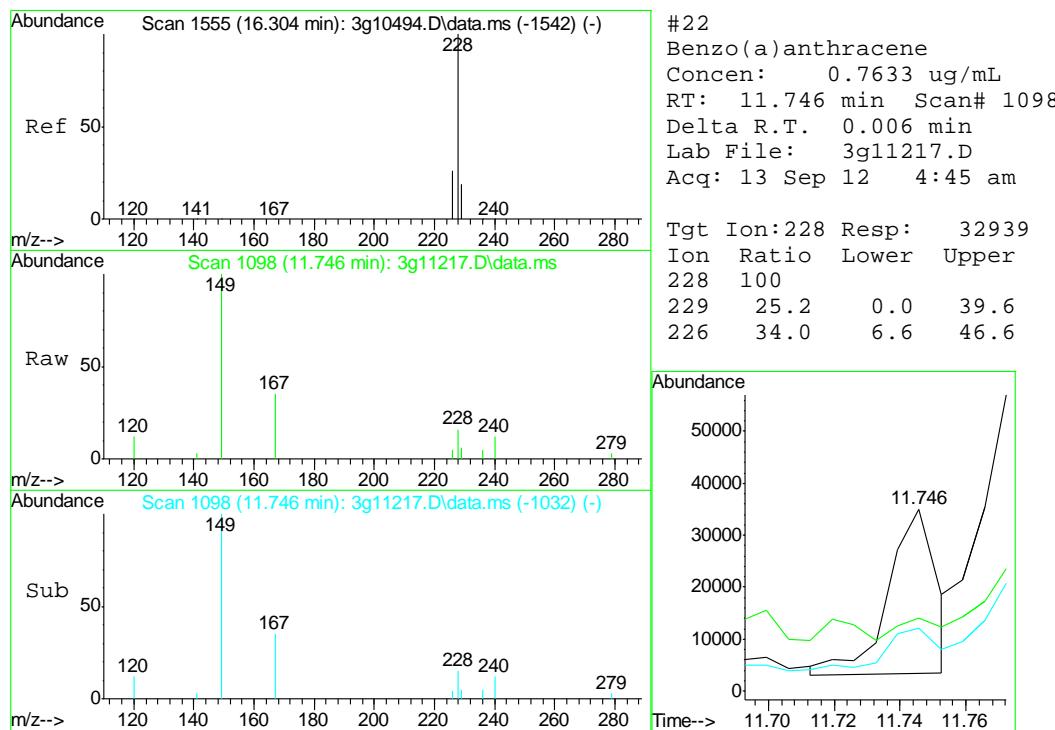
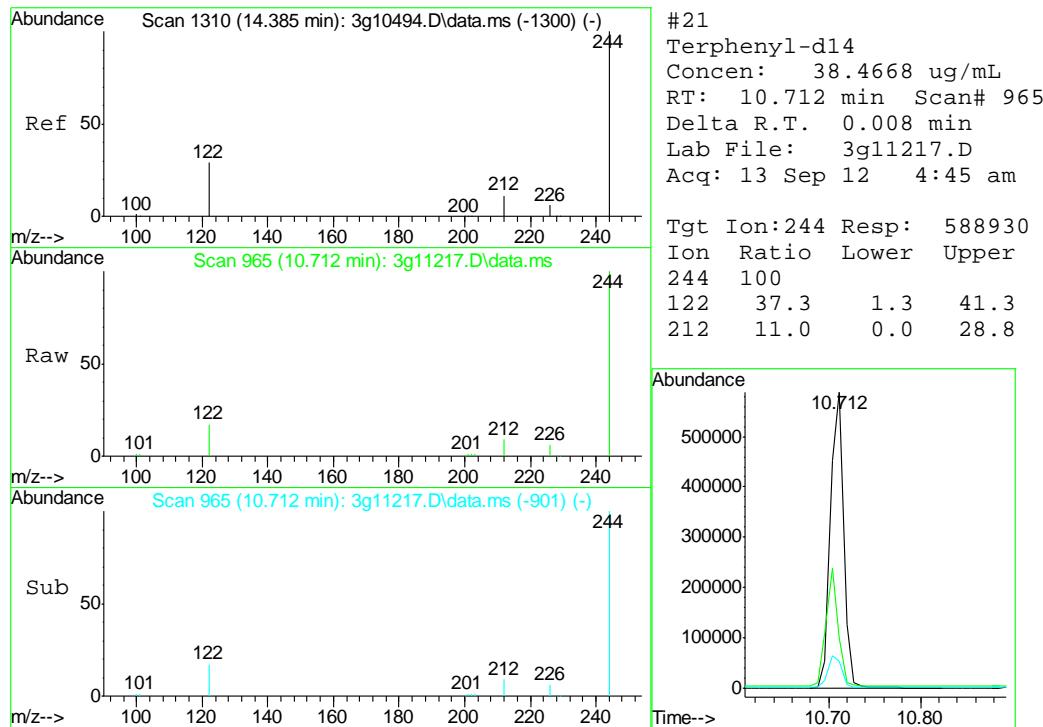


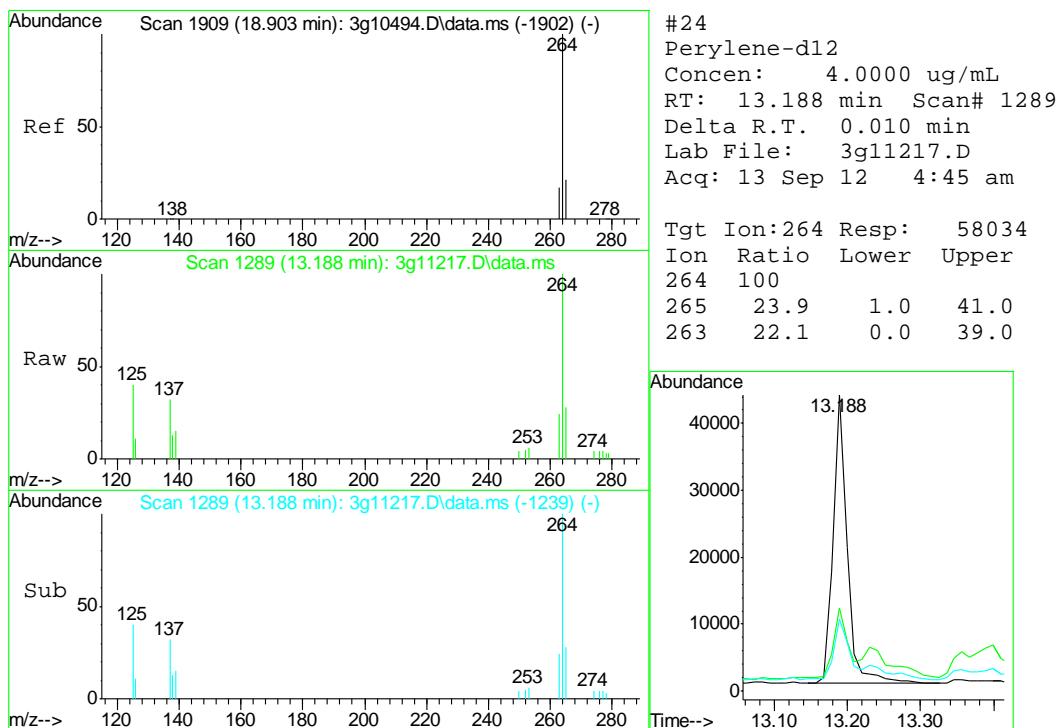
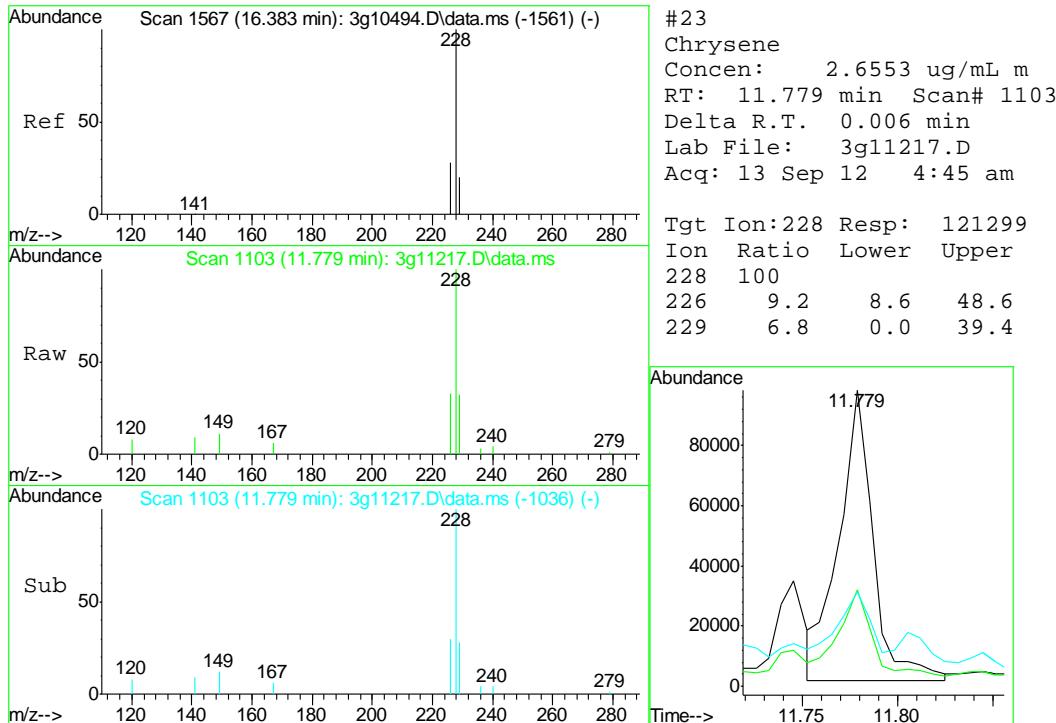


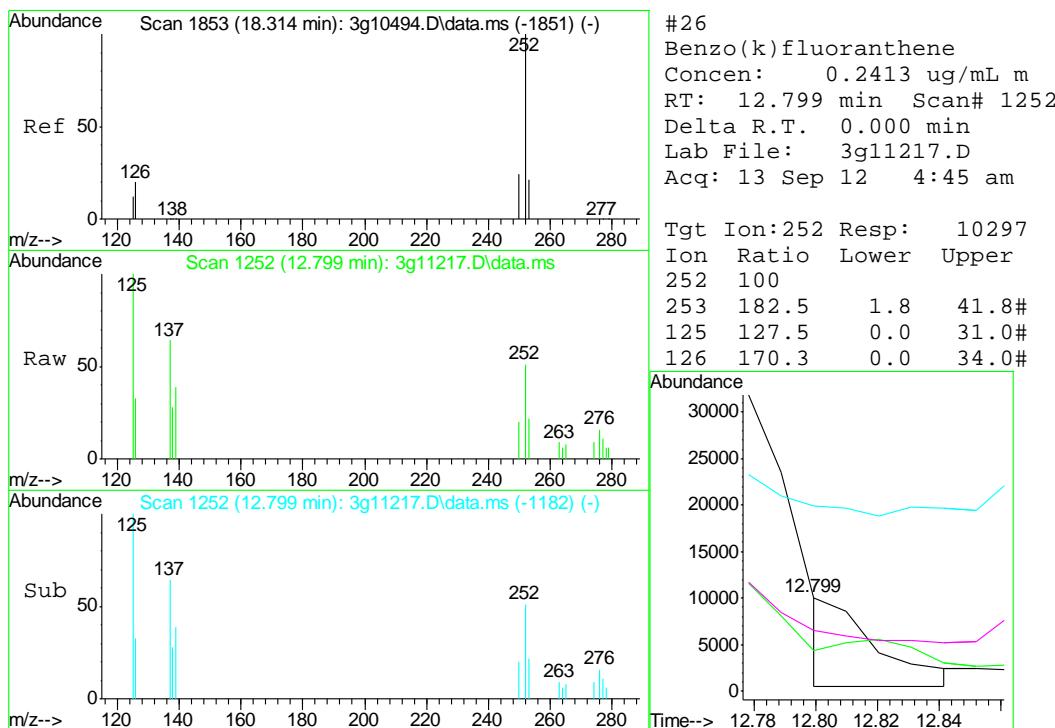
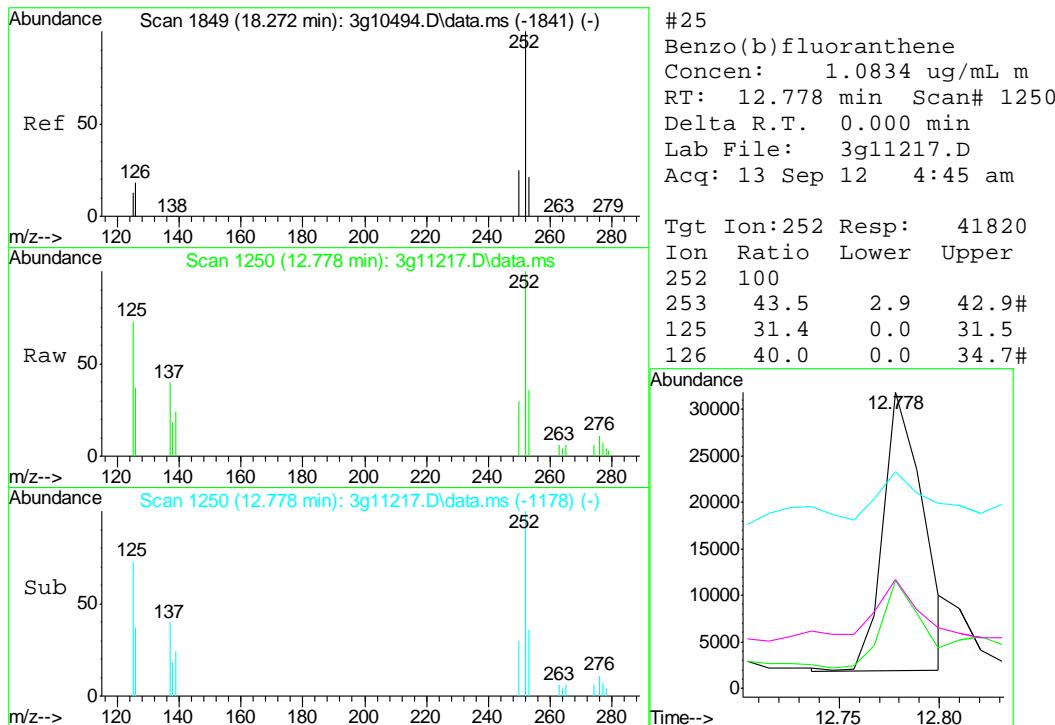


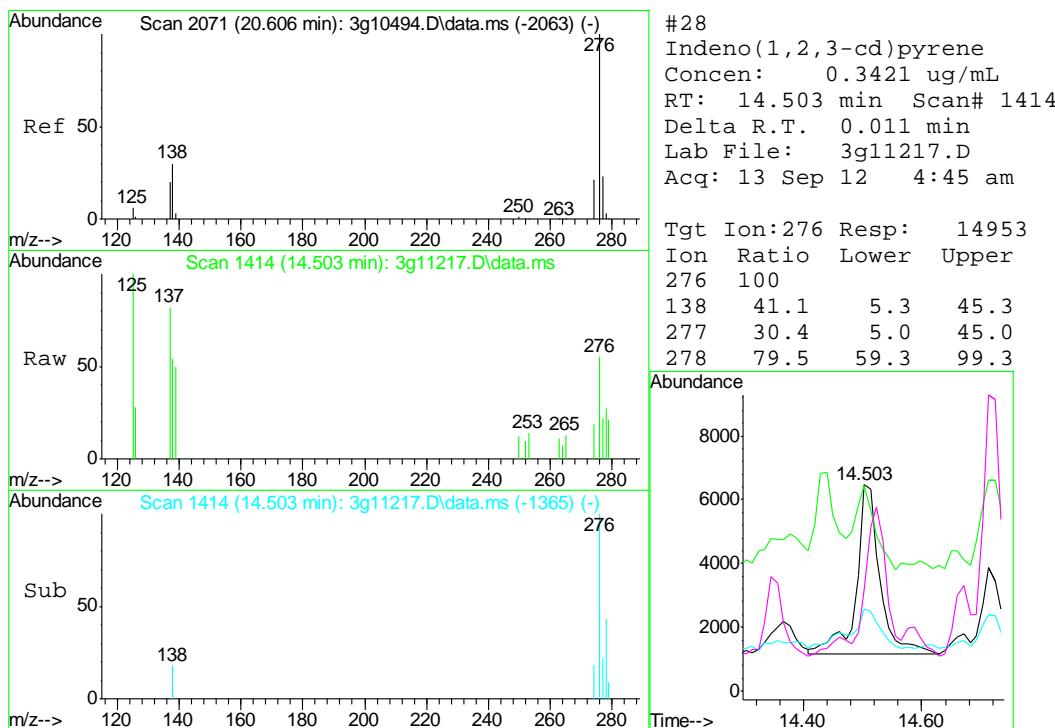
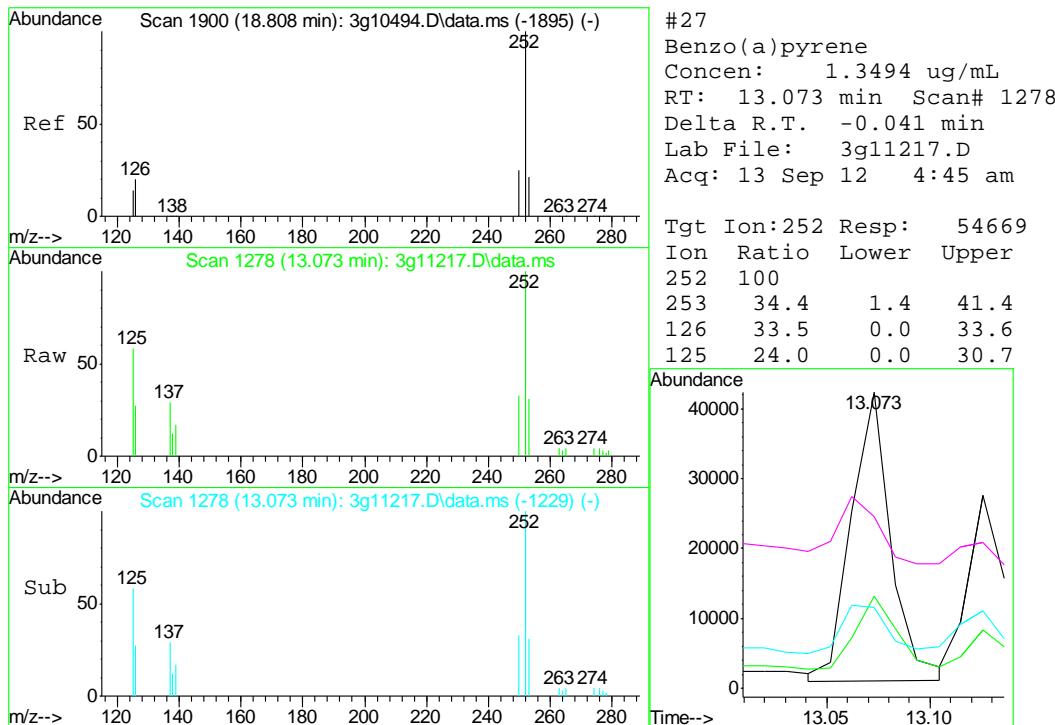


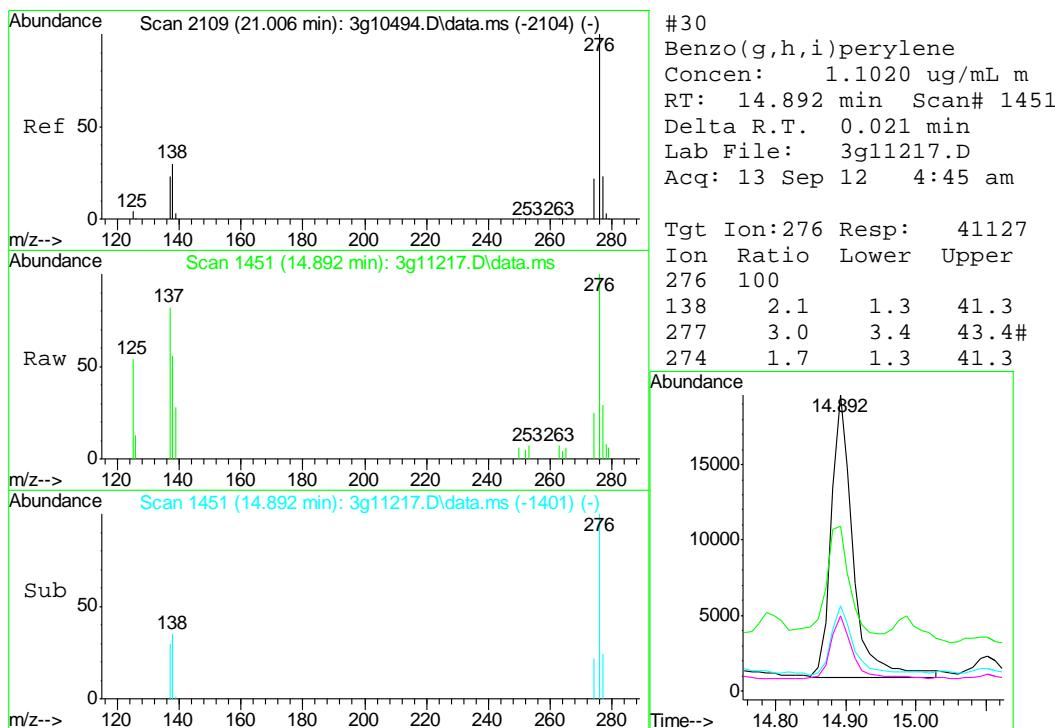
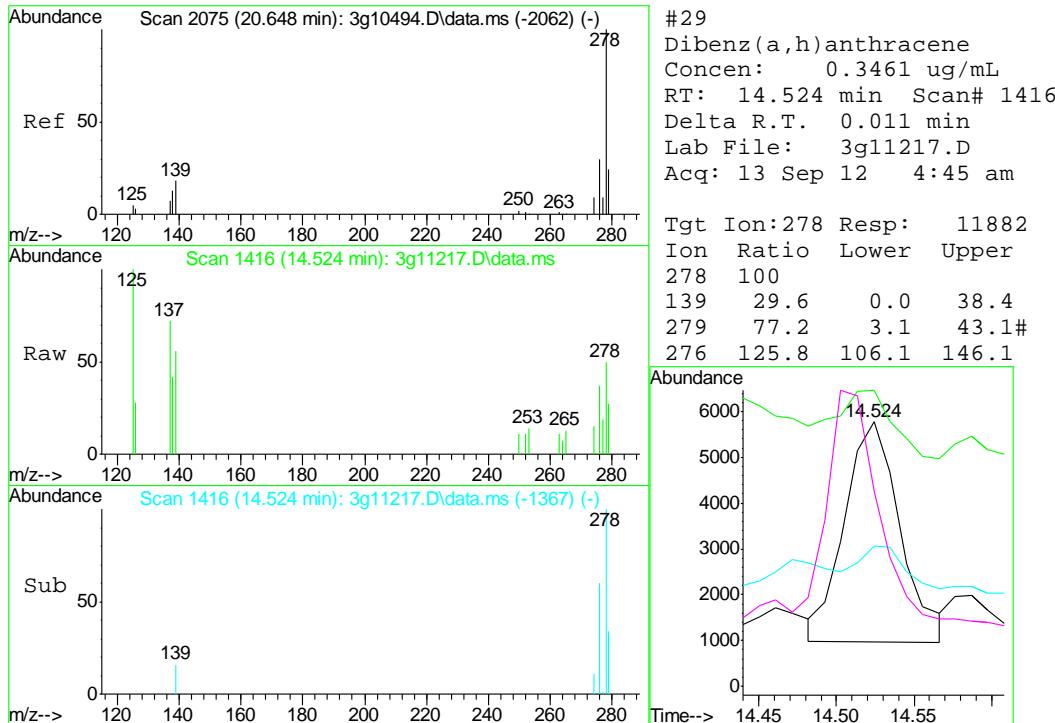












**Manual Integrations
APPROVED
(compounds with "m" flag)**
**Judy Nelson
09/13/12 14:04**

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\091112\
Data File : 3g11149.D
Acq On : 11 Sep 2012 7:39 pm
Operator : DONC
Sample : OP6602-MB
Misc : OP6602,E3G518,30.00,,,1,1
ALS Vial : 20 Sample Multiplier: 1

Quant Time: Sep 12 14:01:43 2012
Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
Quant Title : PAHSIM BASE
QLast Update : Thu Sep 06 09:42:23 2012
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.922	136	149108	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.640	164	88112	4.0000	ug/mL	0.00
15) Phenanthrene-d10	9.121	188	148358	4.0000	ug/mL	0.00
19) Chrysene-d12	11.759	240	119708	4.0000	ug/mL	0.00
24) Perylene-d12	13.189	264	64389	4.0000	ug/mL	0.01

System Monitoring Compounds

2) Nitrobenzene-d5	5.236	82	705385	48.0824	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	= 96.16%	
7) 2-Fluorobiphenyl	6.978	172	1715263	46.7981	ug/mL	0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	= 93.60%	
21) Terphenyl-d14	10.712	244	853237	47.3048	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 94.60%	

Target Compounds

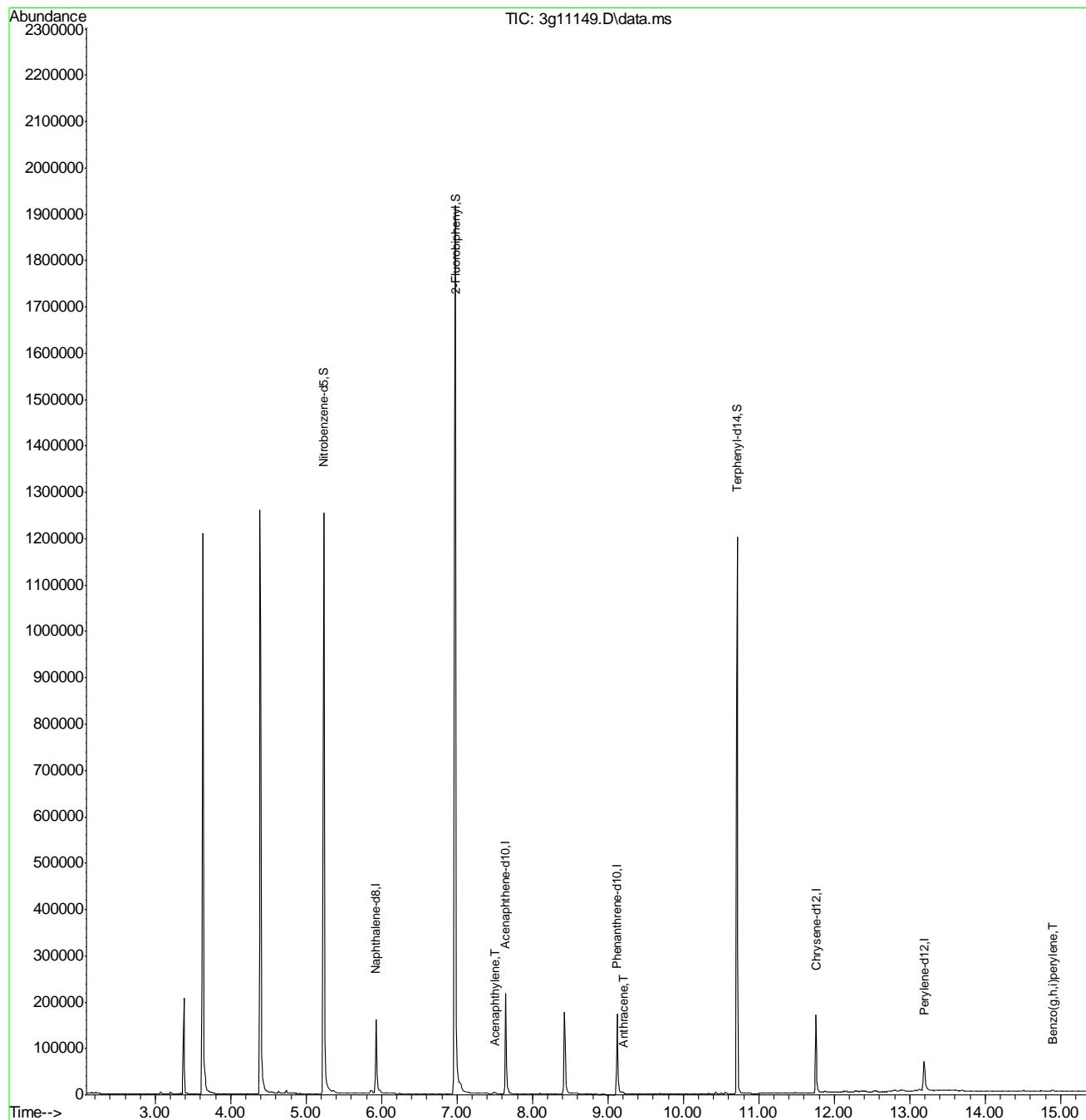
				Qvalue
3) N-Nitrosodimethylamine	2.625	74	12	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.947	128	211	N.D.
8) 2-Methylnaphthalene	6.632	142	159	N.D.
9) 1-Methylnaphthalene	6.719	142	210	N.D.
10) Acenaphthylene	7.498	152	4998m	0.1056 ug/mL
11) Acenaphthene	7.640	154	763	N.D.
12) Dibenzofuran	7.959	168	504	N.D.
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	0.000	178	0	N.D. d
17) Anthracene	9.200	178	3285m	0.0599 ug/mL
18) Fluoranthene	10.332	202	1940	N.D.
20) Pyrene	10.561	202	2753	N.D.
22) Benzo(a)anthracene	0.000	228	0	N.D. d
23) Chrysene	0.000	228	0	N.D. d
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d
26) Benzo(k)fluoranthene	0.000	252	0	N.D. d
27) Benzo(a)pyrene	13.125	252	1631	N.D.
28) Indeno(1,2,3-cd)pyrene	14.514	276	2365	N.D.
29) Dibenz(a,h)anthracene	14.524	278	420	N.D.
30) Benzo(g,h,i)perylene	14.892	276	3754m	0.0907 ug/mL

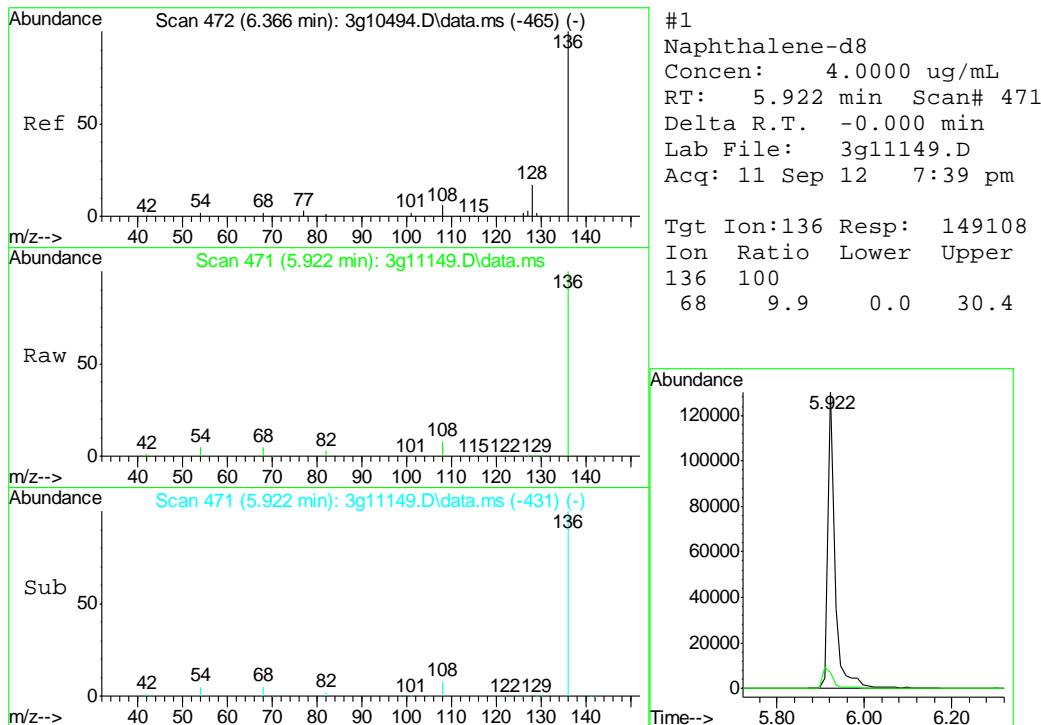
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\091112\
 Data File : 3g11149.D
 Acq On : 11 Sep 2012 7:39 pm
 Operator : DONC
 Sample : OP6602-MB
 Misc : OP6602,E3G518,30.00,,,1,1
 ALS Vial : 20 Sample Multiplier: 1

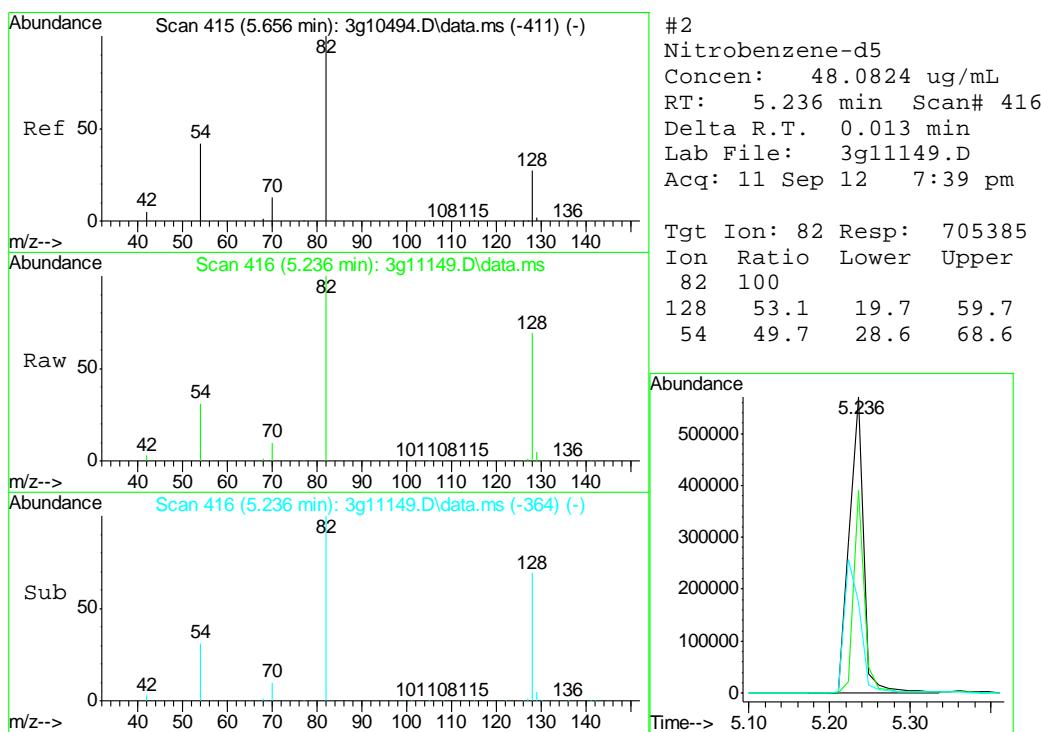
Quant Time: Sep 12 14:01:43 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G511.M
 Quant Title : PAHSIM BASE
 QLast Update : Thu Sep 06 09:42:23 2012
 Response via : Initial Calibration

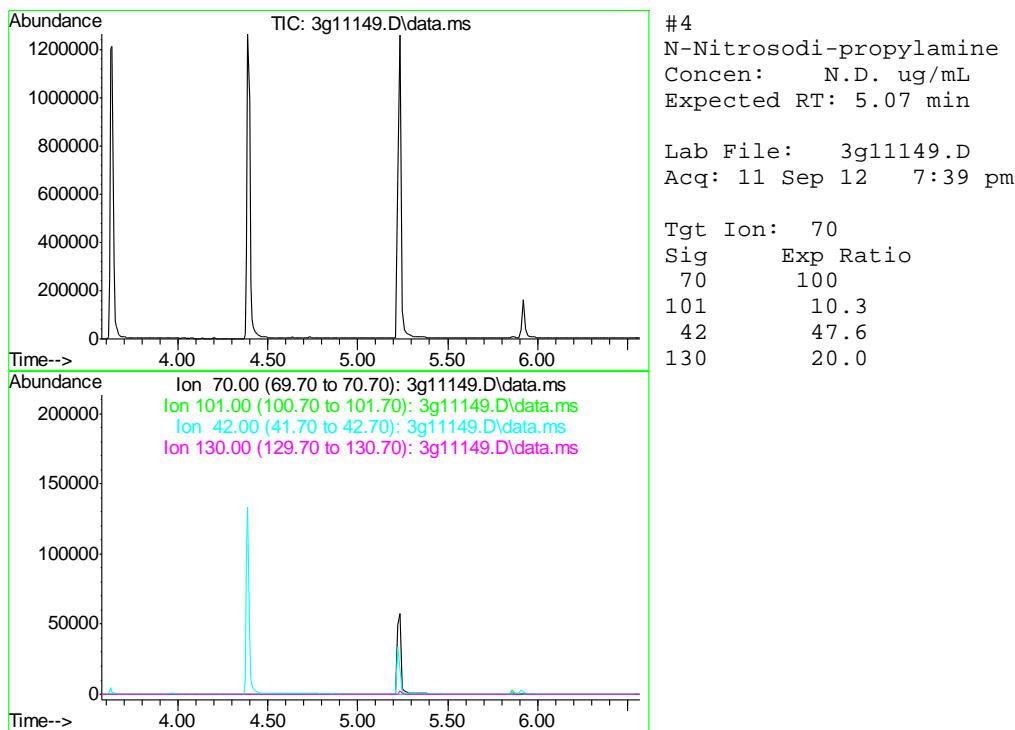
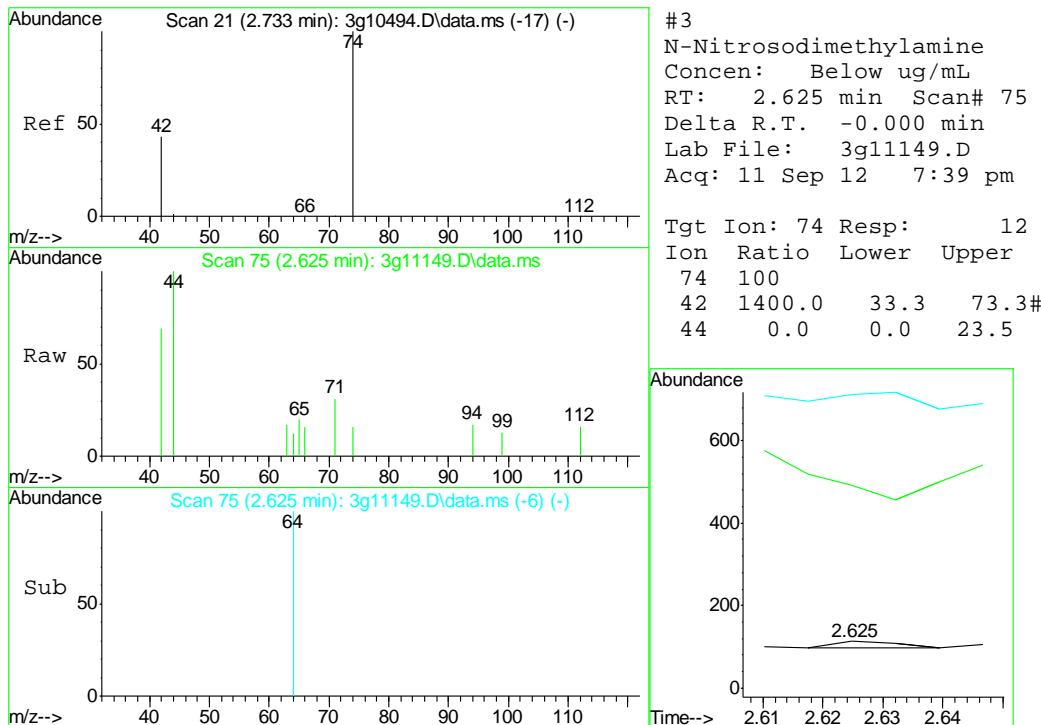


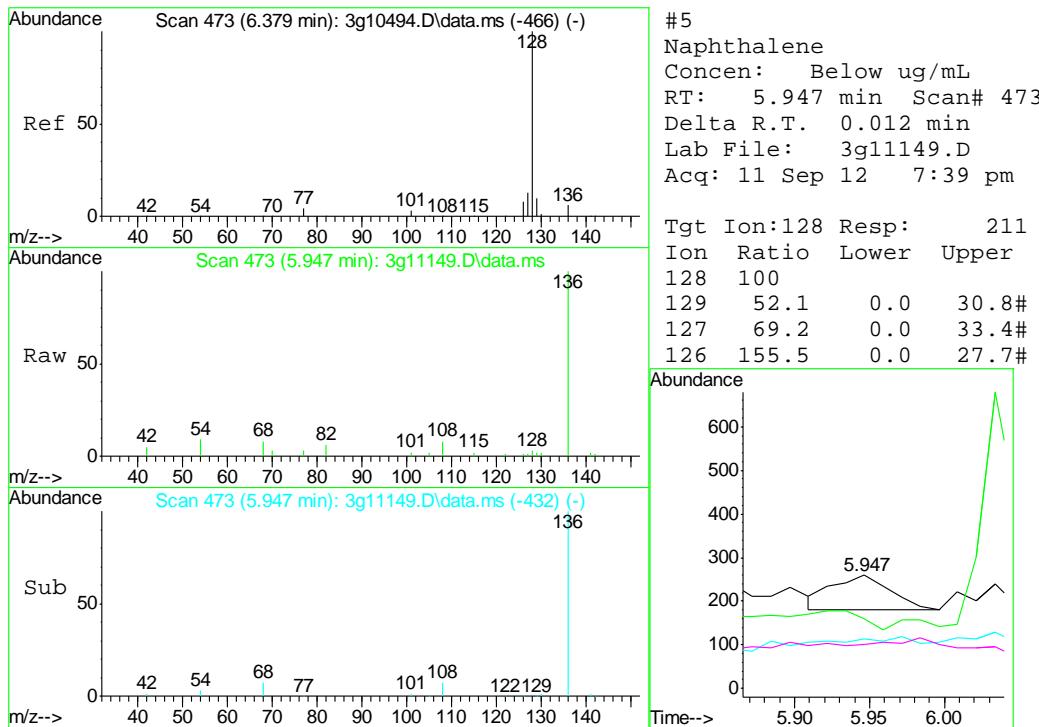


9.2.1

9

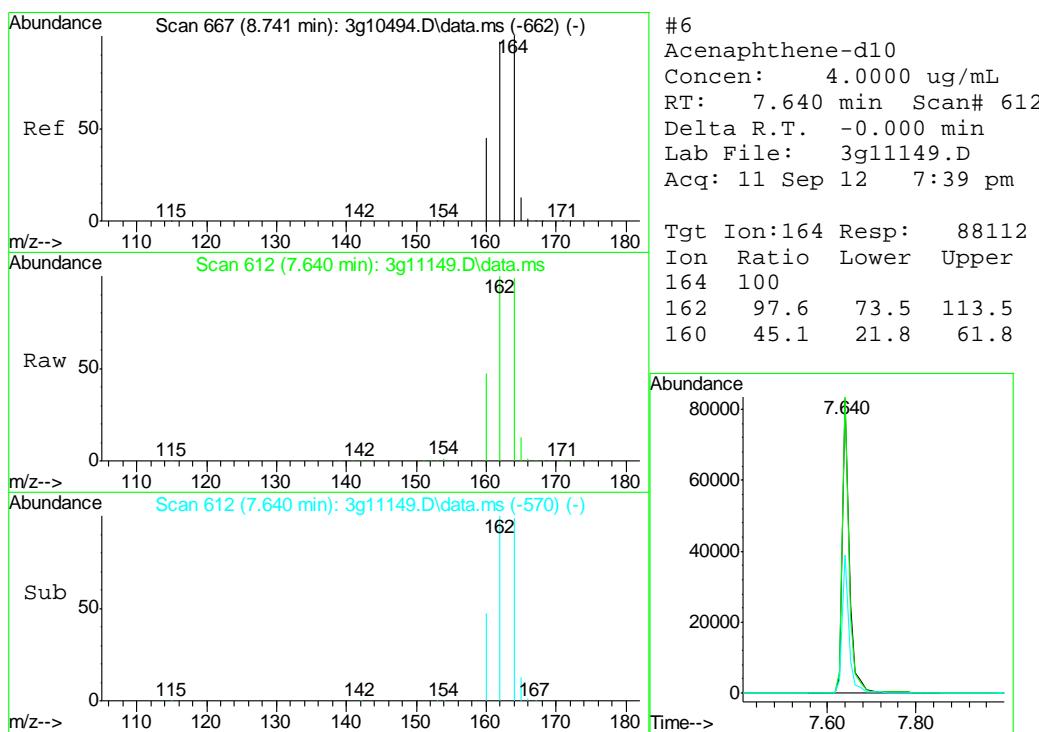


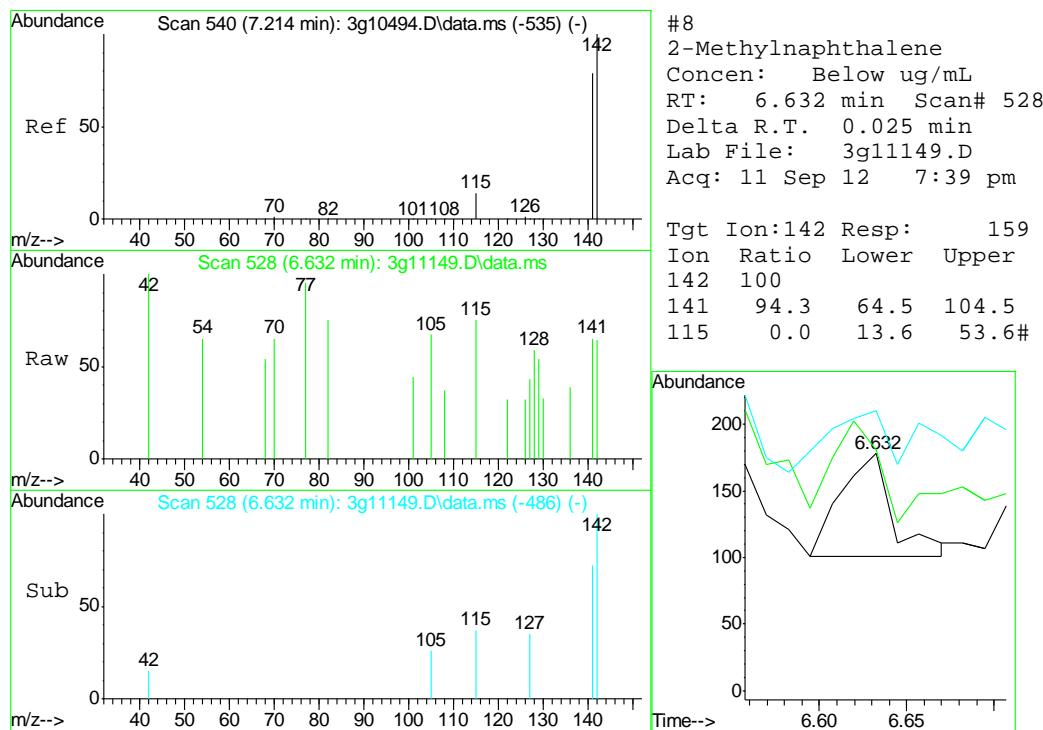
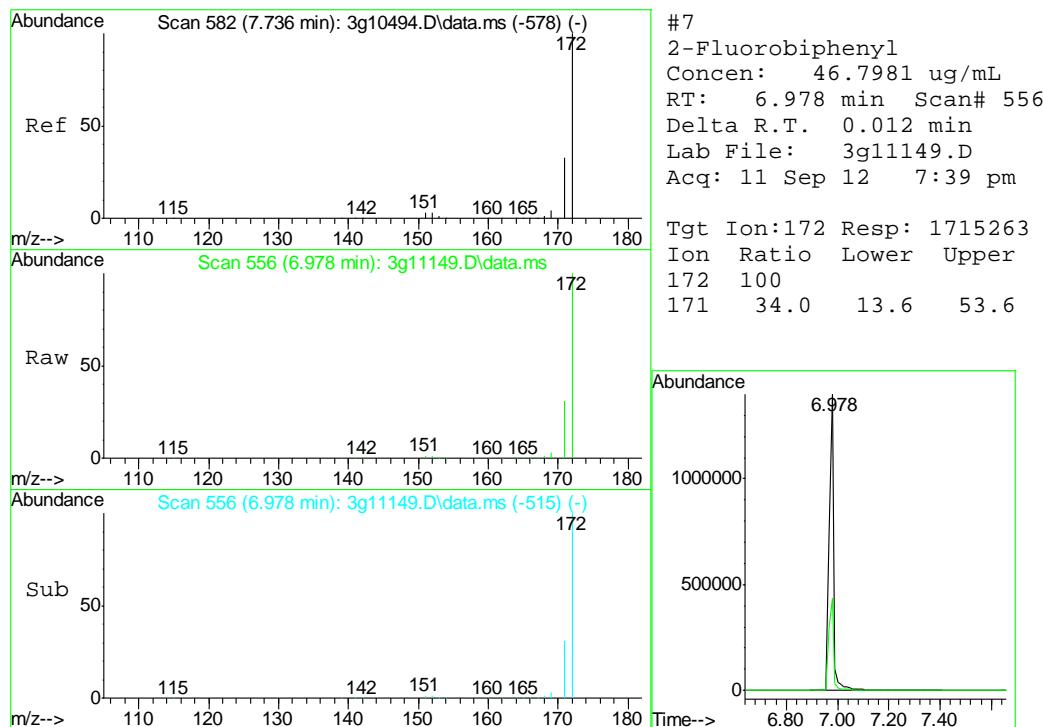


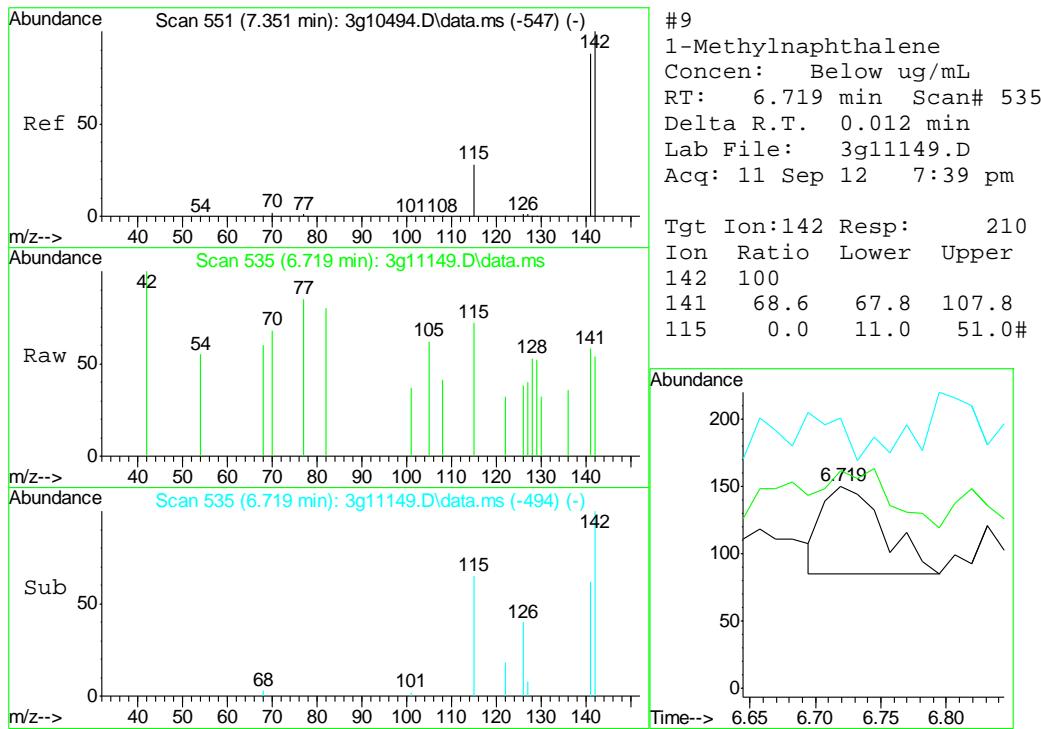


9.2.1

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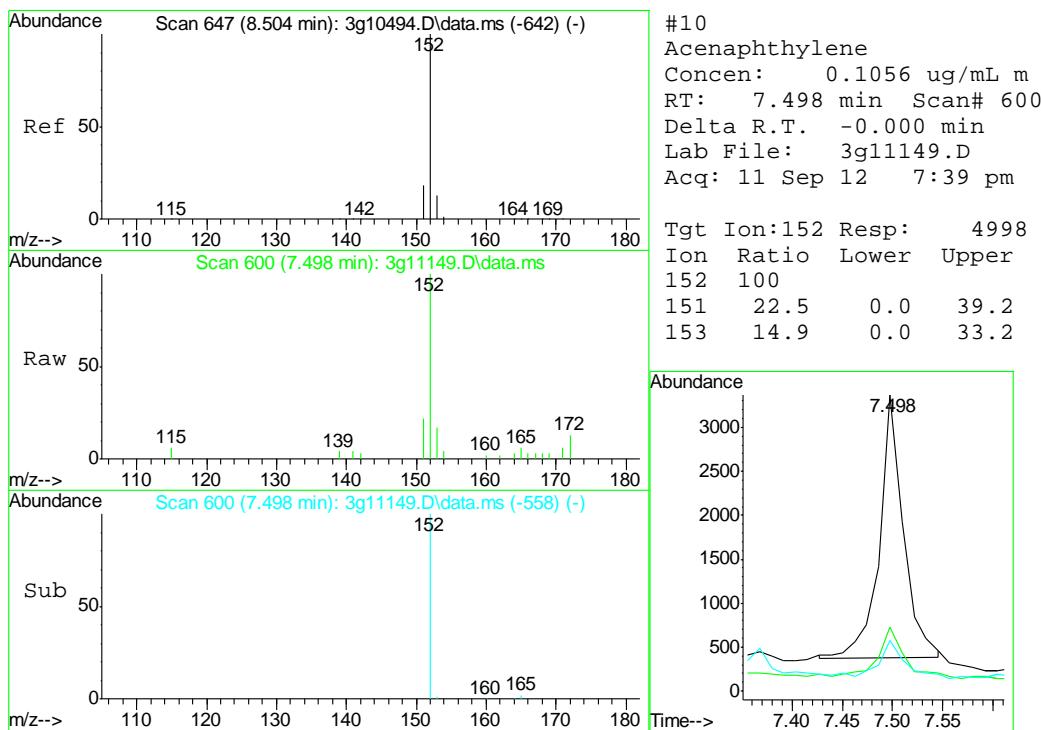


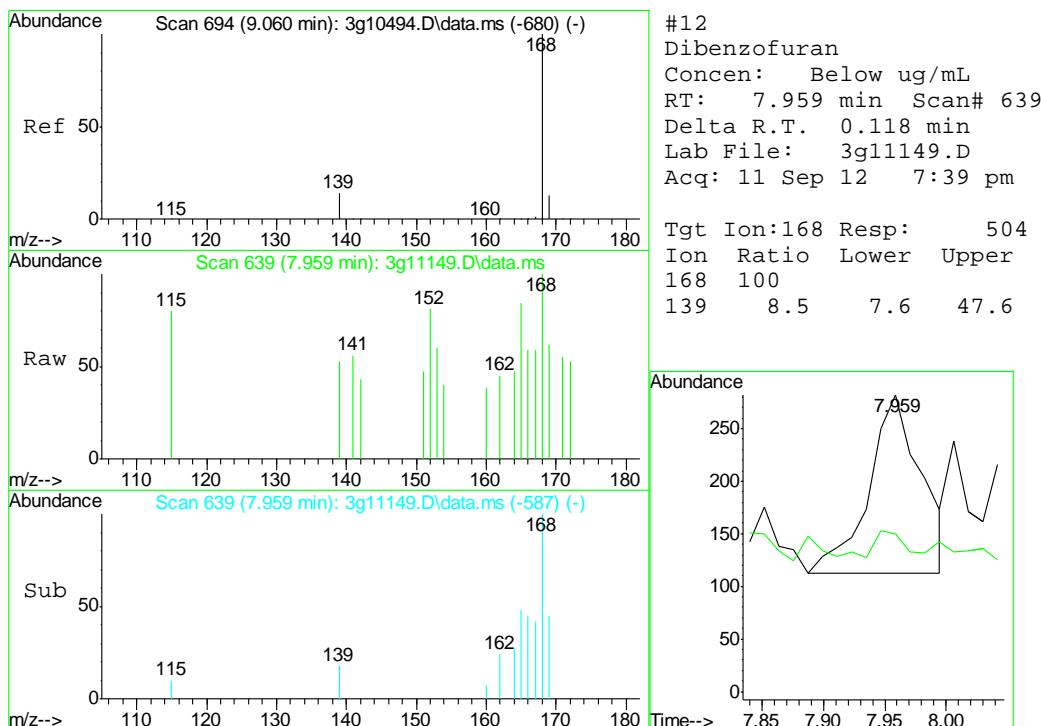
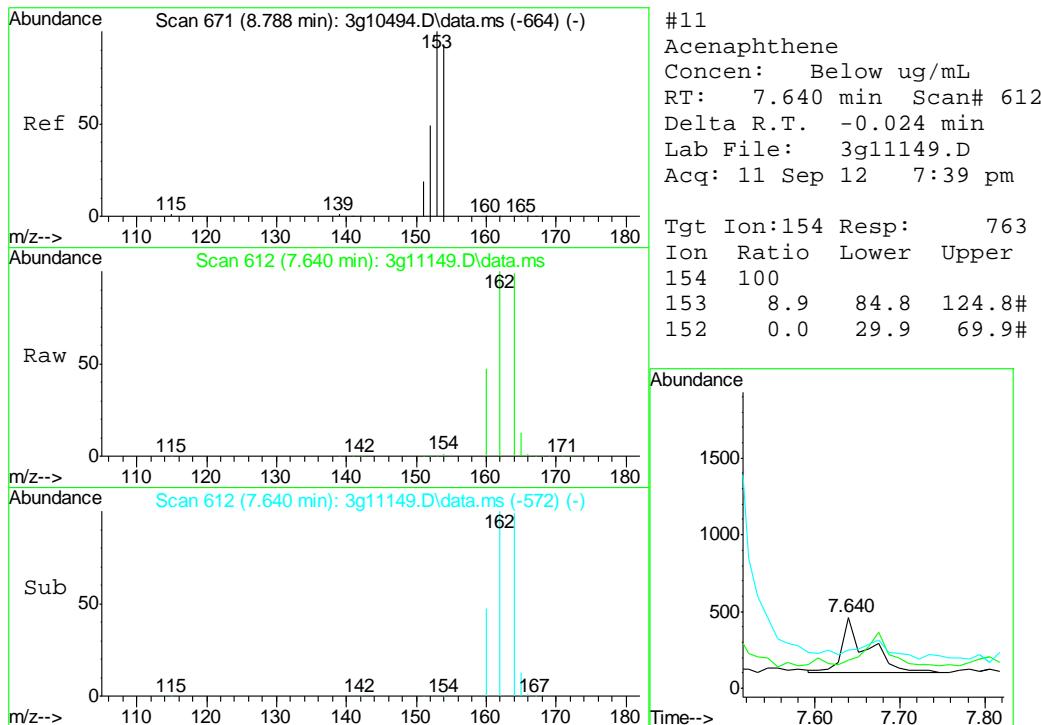


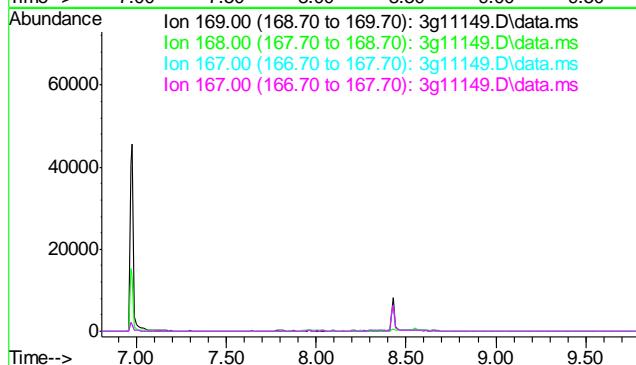
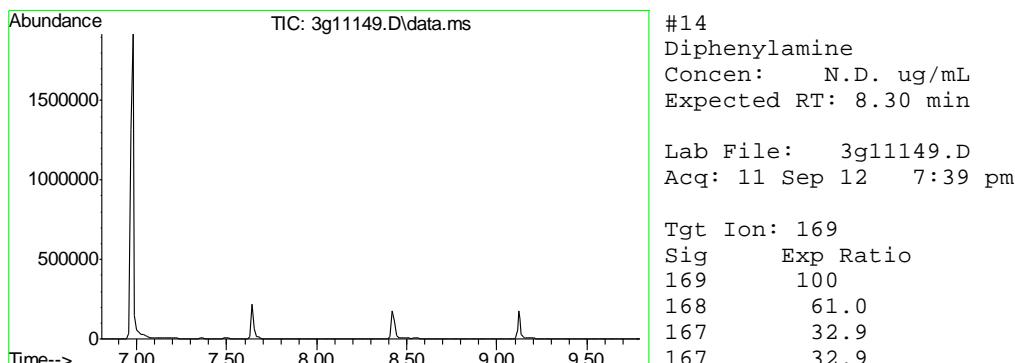
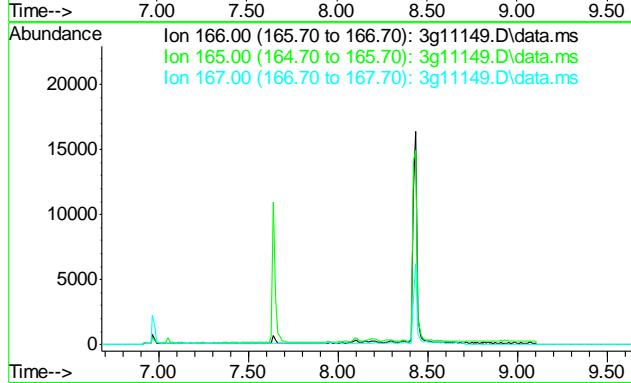
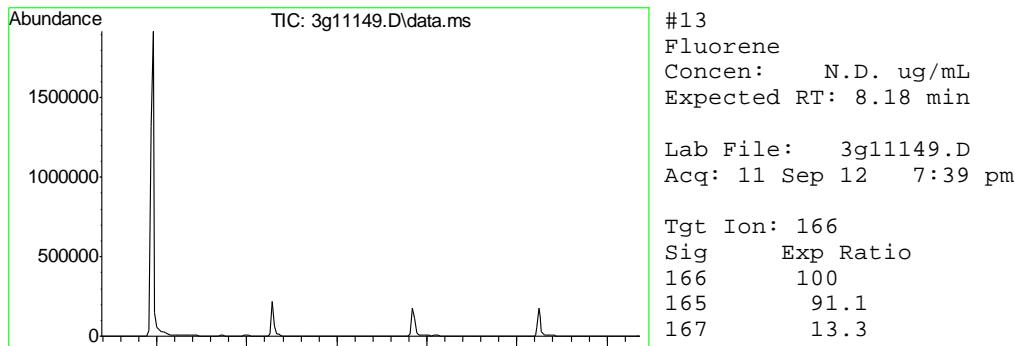


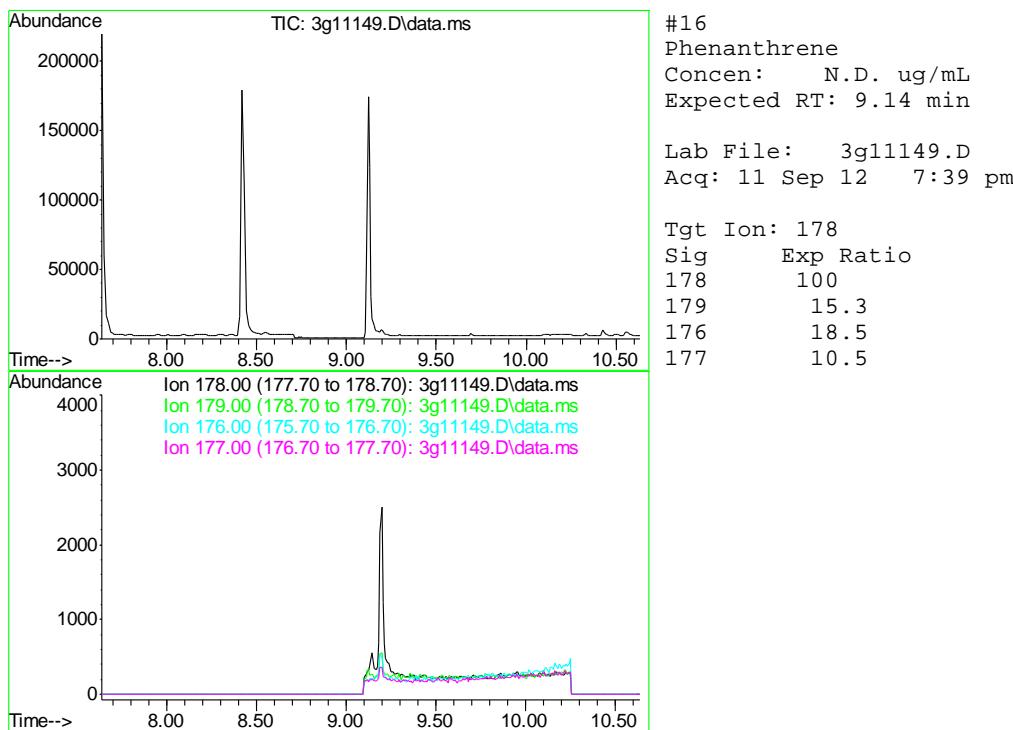
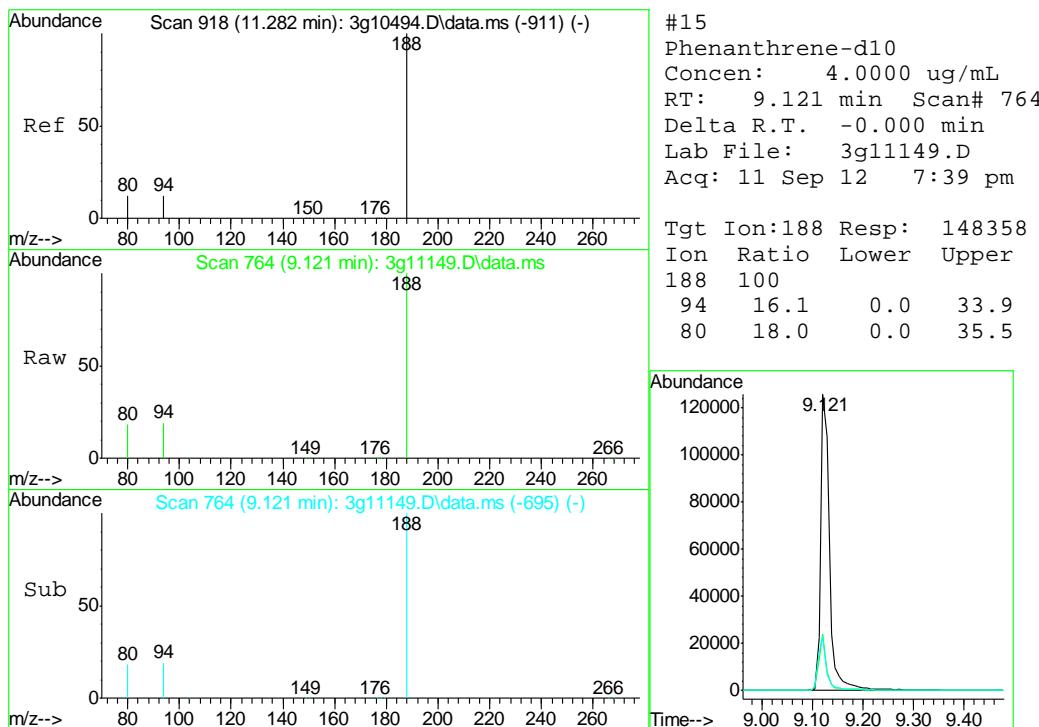
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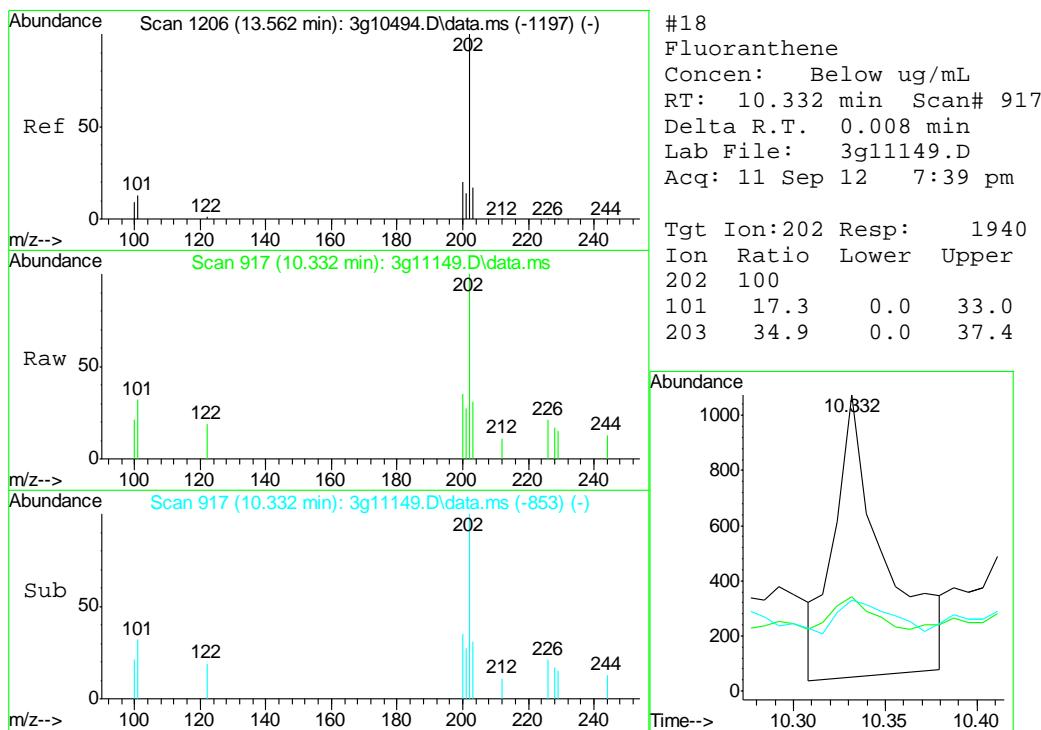
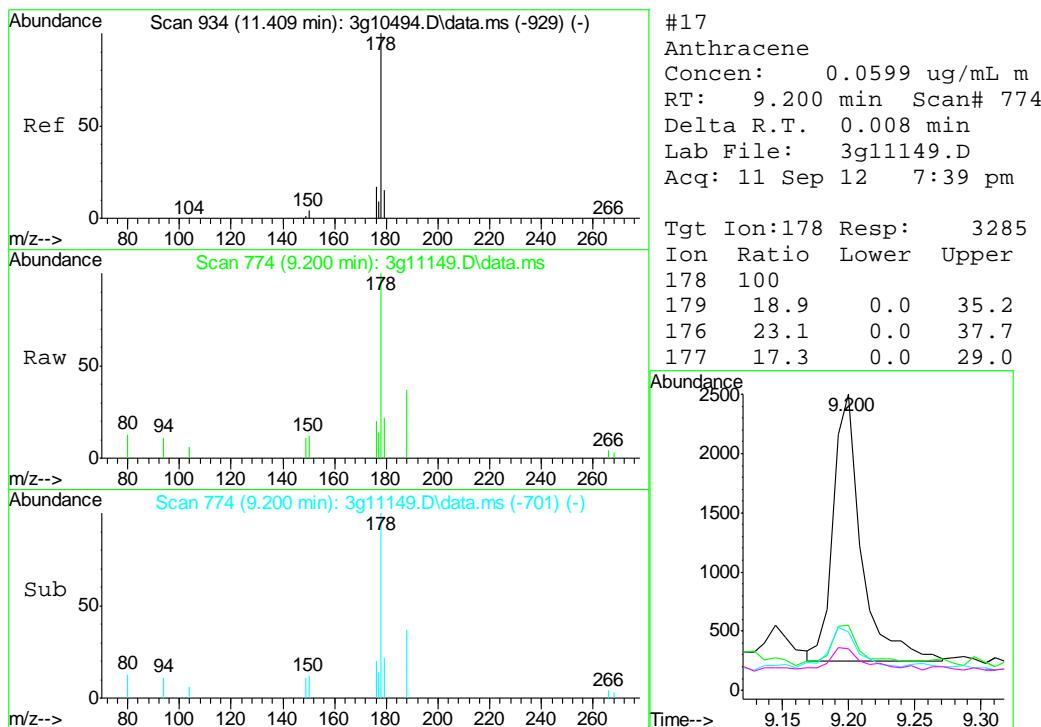
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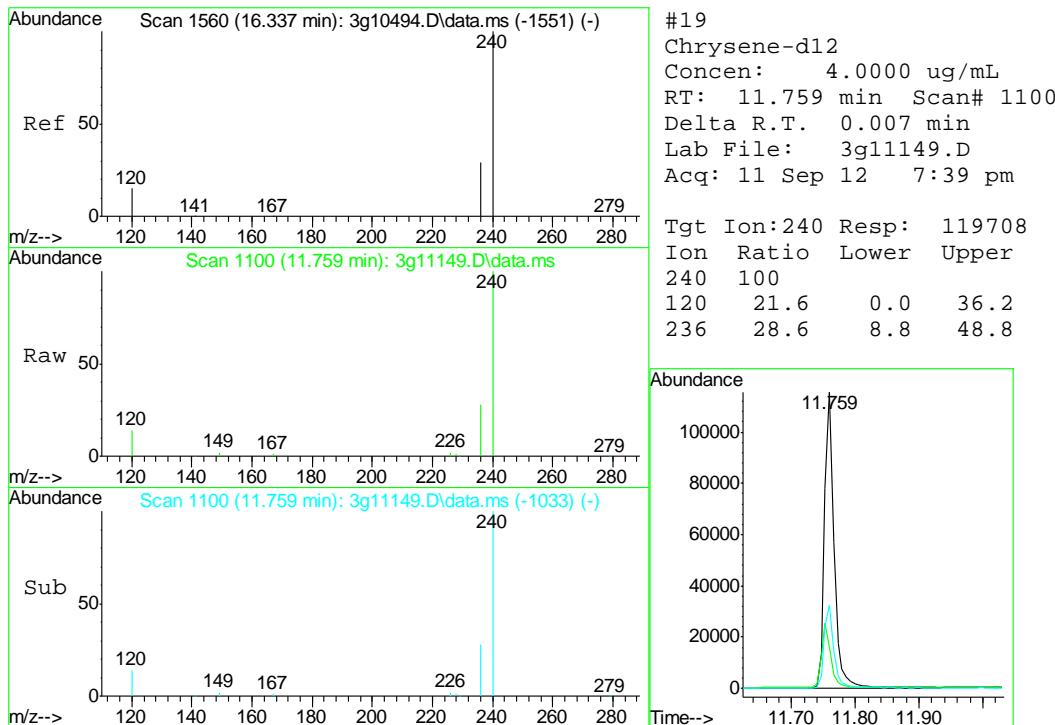






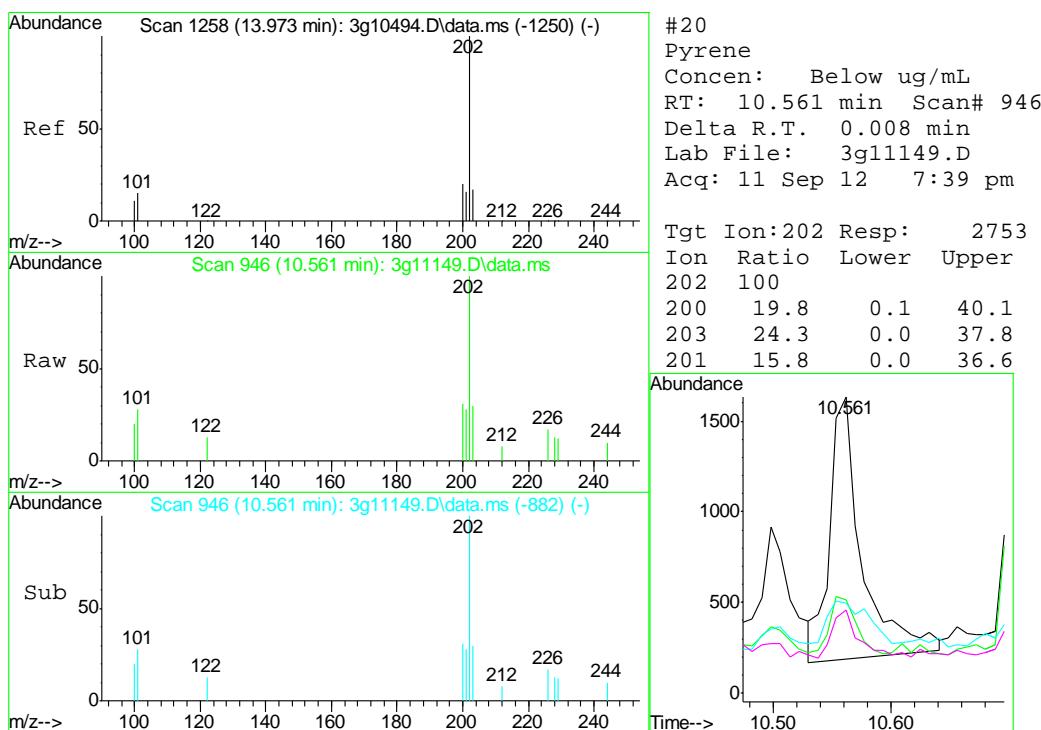


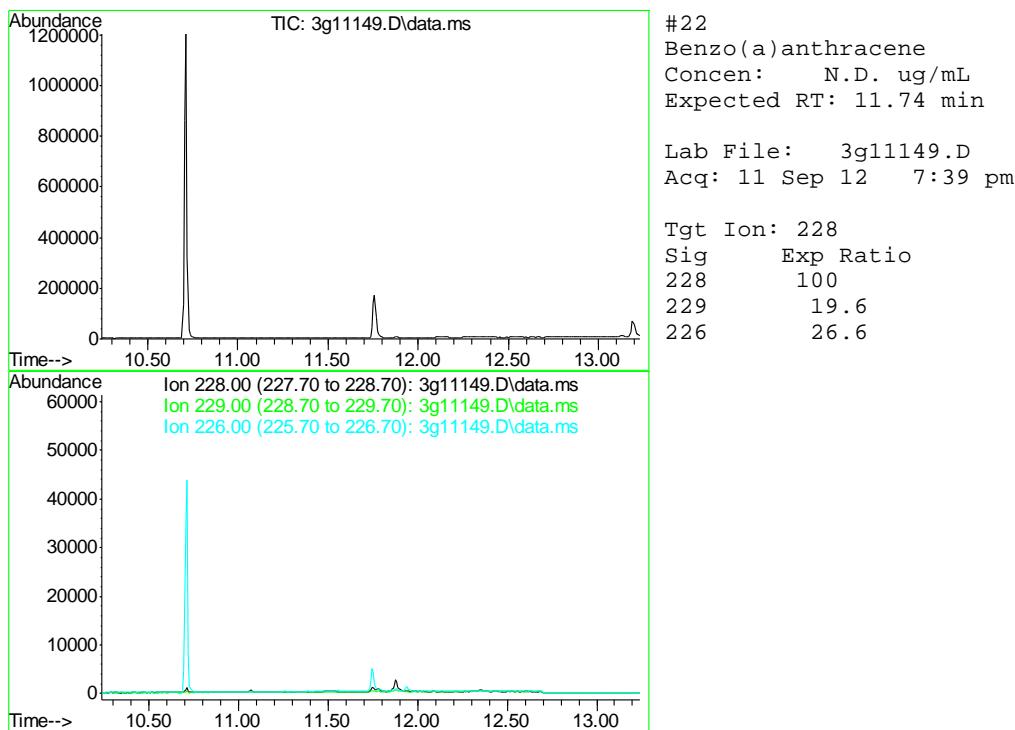
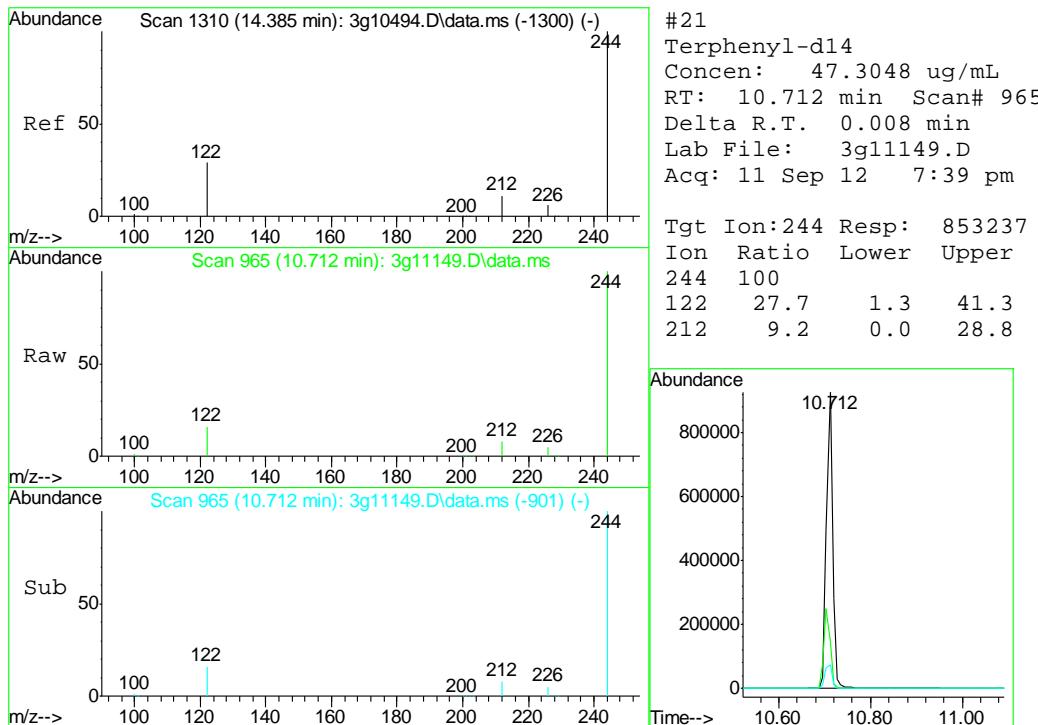


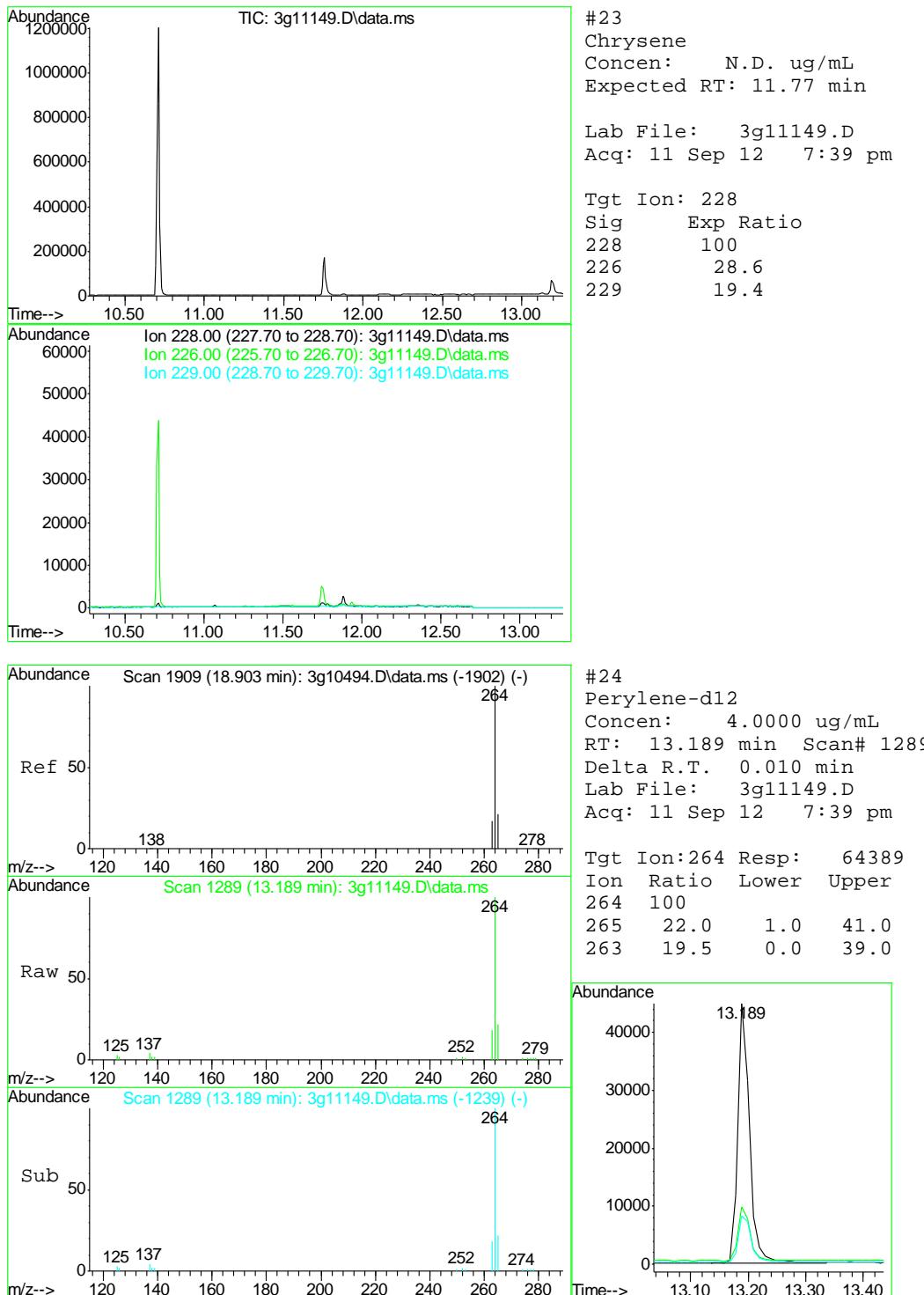


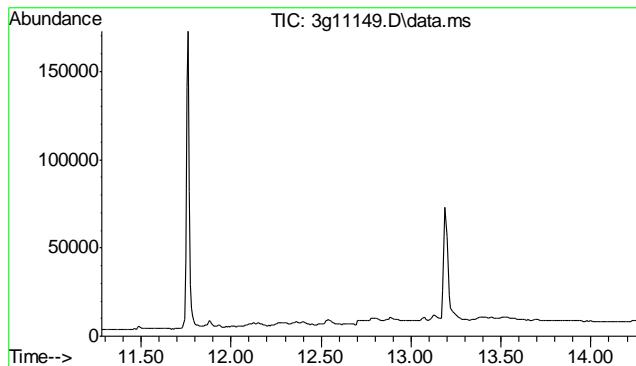
9.2.1

9





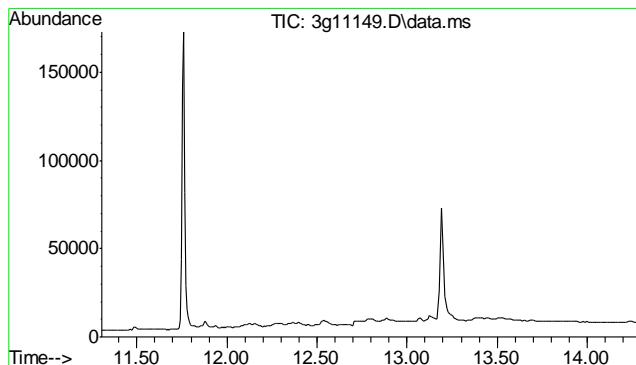
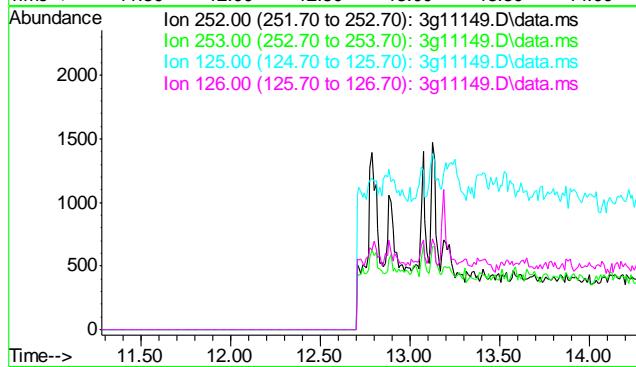




#25
Benzo(b)fluoranthene
Concen: N.D. ug/mL
Expected RT: 12.78 min

Lab File: 3g11149.D
Acq: 11 Sep 12 7:39 pm

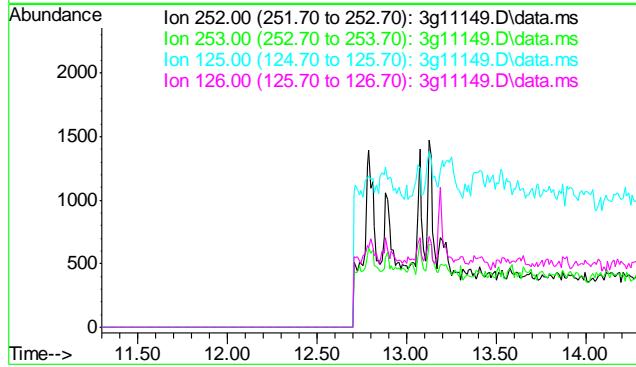
Tgt Ion: 252
Sig Exp Ratio
252 100
253 22.9
125 11.5
126 14.7

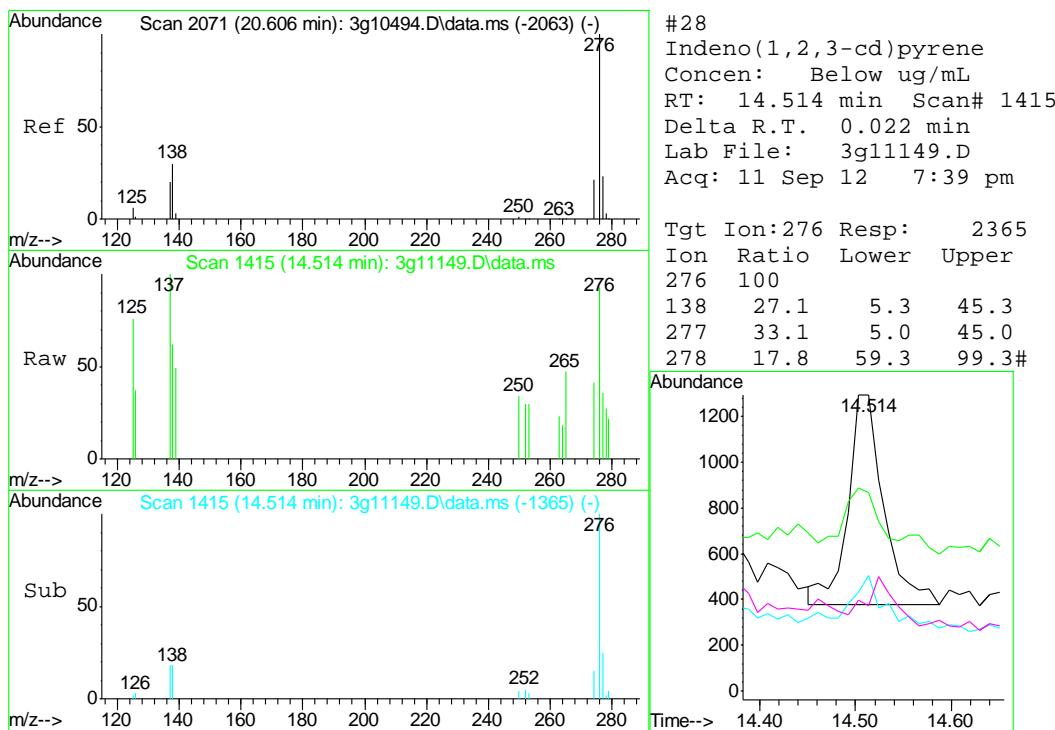
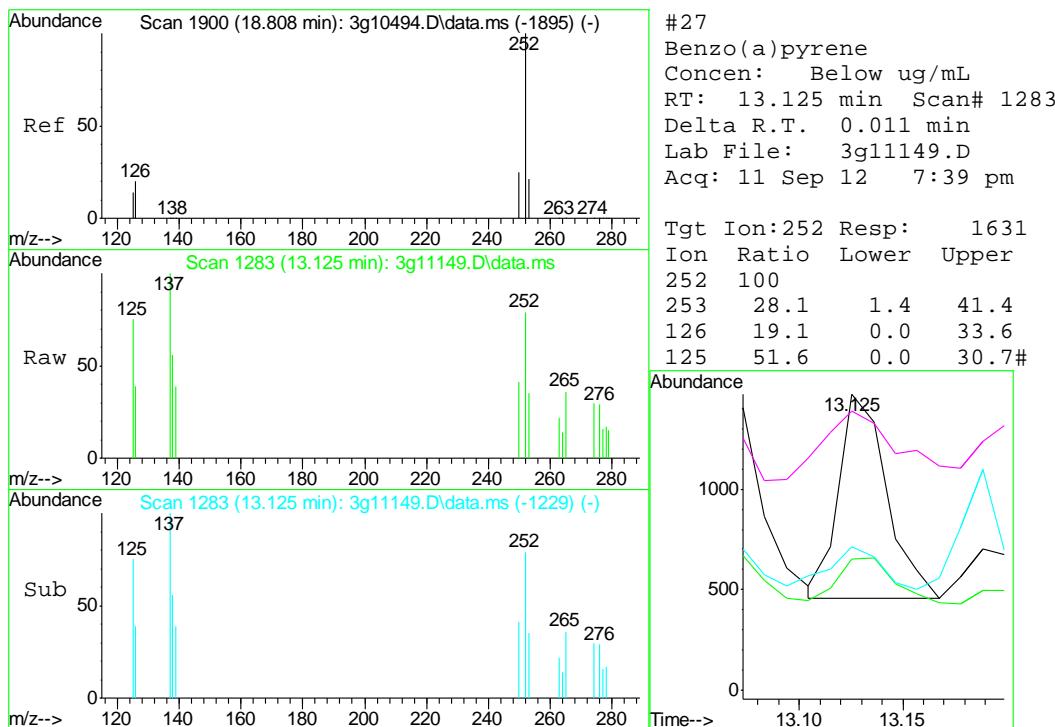


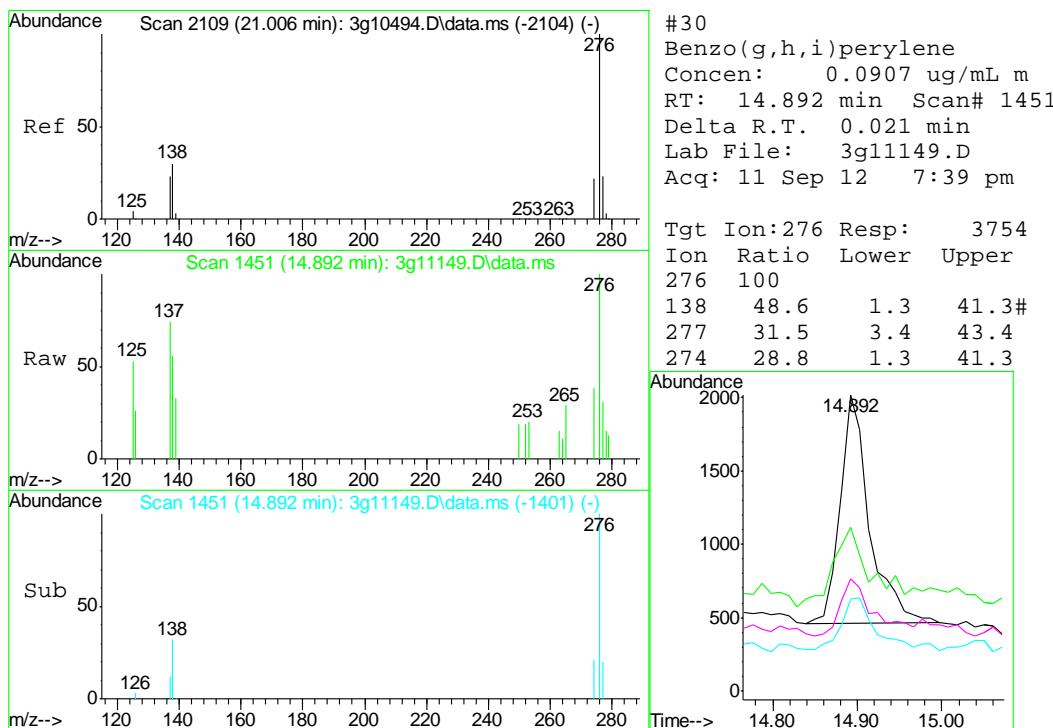
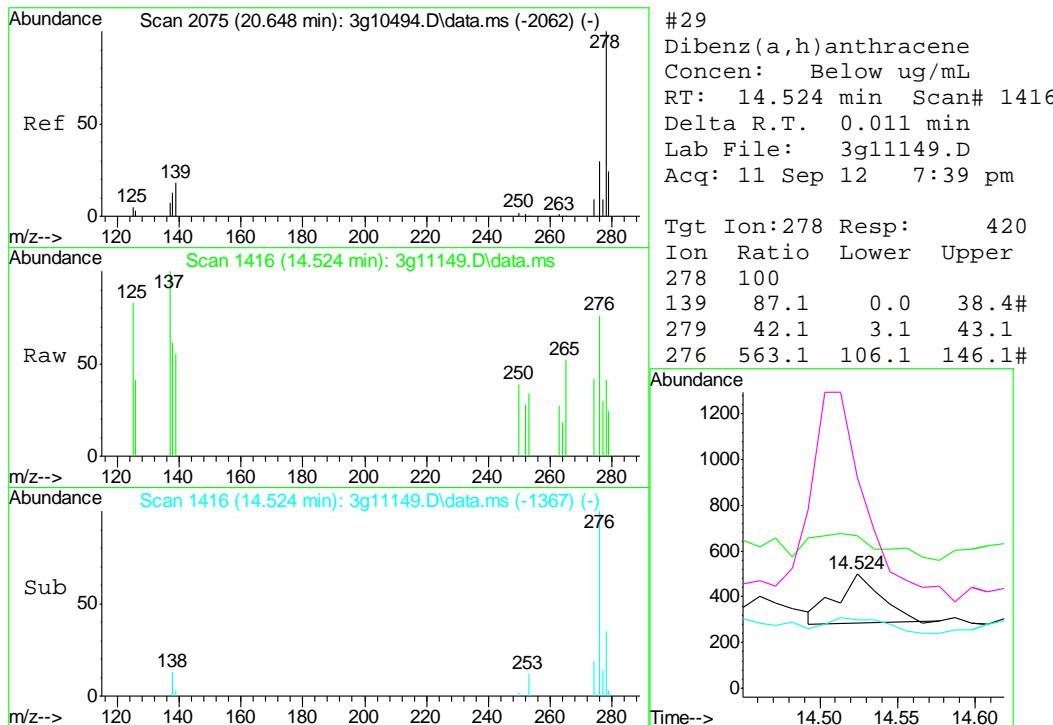
#26
Benzo(k)fluoranthene
Concen: N.D. ug/mL
Expected RT: 12.80 min

Lab File: 3g11149.D
Acq: 11 Sep 12 7:39 pm

Tgt Ion: 252
Sig Exp Ratio
252 100
253 21.8
125 11.0
126 14.0









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: D38480
Account: XTOKWR XTO Energy
Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB958-MB	GB17458.D	1	09/10/12	SK	n/a	n/a	GGB958

The QC reported here applies to the following samples:

Method: SW846 8015B

D38480-1, D38480-2, D38480-3

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	100% 60-140%

10.1.1

10

Blank Spike Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB958-BS	GB17459.D	1	09/10/12	SK	n/a	n/a	GGB958

The QC reported here applies to the following samples:

Method: SW846 8015B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	126	115	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	107%	60-140%

10.2.1

10

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D38454-1MS	GB17461.D	1	09/10/12	SK	n/a	n/a	GGB958
D38454-1MSD	GB17462.D	1	09/10/12	SK	n/a	n/a	GGB958
D38454-1	GB17460.D	1	09/10/12	SK	n/a	n/a	GGB958

The QC reported here applies to the following samples:

Method: SW846 8015B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	D38454-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		168	194	115	197	117	2	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D38454-1	Limits
120-82-1	1,2,4-Trichlorobenzene	106%	106%	94%	60-140%

* = Outside of Control Limits.

10.3.1
10



GC Volatiles

Raw Data

Judy Nelson
 09/11/12 08:56

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17471.D\FID1A.CH Vial: 17
 Signal #2 : Y:\1\DATA\091012\GB17471.D\FID2B.CH
 Acq On : 10 Sep 2012 7:14 pm Operator: StephK
 Sample : D38480-1, 50X Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.064,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 11 08:17:16 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Mon Sep 10 11:02:56 2012
 Response via : Initial Calibration
 DataAcq Meth : TVB4.M

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

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

System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.36	2761265	88.124 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	19380807	119.246 %	

Target Compounds

1) H	TVH-Gasoline	7.23	58975870	0.905	mg/L
4) T	Methyl-t-butyl-ether	2.20	203058	1.564	ug/L
5) T	Benzene	4.10	769084	1.908	ug/L
6) T	Toluene	7.64	14054678	35.467	ug/L
7) T	Ethylbenzene	10.28	3138365	9.278	ug/L
8) T	m,p-Xylene	10.46	13337095	36.167	ug/L
9) T	o-Xylene	10.96	1085060	3.304	ug/L
11) T	Naphthalene	14.58	7766841	39.364	ug/L

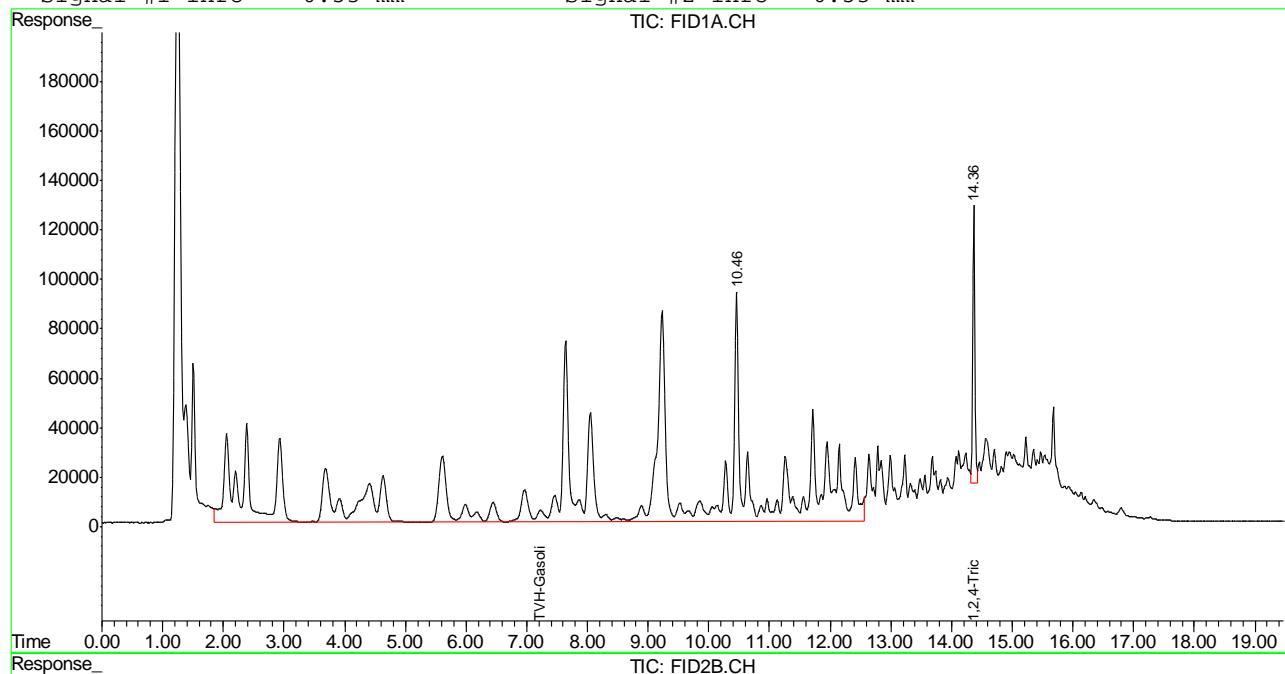
(f)=RT Delta > 1/2 Window (m)=manual int.
 GB17471.D TB868GB868SOIL.M Tue Sep 11 08:26:13 2012 GC

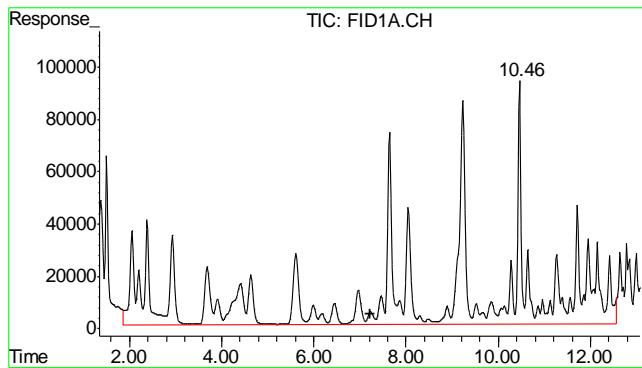
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17471.D\FID1A.CH Vial: 17
 Signal #2 : Y:\1\DATA\091012\GB17471.D\FID2B.CH
 Acq On : 10 Sep 2012 7:14 pm Operator: StephK
 Sample : D38480-1, 50X Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.064,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 11 7:28 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Mon Sep 10 11:02:56 2012
 Response via : Multiple Level Calibration
 DataAcq Meth : TVB4.M

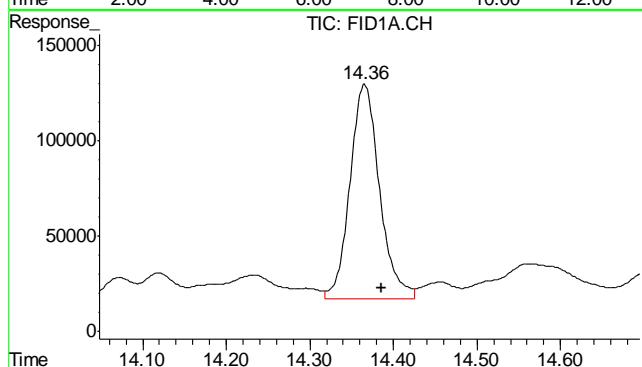
Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





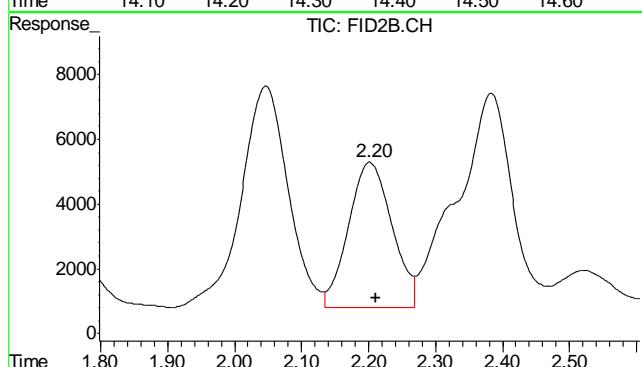
#1 TVH-Gasoline

R.T.: 7.230 min
Delta R.T.: 0.000 min
Response: 58975870
Conc: 0.90 mg/L m



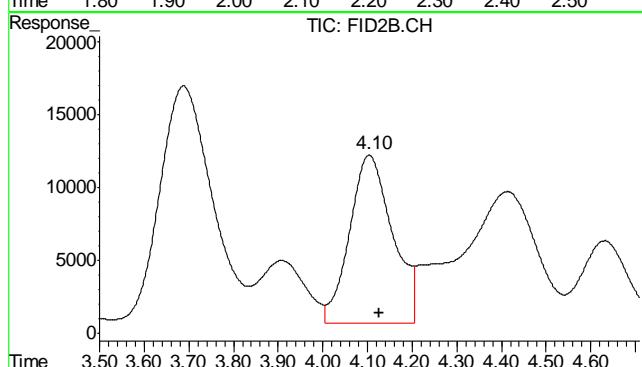
#2 1,2,4-Trichlorobenzene

R.T.: 14.365 min
Delta R.T.: -0.021 min
Response: 2761265
Conc: 88.12 % m



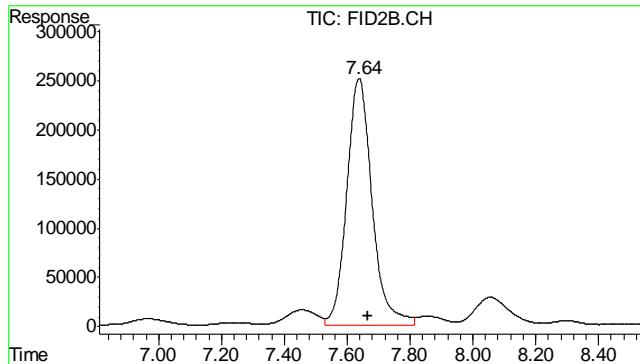
#4 Methyl-t-butyl-ether

R.T.: 2.202 min
Delta R.T.: -0.009 min
Response: 203058
Conc: 1.56 ug/L



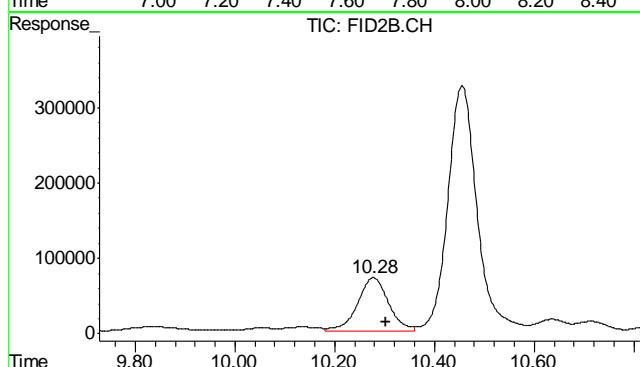
#5 Benzene

R.T.: 4.104 min
Delta R.T.: -0.023 min
Response: 769084
Conc: 1.91 ug/L



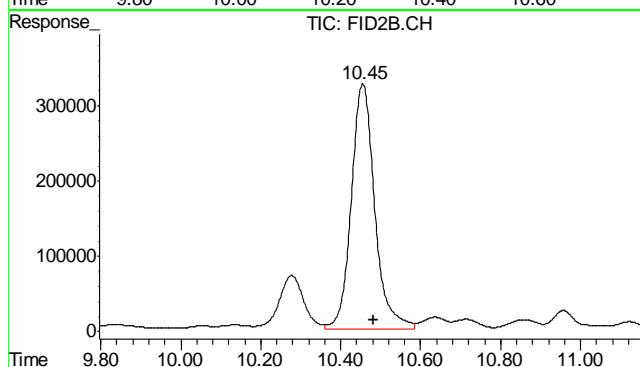
#6 Toluene

R.T.: 7.638 min
 Delta R.T.: -0.029 min
 Response: 14054678
 Conc: 35.47 ug/L



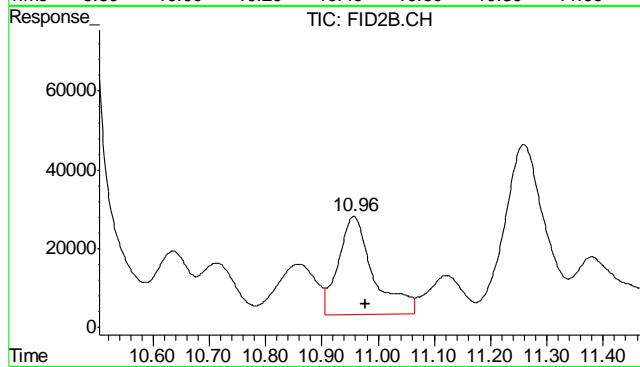
#7 Ethylbenzene

R.T.: 10.278 min
 Delta R.T.: -0.024 min
 Response: 3138365
 Conc: 9.28 ug/L



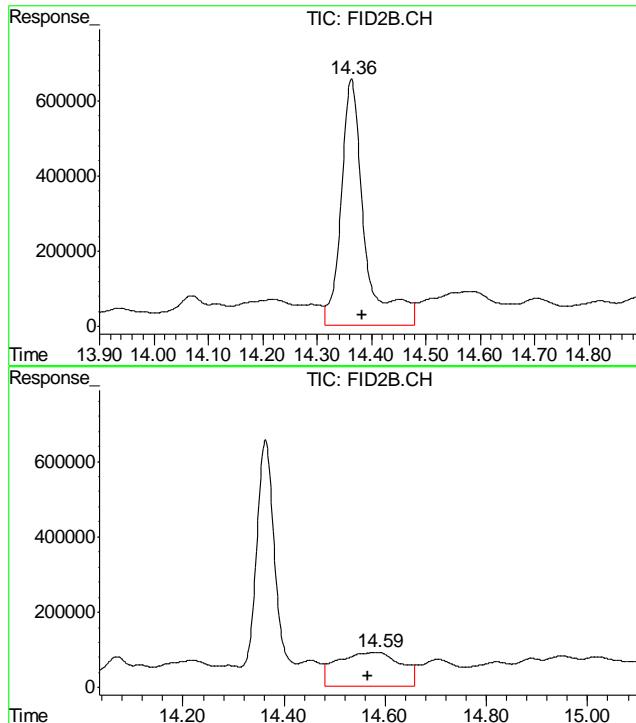
#8 m,p-Xylene

R.T.: 10.455 min
 Delta R.T.: -0.026 min
 Response: 13337095
 Conc: 36.17 ug/L



#9 o-Xylene

R.T.: 10.957 min
 Delta R.T.: -0.021 min
 Response: 1085060
 Conc: 3.30 ug/L



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

R.T.: 14.363 min
Delta R.T.: -0.020 min
Response: 19380807
Conc: 119.25 %

#11 Naphthalene

R.T.: 14.580 min
Delta R.T.: 0.014 min
Response: 7766841
Conc: 39.36 ug/L

11.1.1

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson
09/11/12 08:56

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17472.D\FID1A.CH Vial: 18
 Signal #2 : Y:\1\DATA\091012\GB17472.D\FID2B.CH
 Acq On : 10 Sep 2012 7:49 pm Operator: StephK
 Sample : D38480-2, 50X Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.062,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 11 08:17:20 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Mon Sep 10 11:02:56 2012
 Response via : Initial Calibration
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.36	2918282	93.135 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	16620672	102.264 %	

Target Compounds

1) H	TVH-Gasoline	7.23	23151874	0.335 mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	4.09	3250042	8.065 ug/L
6) T	Toluene	7.63	6638462	16.752 ug/L
7) T	Ethylbenzene	10.27	1068341	3.158 ug/L
8) T	m,p-Xylene	10.45	5433293	14.512 ug/L
9) T	o-Xylene	10.95	1023567	3.117 ug/L
11) T	Naphthalene	14.55	6179507	31.319 ug/L

11.1.2
11

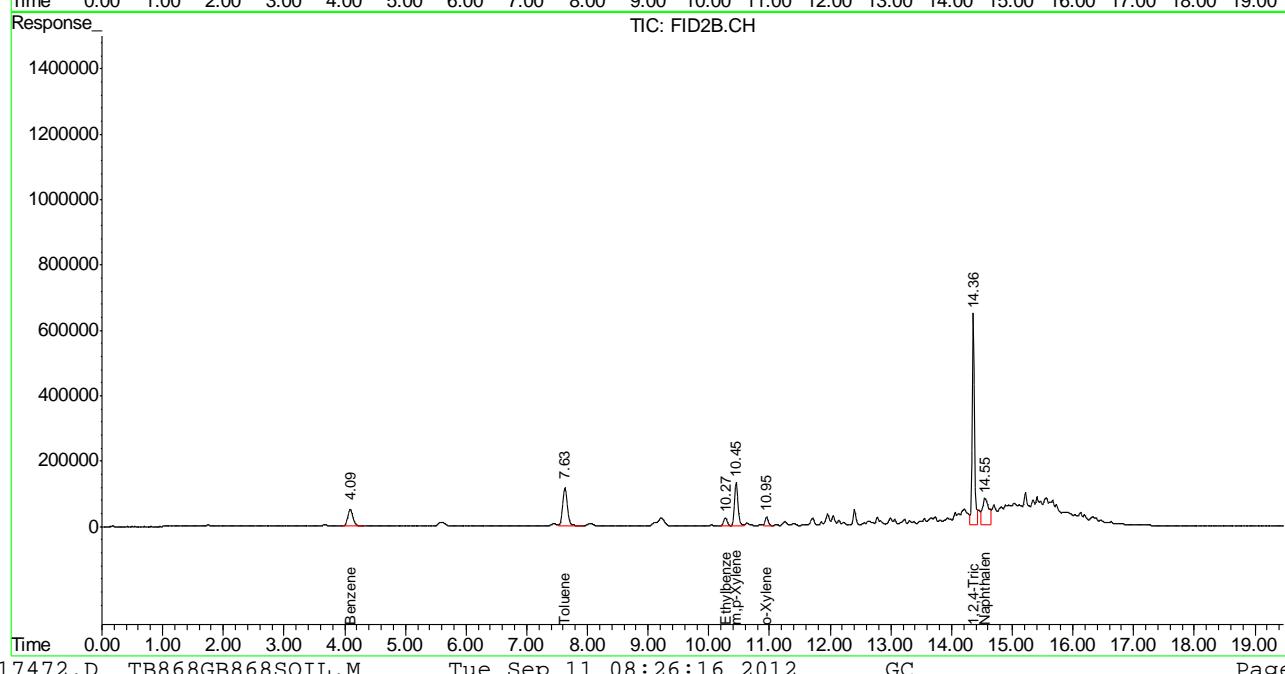
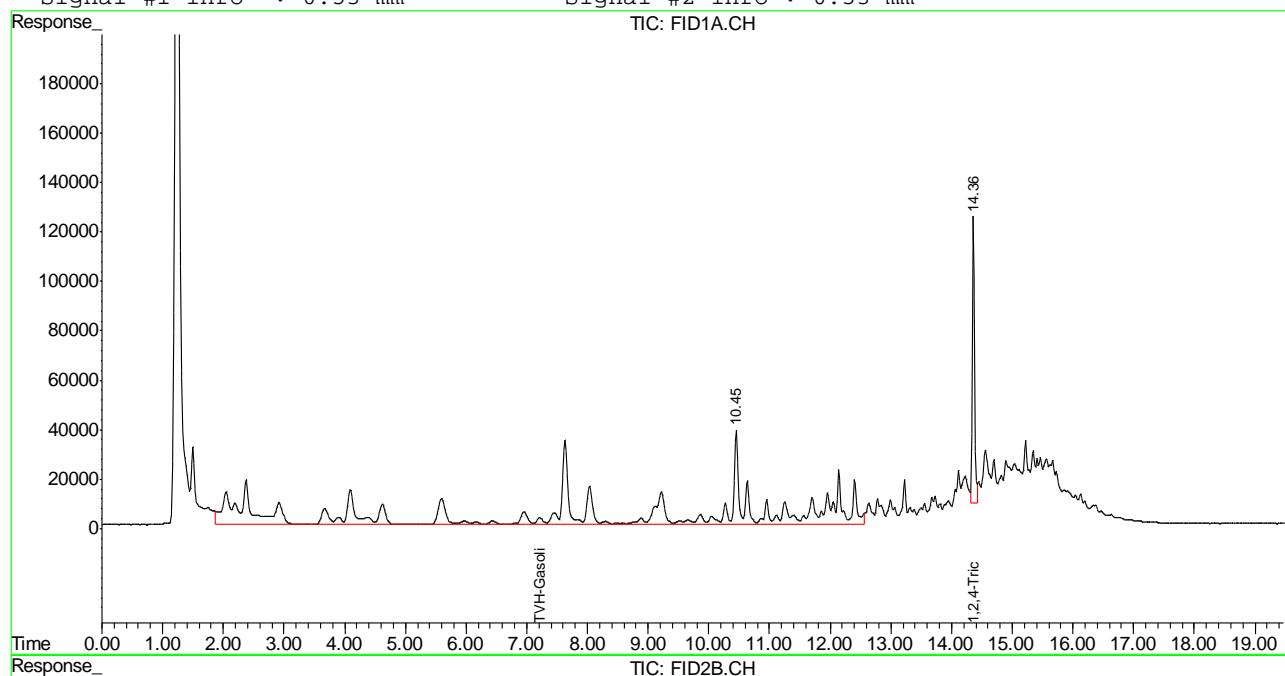
(f)=RT Delta > 1/2 Window (m)=manual int.
 GB17472.D TB868GB868SOIL.M Tue Sep 11 08:26:16 2012 GC

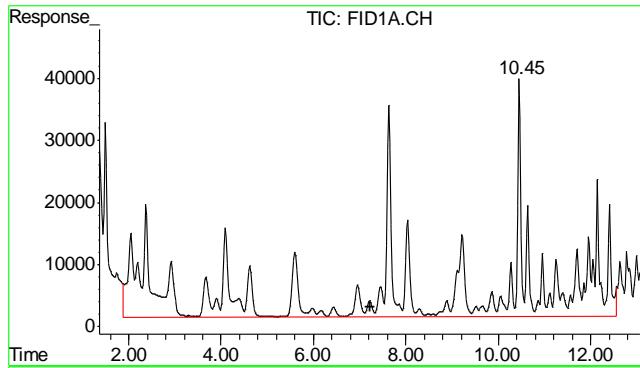
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17472.D\FID1A.CH Vial: 18
 Signal #2 : Y:\1\DATA\091012\GB17472.D\FID2B.CH
 Acq On : 10 Sep 2012 7:49 pm Operator: StephK
 Sample : D38480-2, 50X Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.062,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 11 7:28 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Mon Sep 10 11:02:56 2012
 Response via : Multiple Level Calibration
 DataAcq Meth : TVB4.M

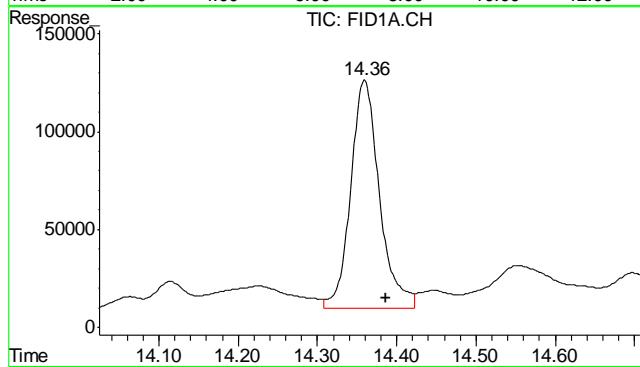
Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





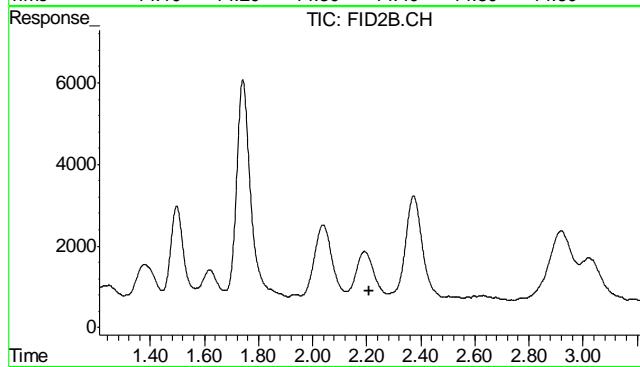
#1 TVH-Gasoline

R.T.: 7.230 min
Delta R.T.: 0.000 min
Response: 23151874
Conc: 0.33 mg/L m



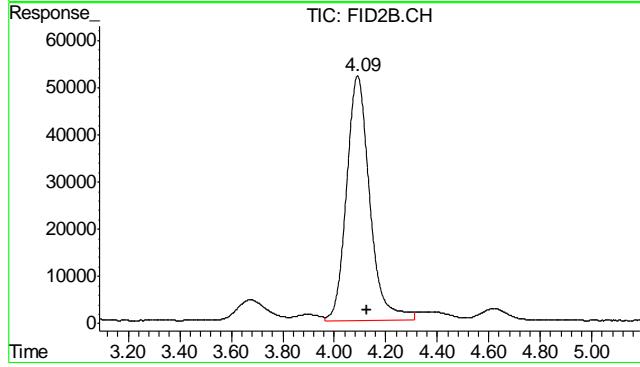
#2 1,2,4-Trichlorobenzene

R.T.: 14.359 min
Delta R.T.: -0.027 min
Response: 2918282
Conc: 93.13 % m



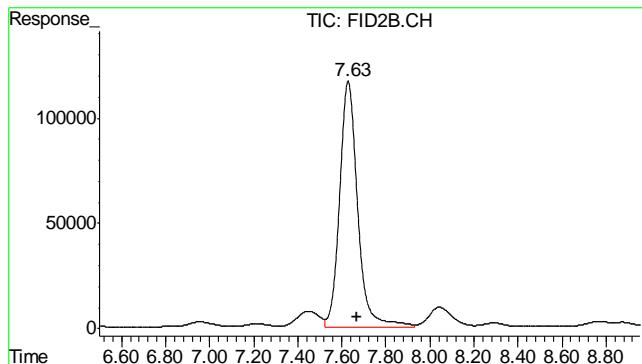
#4 Methyl-t-butyl-ether

R.T.: 0.000 min
Exp R.T.: 2.211 min
Response: 0
Conc: N.D.



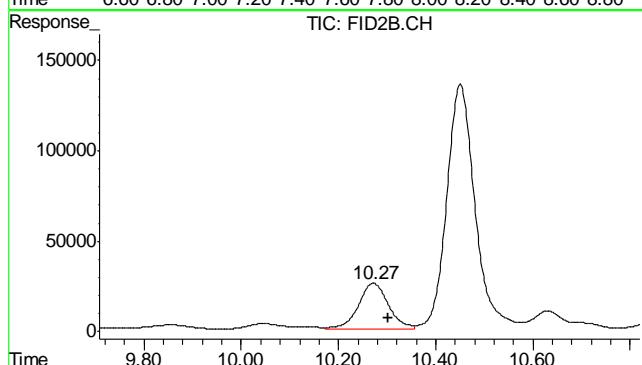
#5 Benzene

R.T.: 4.091 min
Delta R.T.: -0.036 min
Response: 3250042
Conc: 8.06 ug/L



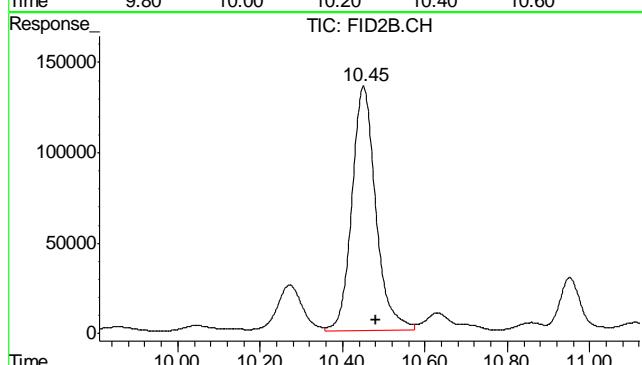
#6 Toluene

R.T.: 7.629 min
Delta R.T.: -0.037 min
Response: 6638462
Conc: 16.75 ug/L



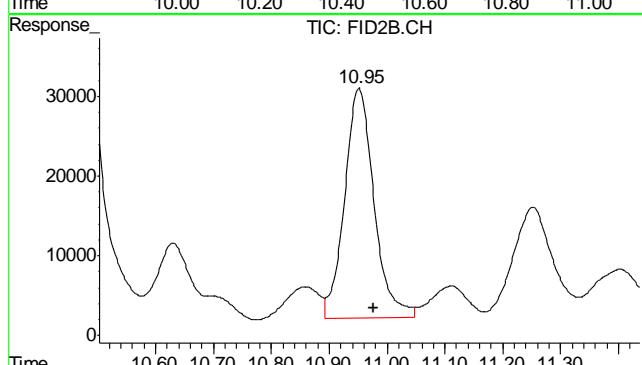
#7 Ethylbenzene

R.T.: 10.272 min
Delta R.T.: -0.030 min
Response: 1068341
Conc: 3.16 ug/L



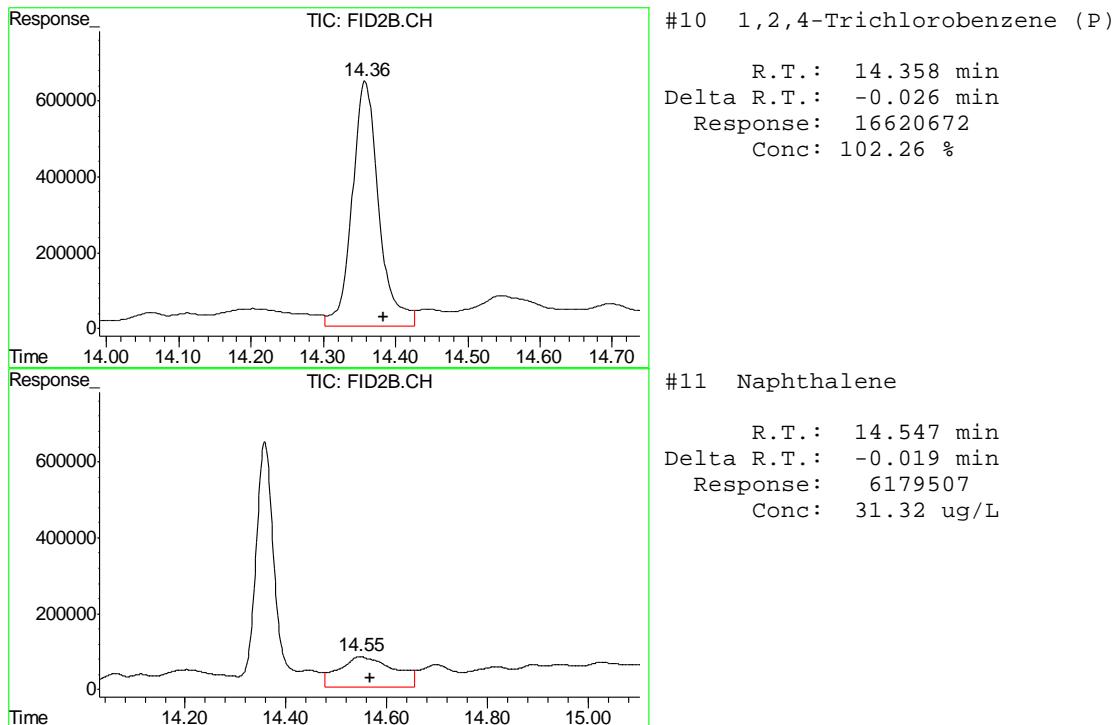
#8 m,p-Xylene

R.T.: 10.451 min
Delta R.T.: -0.031 min
Response: 5433293
Conc: 14.51 ug/L



#9 o-Xylene

R.T.: 10.952 min
Delta R.T.: -0.026 min
Response: 1023567
Conc: 3.12 ug/L



11.1.2

11

Judy Nelson
 09/11/12 08:56

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17474.D\FID1A.CH Vial: 20
 Signal #2 : Y:\1\DATA\091012\GB17474.D\FID2B.CH
 Acq On : 10 Sep 2012 8:59 pm Operator: StephK
 Sample : D38480-3, 50X Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.078,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 11 08:18:01 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Tue Sep 11 08:17:43 2012
 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	2772102	88.469 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.37	19645576	120.875 %	

Target Compounds

1) H	TVH-Gasoline	7.23	41048209	0.620 mg/L
4) T	Methyl-t-butyl-ether	2.21	90659	0.698 ug/L
5) T	Benzene	4.11	7018387	17.416 ug/L
6) T	Toluene	7.64	16859988	42.546 ug/L
7) T	Ethylbenzene	10.28	1077039	3.184 ug/L
8) T	m,p-Xylene	10.46	15160284	41.162 ug/L
9) T	o-Xylene	10.96	1645166	5.010 ug/L
11) T	Naphthalene	14.57	13438626	68.110 ug/L

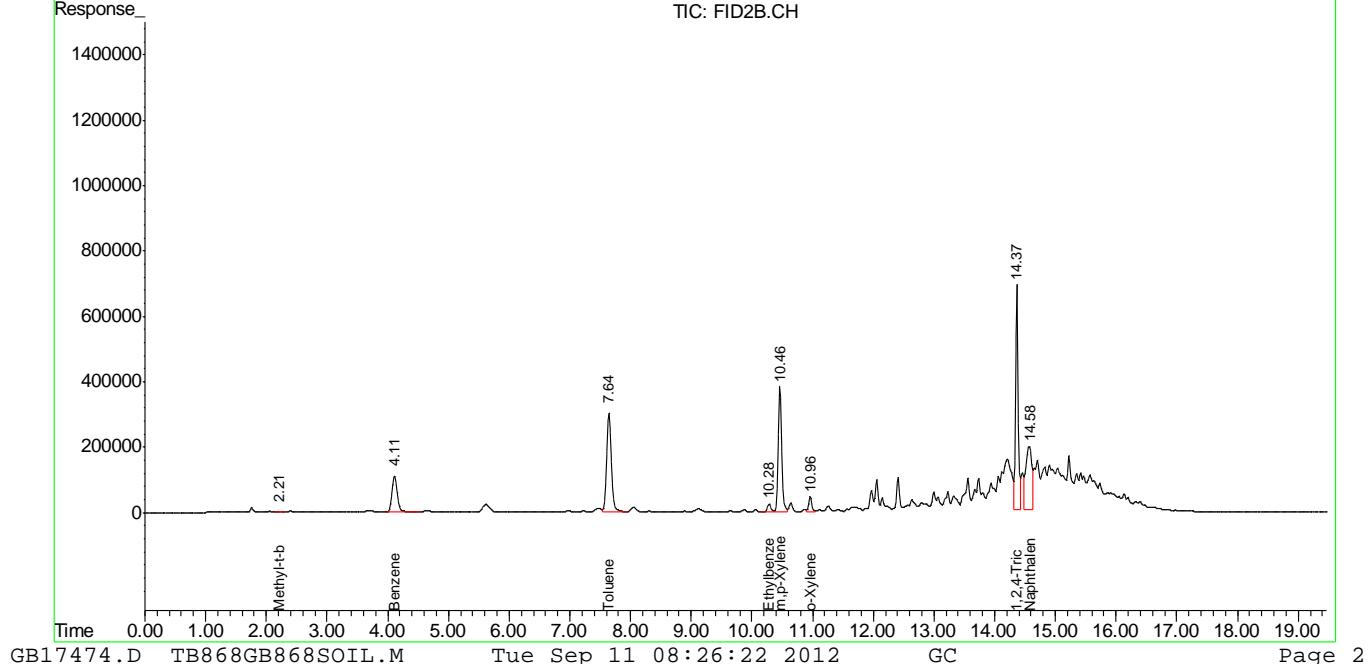
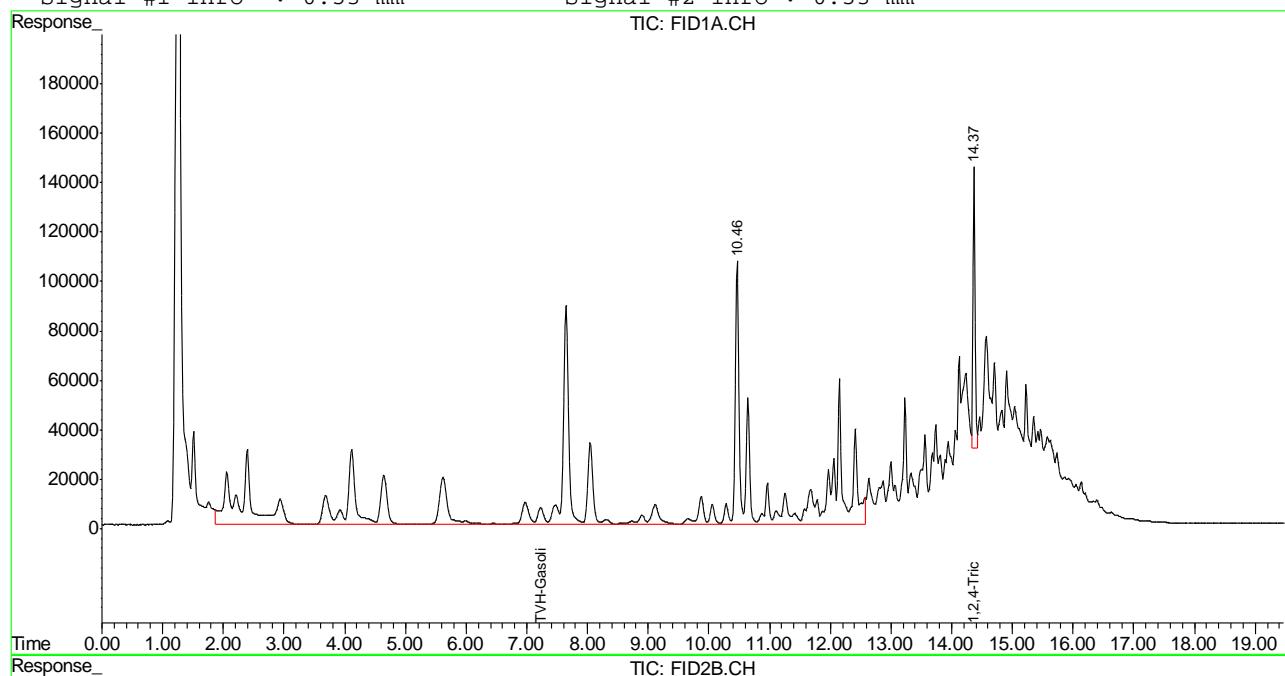
(f)=RT Delta > 1/2 Window (m)=manual int.
 GB17474.D TB868GB868SOIL.M Tue Sep 11 08:26:22 2012 GC

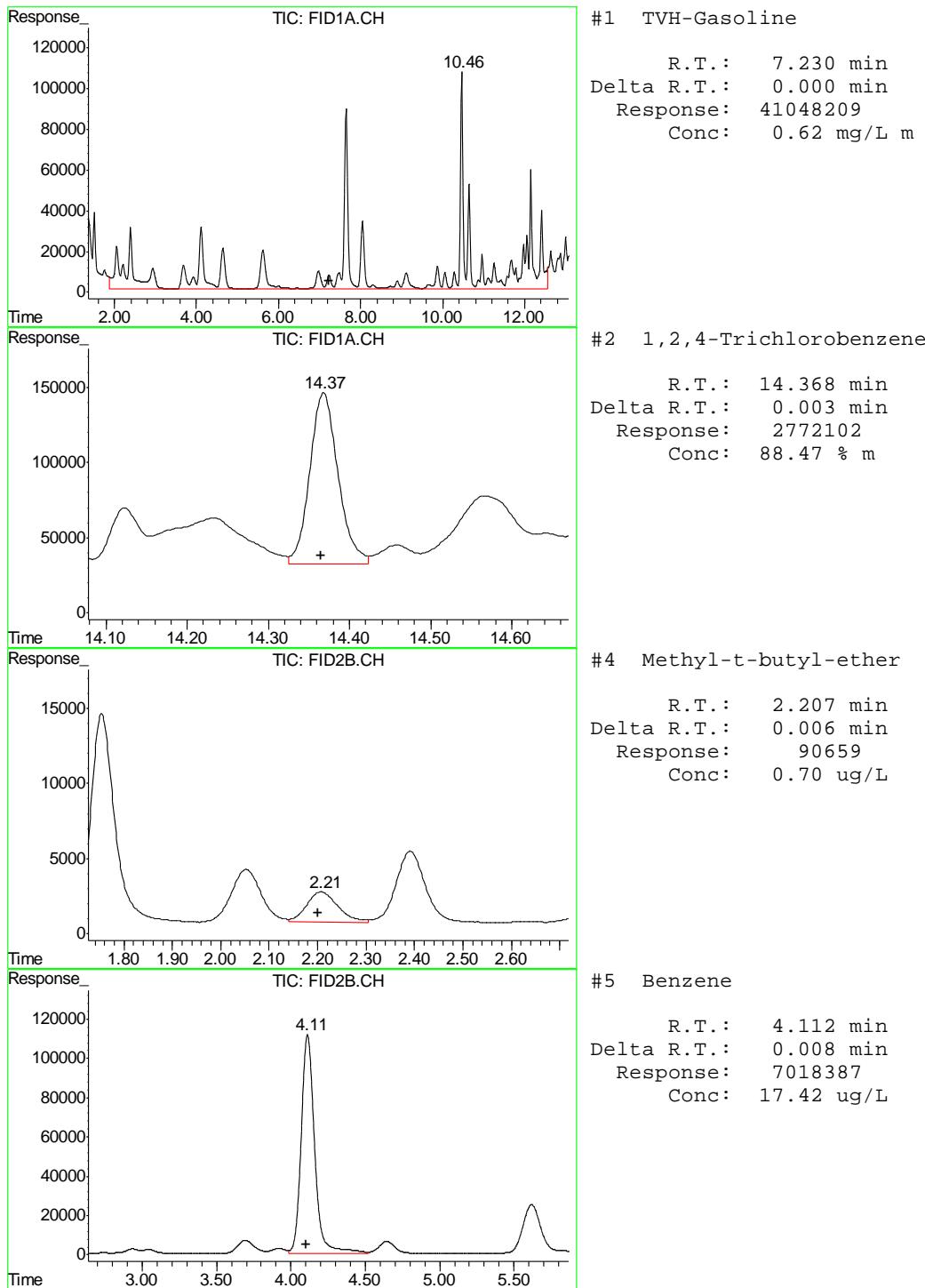
Quantitation Report (QT Reviewed)

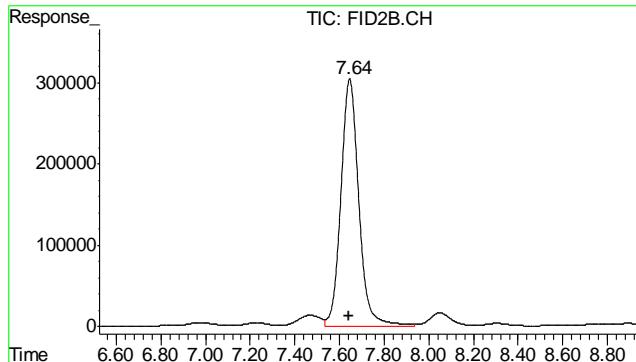
Signal #1 : Y:\1\DATA\091012\GB17474.D\FID1A.CH Vial: 20
 Signal #2 : Y:\1\DATA\091012\GB17474.D\FID2B.CH
 Acq On : 10 Sep 2012 8:59 pm Operator: StephK
 Sample : D38480-3, 50X Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.078,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 11 7:29 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Tue Sep 11 08:17:43 2012
 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

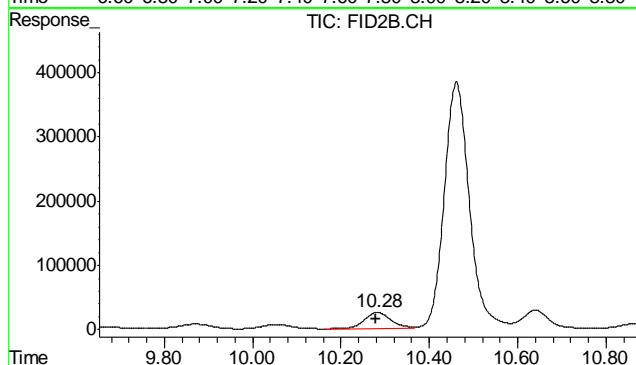






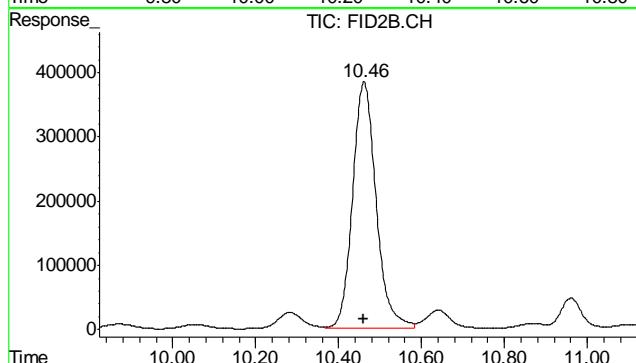
#6 Toluene

R.T.: 7.645 min
Delta R.T.: 0.003 min
Response: 16859988
Conc: 42.55 ug/L



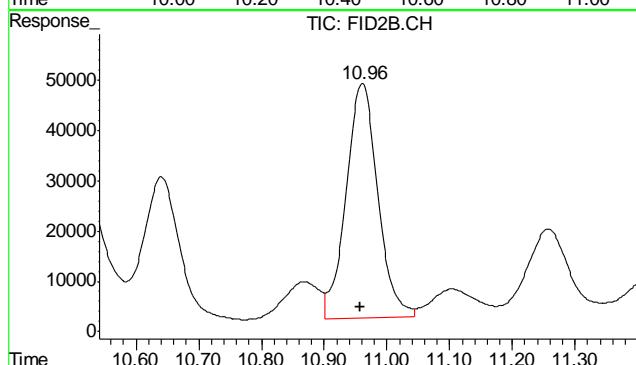
#7 Ethylbenzene

R.T.: 10.282 min
Delta R.T.: 0.004 min
Response: 1077039
Conc: 3.18 ug/L



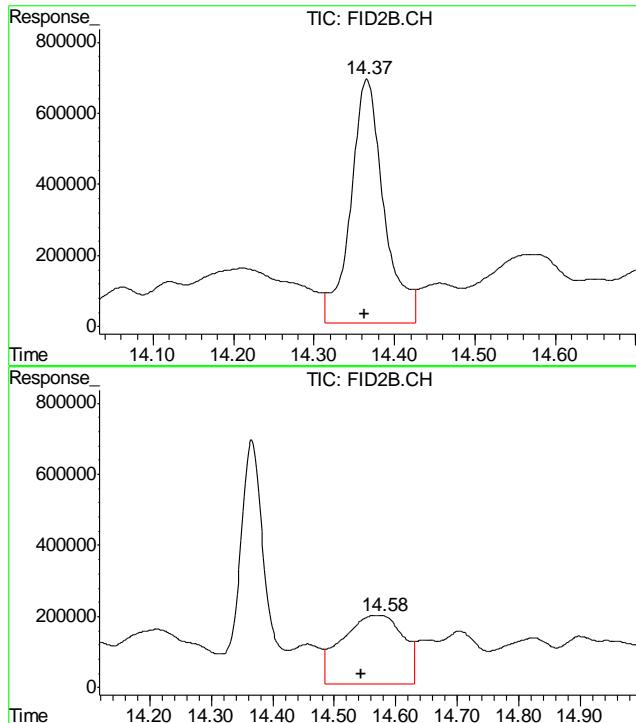
#8 m,p-Xylene

R.T.: 10.461 min
Delta R.T.: 0.002 min
Response: 15160284
Conc: 41.16 ug/L



#9 o-Xylene

R.T.: 10.961 min
Delta R.T.: 0.003 min
Response: 1645166
Conc: 5.01 ug/L



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

R.T.: 14.366 min
 Delta R.T.: 0.003 min
 Response: 19645576
 Conc: 120.88 %

#11 Naphthalene

R.T.: 14.572 min
 Delta R.T.: 0.028 min
 Response: 13438626
 Conc: 68.11 ug/L

11.1.3

11

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17458.D\FID1A.CH Vial: 4
 Signal #2 : Y:\1\DATA\091012\GB17458.D\FID2B.CH
 Acq On : 10 Sep 2012 11:26 am Operator: StephK
 Sample : MB Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 10 12:29:53 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Mon Sep 10 11:02:56 2012
 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	3120564	99.590	%
10) S 1,2,4-Trichlorobenzene (P)	14.37	16470036	101.337	%

Target Compounds

1) H TVH-Gasoline	7.23	3428685	<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	125866	0.318	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	194561	0.986	ug/L

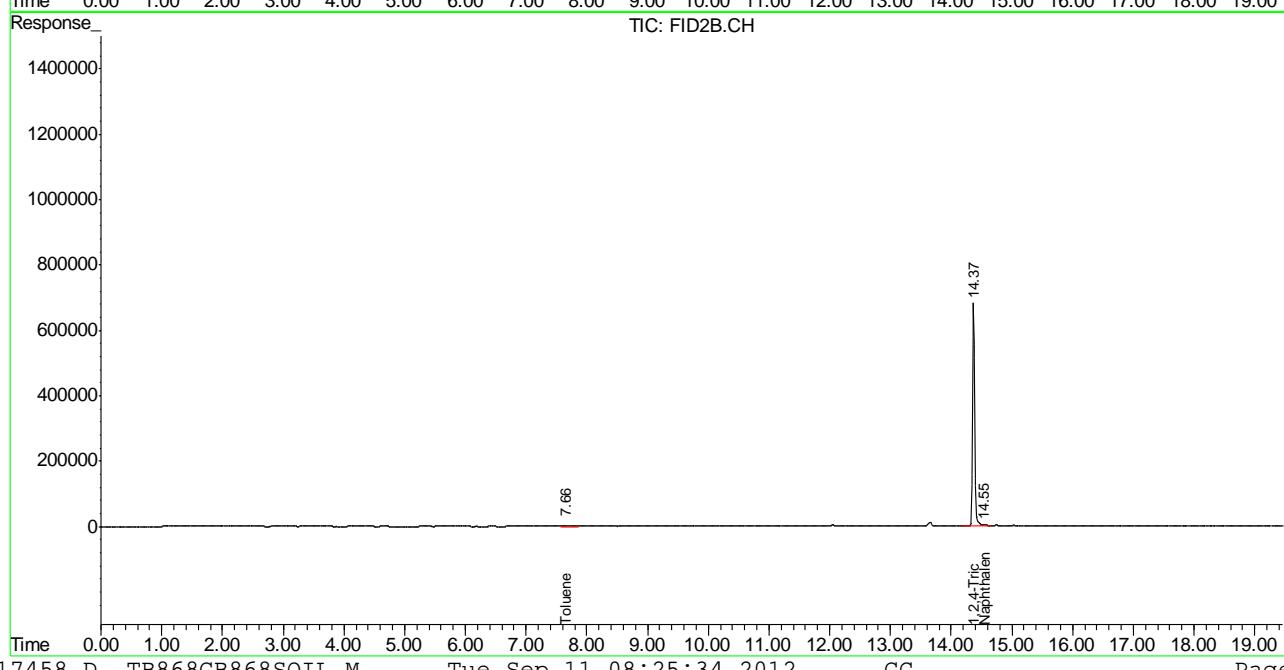
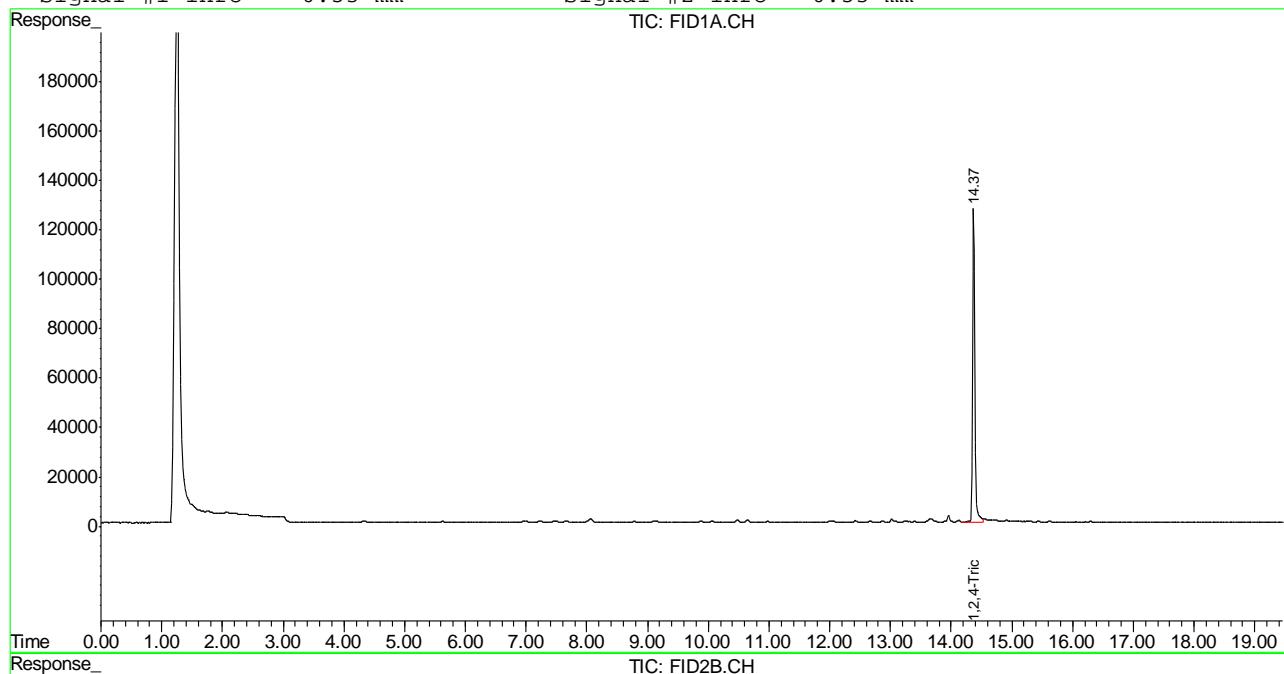
(f)=RT Delta > 1/2 Window (m)=manual int.
 GB17458.D TB868GB868SOIL.M Tue Sep 11 08:25:34 2012 GC

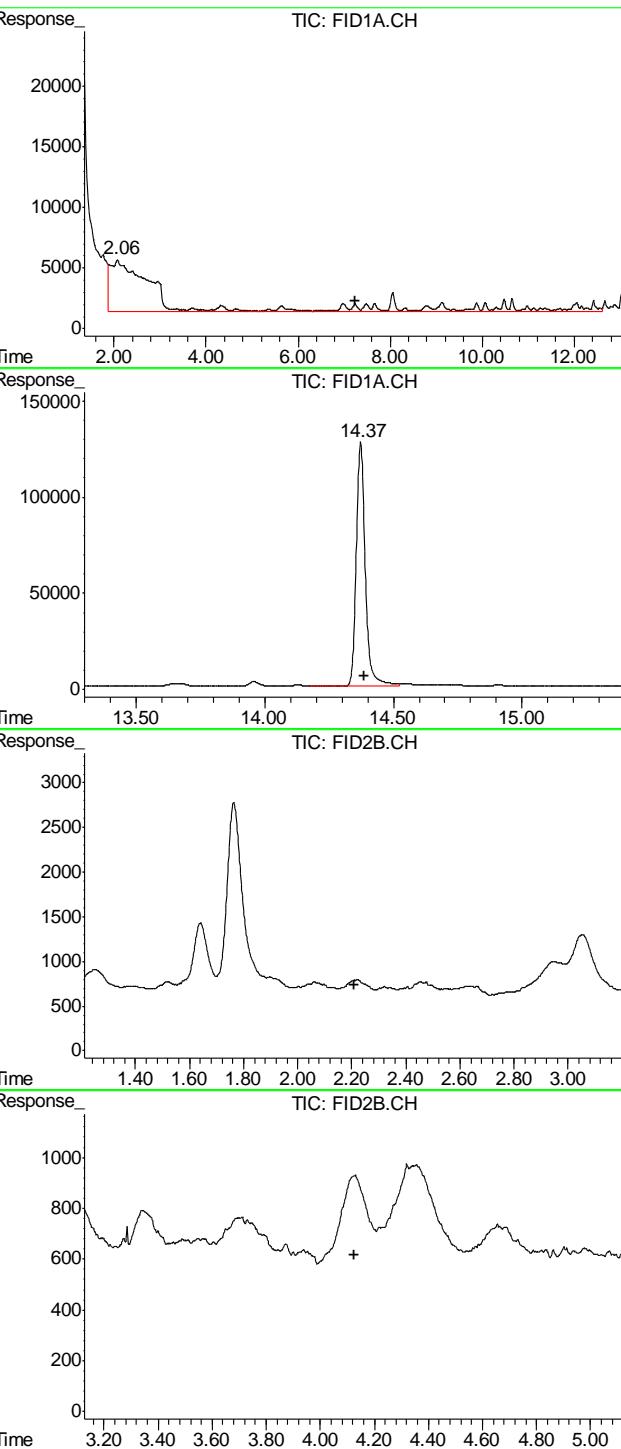
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\091012\GB17458.D\FID1A.CH Vial: 4
 Signal #2 : Y:\1\DATA\091012\GB17458.D\FID2B.CH
 Acq On : 10 Sep 2012 11:26 am Operator: StephK
 Sample : MB Inst : GC/MS Ins
 Misc : GC3092,GGB958,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Sep 10 11:44 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Mon Sep 10 11:02:56 2012
 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: 3428685
 Conc: N.D.

#2 1,2,4-Trichlorobenzene

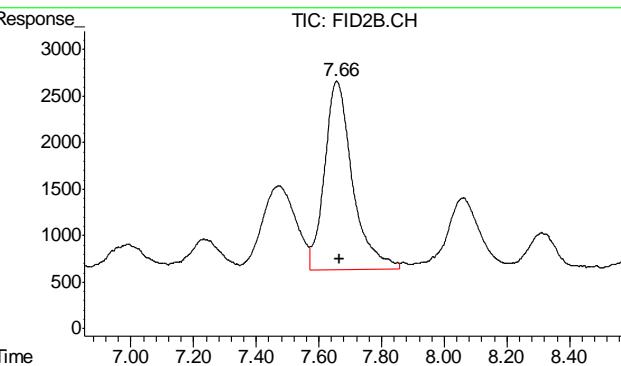
R.T.: 14.372 min
 Delta R.T.: -0.014 min
 Response: 3120564
 Conc: 99.59 %

#4 Methyl-t-butyl-ether

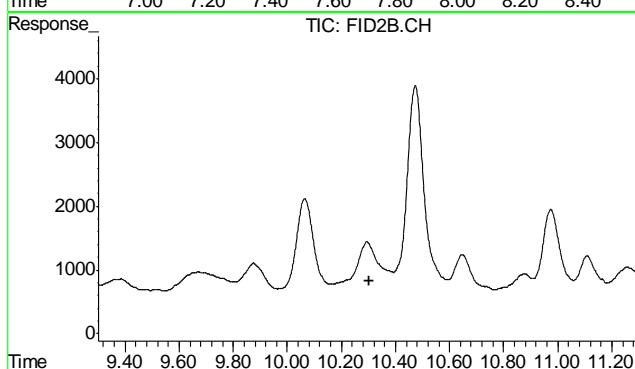
R.T.: 0.000 min
 Exp R.T. : 2.211 min
 Response: 0
 Conc: N.D.

#5 Benzene

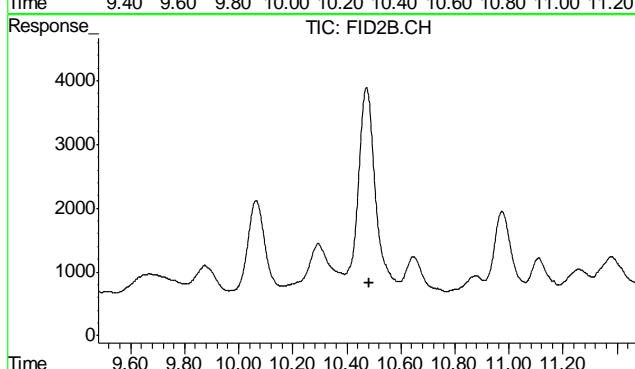
R.T.: 0.000 min
 Exp R.T. : 4.127 min
 Response: 0
 Conc: N.D.



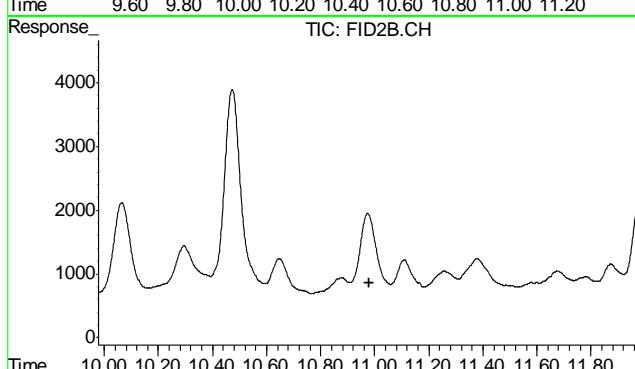
#6 Toluene
R.T.: 7.657 min
Delta R.T.: -0.009 min
Response: 125866
Conc: 0.32 ug/L



#7 Ethylbenzene
R.T.: 0.000 min
Exp R.T. : 10.302 min
Response: 0
Conc: N.D.



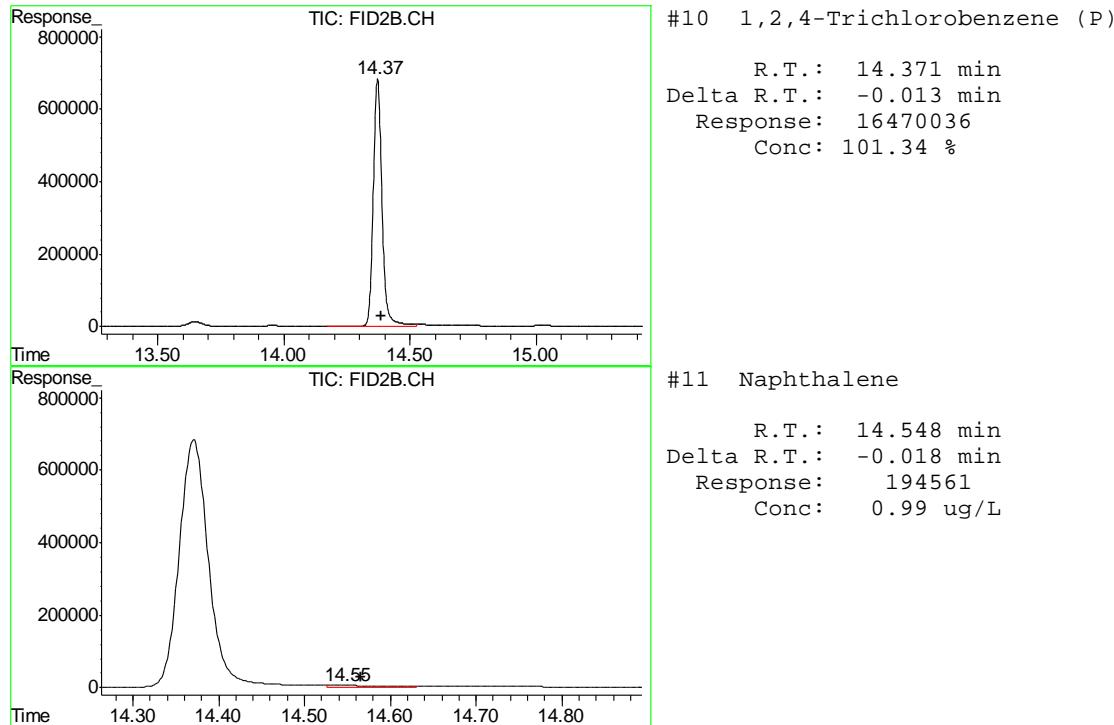
#8 m,p-Xylene
R.T.: 0.000 min
Exp R.T. : 10.482 min
Response: 0
Conc: N.D.



#9 o-Xylene
R.T.: 0.000 min
Exp R.T. : 10.978 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: D38480
Account: XTOKWR XTO Energy
Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6603-MB	FD17336.D	1	09/12/12	AW	09/11/12	OP6603	GFD891

The QC reported here applies to the following samples:

Method: SW846-8015B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	ND	13	8.7	mg/kg	

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	91% 43-136%

Blank Spike Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6603-BS	FD17338.D	1	09/12/12	AW	09/11/12	OP6603	GFD891

The QC reported here applies to the following samples:

Method: SW846-8015B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	555	83	58-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	86%	43-136%

* = Outside of Control Limits.

12.2.1

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Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D38480

Account: XTOKWR XTO Energy

Project: PCU 197-36A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6603-MS	FD17340.D	1	09/12/12	AW	09/11/12	OP6603	GFD891
OP6603-MSD	FD17342.D	1	09/12/12	AW	09/11/12	OP6603	GFD891
D38483-1	FD17344.D	1	09/12/12	AW	09/11/12	OP6603	GFD891

The QC reported here applies to the following samples:

Method: SW846-8015B

D38480-1, D38480-2, D38480-3

CAS No.	Compound	D38483-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	213		739	766	75	753	73	2	20-183/43
CAS No.	Surrogate Recoveries	MS		MSD		D38483-1		Limits		
84-15-1	o-Terphenyl	73%		67%		70%		43-136%		

* = Outside of Control Limits.

12.3.1
12



GC Semi-volatiles

Raw Data

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Nelson
09/13/12 15:50

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17346.D Vi
 Acq On : 9-12-2012 04:08:23 PM Operator: alexwl
 Sample : D38480-1 Inst : FID5
 Misc : OP6603,GFD891,30.10,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 09:06:27 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Initial Calibration
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	9.31	29405007	735.255 mg/L m
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	7.16	328605542	10065.119 mg/L

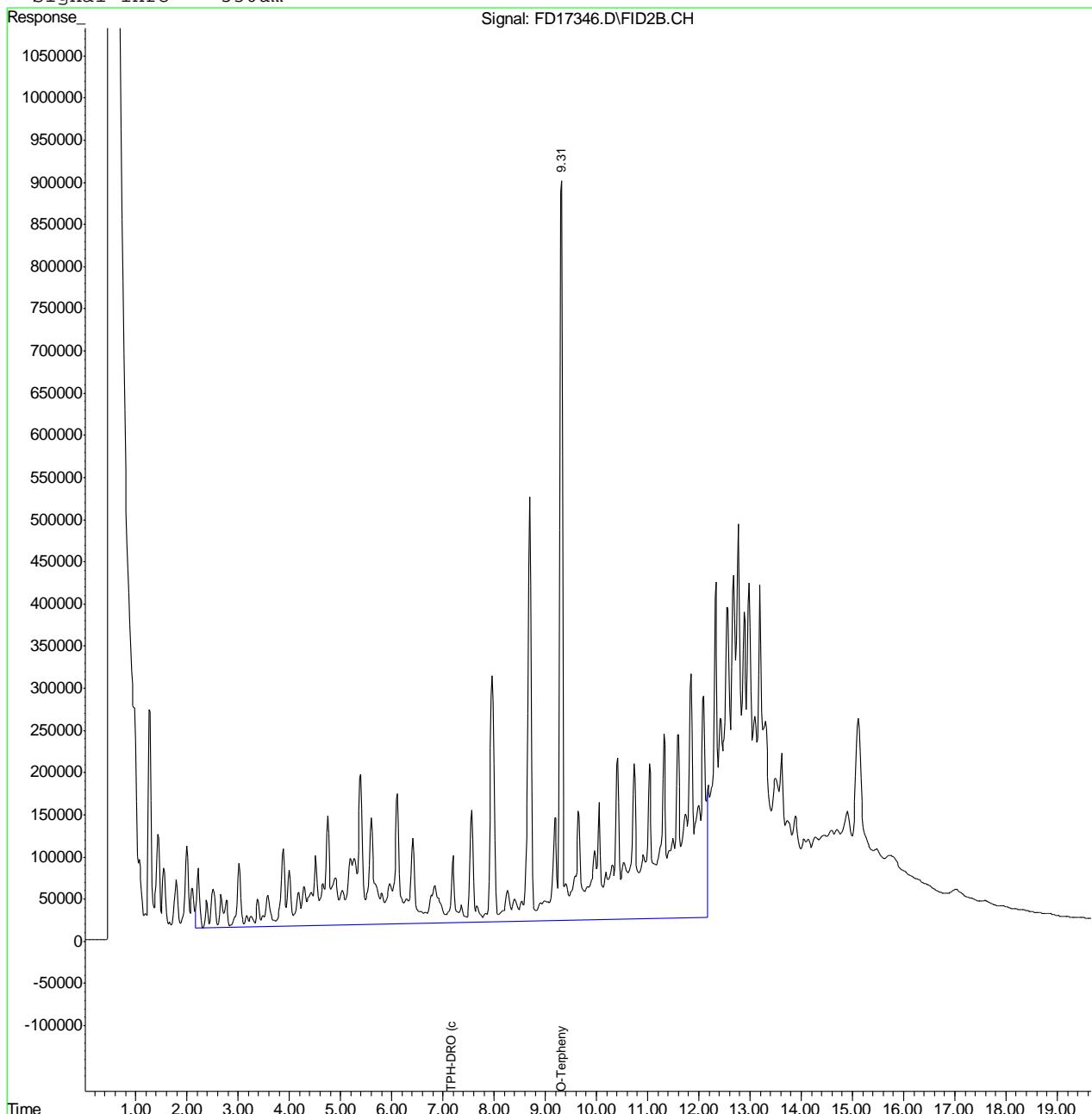
(f)=RT Delta > 1/2 Window (m)=manual int.
 FD17346.D DRO-GFD889R.M Thu Sep 13 09:41:25 2012 GC

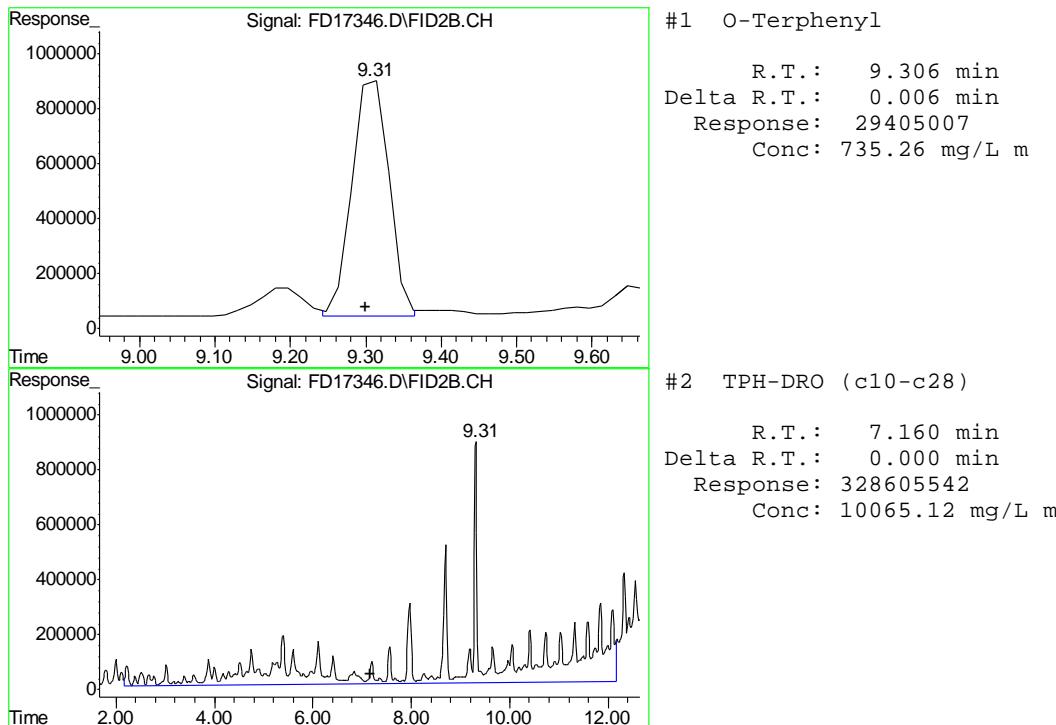
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17346.D Vial: 58
 Acq On : 9-12-2012 04:08:23 PM Operator: alexwl
 Sample : D38480-1 Inst : FID5
 Misc : OP6603,GFD891,30.10,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 9:06 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Multiple Level Calibration
 DataAcq Meth : DRODUAL.M

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





Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17348.D Vial: 59
 Acq On : 9-12-2012 04:34:24 PM Operator: alexwl
 Sample : D38480-2 Inst : FID5
 Misc : OP6603,GFD891,30.05,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 09:07:00 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Initial Calibration
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	9.31	30713320	767.969 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	7.16	93661576	2868.835 mg/L

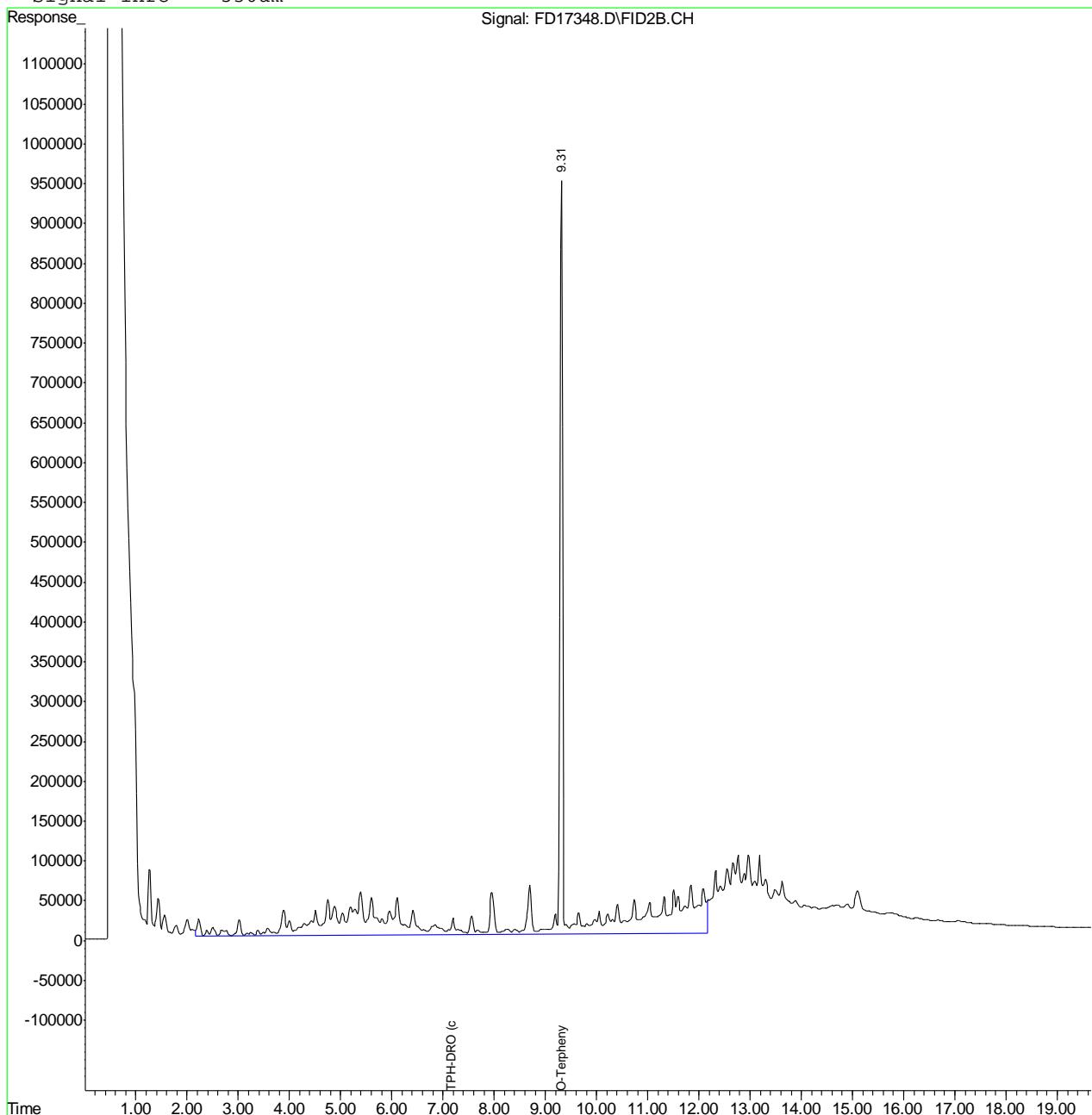
(f)=RT Delta > 1/2 Window (m)=manual int.
 FD17348.D DRO-GFD889R.M Thu Sep 13 09:41:26 2012 GC

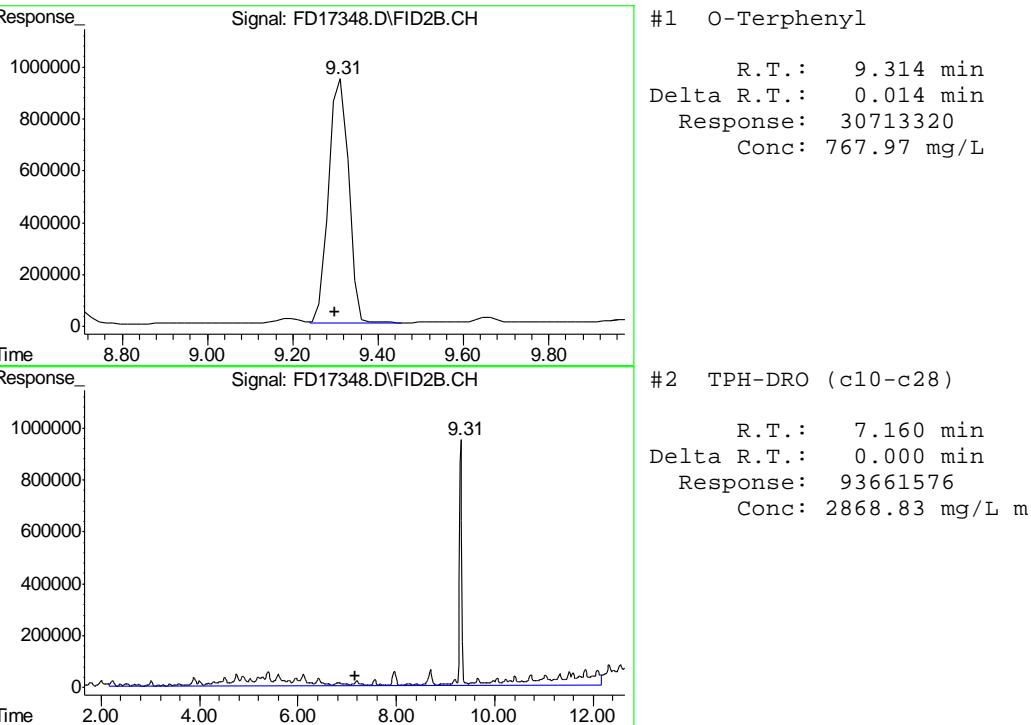
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17348.D Vial: 59
 Acq On : 9-12-2012 04:34:24 PM Operator: alexwl
 Sample : D38480-2 Inst : FID5
 Misc : OP6603,GFD891,30.05,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 9:07 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Multiple Level Calibration
 DataAcq Meth : DRODUAL.M

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





Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17350.D Vial: 60
 Acq On : 9-12-2012 05:00:40 PM Operator: alexwl
 Sample : D38480-3 Inst : FID5
 Misc : OP6603,GFD891,30.11,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 09:07:19 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Initial Calibration
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	9.32	36257724	906.604 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	7.16	174254545	5337.380 mg/L

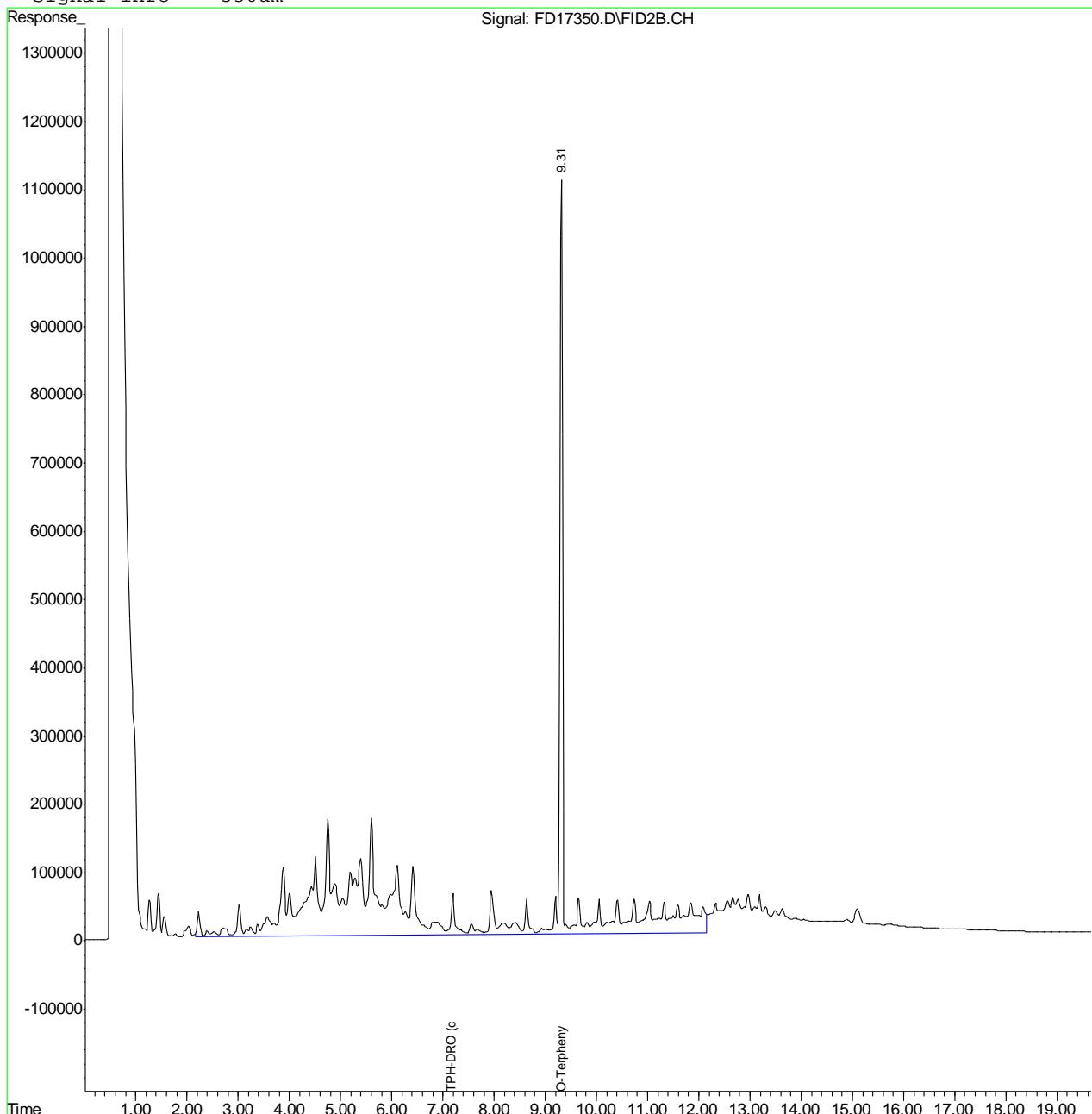
(f)=RT Delta > 1/2 Window (m)=manual int.
 FD17350.D DRO-GFD889R.M Thu Sep 13 09:41:27 2012 GC

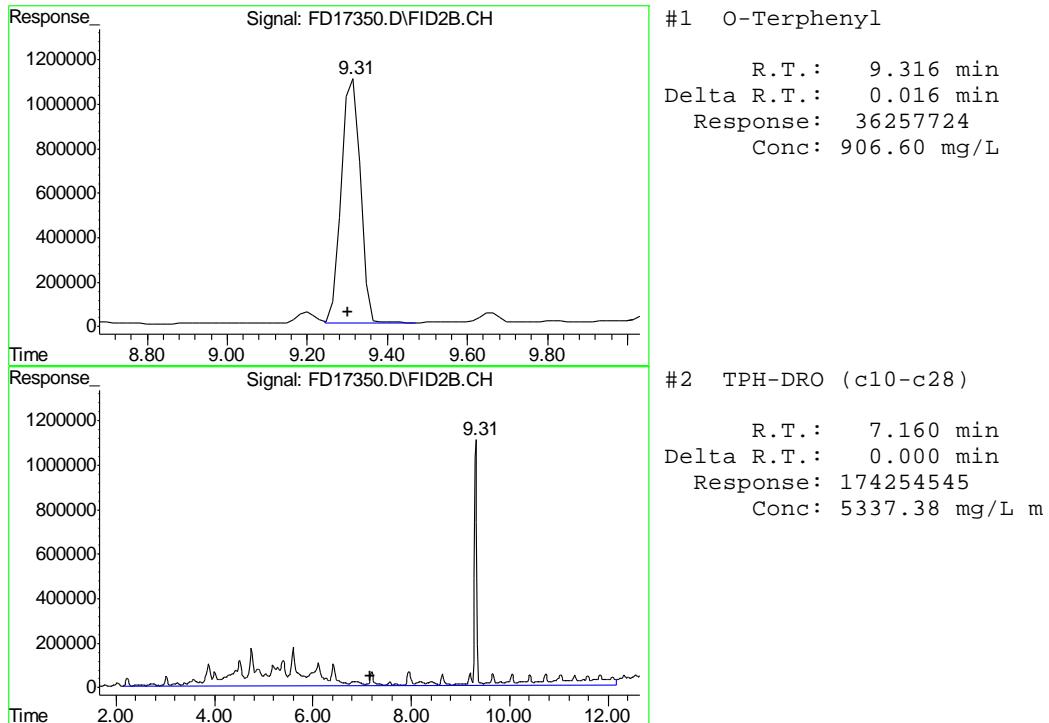
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17350.D Vial: 60
 Acq On : 9-12-2012 05:00:40 PM Operator: alexwl
 Sample : D38480-3 Inst : FID5
 Misc : OP6603,GFD891,30.11,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 9:07 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Multiple Level Calibration
 DataAcq Meth : DRODUAL.M

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



13.1.3
13

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17336.D Vial: 53
 Acq On : 9-12-2012 01:58:44 PM Operator: alexwl
 Sample : OP6603-MB Inst : FID5
 Misc : OP6603,GFD891,30.00,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 09:04:14 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Initial Calibration
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	9.32	36449548	911.400 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	7.16	588395	18.022 mg/L

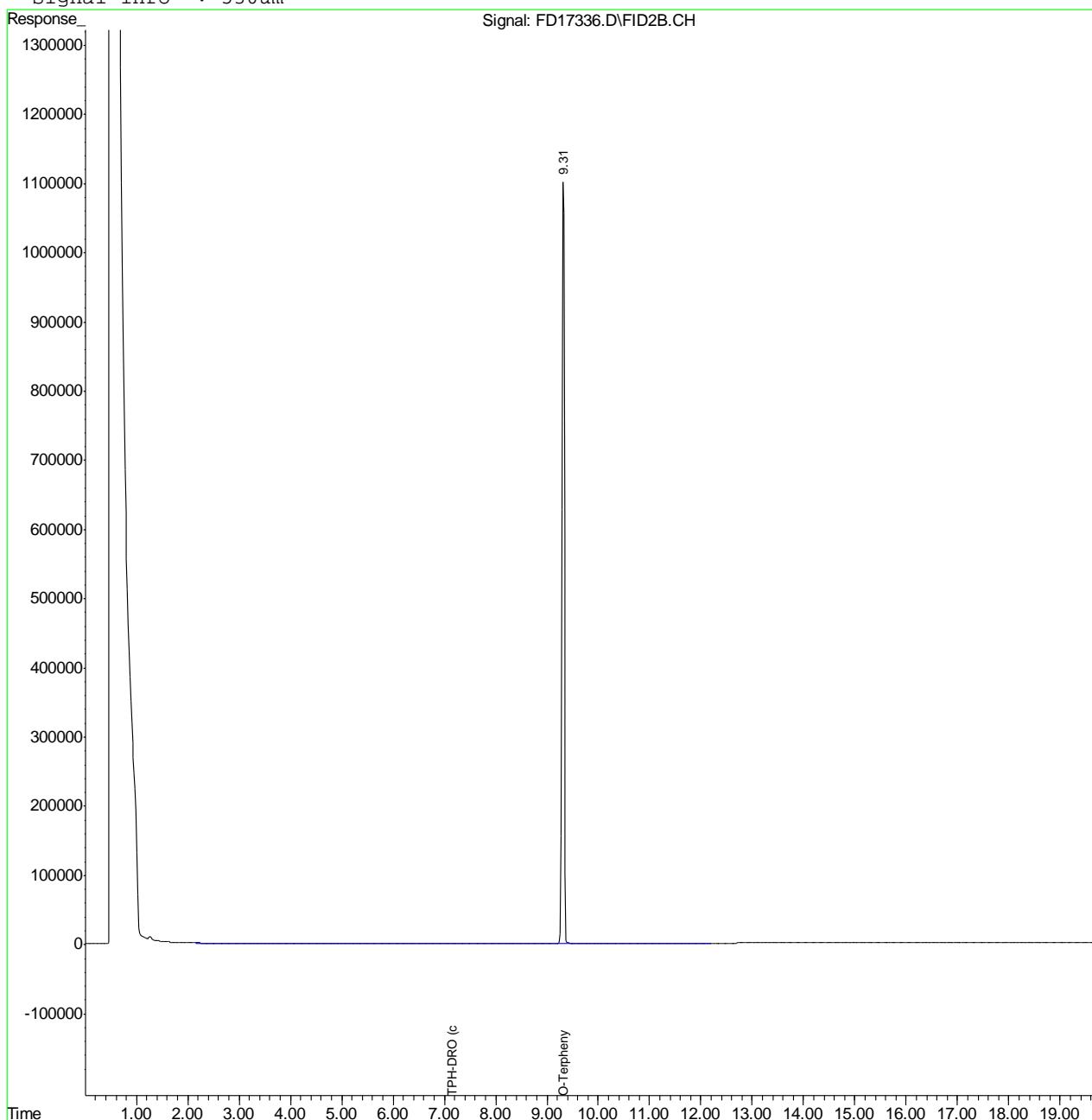
(f)=RT Delta > 1/2 Window (m)=manual int.
 FD17336.D DRO-GFD889R.M Thu Sep 13 09:41:20 2012 GC

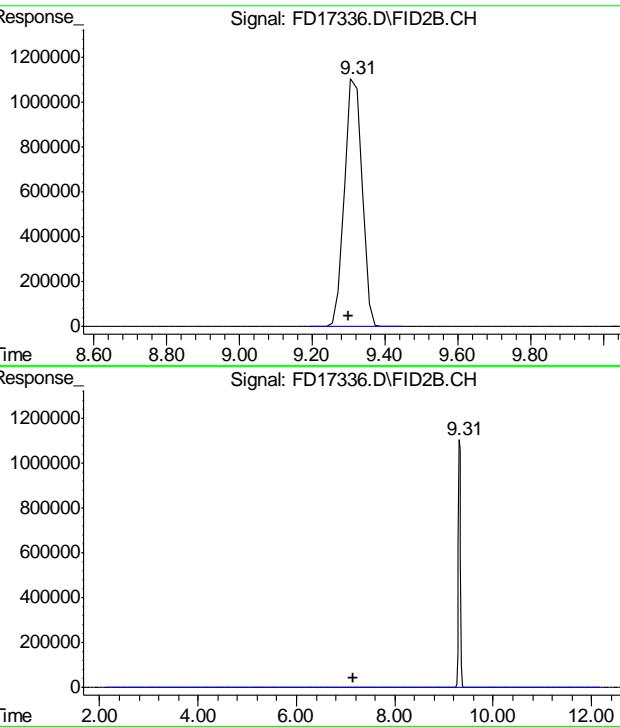
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\SEPTEMBER\FD091212.SEC\FD17336.D Vial: 53
 Acq On : 9-12-2012 01:58:44 PM Operator: alexwl
 Sample : OP6603-MB Inst : FID5
 Misc : OP6603,GFD891,30.00,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Sep 13 9:04 2012 Quant Results File: DRO-GFD889R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD889R.M (Chemstation Integrator)
 Title : 8015B TEH
 Last Update : Wed Sep 12 10:18:23 2012
 Response via : Multiple Level Calibration
 DataAcq Meth : DRODUAL.M

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





#1 O-Terphenyl
 R.T.: 9.320 min
 Delta R.T.: 0.020 min
 Response: 36449548
 Conc: 911.40 mg/L

#2 TPH-DRO (c10-c28)
 R.T.: 7.160 min
 Delta R.T.: 0.000 min
 Response: 588395
 Conc: 18.02 mg/L



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: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8357
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

09/11/12

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.0009	-0.0022	<0.10

Associated samples MP8357: D38480-1, D38480-2, D38480-3

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: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8357
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 09/11/12

Metal	D38479-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.025	0.57	0.55	99.1 75-125

Associated samples MP8357: D38480-1, D38480-2, D38480-3

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: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8357
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

09/11/12

Metal	D38479-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.025	0.53	0.509	99.3	7.3	

Associated samples MP8357: D38480-1, D38480-2, D38480-3

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: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8357
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 09/11/12

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

Associated samples MP8357: D38480-1, D38480-2, D38480-3

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8358
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

09/11/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	2.1	.57		
Antimony	3.0	.36	.12		
Arsenic	2.5	.54	.56		
Barium	1.0	.08	.11	0.010	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.43	.06		
Cadmium	1.0	.06	.036	-0.010	<1.0
Calcium	40	.84	9		
Chromium	1.0	.03	.03	-0.010	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.0	<1.0
Iron	7.0	.19	.87		
Lead	5.0	.24	.24	-0.030	<5.0
Lithium	0.20	.28	.054		
Magnesium	20	2.2	.98		
Manganese	0.50	.12	.022		
Molybdenum	1.0	.21	.08		
Nickel	3.0	.05	.026	0.030	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	15	7		
Selenium	5.0	.61	.36	0.35	<5.0
Silicon	5.0	.65	.37		
Silver	3.0	.05	.06	0.0	<3.0
Sodium	40	2.1	1.9		
Strontium	5.0	.02	.017		
Thallium	1.0	.29	.53		
Tin	5.0	1.2	2		
Titanium	1.0	.01	.038		
Uranium	5.0	.46	.26		
Vanadium	1.0	.03	.036		
Zinc	3.0	.08	.37	-0.070	<3.0

Associated samples MP8358: D38480-1, D38480-2, D38480-3

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8358
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8358
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 09/11/12

Metal	D38480-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	1270	1790	240	216.8(a) 75-125
Beryllium				
Boron				
Cadmium	0.43	62.1	60	102.8 75-125
Calcium				
Chromium	28.9	85.4	60	94.2 75-125
Cobalt				
Copper	28.8	106	60	128.7N(b) 75-125
Iron				
Lead	15.0	135	120	100.0 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	22.4	78.1	60	92.9 75-125
Phosphorus	anr			
Potassium				
Selenium	1.5	118	120	97.1 75-125
Silicon				
Silver	0.30	25.7	24	105.9 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	51.8	111	60	98.7 75-125

Associated samples MP8358: D38480-1, D38480-2, D38480-3

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8358
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8358
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

09/11/12

Metal	D38480-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	anr					
Barium	1270	2280	254	397.0 (a)	24.1 (b)	20
Beryllium						
Boron						
Cadmium	0.43	65.6	63.6	102.5	5.5	20
Calcium						
Chromium	28.9	86.5	63.6	90.6	1.3	20
Cobalt						
Copper	28.8	102	63.6	115.1	3.8	20
Iron						
Lead	15.0	138	127	96.7	2.2	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	22.4	76.2	63.6	84.6	2.5	20
Phosphorus	anr					
Potassium						
Selenium	1.5	124	127	96.3	5.0	20
Silicon						
Silver	0.30	26.9	25.4	104.5	4.6	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	51.8	110	63.6	91.5	0.9	20

Associated samples MP8358: D38480-1, D38480-2, D38480-3

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8358
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8358
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 09/11/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	221	200	110.5	80-120
Beryllium				
Boron				
Cadmium	50.7	50	101.4	80-120
Calcium				
Chromium	51.0	50	102.0	80-120
Cobalt				
Copper	51.6	50	103.2	80-120
Iron				
Lead	100	100	100.0	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	48.5	50	97.0	80-120
Phosphorus	anr			
Potassium				
Selenium	94.7	100	94.7	80-120
Silicon				
Silver	20.4	20	102.0	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	50.4	50	100.8	80-120

Associated samples MP8358: D38480-1, D38480-2, D38480-3

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8358
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

14.2.3
14

SERIAL DILUTION RESULTS SUMMARY

Login Number: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8358
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: ug/l

Prep Date: 09/11/12

Metal	D38480-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	10100	10400	3.5	0-10
Beryllium				
Boron				
Cadmium	3.40	0.00	100.0(a)	0-10
Calcium				
Chromium	230	244	6.3	0-10
Cobalt				
Copper	228	200	12.7*(b)	0-10
Iron				
Lead	119	124	4.0	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	178	191	6.9	0-10
Phosphorus	anr			
Potassium				
Selenium	11.7	0.00	100.0(a)	0-10
Silicon				
Silver	2.40	4.00	66.7 (a)	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	411	439	6.7	0-10

Associated samples MP8358: D38480-1, D38480-2, D38480-3

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8358
Matrix Type: SOLID

Methods: SW846 6010C
Units: ug/l

Prep Date:

Metal

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8359
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date:

09/11/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	0.0047	<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 MP8359: D38480-1, D38480-2, D38480-3

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: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8359
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 09/11/12

Metal	D38480-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	16.8	168	120	126.1N(a) 75-125
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8359: D38480-1, D38480-2, D38480-3

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

(a) Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8359
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

09/11/12

Metal	D38480-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	16.8	172	127	122.0	2.4	20
Barium						
Beryllium						
Boron						
Cadmium						
Calcium						
Chromium						
Cobalt						
Copper						
Iron						
Lead						
Magnesium						
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium						
Silver						
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP8359: D38480-1, D38480-2, D38480-3

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: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8359
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 09/11/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	105	100	105.0	80-120
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8359: D38480-1, D38480-2, D38480-3

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D38480
 Account: XTOKWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8359
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: ug/l

Prep Date: 09/11/12

Metal	D38480-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	134	137	2.3	0-10
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8359: D38480-1, D38480-2, D38480-3

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: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8372
Matrix Type: AQUEOUS

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

Prep Date:

09/11/12

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	930	<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	37.5	<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	-71	<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 MP8372: D38480-1A, D38480-2A, D38480-3A

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8372
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: D38480
 Account: XTOKWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8372
 Matrix Type: AQUEOUS

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

Prep Date:

09/11/12

Metal	D38518-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	129000	256000	125000	101.6
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	78.5	125000	125000	99.9
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	2590000	2490000	125000	-80.0(a)
Strontium				75-125
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8372: D38480-1A, D38480-2A, D38480-3A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8372
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 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: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8372
 Matrix Type: AQUEOUS

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

Prep Date: 09/11/12

Metal	D38518-1A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	129000	247000	125000	94.4	3.6
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	78.5	125000	125000	99.9	0.0
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	2590000	2290000	125000	-240.0 (a)	8.4
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP8372: D38480-1A, D38480-2A, D38480-3A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D38480
Account: XTOKRWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8372
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 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: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8372
 Matrix Type: AQUEOUS

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

Prep Date: 09/11/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	132000	125000	105.6	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	122000	125000	97.6	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	125000	125000	100.0	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8372: D38480-1A, D38480-2A, D38480-3A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8372
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

14.4.3
14

SERIAL DILUTION RESULTS SUMMARY

Login Number: D38480
 Account: XTOKRWR - XTO Energy
 Project: PCU 197-36A

QC Batch ID: MP8372
 Matrix Type: AQUEOUS

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

Prep Date:

09/11/12

Metal	D38518-1A	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	25700	26000		1.0	0-10
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	15.7	0.00	100.0(a)	0-10	
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	517000	527000		1.9	0-10
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP8372: D38480-1A, D38480-2A, D38480-3A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

QC Batch ID: MP8372
Matrix Type: AQUEOUS

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

Prep Date:

Metal

- (anr) Analyte not requested
(a) Percent difference acceptable due to low initial sample concentration (< 50 times IDL).



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: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP8138/GN16704	1.0	0.0	mg/kg	60.7	67.2	111.0	80-120%
Specific Conductivity	GP8183/GN16757	1.0	<1.0	umhos/cm	9989	9910	99.2	90-110%
pH	GN16667			su	8.00	8.01	100.1	99.3-100.7%

Associated Samples:

Batch GP8138: D38480-1, D38480-2, D38480-3

Batch GP8183: D38480-1, D38480-2, D38480-3

Batch GN16667: D38480-1, D38480-2, D38480-3

(*) Outside of QC limits

DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP8138/GN16704 GN16683	D38513-1 D38518-2	mg/kg mv	0.0 141	0.0 143	22.4(a) 1.4	0-20% 0-20%

Associated Samples:

Batch GP8138: D38480-1, D38480-2, D38480-3

Batch GN16683: D38480-1, D38480-2, D38480-3

(*) Outside of QC limits

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

MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP8138/GN16704	D38513-1	mg/kg	0.0	40	40.5	101.0	75-125%

Associated Samples:

Batch GP8138: D38480-1, D38480-2, D38480-3

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D38480
Account: XTOKWR - XTO Energy
Project: PCU 197-36A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP8138/GN16704	D38513-1	mg/kg	0.0	40	41.0	1.2	

Associated Samples:

Batch GP8138: D38480-1, D38480-2, D38480-3

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

15.4

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