

Quantitation Report (Qedit)

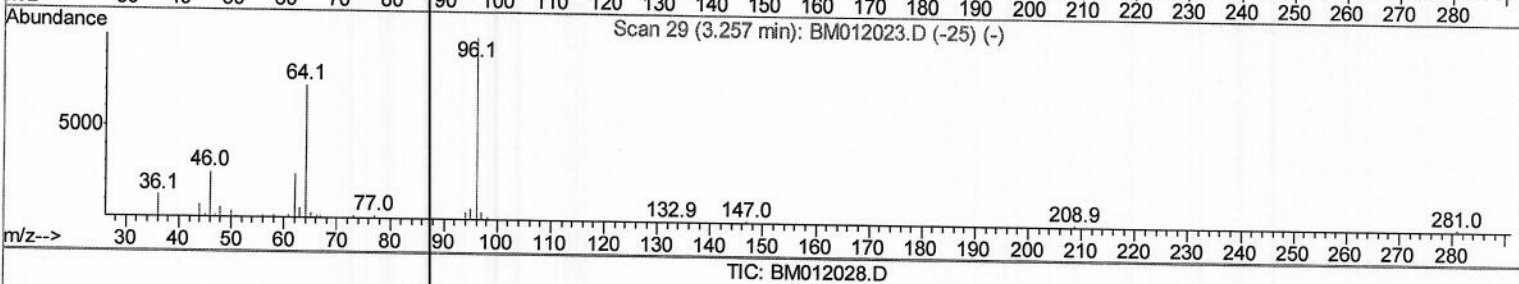
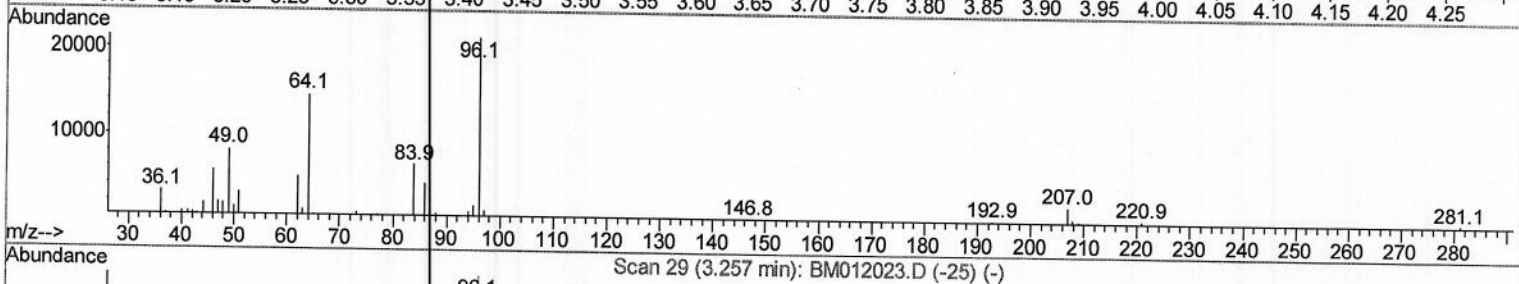
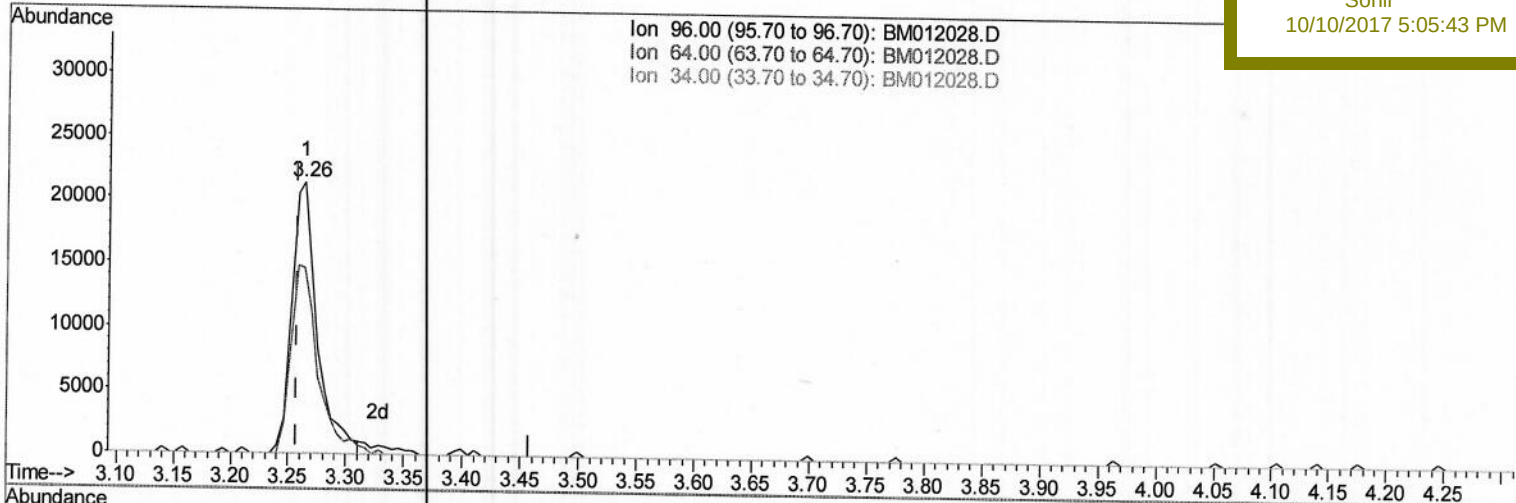
Data Path : Z:\HPCHEM1\BNA M\DATA\BM100917\
 Data File : BM012028.D
 Acq On : 10 Oct 2017 00:27
 Operator : SJ/JU
 Sample : I5666-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 10 02:38:08 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 02:27:01 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 H0AD0

Manual Integrations
 APPROVED

Sohil
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(3) 1,4-Dioxane-d8 (S)
 3.263min (+0.006) 7.18ng/uL
 response 32591

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	71.89
34.00	0.00	0.00
0.00	0.00	0.00

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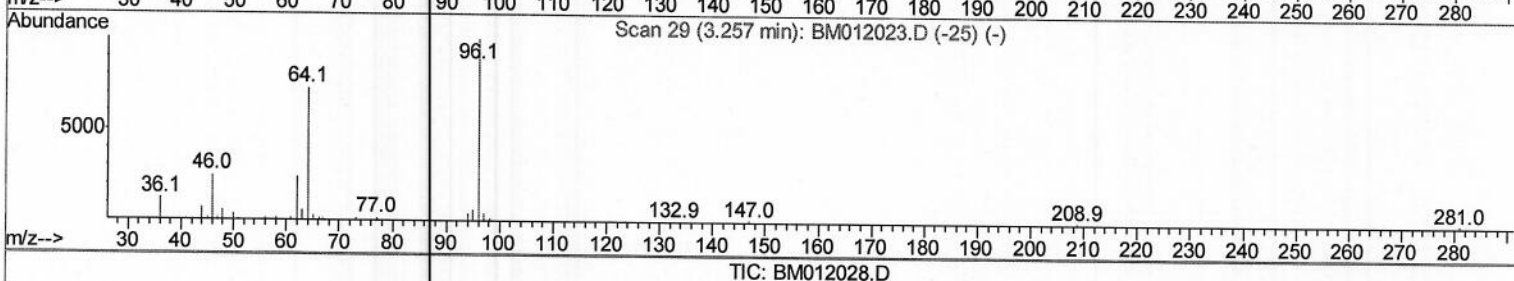
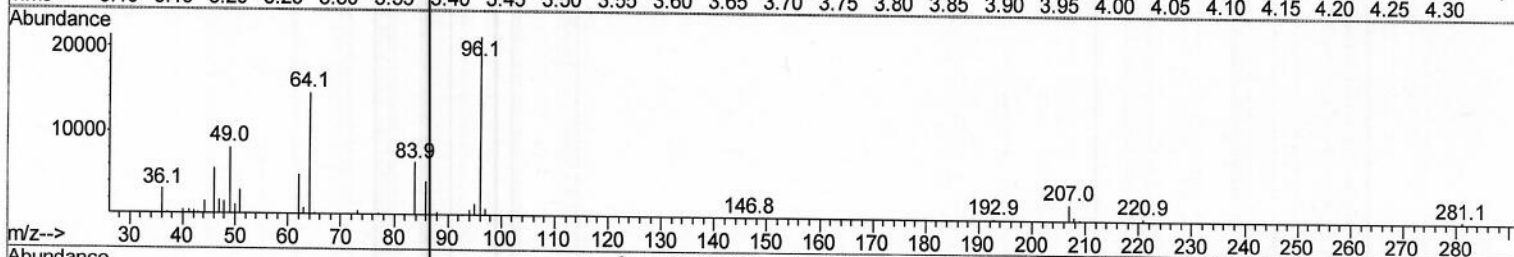
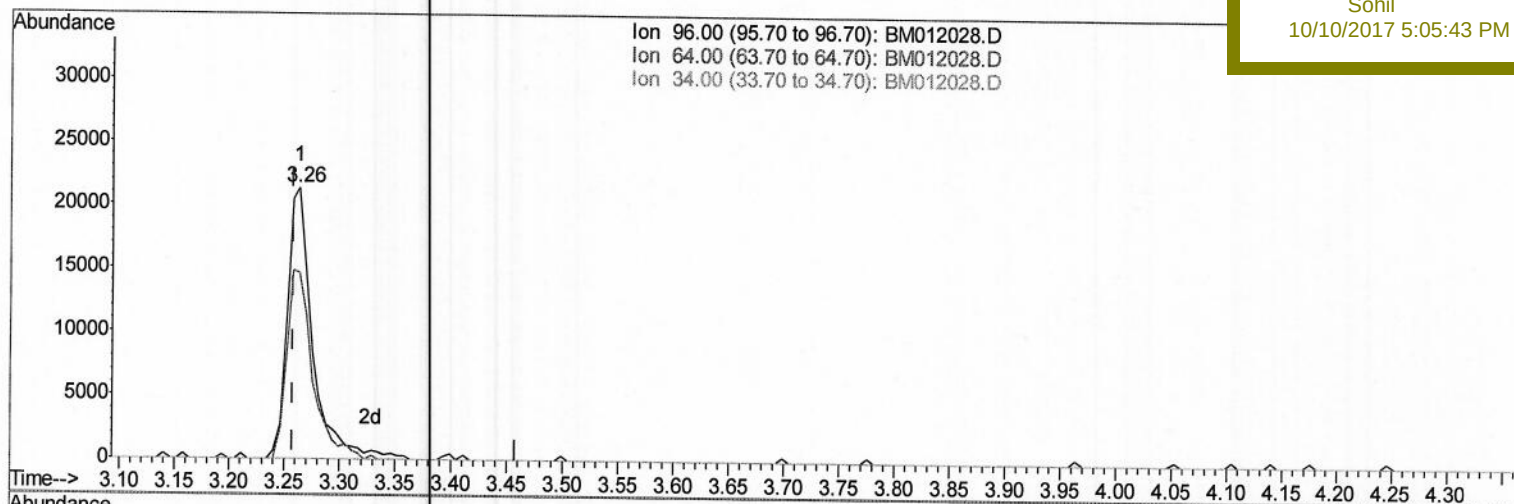
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(3) 1,4-Dioxane-d8 (S)

3.263min (+0.006) 7.50ng/uL m

response 34027

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	68.86
34.00	0.00	0.00
0.00	0.00	0.00

SJ
 10/17/17

Quantitation Report (Qedit)

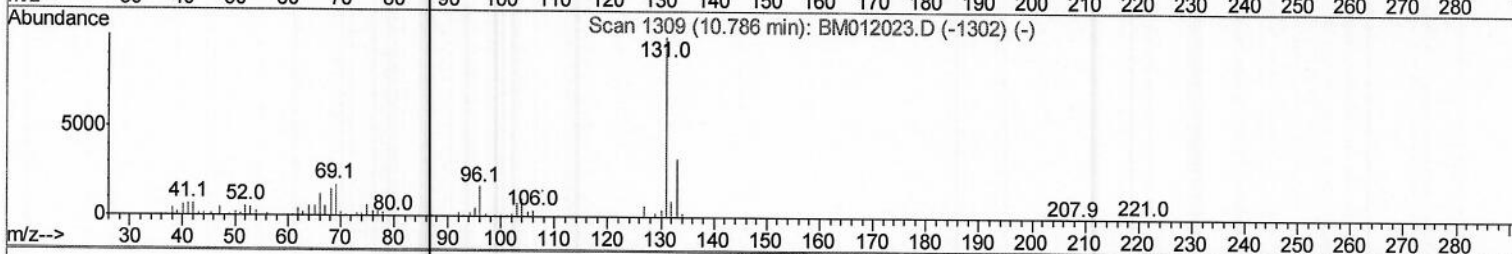
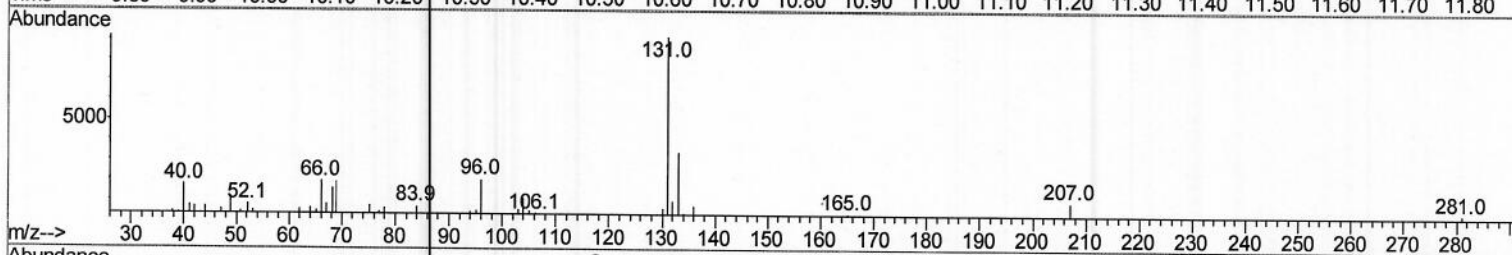
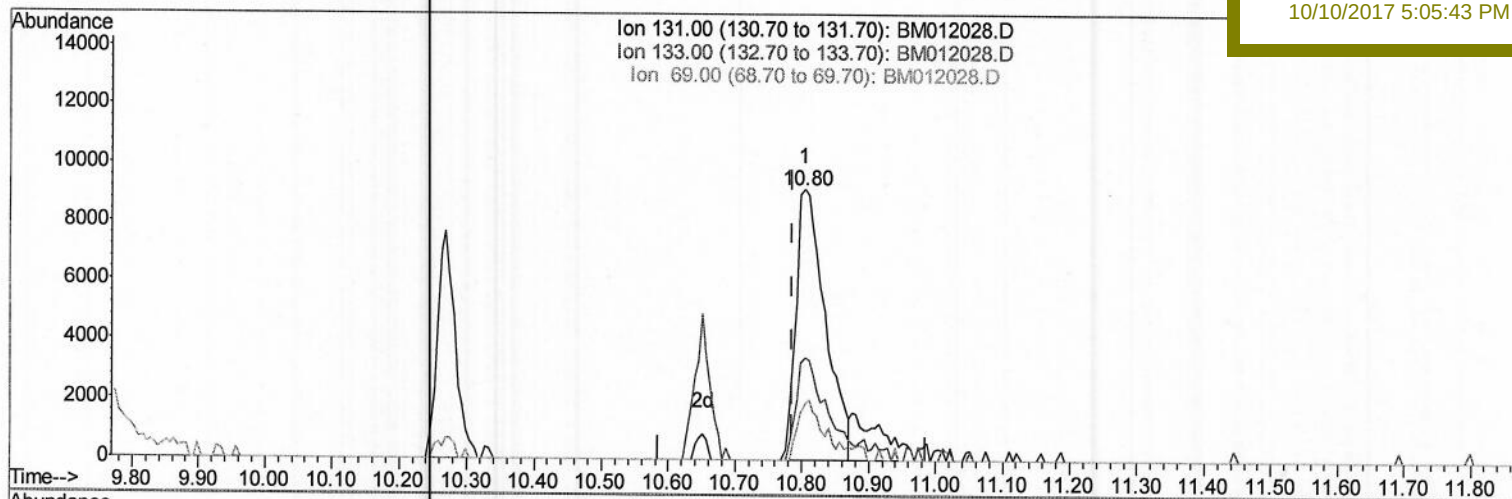
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 Acq On : 10 Oct 2017 00:27
 Operator : SJ/JU
 Sample : I5666-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 10 02:53:41 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
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TIC: BM012028.D

(29) 4-Chloroaniline-d4 (S)
 10.804min (+0.018) 1.66ng/ul
 response 27683

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	37.19
69.00	24.30	20.38
0.00	0.00	0.00

Quantitation Report (Qedit)

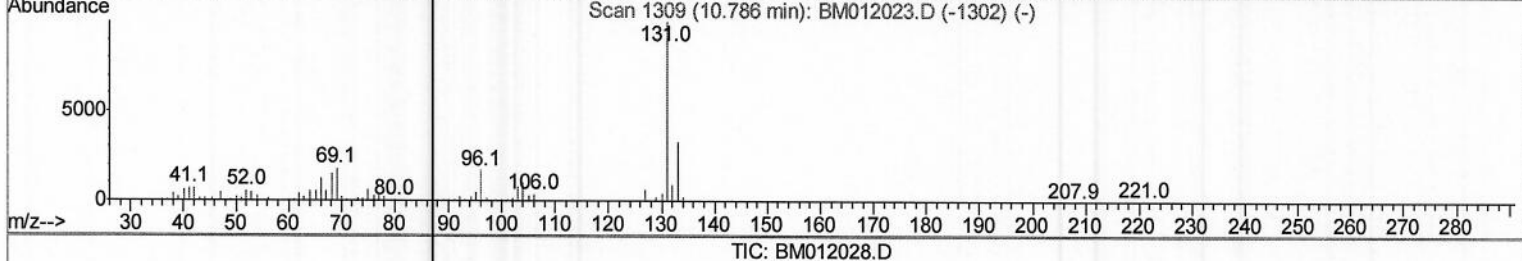
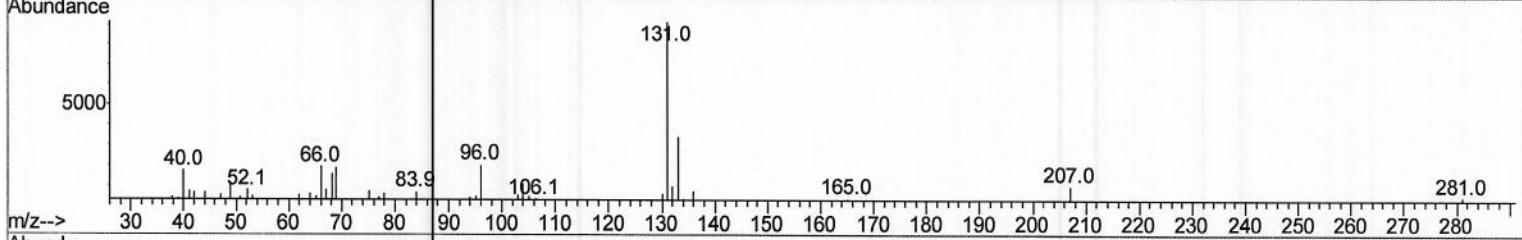
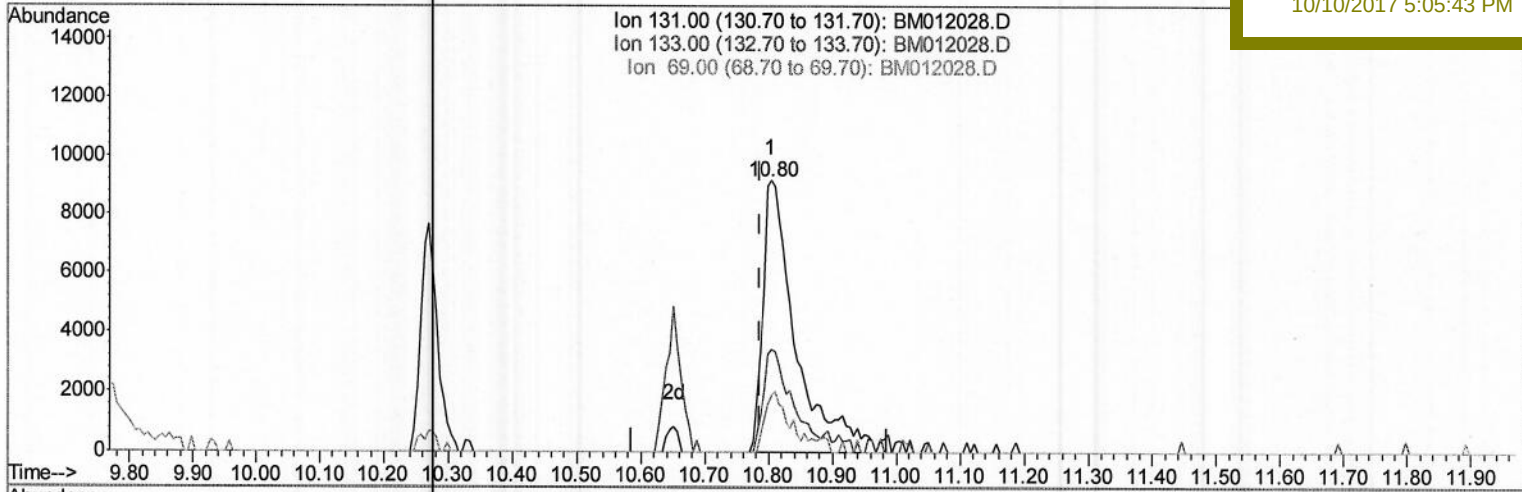
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 Operator : SJ/JU
 Sample : I5666-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

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Instrument :
 BNA_M
 ClientSampled :
 H0AD0

Manual Integrations
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TIC: BM012028.D

(29) 4-Chloroaniline-d4 (S)

10.804min (+0.018) 1.97ng/ul m

response 32842

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	37.19
69.00	24.30	20.38
0.00	0.00	0.00

> SJ
 10/17/17

Quantitation Report (Qedit)

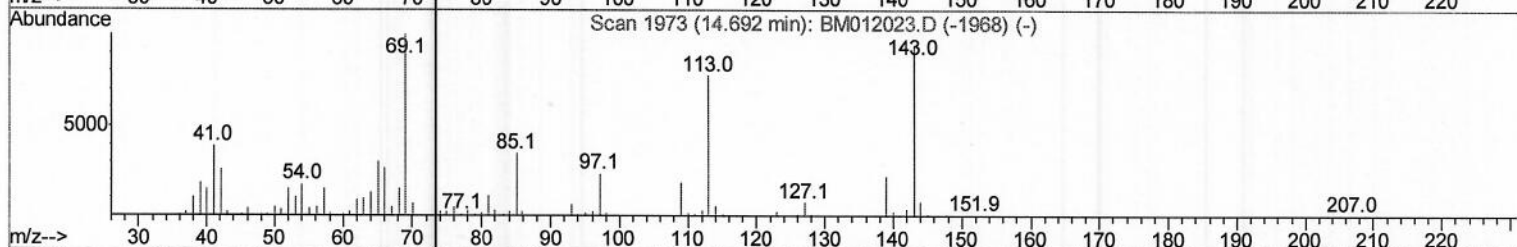
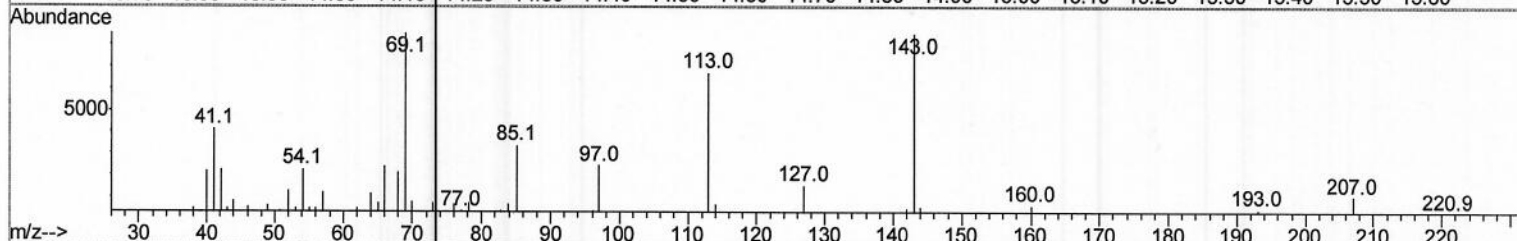
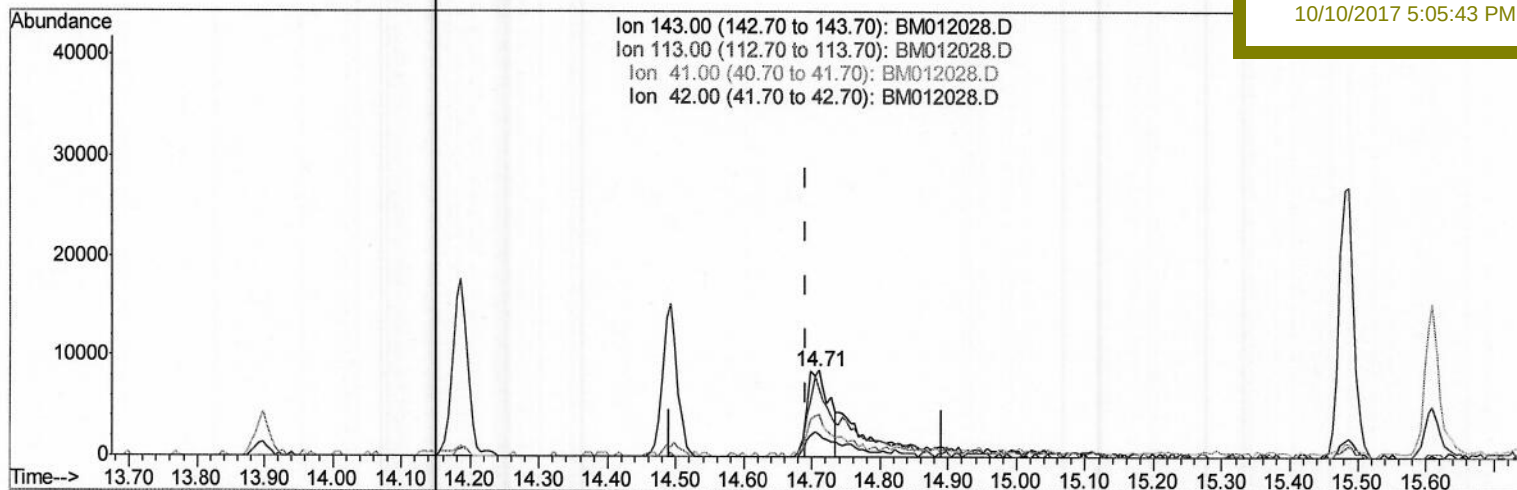
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Instrument :
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 ClientSampleId :
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Manual Integrations
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TIC: BM012028.D

(51) 4-Nitrophenol-d4 (S)

14.710min (+0.018) 2.13ng/ul

response 19074

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	79.20#
41.00	34.80	48.43#
42.00	20.50	26.18#

Quantitation Report (Qedit)

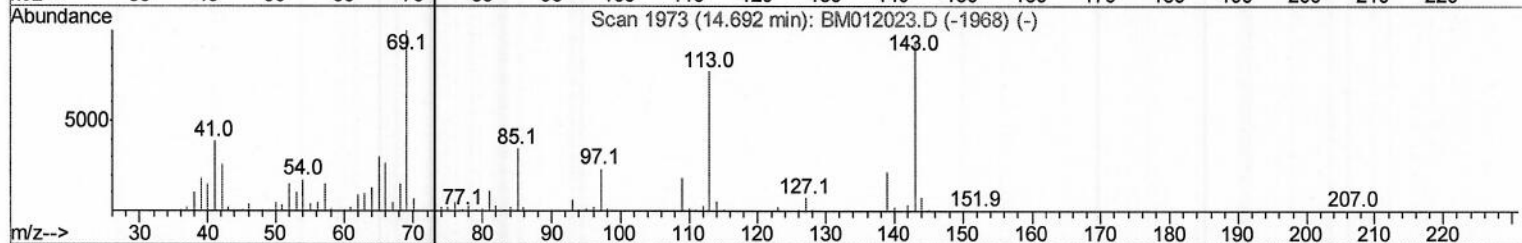
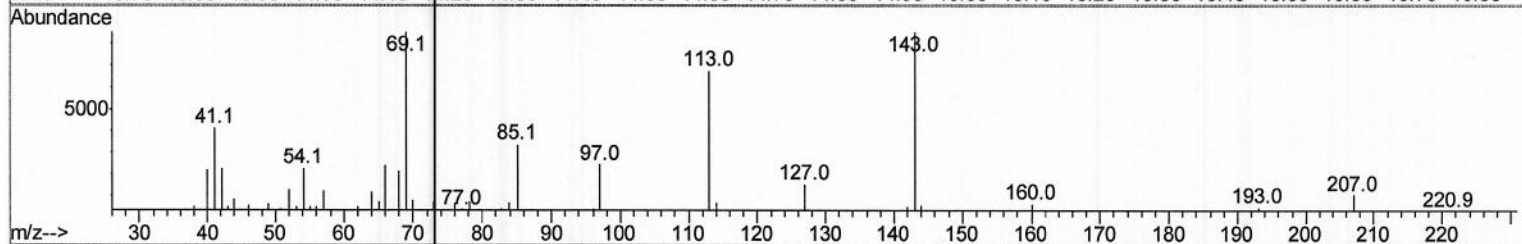
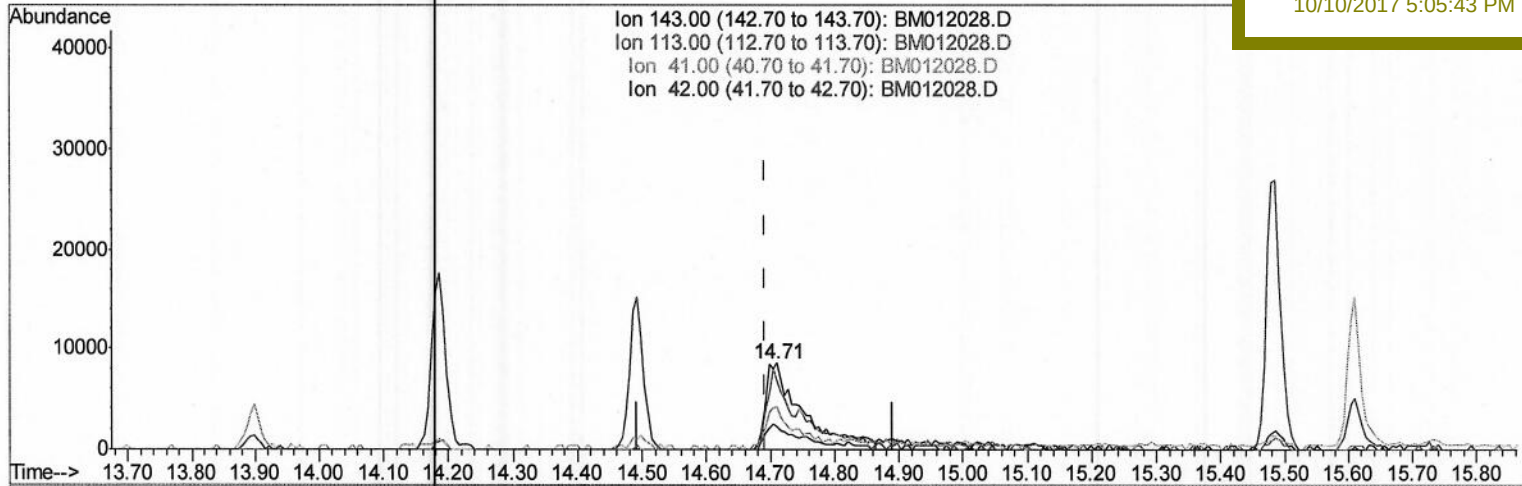
Data Path : Z:\HPCHEM1\BNA M\DATA\BM100917\
 Data File : BM012028.D
 Acq On : 10 Oct 2017 00:27
 Operator : SJ/JU
 Sample : I5666-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 10 02:54:34 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 02:27:01 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleID :
 H0AD0

Manual Integrations
 APPROVED

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TIC: BM012028.D

(51) 4-Nitrophenol-d4 (S)

14.710min (+0.018) 3.83ng/ul m

response 34265

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	79.20#
41.00	34.80	48.43#
42.00	20.50	26.18#

> SJ
 10/17/17

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100917\
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 Acq On : 10 Oct 2017 00:27
 Operator : SJ/JU
 Sample : I5666-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Time: Oct 10 02:55:23 2017
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	214610	20.00	ng/ul	0.00
18) Naphthalene-d8	10.65	136	940712	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.49	164	590179	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.23	188	1427486	20.00	ng/ul	0.00
75) Chrysene-d12	21.41	240	1692582	20.00	ng/ul	0.00
83) Perylene-d12	23.73	264	1794645	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.26	96	34027m	7.50	ng/uL	0.00
5) Phenol-d5	7.00	99	121919	6.60	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.17	67	319350	29.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.37	132	364674	24.62	ng/ul	0.00
13) 4-Methylphenol-d8	8.55	113	246409	15.44	ng/ul	0.00
19) Nitrobenzene-d5	9.01	128	211498	31.57	ng/ul	0.00
22) 2-Nitrophenol-d4	9.73	143	240805	31.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.27	165	438471	28.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.80	131	32842m	1.97	ng/ul	0.02
43) Dimethylphthalate-d6	13.90	166	1699363	33.22	ng/ul	0.00
46) Acenaphthylene-d8	14.19	160	1980401	32.95	ng/ul	0.00
51) 4-Nitrophenol-d4	14.71	143	34265m	3.83	ng/ul	0.02
57) Fluorene-d10	15.49	176	1434642	31.78	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.61	200	254797	26.53	ng/ul	0.00
70) Anthracene-d10	17.33	188	2369823	33.69	ng/ul	0.00
76) Pyrene-d10	19.63	212	2918739	38.23	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.58	264	2928053	33.92	ng/ul	0.00

Target Compounds Ovalue

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

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