

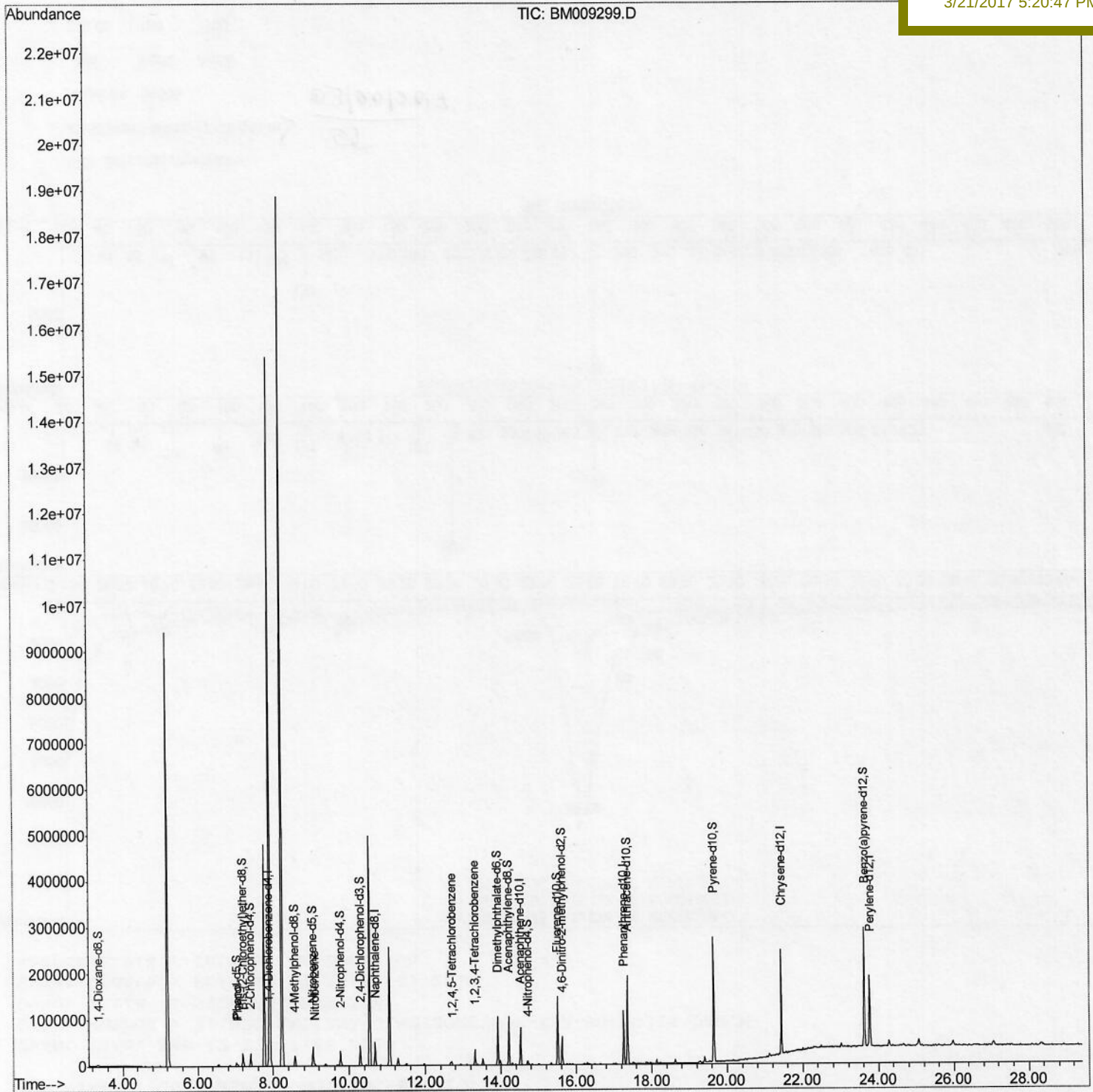
Data Path : Z:\HPCHEM1\BNA_M\Data\BM032017\
Data File : BM009299.D
Acq On : 20 Mar 2017 17:34
Operator : SJ/MA
Sample : I2303-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Mar 20 18:20:50 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031317MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 20 13:19:54 2017
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
C0BH2

Manual Integrations
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Quantitation Report (Qedit)

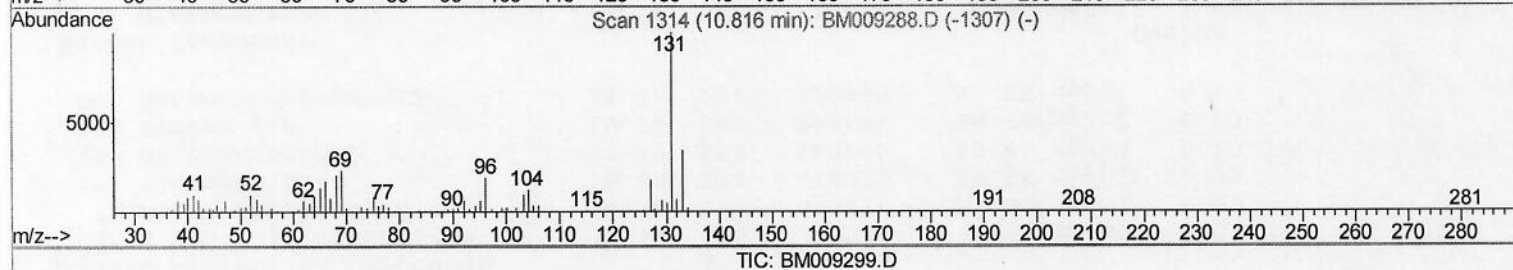
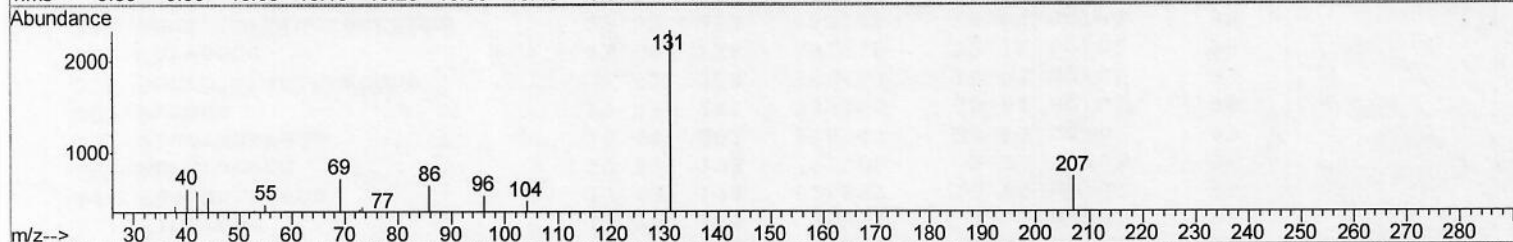
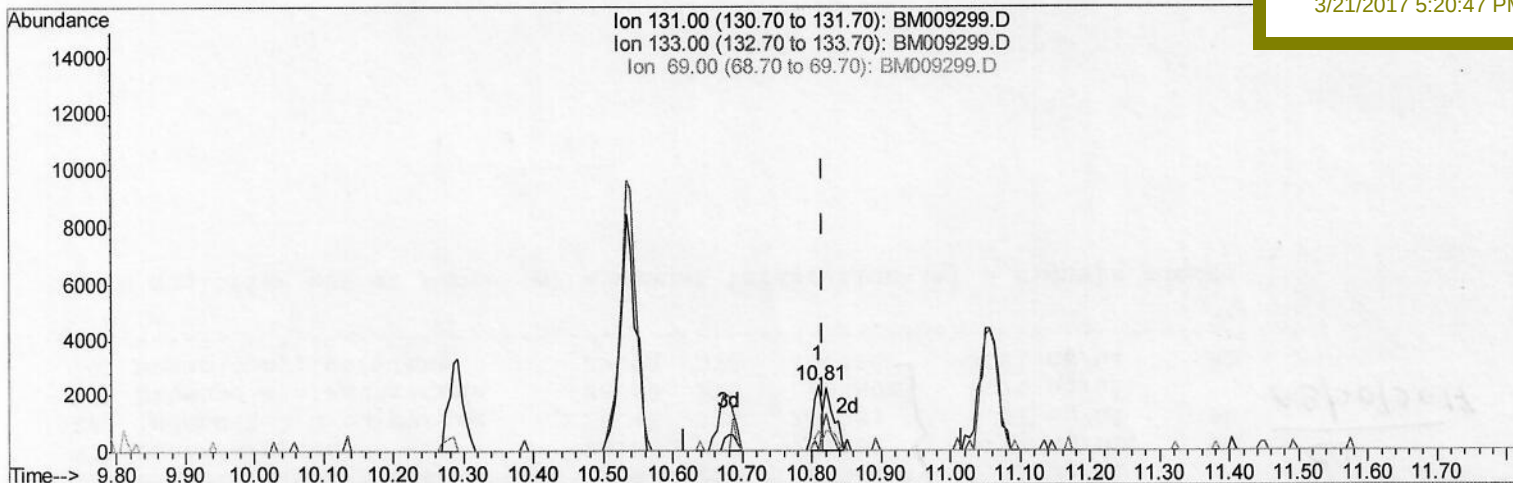
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Manual Integrations
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(29) 4-Chloroaniline-d4 (S)

10.810min (-0.006) 0.29ng/ul

response 2100

Ion	Exp%	Act%
131.00	100	100
133.00	32.10	0.00#
69.00	16.20	30.42#
0.00	0.00	0.00

Quantitation Report (Qedit)

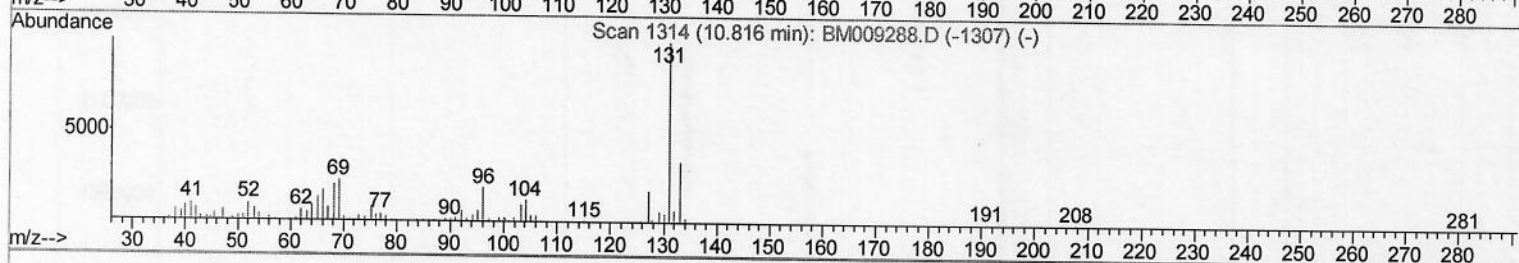
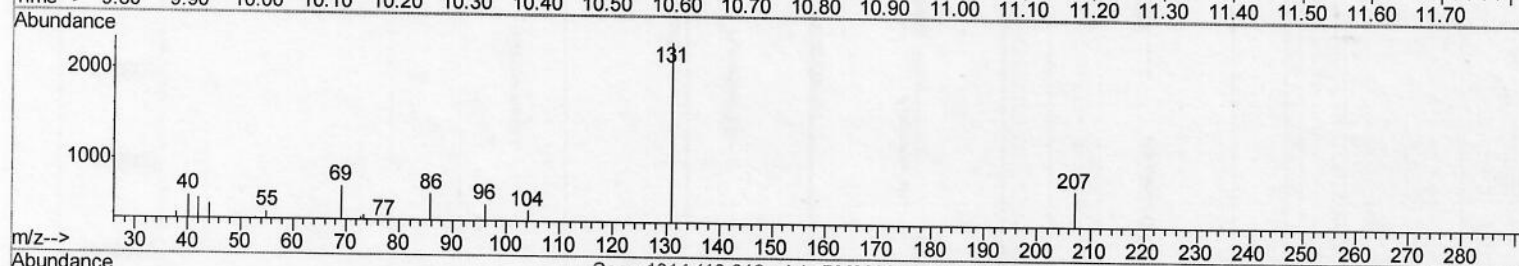
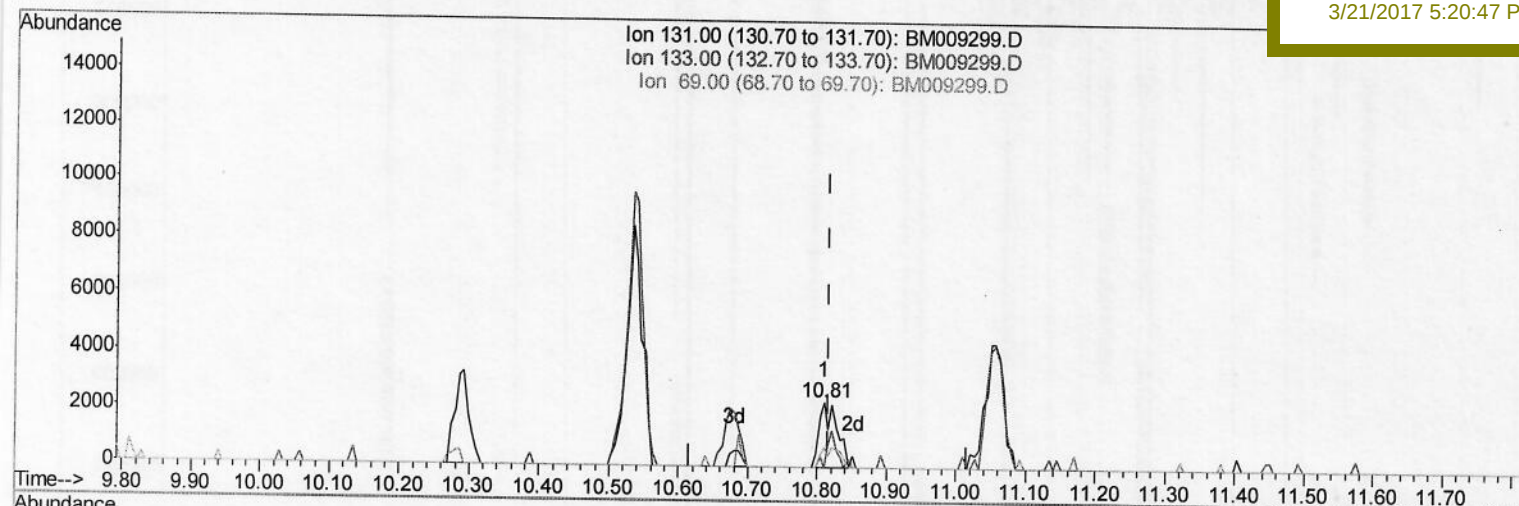
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Manual Integrations
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TIC: BM009299.D

(29) 4-Chloroaniline-d4 (S)

10.810min (-0.006) 0.56ng/ul m>

response 4088

Ion	Exp%	Act%
131.00	100	100
133.00	32.10	0.00#
69.00	16.20	30.42#
0.00	0.00	0.00

55
03/27/2017

Quantitation Report (Qedit)

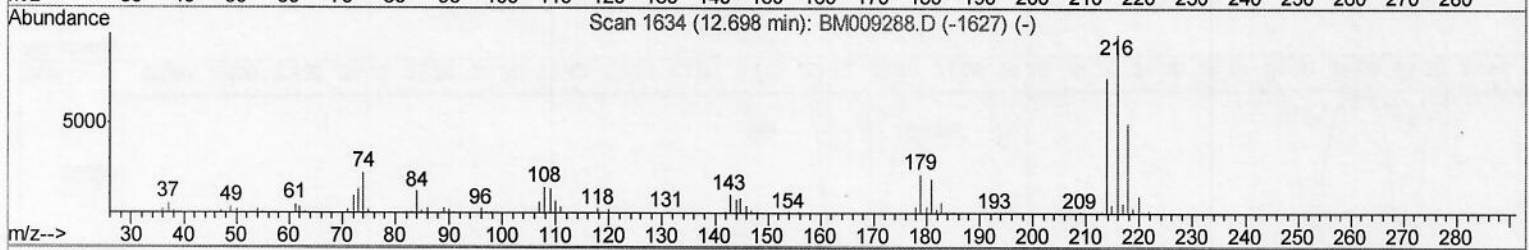
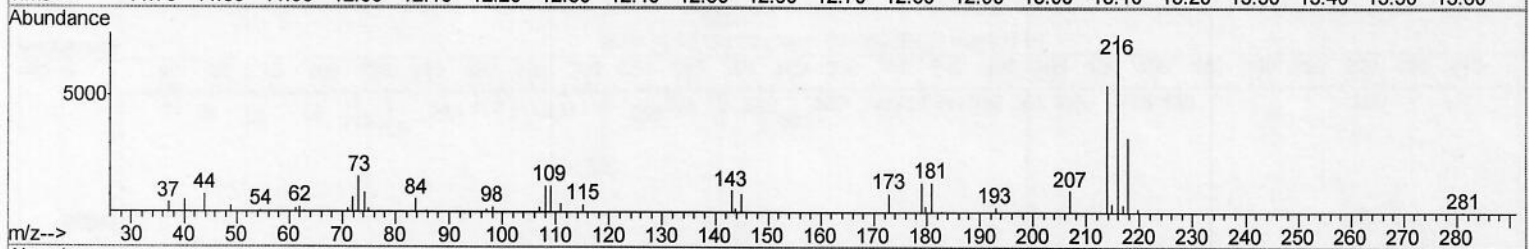
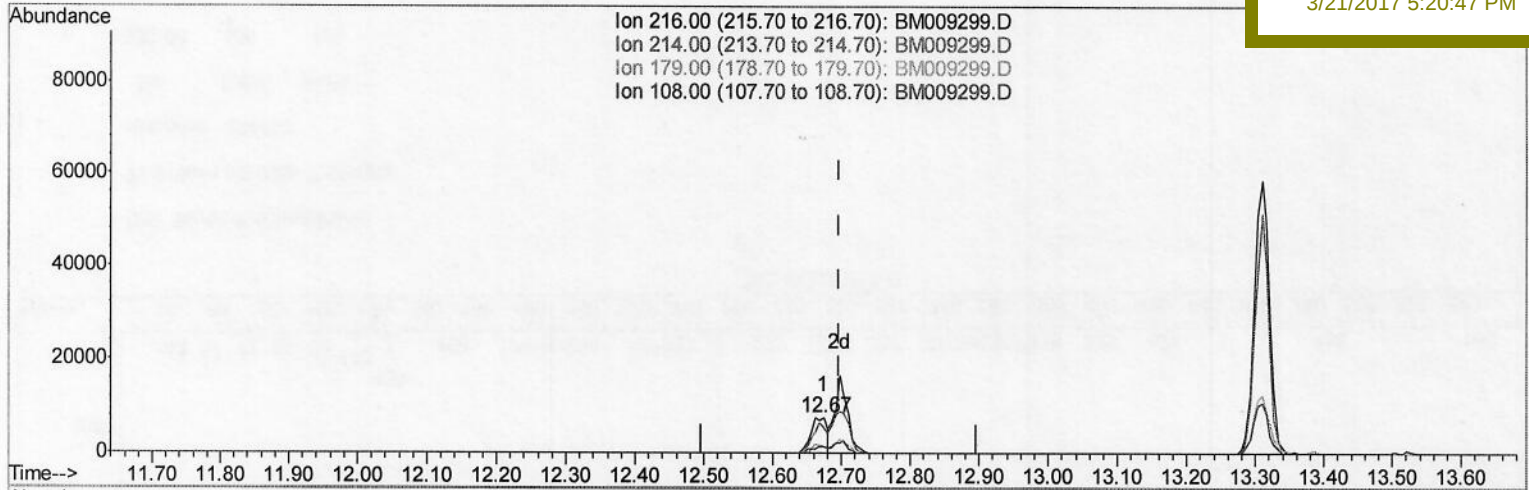
Data Path : Z:\HPCHEM1\BNA_M\Data\BM032017\
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ClientSampleId :
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Manual Integrations
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TIC: BM009299.D

(36) 1,2,4,5-Tetrachlorobenzene

12.674min (-0.024) 1.73ng/ul

response 12235

Ion	Exp%	Act%
216.00	100	100
214.00	78.10	72.48
179.00	22.70	19.51
108.00	17.70	17.70

Quantitation Report (Qedit)

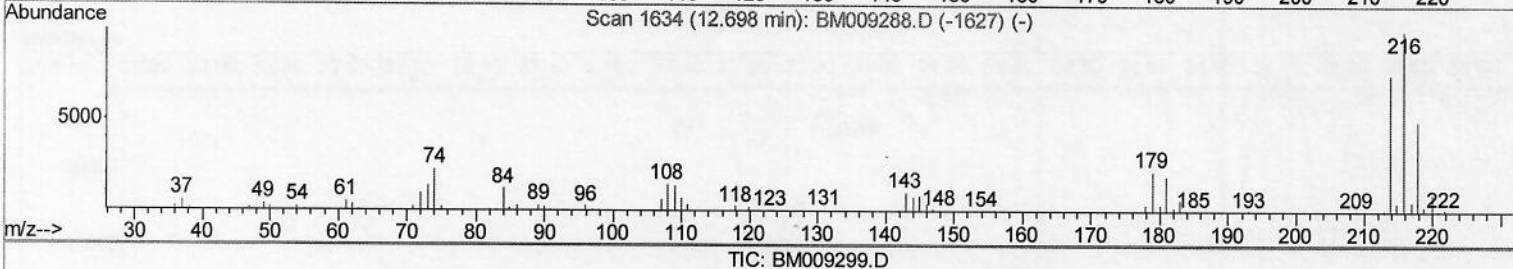
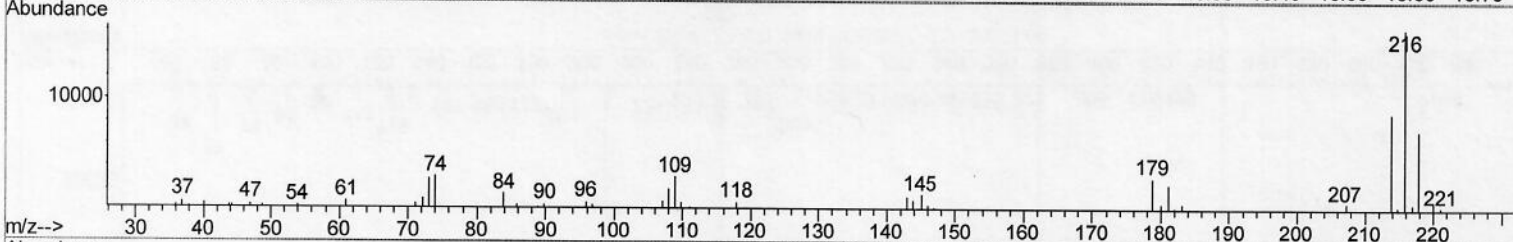
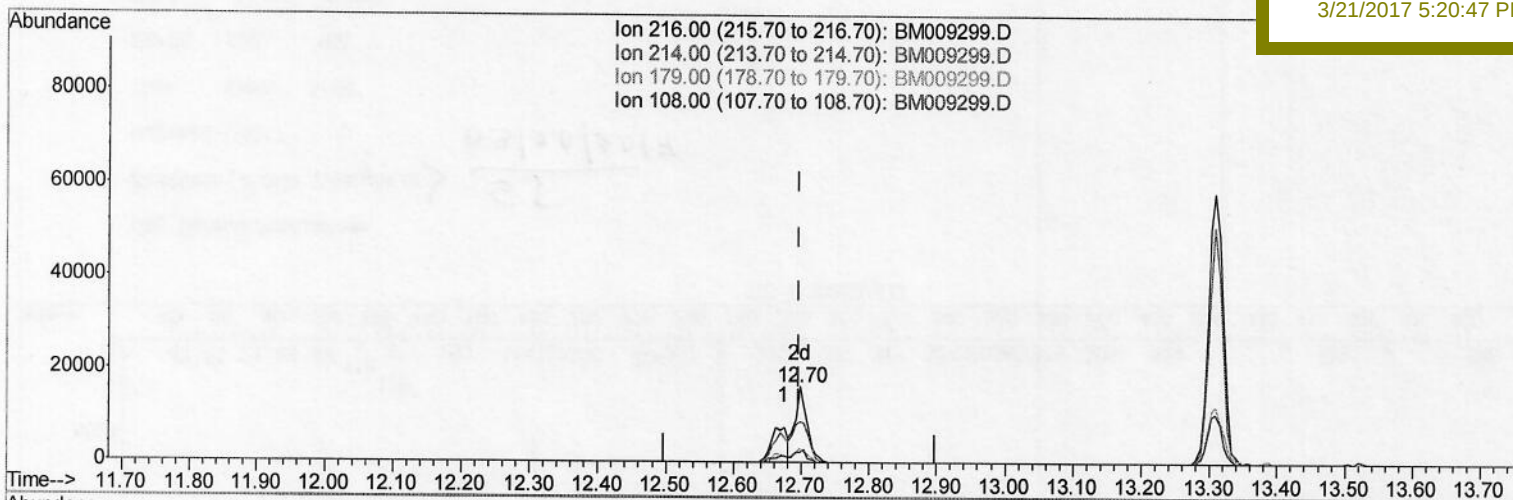
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 Data File : BM009299.D
 Acq On : 20 Mar 2017 17:34
 Operator : SJ/MA
 Sample : I2303-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BH2

Quant Time: Mar 20 18:18:42 2017
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Manual Integrations
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TIC: BM009299.D

(36) 1,2,4,5-Tetrachlorobenzene

12.698min (-0.000) 2.92ng/ul m

response 20691

55
 03/27/2017

Ion	Exp%	Act%
216.00	100	100
214.00	78.10	54.45#
179.00	22.70	18.75
108.00	17.70	12.65#

Data Path : Z:\HPCHEM1\BNA_M\Data\BM032017\
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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	80276	20.00	ng/ul	0.00
18) Naphthalene-d8	10.67	136	360614	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	235812	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.26	188	628137	20.00	ng/ul	0.00
77) Chrysene-d12	21.43	240	1024118	20.00	ng/ul	0.00
85) Perylene-d12	23.77	264	1121679	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.35	96	9688	5.32	ng/uL	0.00
5) Phenol-d5	7.03	99	34532	3.98	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.20	67	130471	23.07	ng/ul	0.00
9) 2-Chlorophenol-d4	7.40	132	97319	18.34	ng/ul	0.00
13) 4-Methylphenol-d8	8.57	113	68991	9.96	ng/ul	0.00
19) Nitrobenzene-d5	9.04	128	69671	27.05	ng/ul	0.00
22) 2-Nitrophenol-d4	9.76	143	73696	25.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.29	165	135902	22.26	ng/ul	0.00
29) 4-Chloroaniline-d4	10.81	131	4088m	0.56	ng/ul	0.00
43) Dimethylphthalate-d6	13.92	166	577400	32.24	ng/ul	0.00
46) Acenaphthylene-d8	14.21	160	663117	32.20	ng/ul	0.00
51) 4-Nitrophenol-d4	14.71	143	17307	4.87	ng/ul	0.00
57) Fluorene-d10	15.51	176	527073	32.98	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	98124	26.11	ng/ul	0.00
70) Anthracene-d10	17.36	188	1001021	33.67	ng/ul	0.00
78) Pyrene-d10	19.65	212	1316973	31.86	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.62	264	1716000	33.44	ng/ul	0.00

Target Compounds

					Qvalue	
6) Phenol	7.06	94	14240	1.54	ng/ul#	89
20) Nitrobenzene	9.09	77	20903	2.43	ng/ul#	80
36) 1,2,4,5-Tetrachlorobenzene	12.70	216	20691m	2.92	ng/ul	
72) 1,2,3,4-Tetrachlorobenzene	13.31	216	88514	11.99	ng/uL	90

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

SJ
 03/27/2017

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