

Quantitation Report (Qedit)

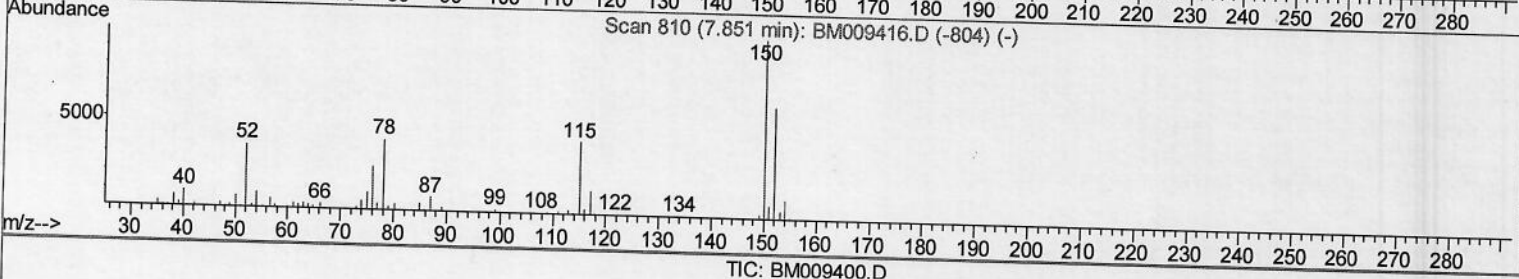
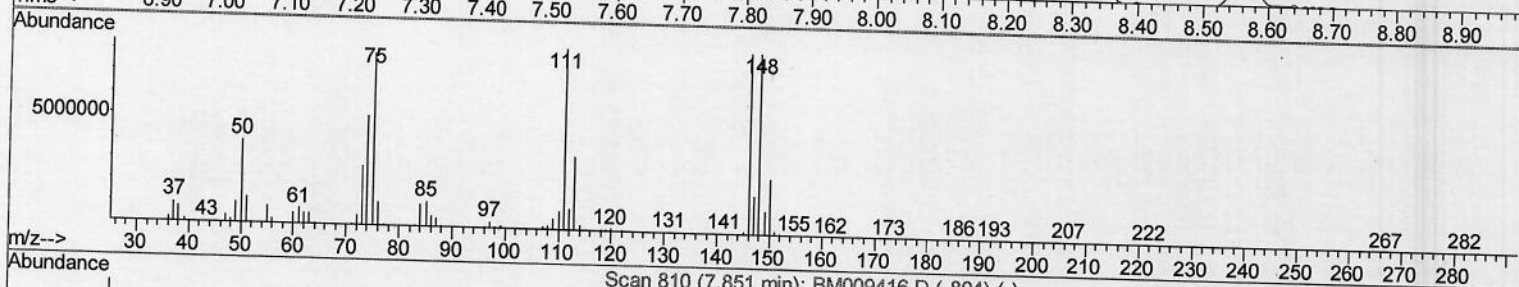
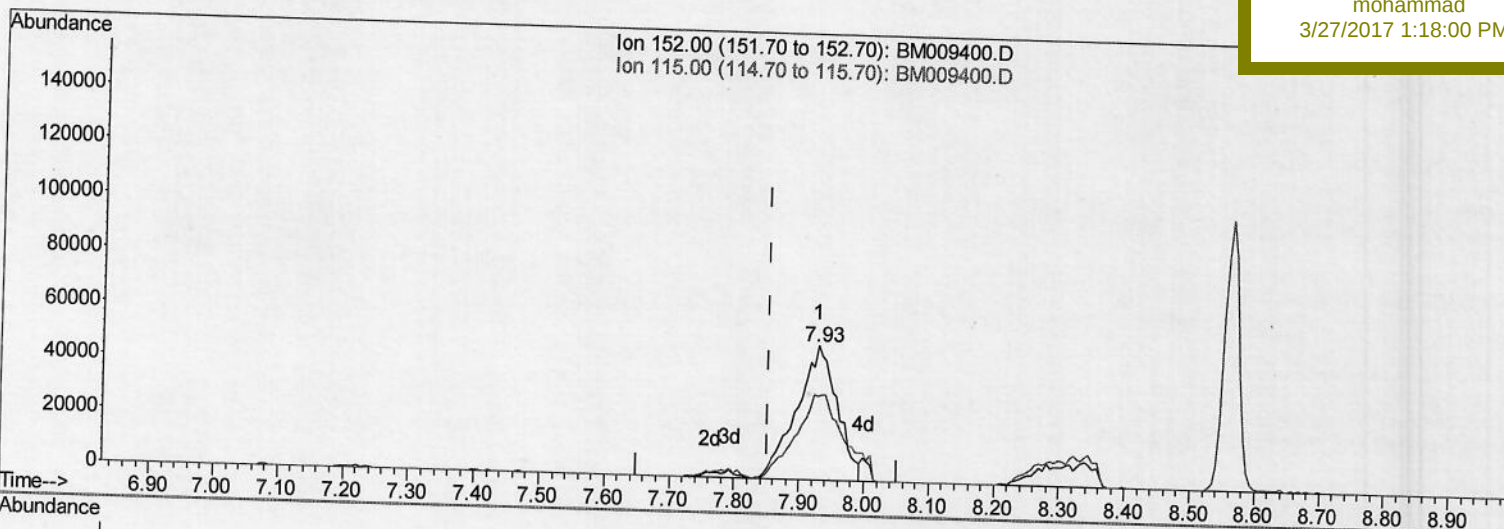
Data Path : Z:\HPCHEM1\BNA M\Data\BM032417\
 Data File : BM009400.D
 Acq On : 25 Mar 2017 03:13
 Operator : SJ/MA
 Sample : I2358-10
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BK5

Manual Integrations
 APPROVED

mohammad
 3/27/2017 1:18:00 PM

Quant Time: Apr 01 04:41:11 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM032317MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 28 06:57:41 2017
 Response via : Initial Calibration



(1) 1,4-Dichlorobenzene-d4 (I)

7.928min (+0.076) 20.00ng/ul

response 215920

Ion	Exp%	Act%
152.00	100	100
115.00	63.00	62.15
0.00	0.00	0.00
0.00	0.00	0.00

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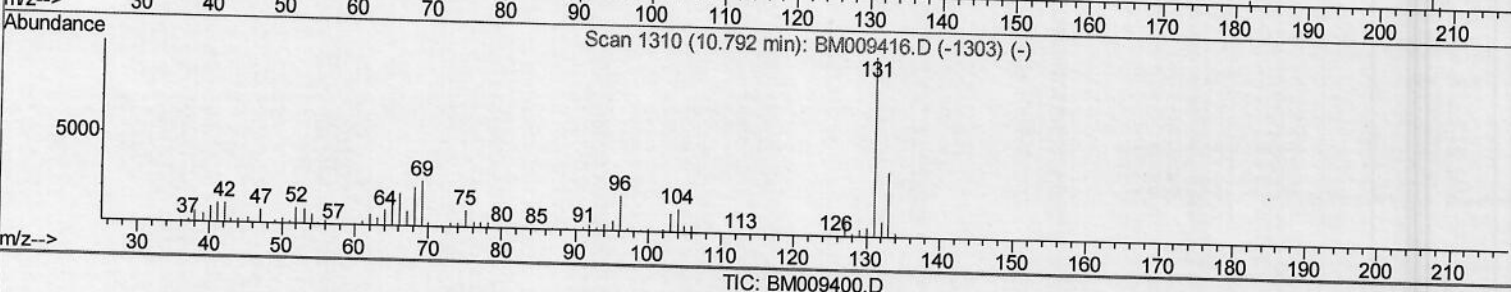
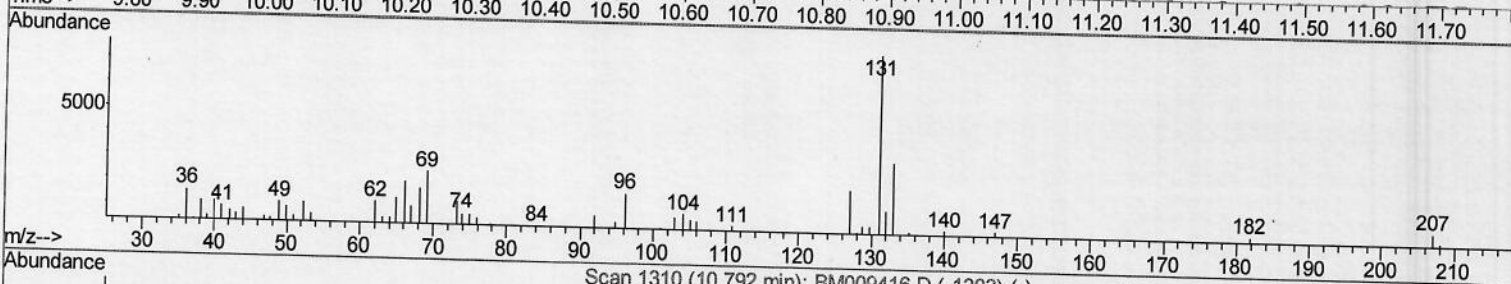
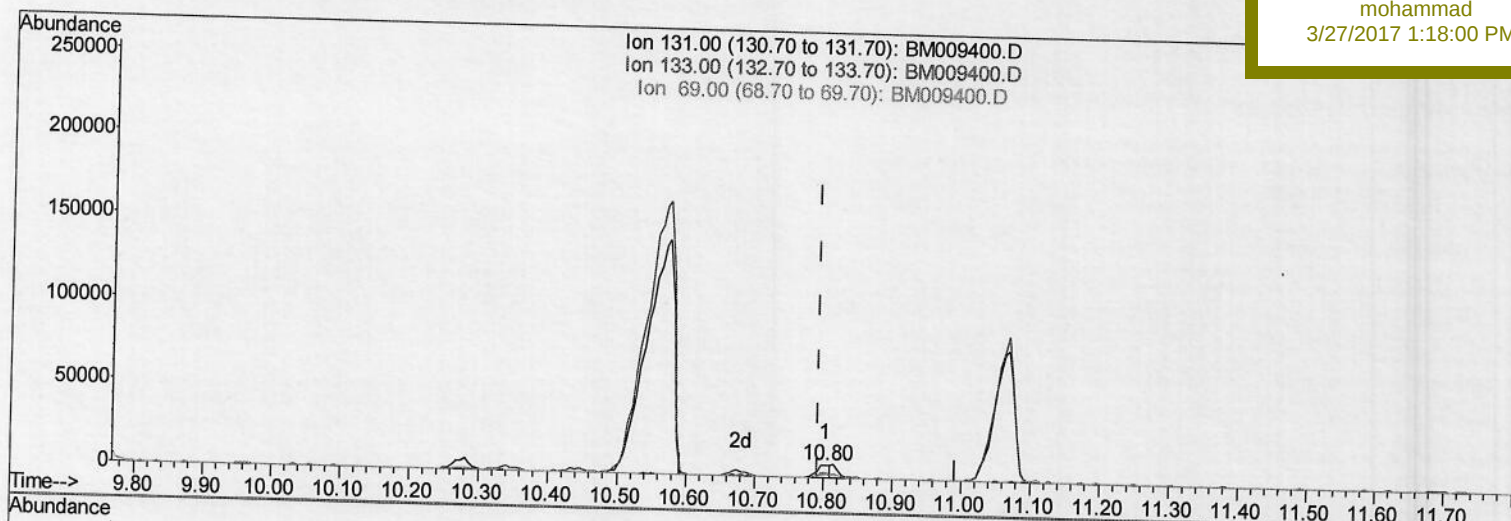
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(29) 4-Chloroaniline-d4 (S)
 10.804min (+0.012) 0.89ng/ul
 response 10347

Ion	Exp%	Act%
131.00	100	100
133.00	32.10	42.22#
69.00	16.20	32.41#
0.00	0.00	0.00

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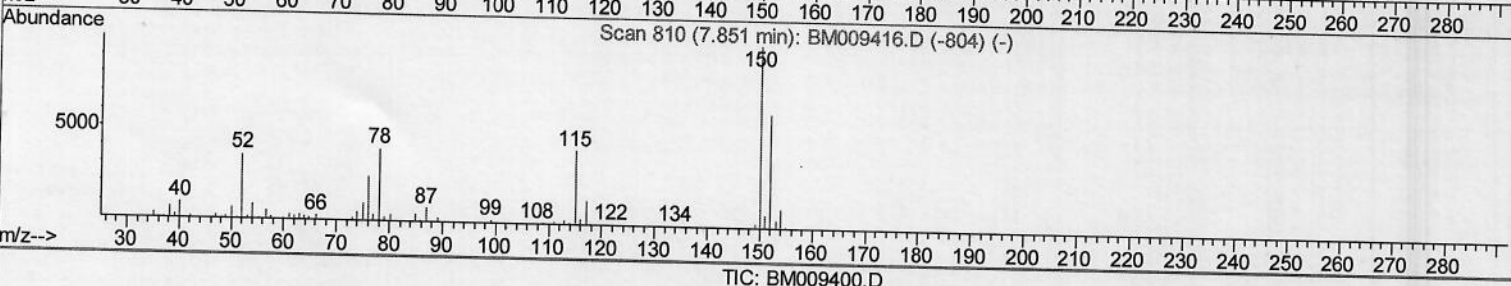
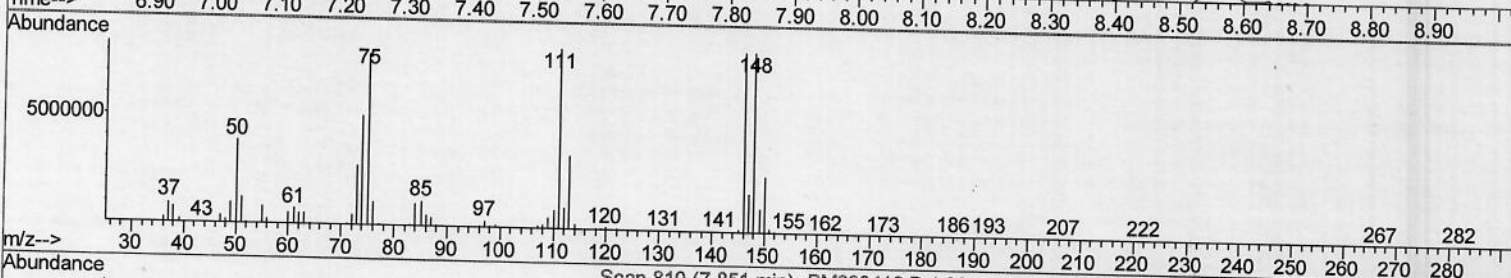
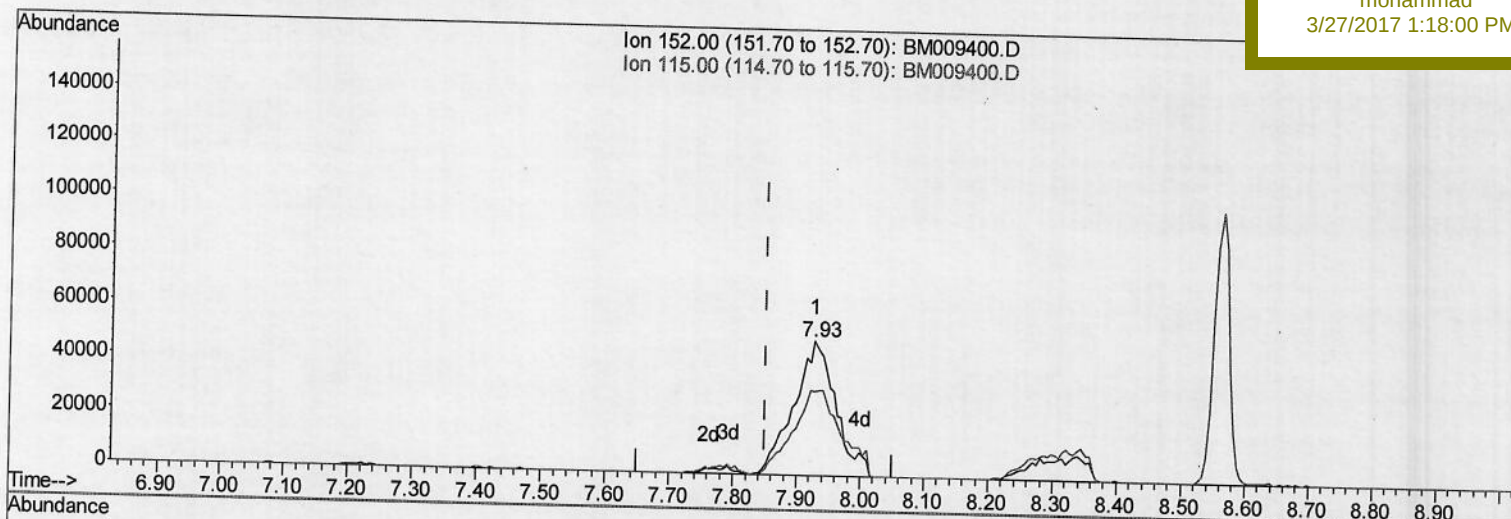
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Quant Time: Mar 27 10:50:37 2017
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TIC: BM009400.D

(1) 1,4-Dichlorobenzene-d4 (I)

7.928min (+0.077) 20.00ng/ul m

response 223818

Ion Exp% Act%

152.00 100 100

115.00 63.00 62.15

0.00 0.00 0.00

0.00 0.00 0.00

26.1
 03/31/17

Quantitation Report (Qedit)

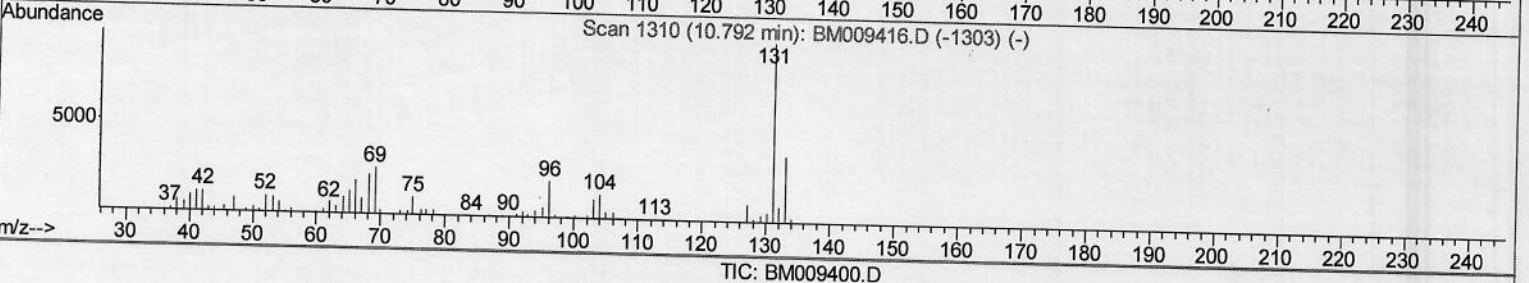
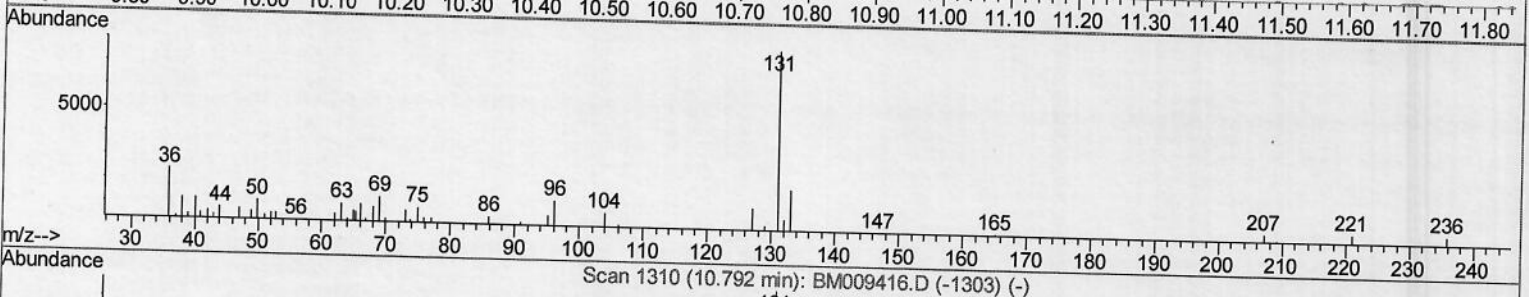
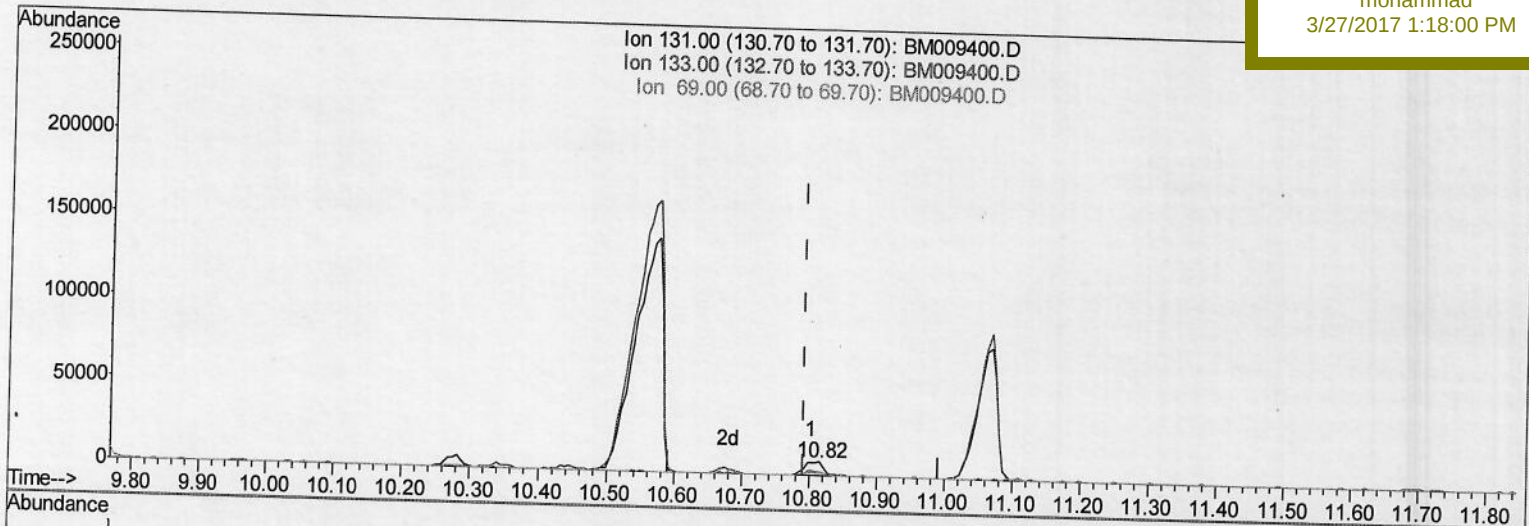
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(29) 4-Chloroaniline-d4 (S)

10.816min (+0.024) 1.36ng/ul m

response 15929

Ion	Exp%	Act%
131.00	100	100
133.00	32.10	26.01
69.00	16.20	17.27
0.00	0.00	0.00

5.1
 03/31/17

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	223818m	20.00	ng/ul	0.08
18) Naphthalene-d8	10.67	136	645948	20.00	ng/ul	0.02
35) Acenaphthene-d10	14.50	164	365197	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.24	188	804664	20.00	ng/ul	0.00
77) Chrysene-d12	21.42	240	1038125	20.00	ng/ul	0.00
85) Perylene-d12	23.74	264	964140	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	17448	2.97	ng/uL	0.00
5) Phenol-d5	7.02	99	83511	4.13	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.22	67	293488	20.34	ng/ul	0.03
9) 2-Chlorophenol-d4	7.40	132	231916	17.02	ng/ul	0.02
13) 4-Methylphenol-d8	8.56	113	171321	11.34	ng/ul	0.01
19) Nitrobenzene-d5	9.03	128	155640	32.78	ng/ul	0.01
22) 2-Nitrophenol-d4	9.75	143	173948	33.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.27	165	299405	29.56	ng/ul	0.00
29) 4-Chloroaniline-d4	10.82	131	15929m	1.36	ng/ul	0.02
43) Dimethylphthalate-d6	13.90	166	1180483	45.96	ng/ul	0.00
46) Acenaphthylene-d8	14.19	160	1359225	42.19	ng/ul	0.00
51) 4-Nitrophenol-d4	14.69	143	41258	8.71	ng/ul	0.00
57) Fluorene-d10	15.49	176	1000700	42.47	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.61	200	185146	36.40	ng/ul	0.00
70) Anthracene-d10	17.34	188	1631867	42.52	ng/ul	0.00
78) Pyrene-d10	19.63	212	1929243	43.84	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.60	264	1913871	43.10	ng/ul	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.04	94	30746	1.42	ng/ul#	72
10) 2-Chlorophenol	7.43	128	76780	5.29	ng/ul#	77
17) Hexachloroethane	8.95	117	9076	1.30	ng/ul#	80
20) Nitrobenzene	9.07	77	2041405	127.52	ng/ul#	82
23) 2-Nitrophenol	9.77	139	47506	8.72	ng/ul#	67
36) 1,2,4,5-Tetrachlorobenzene	12.68	216	413862	33.73	ng/ul	98
52) 4-Nitrophenol	14.70	109	16596	2.99	ng/ul#	74
72) 1,2,3,4-Tetrachlorobenzene	13.29	216	1524622	128.88	ng/uL	99

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

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