

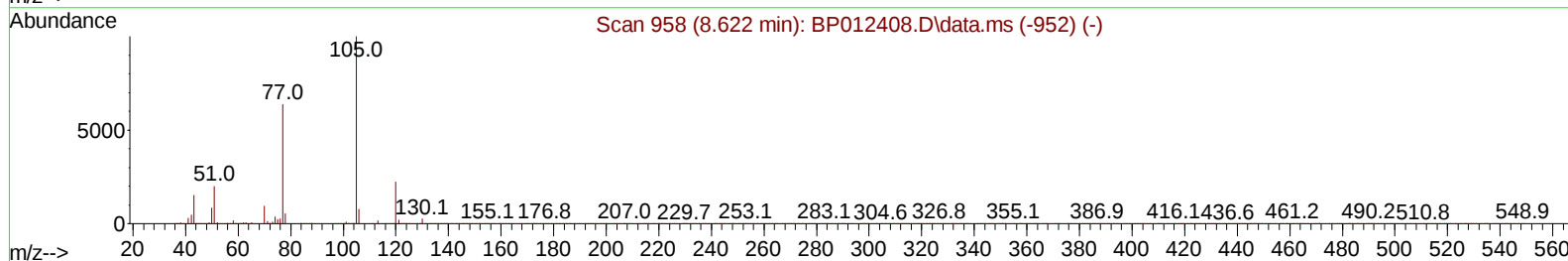
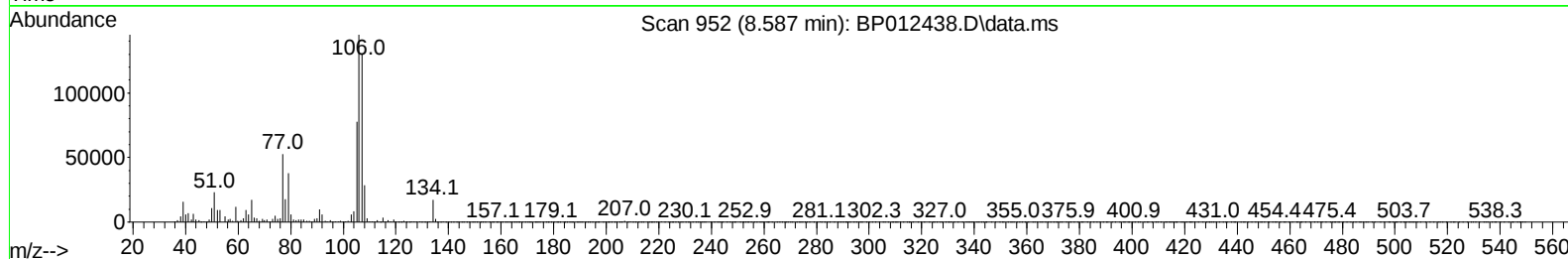
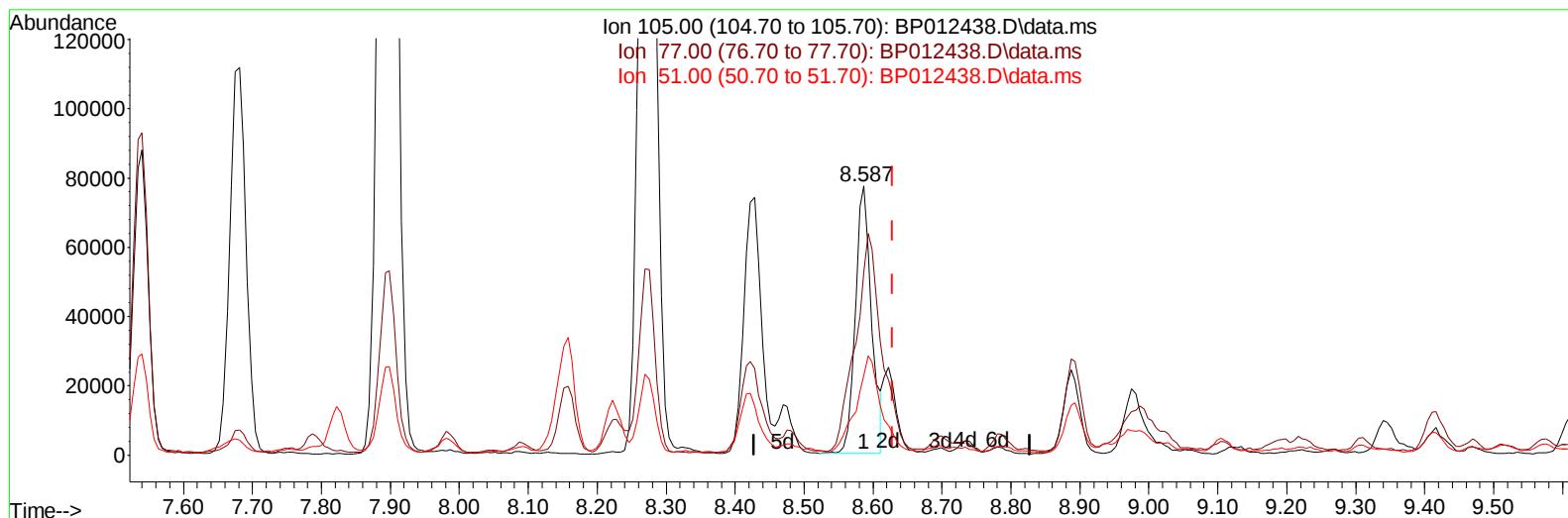
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012438.D
 Acq On : 01 Nov 2022 21:19
 Operator : CG/JU
 Sample : N5300-03
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA80

Manual IntegrationsAPPROVED

Quant Time: Nov 01 23:01:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/02/2022
 Supervised By :mohammad ahmed 11/02/2022



TIC: BP012438.D\data.ms

(16) Acetophenone

8.587min (-0.041) 4.19 ng/u1

response 125485

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	75.40	67.93
51.00	24.90	29.61
0.00	0.00	0.00

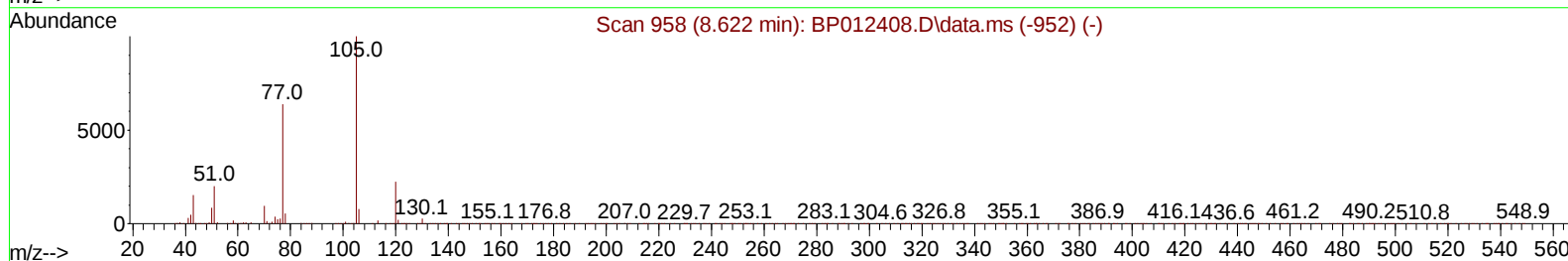
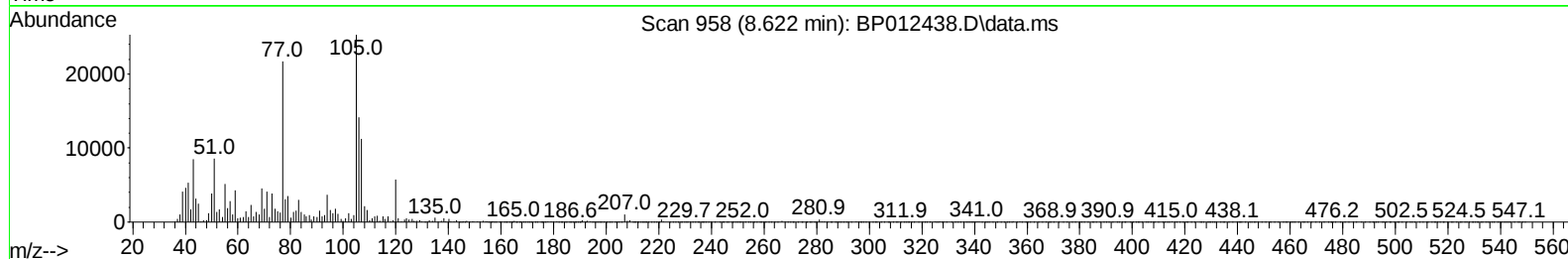
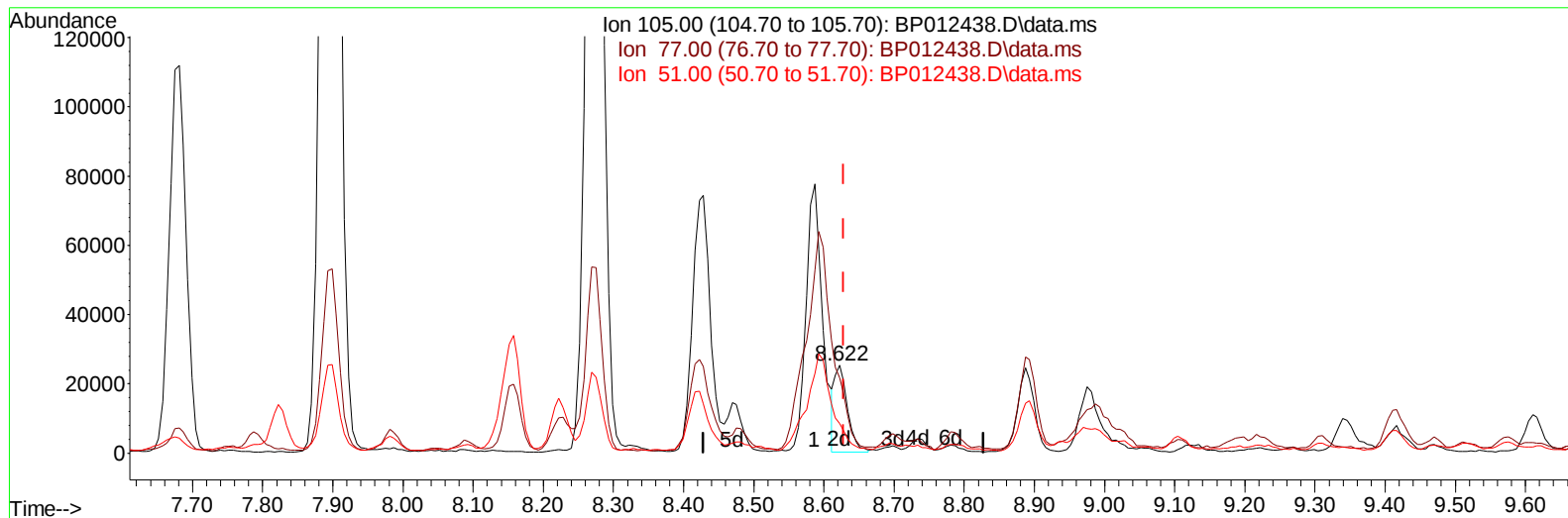
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TIC: BP012438.D\data.ms

(16) Acetophenone

8.622min (-0.006) 1.09 ng/u1 m

response 32533

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	75.40	85.90
51.00	24.90	33.99#
0.00	0.00	0.00

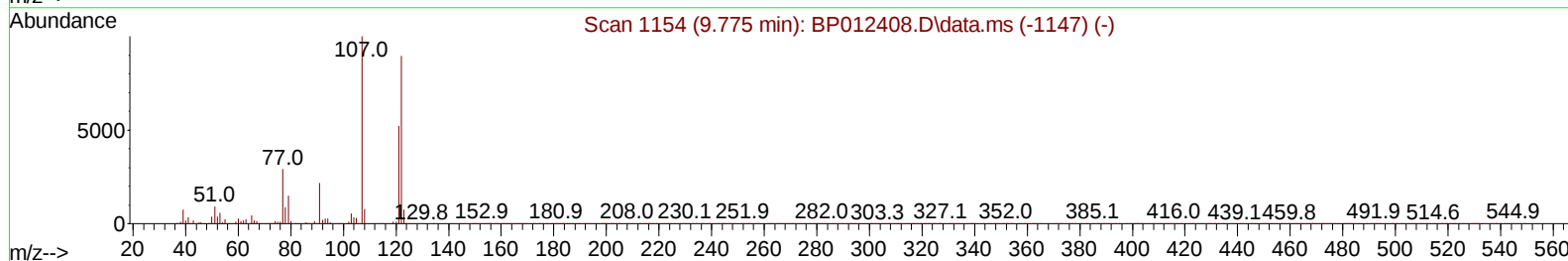
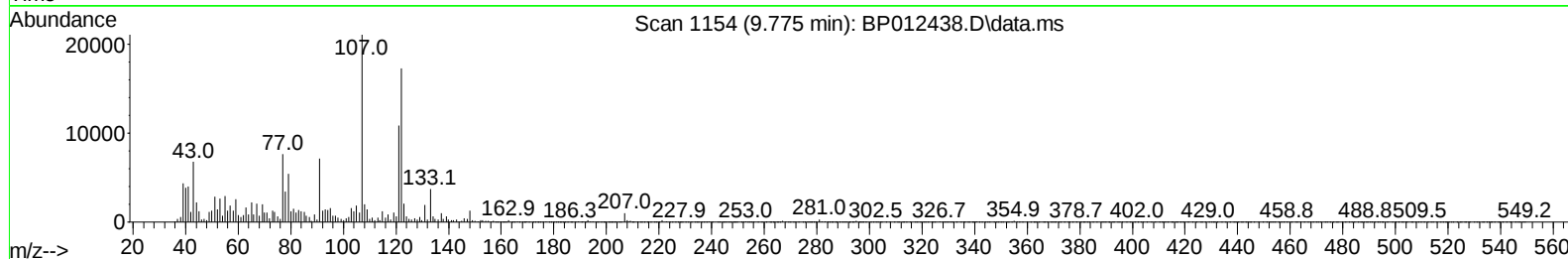
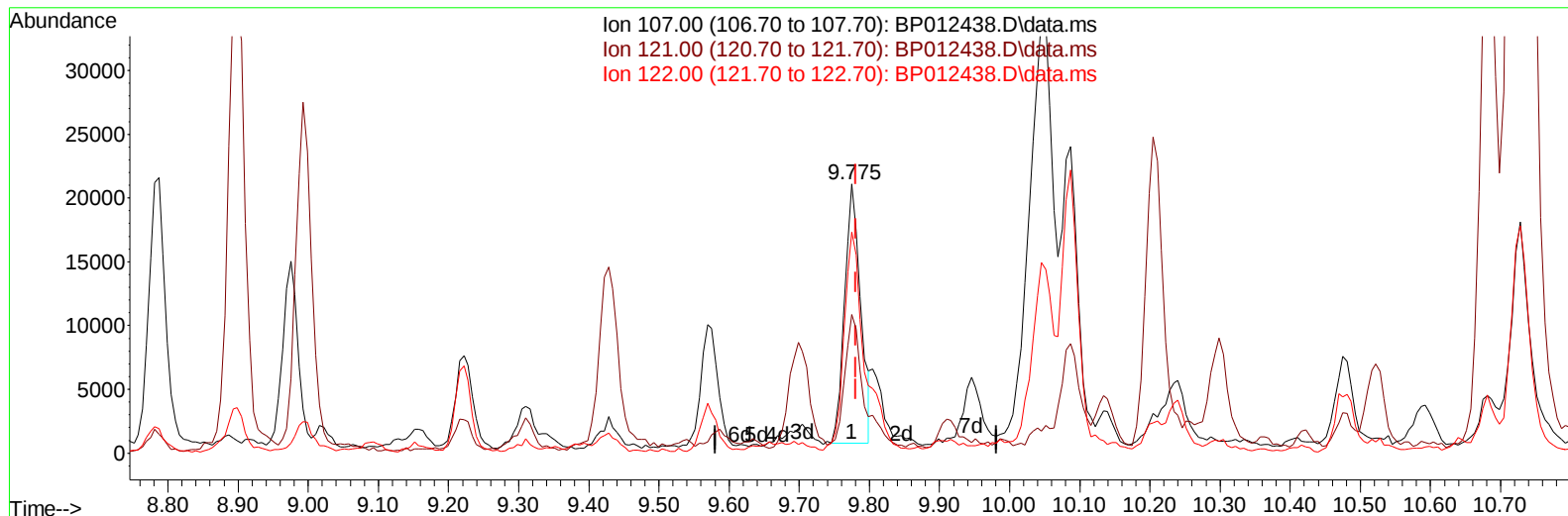
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 Sample : N5300-03
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA80

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TIC: BP012438.D\data.ms

(26) 2,4-Dimethylphenol

9.775min (-0.006) 1.51 ng/u1

response 32446

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	51.10	51.58
122.00	88.20	82.06
0.00	0.00	0.00

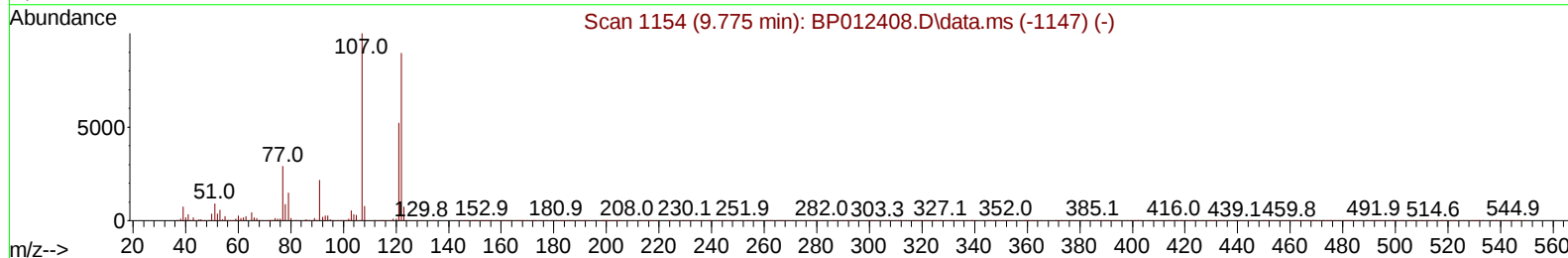
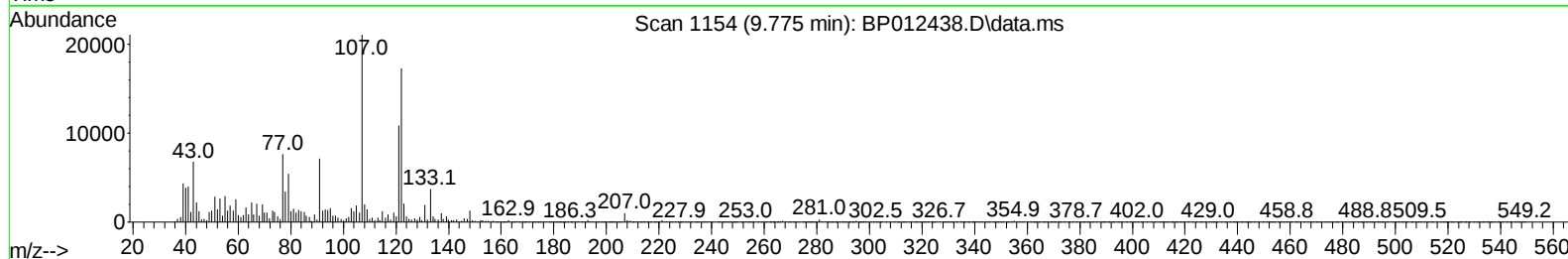
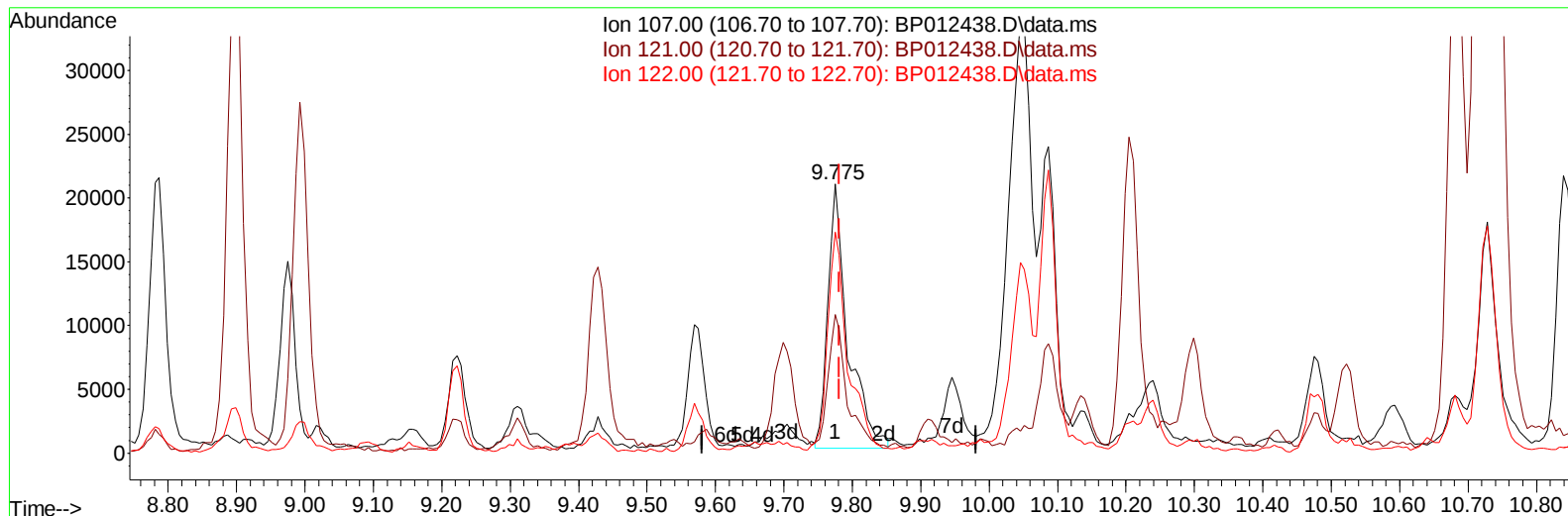
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 Sample : N5300-03
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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TIC: BP012438.D\data.ms

(26) 2,4-Dimethylphenol

9.775min (-0.006) 1.96 ng/ul m

response 42052

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	51.10	51.58
122.00	88.20	82.06
0.00	0.00	0.00

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 Sample : N5300-03
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	272096	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	1212926	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	776069	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1665620	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1761554	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1704730	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	22771	3.370	ng/uL	0.00
4) Pyridine-d5	3.669	84	86982	4.481	ng/ul	0.01
7) Phenol-d5	6.969	99	172729	7.078	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	413811	28.404	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	454042	25.476	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	331383	17.168	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	299074	38.023	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	328803	40.275	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	614811	34.749	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	470826	17.872	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2040287	38.405	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	2284447	37.628	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	73740	7.334	ng/ul	0.02
60) Fluorene-d10	15.439	176	1650645	36.290	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	36894	4.254	ng/ul	0.00
73) Anthracene-d10	17.304	188	2779733	38.732	ng/ul	0.00
81) Pyrene-d10	19.545	212	3251278	36.762	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.533	264	3254554	39.072	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.269	88	51467	7.164	ng/uL	98
8) Phenol	6.999	94	33980	1.339	ng/ul#	84
10) Bis(2-Chloroethyl)ether	7.222	93	27276	1.339	ng/ul	98
16) Acetophenone	8.622	105	32533m	1.085	ng/ul	
18) 4-Methylphenol	8.569	108	133002	6.236	ng/ul	84
26) 2,4-Dimethylphenol	9.775	107	42052m	1.956	ng/ul	
30) Naphthalene	10.640	128	377508	5.873	ng/ul	99
36) 2-Methylnaphthalene	12.257	142	64613	1.466	ng/ul	95
37) 1-Methylnaphthalene	12.475	142	64972	1.476	ng/ul#	94
52) Acenaphthene	14.504	153	53260	1.113	ng/ul	98
56) Dibenzofuran	14.839	168	85216	1.305	ng/ul	89
86) Bis(2-ethylhexyl)phtha...	21.204	149	220803	3.424	ng/ul	99

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

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