

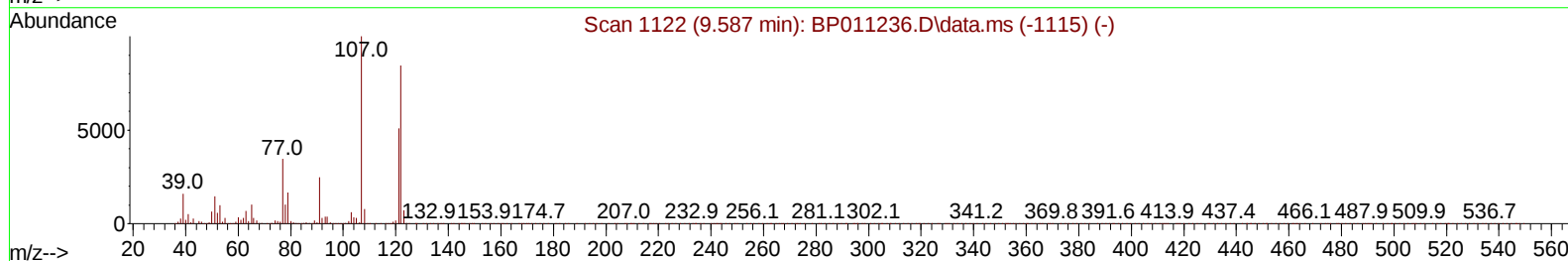
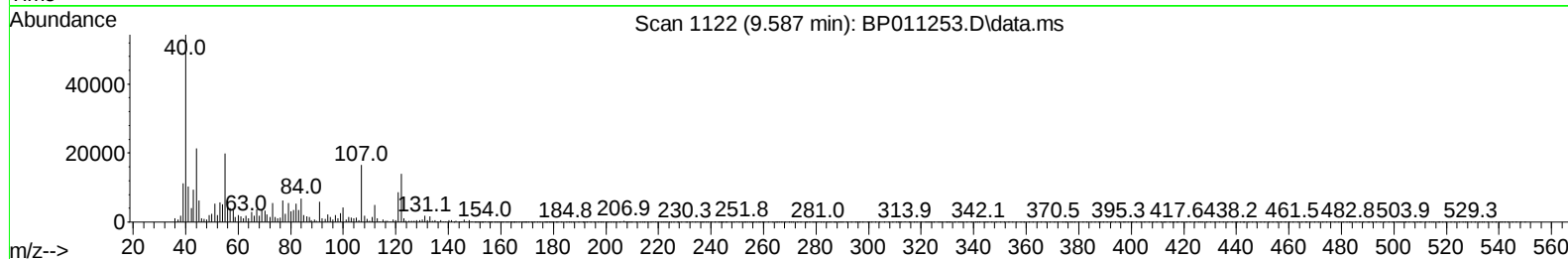
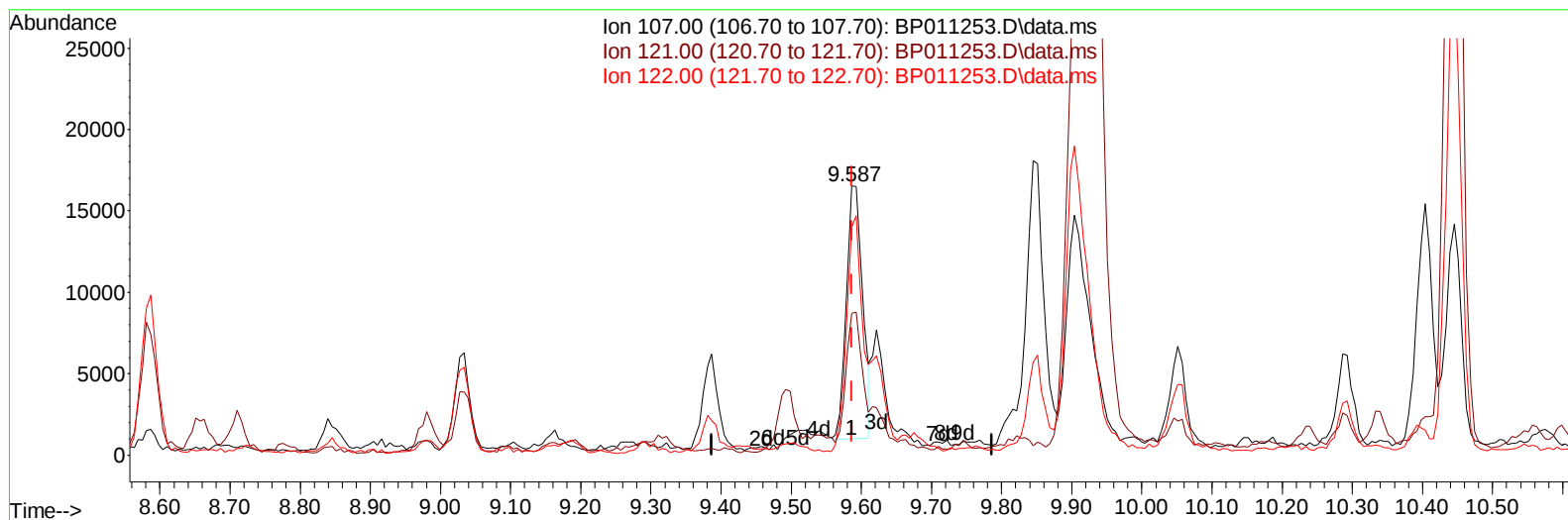
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080322\
 Data File : BP011253.D
 Acq On : 03 Aug 2022 19:51
 Operator : CG/JU
 Sample : N3956-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXF02

Manual IntegrationsAPPROVED

Quant Time: Aug 03 23:01:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/04/2022
 Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011253.D\data.ms

(26) 2,4-Dimethylphenol

9.587min (-0.000) 1.42 ng/u1

response 25301

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	50.40	52.63
122.00	84.60	84.77
0.00	0.00	0.00

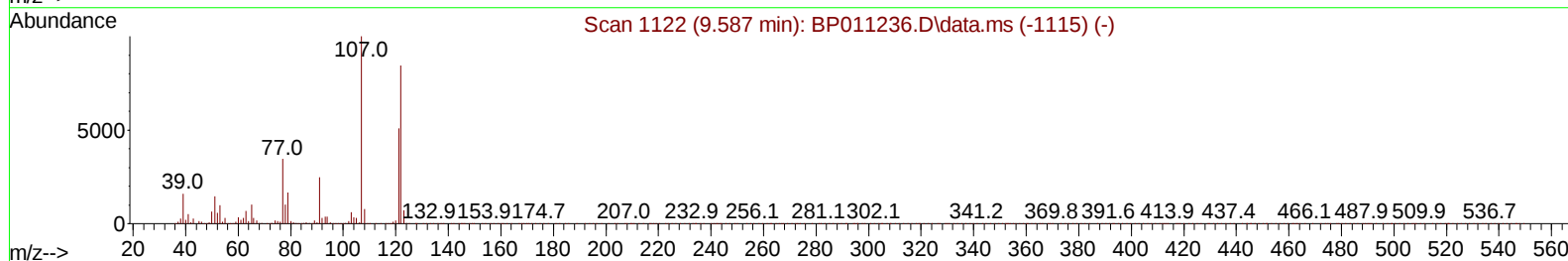
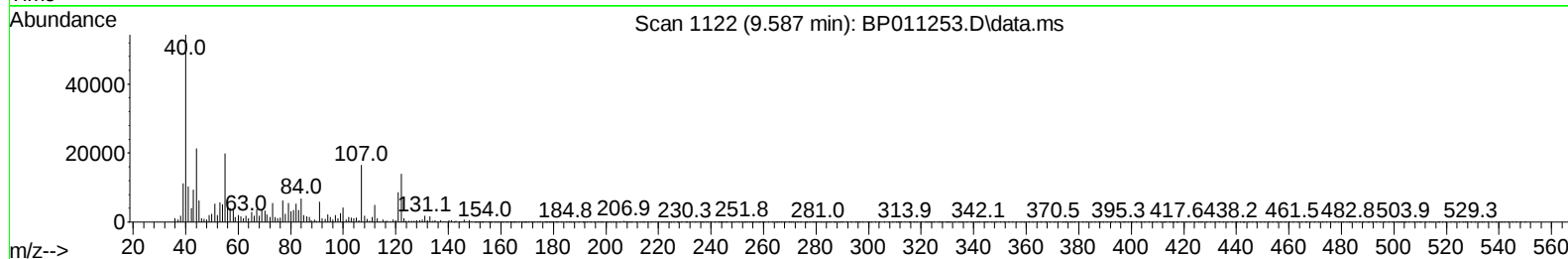
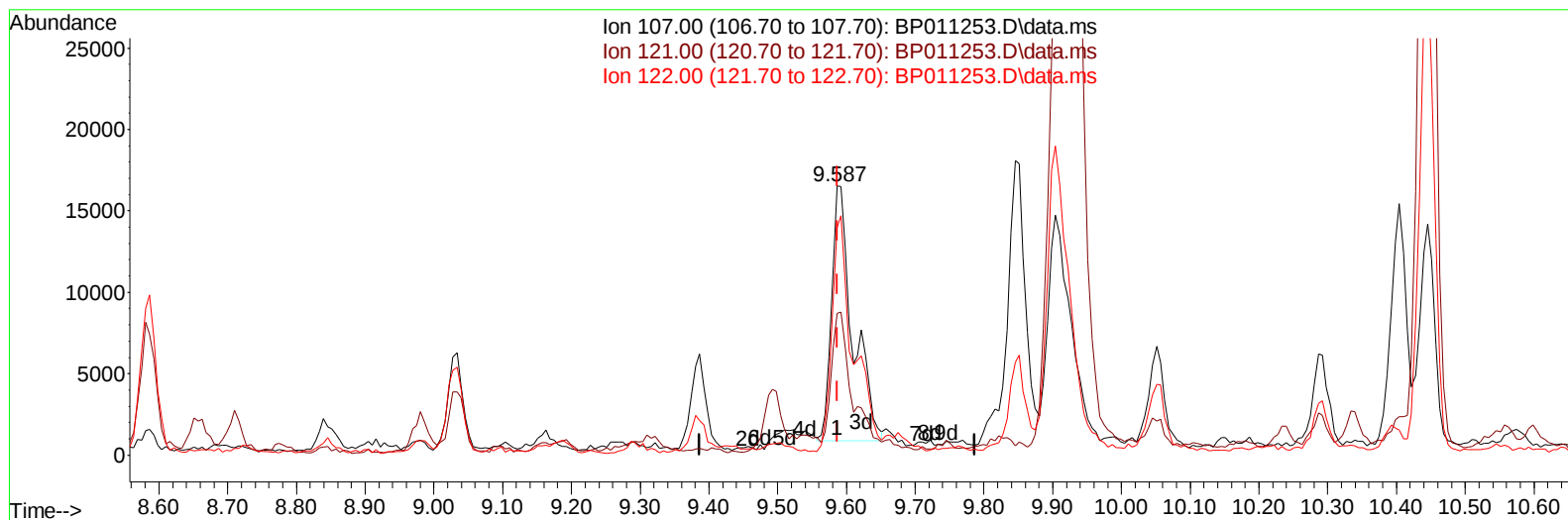
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TIC: BP011253.D\data.ms

(26) 2,4-Dimethylphenol

9.587min (-0.000) 1.89 ng/ul m

response 33764

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	50.40	52.63
122.00	84.60	84.77
0.00	0.00	0.00

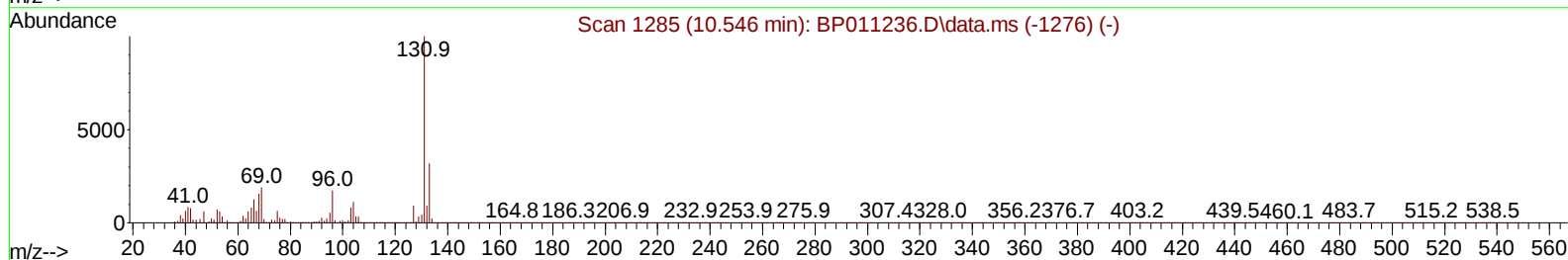
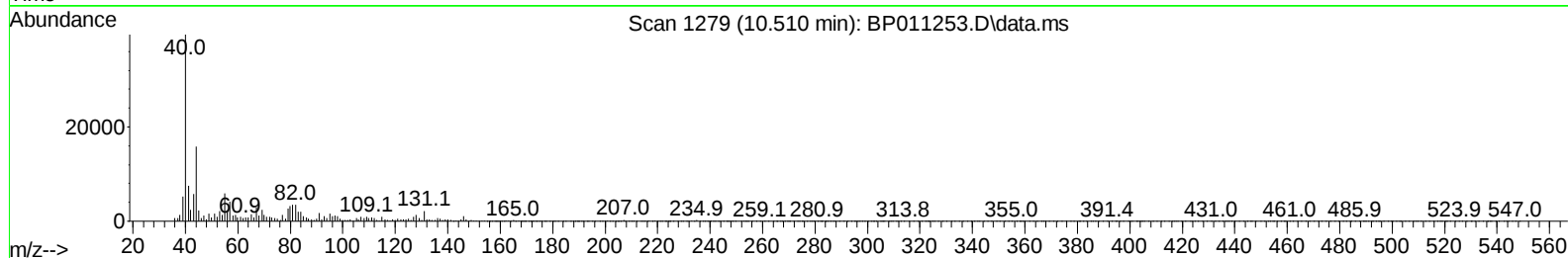
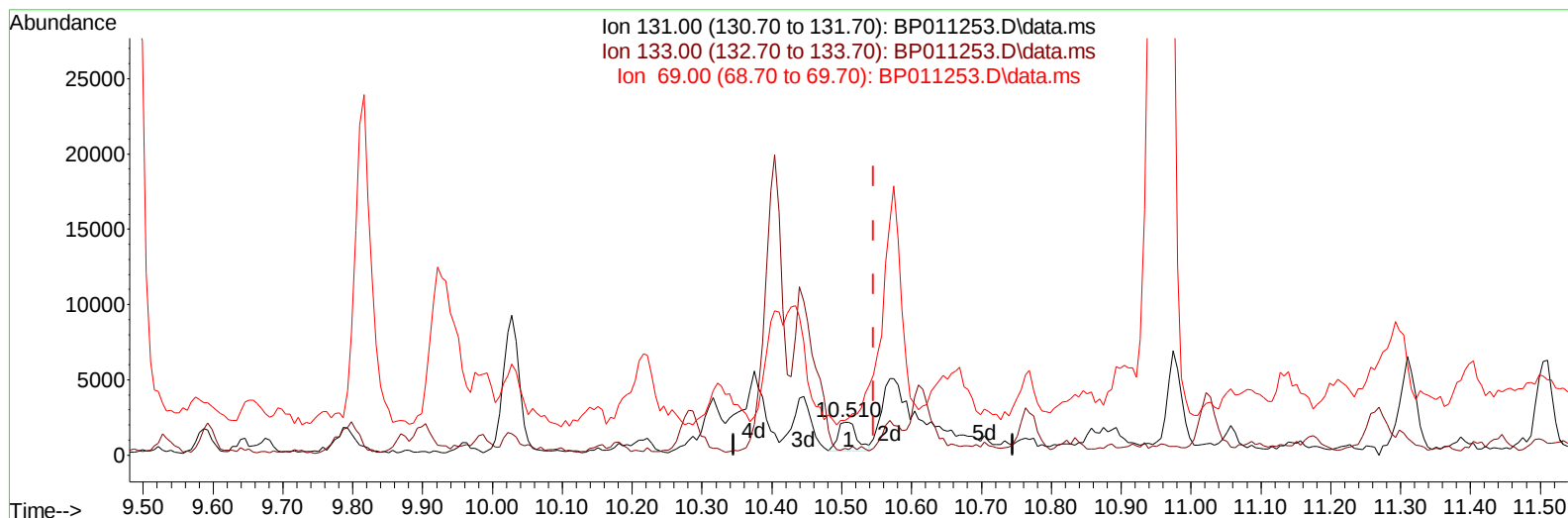
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TIC: BP011253.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.510min (-0.036) 0.18 ng/u1

response 3720

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.00	15.83#
69.00	19.00	113.21#
0.00	0.00	0.00

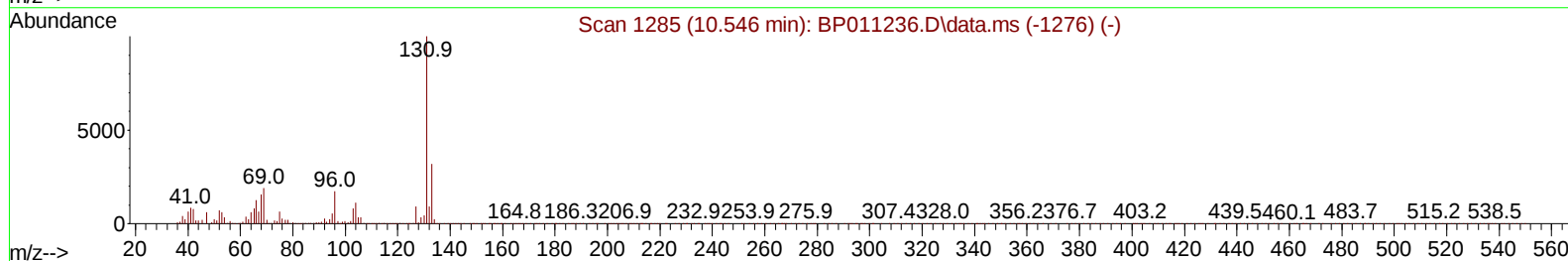
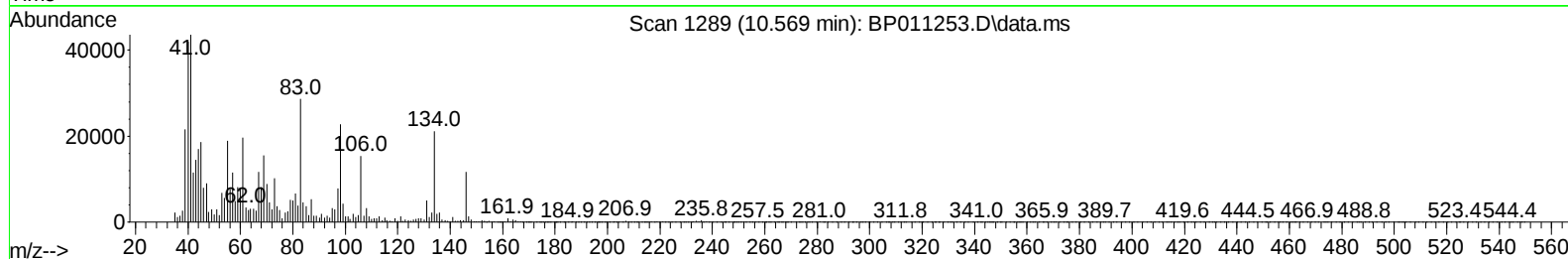
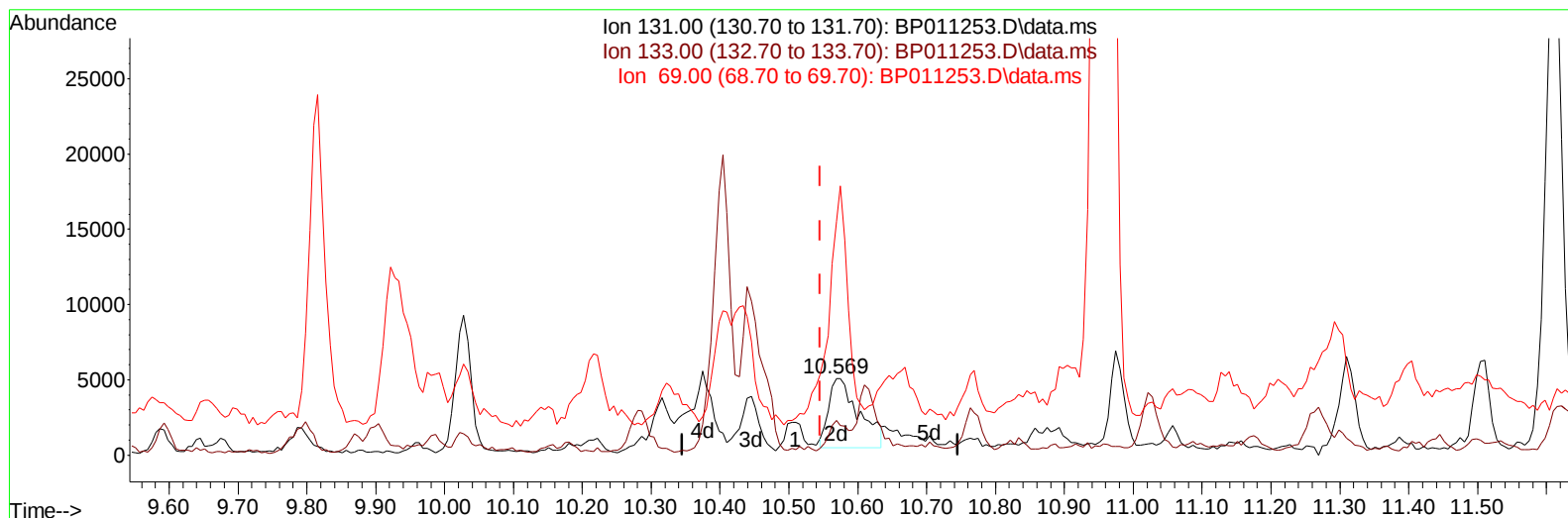
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Instrument :
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TIC: BP011253.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.569min (+ 0.023) 0.70 ng/ul m

response 14160

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.00	44.95#
69.00	19.00	304.16#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.598	152	226955	20.000	ng/ul	0.00
20) Naphthalene-d8	10.392	136	1011044	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	518418	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.045	188	1017649	20.000	ng/ul	0.00
79) Chrysene-d12	21.174	240	976154	20.000	ng/ul	0.00
88) Perylene-d12	23.492	264	1013570	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	7104	2.531	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.793	99	119893	6.339	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.945	67	353110	25.873	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	335249	21.746	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	216062	13.883	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	218328	30.022	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	202205	29.663	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	348049	26.775	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	14160m	0.704	ng/ul	0.02
46) Dimethylphthalate-d6	13.698	166	1124950	31.683	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	1407393	31.504	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	30977	5.130	ng/ul	0.01
60) Fluorene-d10	15.280	176	951418	31.491	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	139610	29.545	ng/ul	0.00
73) Anthracene-d10	17.145	188	1408355	31.489	ng/ul	0.00
81) Pyrene-d10	19.410	212	1545675	31.318	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.345	264	1582392	31.330	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	6.816	94	368758	18.182	ng/ul	99
12) 2-Chlorophenol	7.169	128	227190	14.188	ng/ul#	57
13) 2-Methylphenol	8.063	108	55511	3.617	ng/ul	99
18) 4-Methylphenol	8.392	108	39425	2.396	ng/ul	99
26) 2,4-Dimethylphenol	9.587	107	33764m	1.892	ng/ul	
30) Naphthalene	10.445	128	166217	3.114	ng/ul	99
36) 2-Methylnaphthalene	12.069	142	60909	1.826	ng/ul	98
37) 1-Methylnaphthalene	12.292	142	129273	3.859	ng/ul#	98
52) Acenaphthene	14.339	153	67761	2.075	ng/ul	97
61) Fluorene	15.333	166	49775	1.398	ng/ul#	95
77) Carbazole	17.463	167	53599	1.079	ng/ul#	92

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

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