

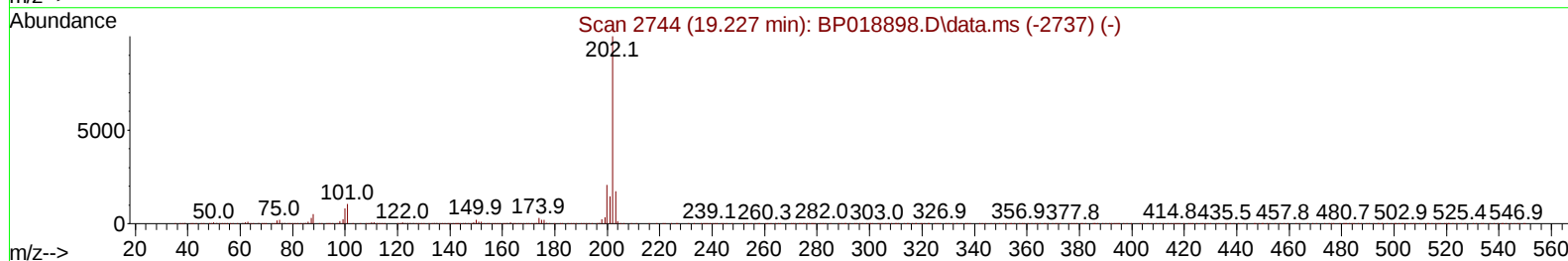
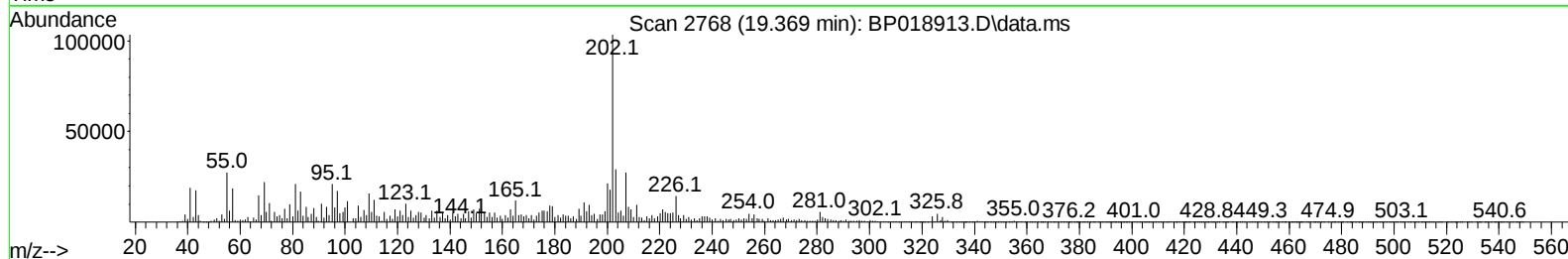
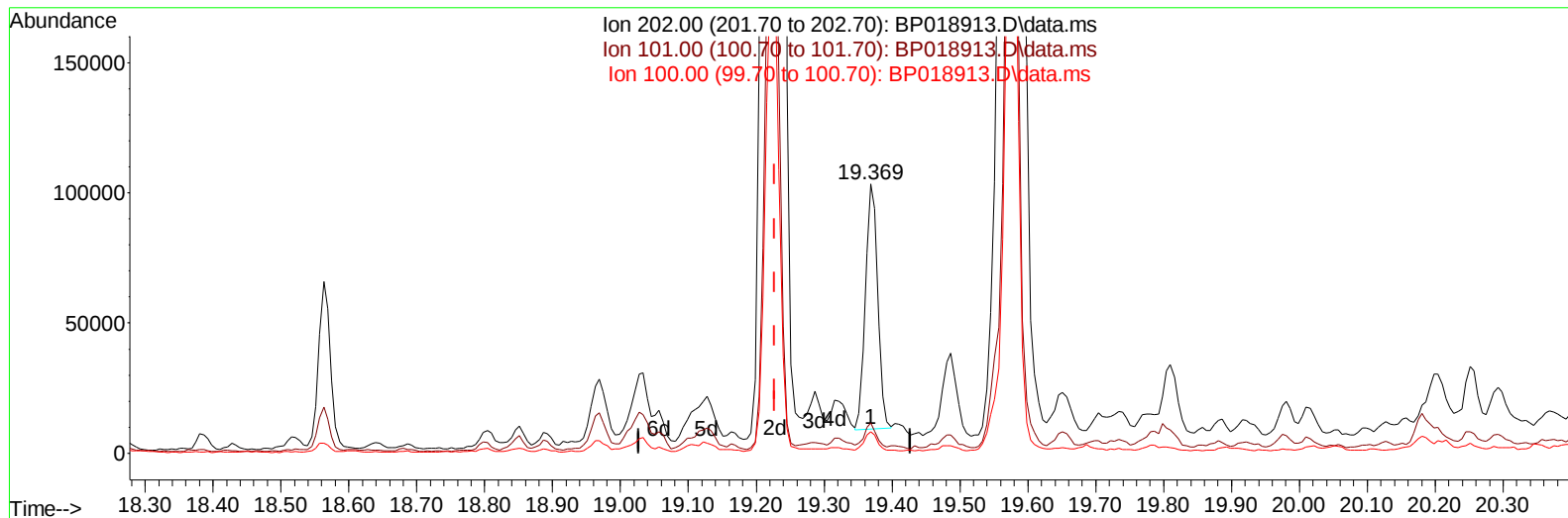
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
 Data File : BP018913.D
 Acq On : 18 Dec 2023 18:43
 Operator : MA/JU
 Sample : 05794-22 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGRG3

Manual IntegrationsAPPROVED

Quant Time: Dec 18 21:42:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 13 23:34:47 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/19/2023
 Supervised By :mohammad ahmed 12/20/2023



TIC: BP018913.D\data.ms

(80) Fluoranthene

19.369min (+ 0.141) 1.22 ng/ul

response 119100

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.40	11.18
100.00	8.90	7.83
0.00	0.00	0.00

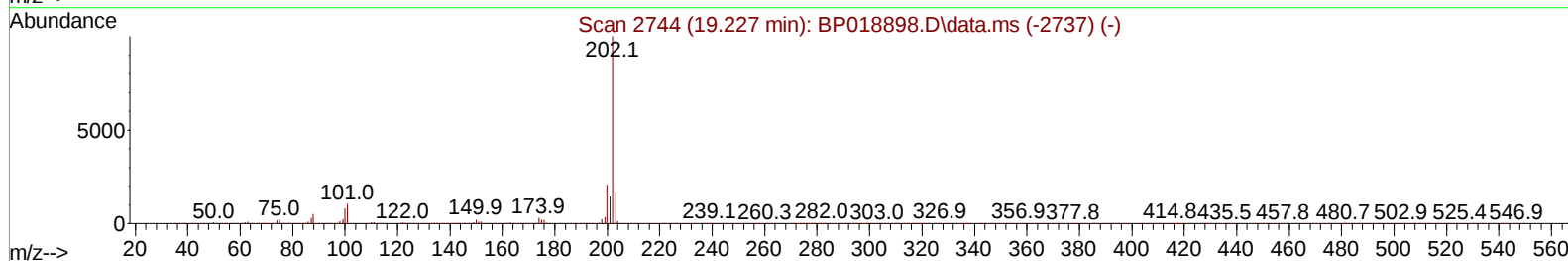
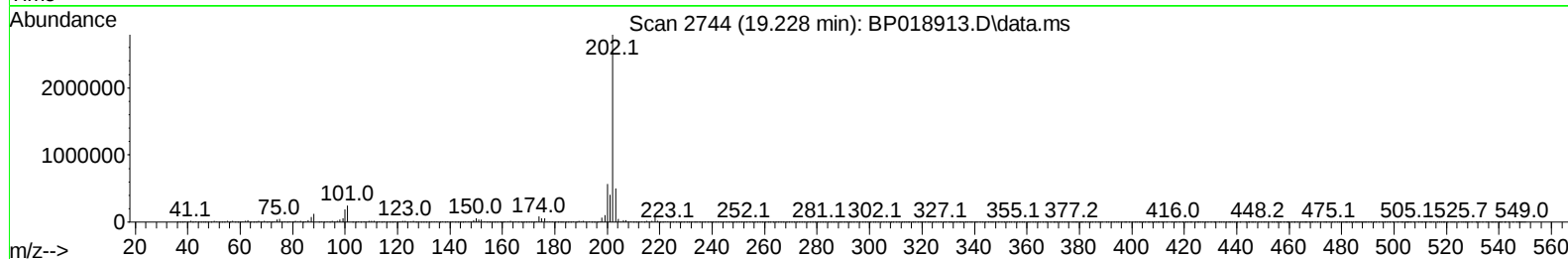
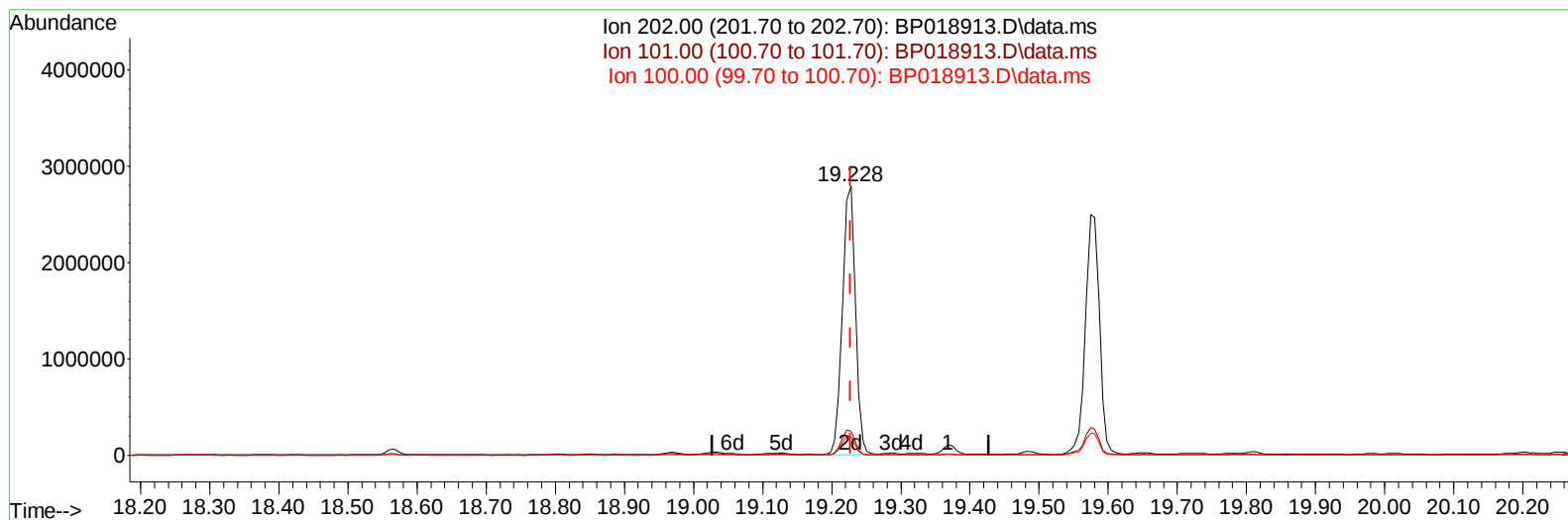
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(80) Fluoranthene

19.228min (+ 0.000) 37.55 ng/ul m

response 3668464

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.40	8.97#
100.00	8.90	6.86#
0.00	0.00	0.00

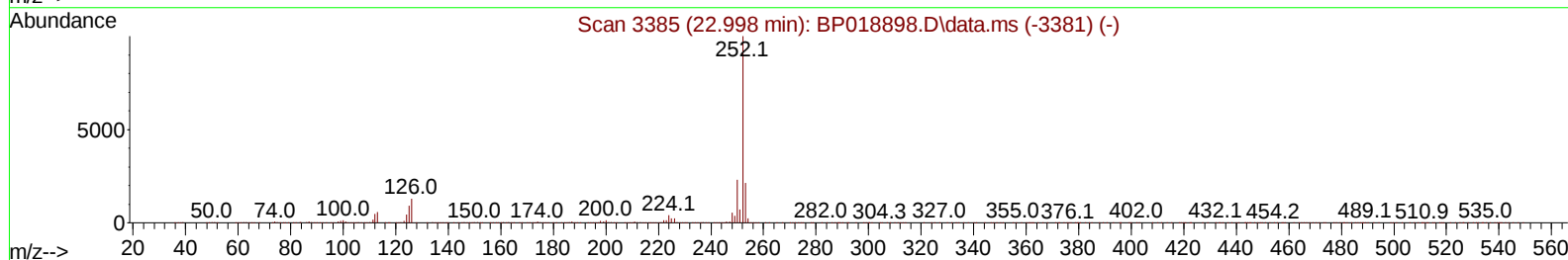
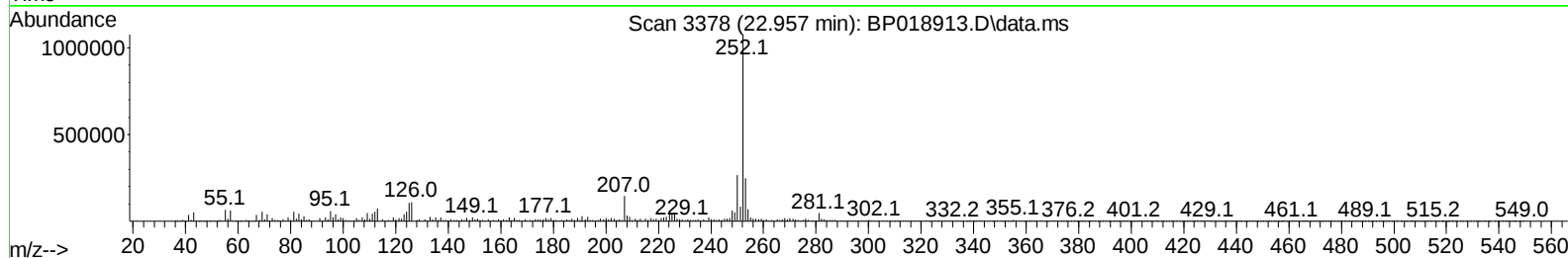
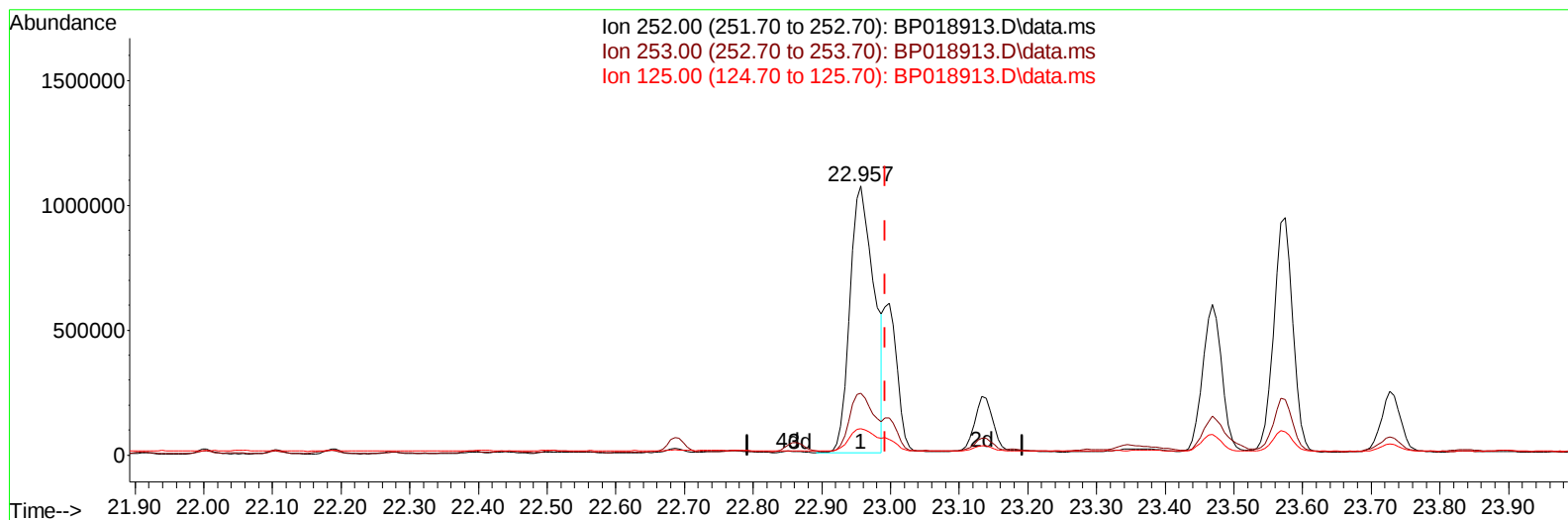
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TIC: BP018913.D\data.ms

(91) Benzo(k)fluoranthene

22.957min (-0.035) 28.13 ng/u1

response 2623597

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.10	23.05
125.00	10.00	9.83
0.00	0.00	0.00

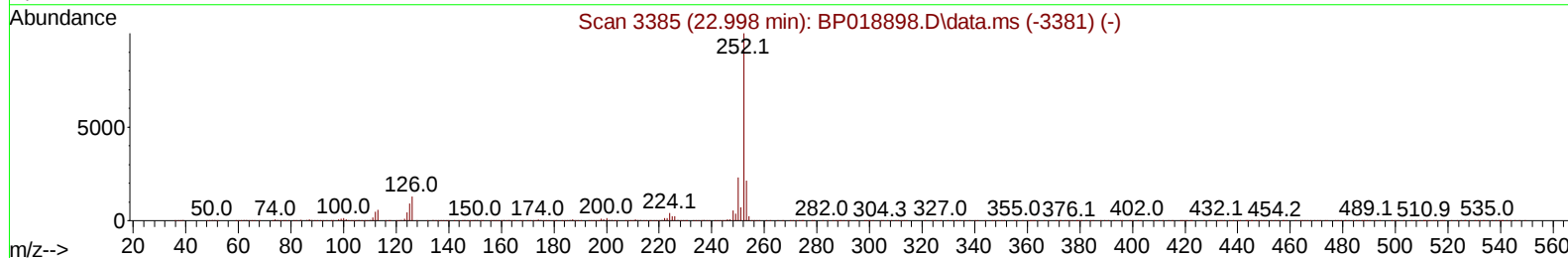
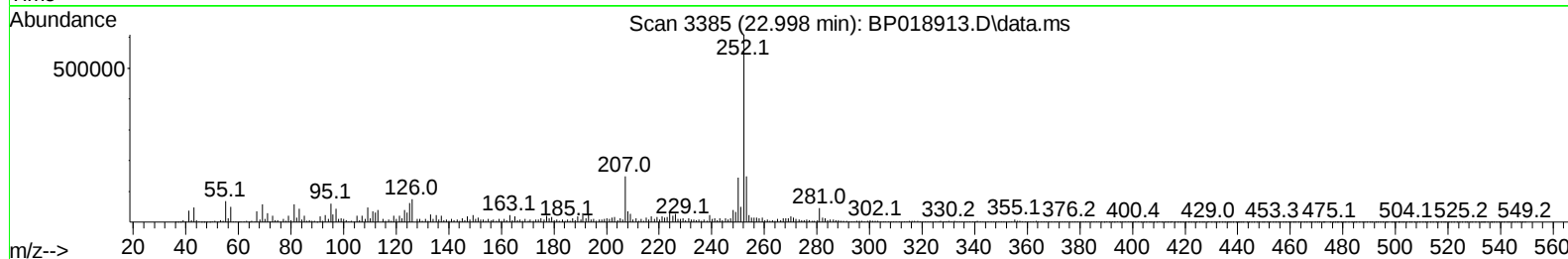
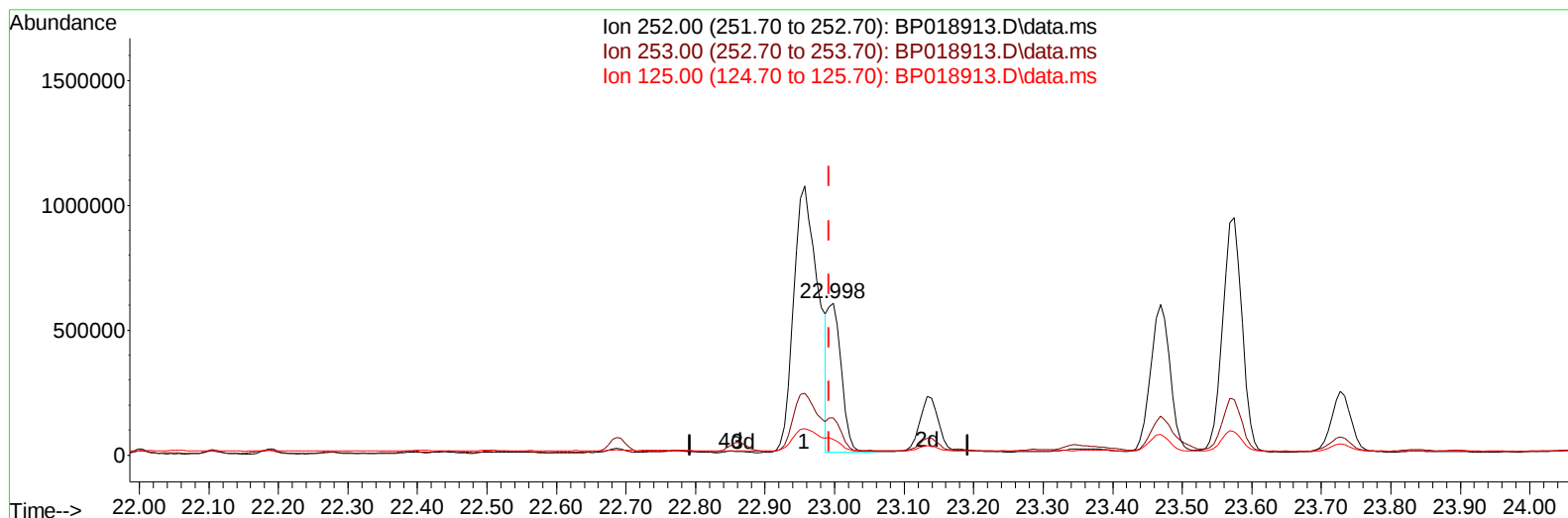
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TIC: BP018913.D\data.ms

(91) Benzo(k)fluoranthene

22.998min (+ 0.006) 8.64 ng/ul m

response 806091

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.10	24.40
125.00	10.00	10.30
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	7.828	152	208379	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	776440	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	491669	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	1121809	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1061724	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1307355	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	7324	1.198	ng/uL	0.00
4) Pyridine-d5	3.687	84	14600	0.891	ng/ul	0.01
7) Phenol-d5	6.999	99	128496	6.145	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.164	67	73986	5.969	ng/ul	0.00
11) 2-Chlorophenol-d4	7.358	132	102565	6.286	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	99843	5.900	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	46971	6.797	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	49624	6.926	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	101117	6.902	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	63866	3.336	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	311870	7.107	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	379010	7.844	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	18056	2.507	ng/ul	0.00
60) Fluorene-d10	15.451	176	303533	8.240	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	12602	1.874	ng/ul	0.00
73) Anthracene-d10	17.310	188	487499	8.266	ng/ul	0.00
81) Pyrene-d10	19.545	212	569105	6.886	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.522	264	575728	7.553	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.034	94	21091	1.028	ng/ul	92
16) Acetophenone	8.652	105	149825	5.940	ng/ul	100
30) Naphthalene	10.675	128	113807	2.416	ng/ul	100
36) 2-Methylnaphthalene	12.281	142	53165	1.593	ng/ul	99
37) 1-Methylnaphthalene	12.499	142	47143	1.399	ng/ul#	96
50) Acenaphthylene	14.181	152	206308	3.798	ng/ul	99
52) Acenaphthene	14.522	153	100818	2.807	ng/ul	97
56) Dibenzofuran	14.857	168	97866	1.960	ng/ul	96
61) Fluorene	15.510	166	164606	4.167	ng/ul	97
69) Hexachlorobenzene	16.510	284	22422	1.418	ng/ul	97
72) Phenanthrene	17.251	178	1885618	27.831	ng/ul	100
74) Anthracene	17.345	178	625111	9.205	ng/ul	99
77) Carbazole	17.616	167	186457	3.445	ng/ul	98
78) Di-n-butylphthalate	18.175	149	119913	1.664	ng/ul	98
80) Fluoranthene	19.228	202	3668464m	37.551	ng/ul	
82) Pyrene	19.575	202	3532357	35.842	ng/ul	94
85) Benzo(a)anthracene	21.286	228	2033275	23.594	ng/ul	97
86) Bis(2-ethylhexyl)phtha...	21.216	149	55507	1.119	ng/ul#	89
87) Chrysene	21.339	228	1984084	25.001	ng/ul	96
90) Benzo(b)fluoranthene	22.957	252	2623597	27.218	ng/ul	98
91) Benzo(k)fluoranthene	22.998	252	806091m	8.643	ng/ul	
93) Benzo(a)pyrene	23.574	252	1829198	20.768	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	26.145	276	1286110	12.221	ng/ul#	93

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
95) Dibenzo(a,h)anthracene	26.157	278		400209	4.572	ng/ul#	74
96) Benzo(g,h,i)perylene	26.910	276		409955	4.841	ng/ul#	92

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

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