

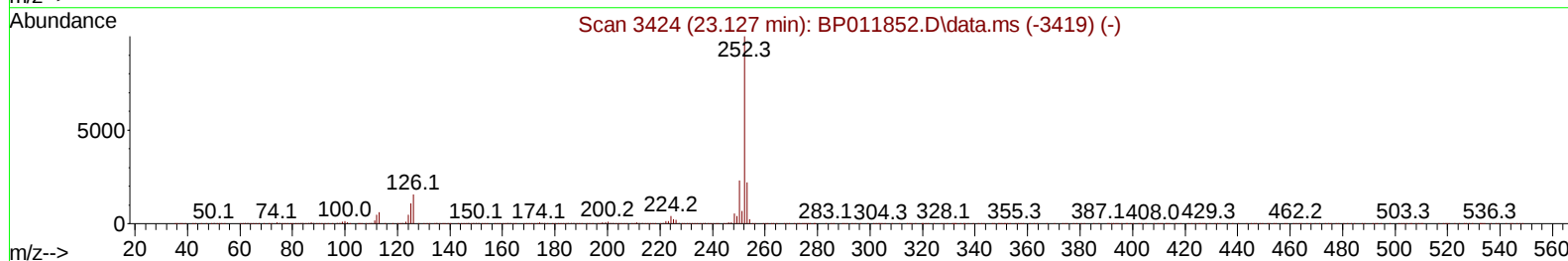
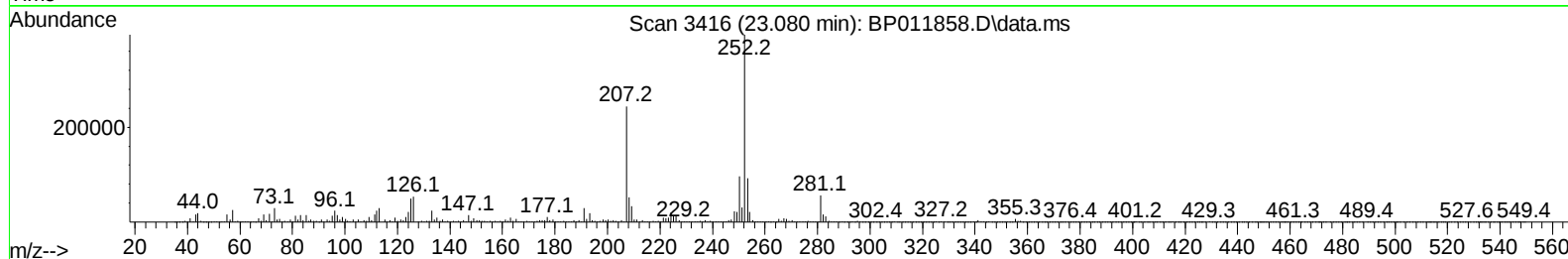
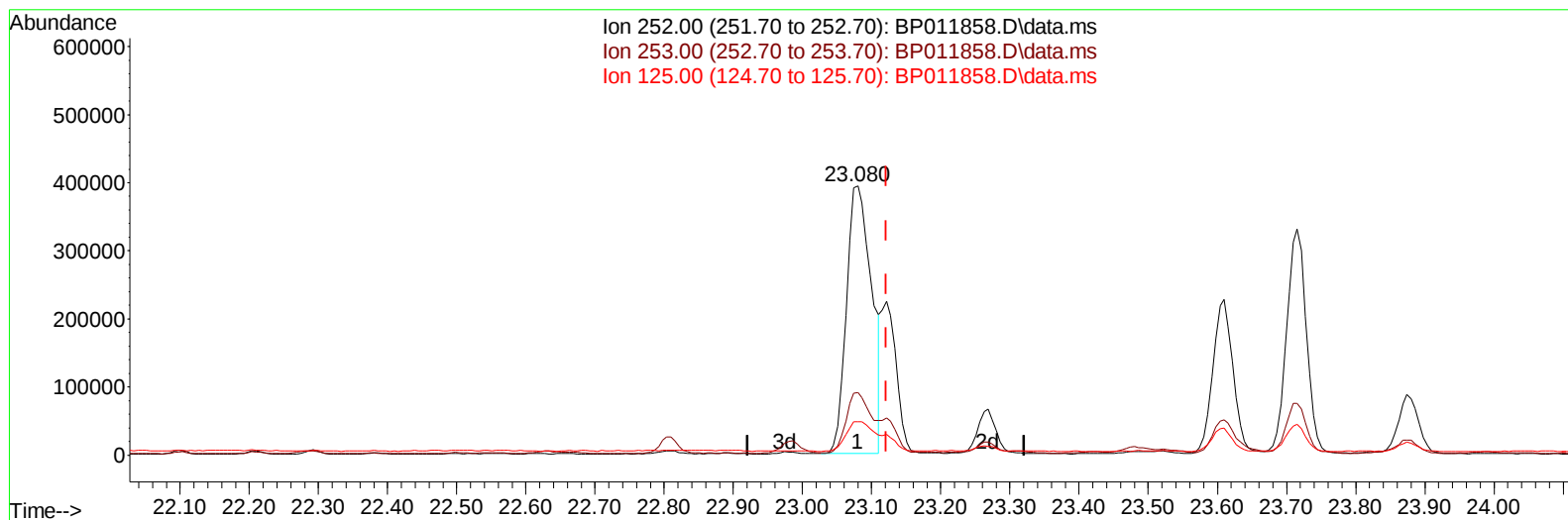
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100122\
 Data File : BP011858.D
 Acq On : 01 Oct 2022 13:01
 Operator : CG/JU
 Sample : N4745-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8N7

Manual IntegrationsAPPROVED

Quant Time: Oct 03 00:59:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/03/2022
 Supervised By :mohammad ahmed 10/04/2022



TIC: BP011858.D\data.ms

(91) Benzo(k)fluoranthene

23.080min (-0.041) 11.36 ng/u1

response 996620

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.50	23.28
125.00	11.50	12.42
0.00	0.00	0.00

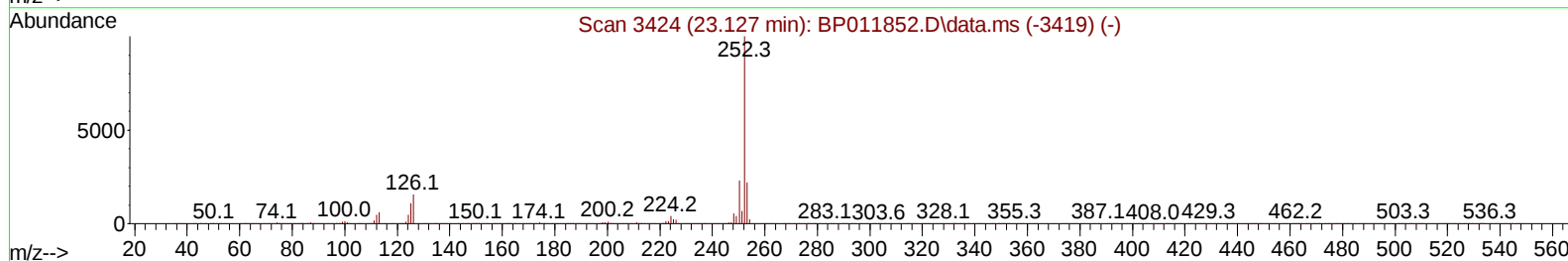
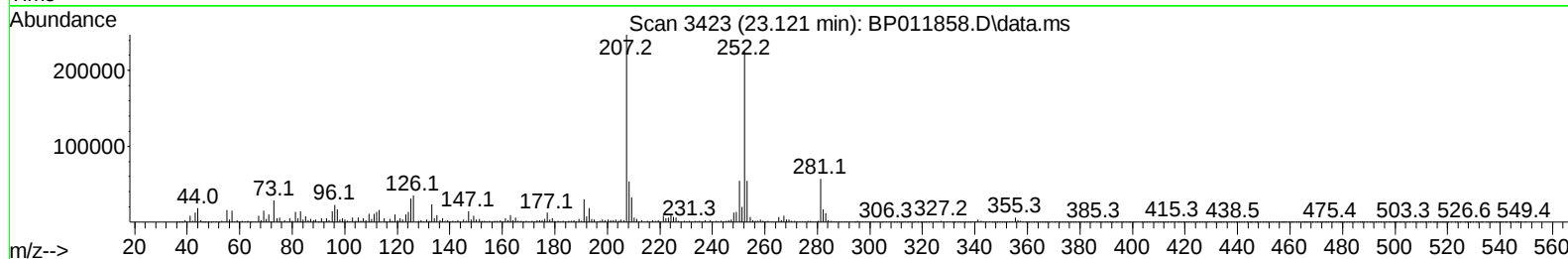
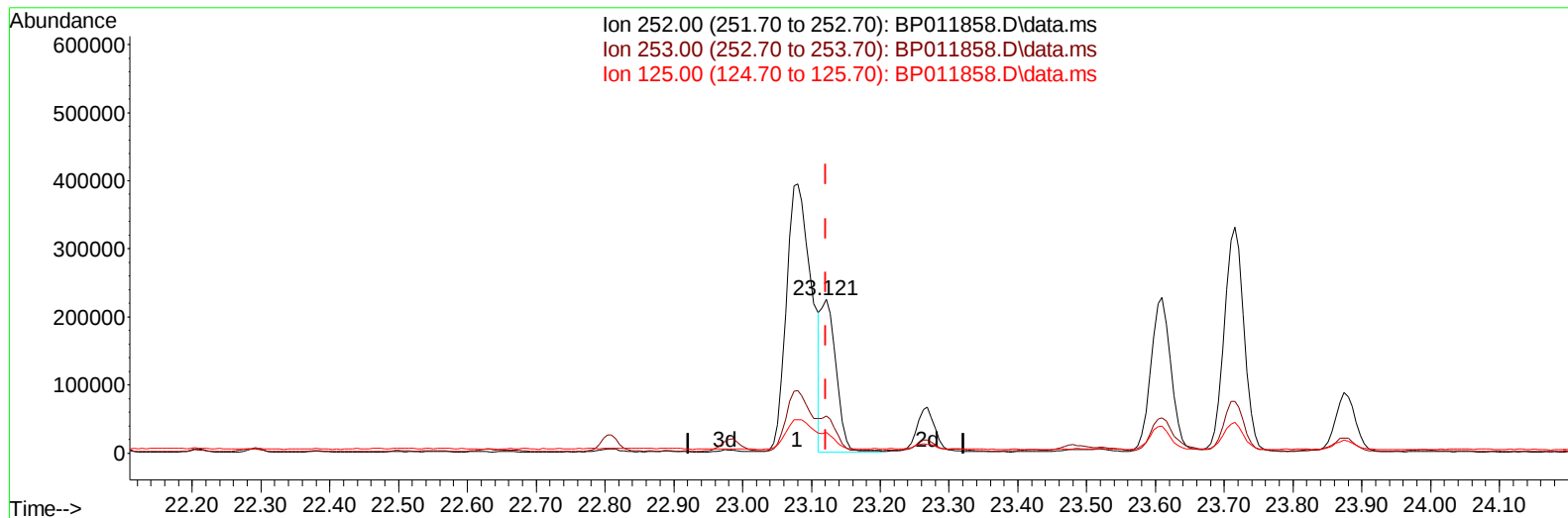
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TIC: BP011858.D\data.ms

(91) Benzo(k)fluoranthene

23.121min (-0.000) 3.89 ng/u1 m

response 341291

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.50	24.03
125.00	11.50	13.56
0.00	0.00	0.00

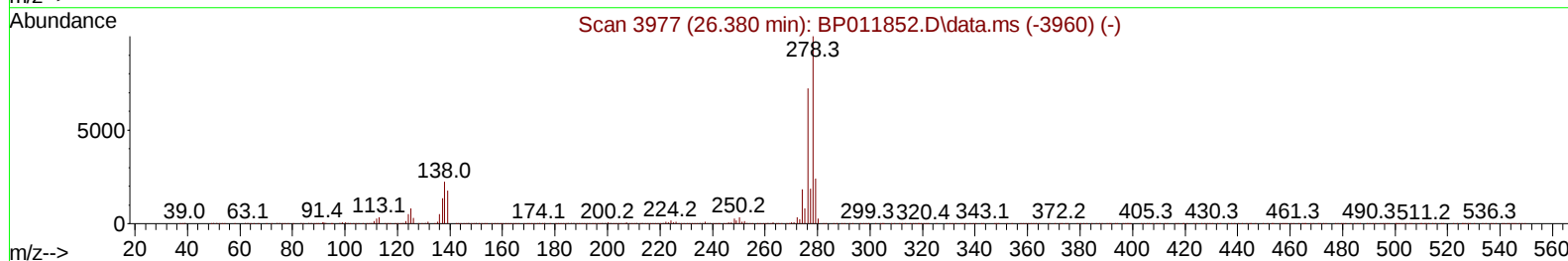
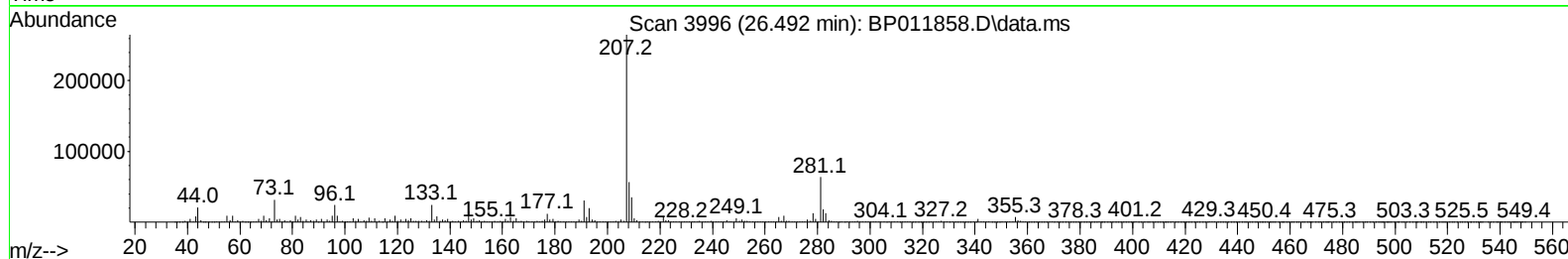
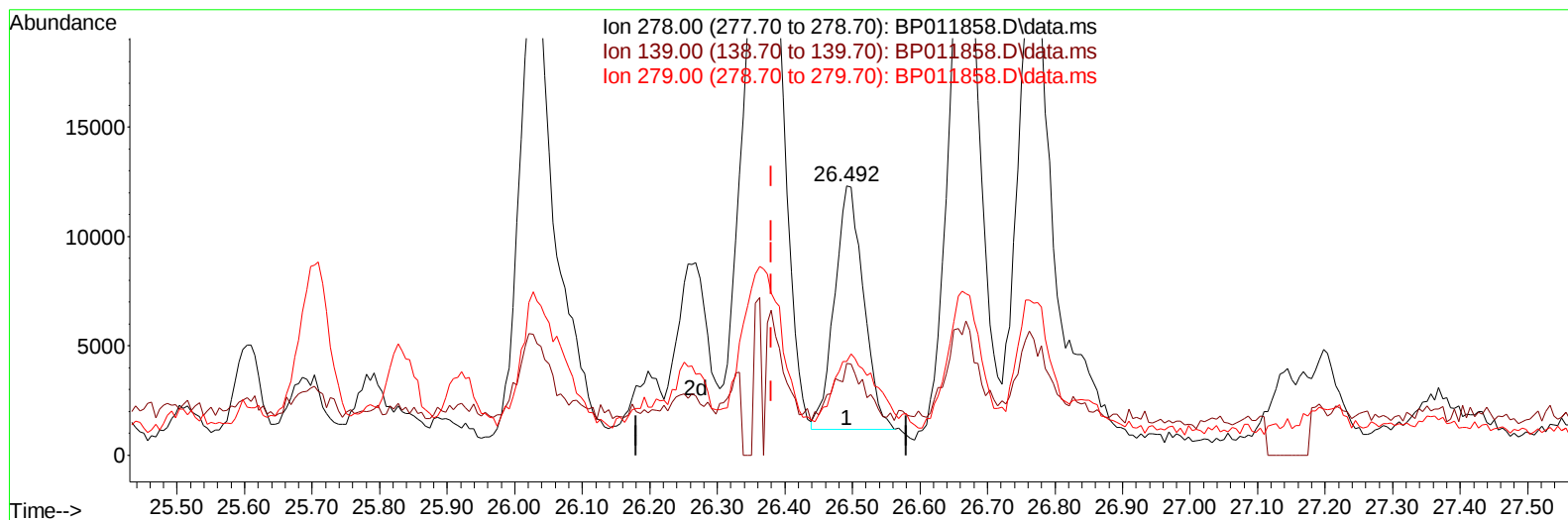
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TIC: BP011858.D\data.ms

(95) Dibenzo(a,h)anthracene

26.492min (+ 0.112) 0.36 ng/u1

response 33218

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	18.80	34.07#
279.00	23.80	34.72#
0.00	0.00	0.00

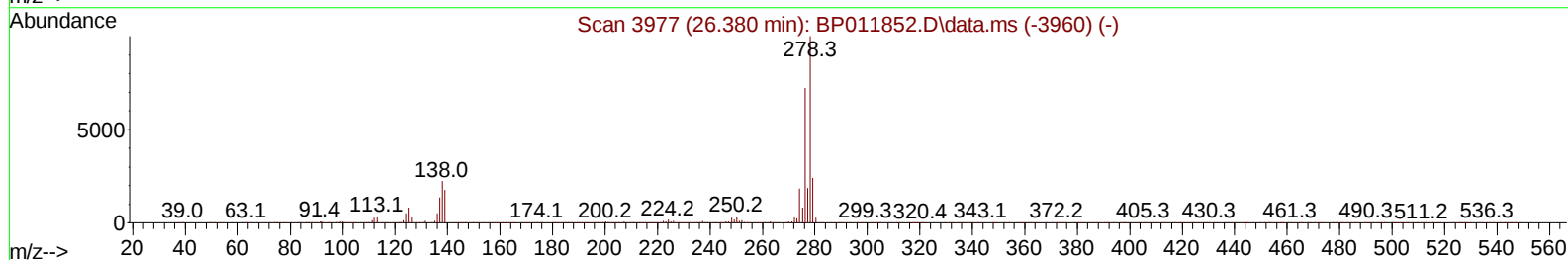
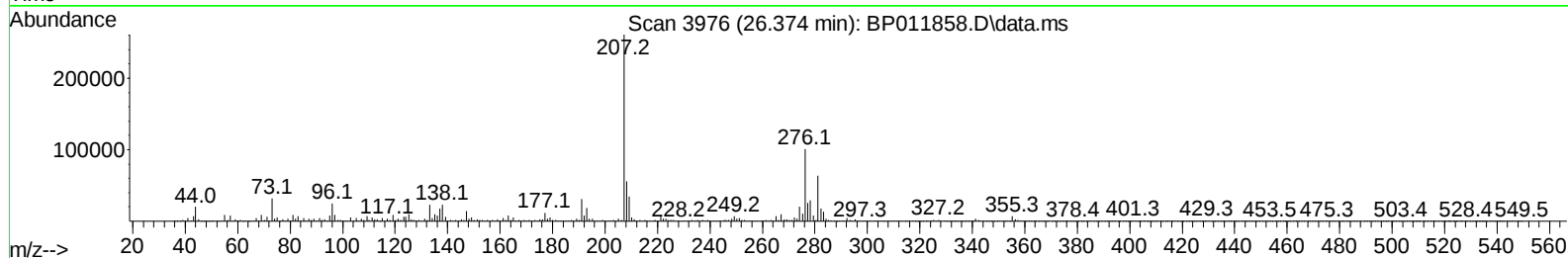
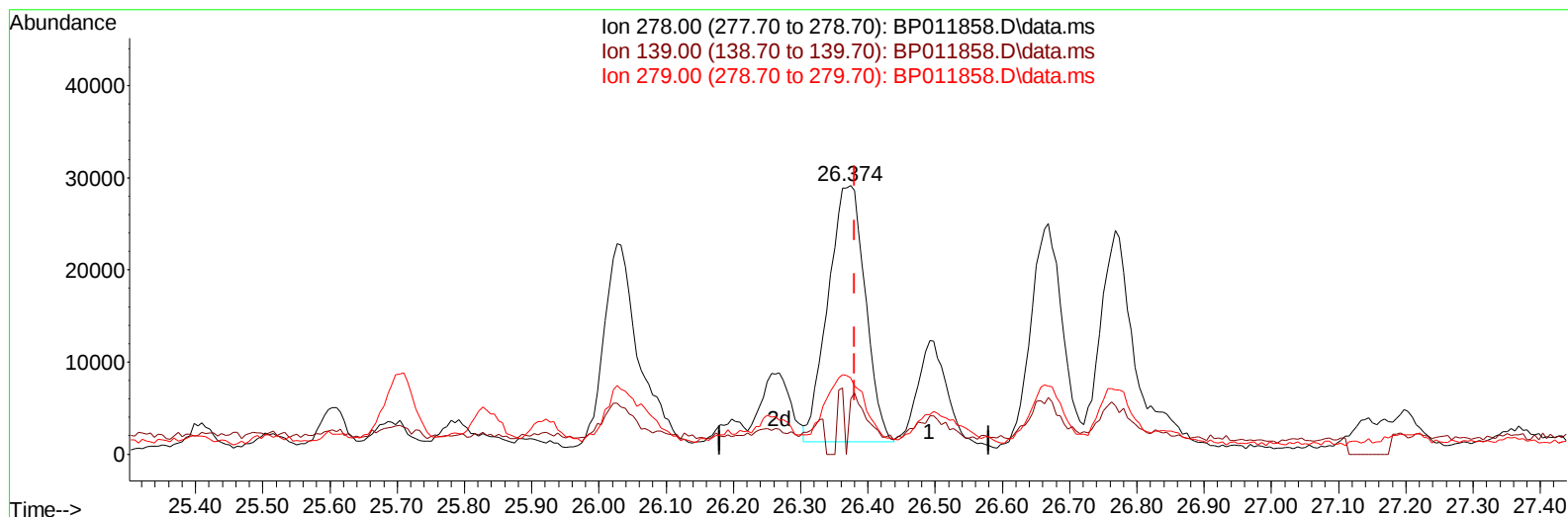
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TIC: BP011858.D\data.ms

(95) Dibenzo(a,h)anthracene

26.374min (-0.006) 1.16 ng/ul m

response 106654

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	18.80	20.72
279.00	23.80	28.47
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.893	152		303951	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136		1377529	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164		900108	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188		1885024	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240		1689523	20.000	ng/ul	0.00
88) Perylene-d12	23.827	264		1427231	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.305	96		21792	3.162	ng/uL	0.00
4) Pyridine-d5	3.734	84		157692	7.811	ng/ul	0.00
7) Phenol-d5	7.057	99		513767	19.579	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67		288473	19.834	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132		442538	21.364	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113		423077	19.504	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128		224375	21.384	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143		248904	22.173	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165		456582	21.989	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131		519868	16.423	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166		1426531	22.162	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160		1667500	22.725	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143		100736	7.558	ng/ul	0.00
60) Fluorene-d10	15.539	176		1309266	23.422	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200		97297	8.528	ng/ul	0.00
73) Anthracene-d10	17.398	188		2045839	23.826	ng/ul	0.00
81) Pyrene-d10	19.633	212		2388882	27.919	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264		1797227	25.155	ng/ul	0.00
Target Compounds							
						Qvalue	
8) Phenol	7.087	94		32796	1.205	ng/ul	95
16) Acetophenone	8.722	105		37058	1.121	ng/ul	99
52) Acenaphthene	14.604	153		61269	1.066	ng/ul	99
61) Fluorene	15.592	166		84392	1.308	ng/ul	100
72) Phenanthrene	17.339	178		1483178	14.317	ng/ul	100
74) Anthracene	17.433	178		303331	2.910	ng/ul	98
77) Carbazole	17.704	167		130803	1.389	ng/ul	98
80) Fluoranthene	19.304	202		2233105	21.326	ng/ul	99
82) Pyrene	19.663	202		1892005	17.354	ng/ul	96
85) Benzo(a)anthracene	21.368	228		914205	8.274	ng/ul	99
87) Chrysene	21.421	228		880887	8.496	ng/ul	97
90) Benzo(b)fluoranthene	23.080	252		996620	10.940	ng/ul	98
91) Benzo(k)fluoranthene	23.121	252		341291m	3.891	ng/ul	
93) Benzo(a)pyrene	23.715	252		652805	8.030	ng/ul	98
94) Indeno(1,2,3-cd)pyrene	26.356	276		426495	4.122	ng/ul	96
95) Dibenzo(a,h)anthracene	26.374	278		106654m	1.160	ng/ul	
96) Benzo(g,h,i)perylene	27.139	276		353111	3.849	ng/ul	97

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

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