

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060324\
 Data File : BN031760.D
 Acq On : 03 Jun 2024 15:17
 Operator : MA/JU
 Sample : P2590-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :

BNA_N

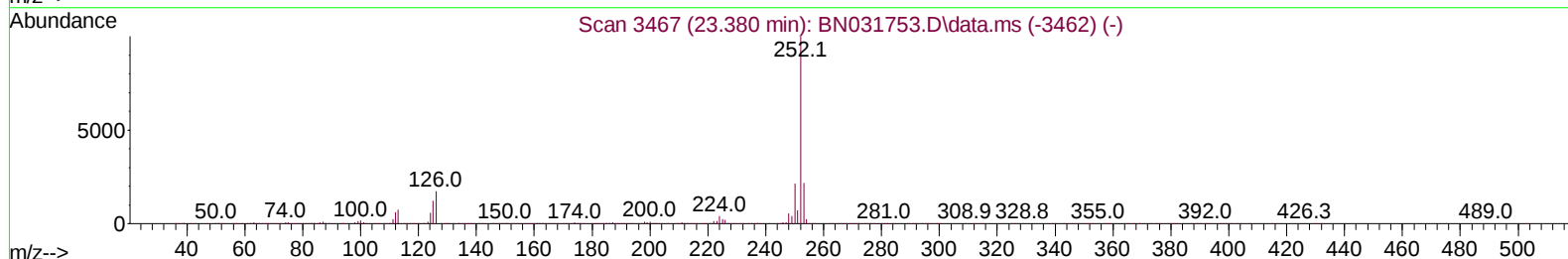
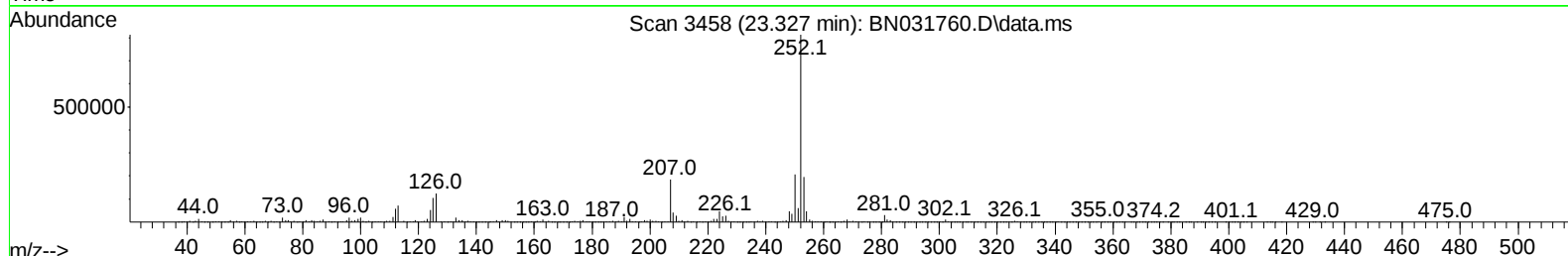
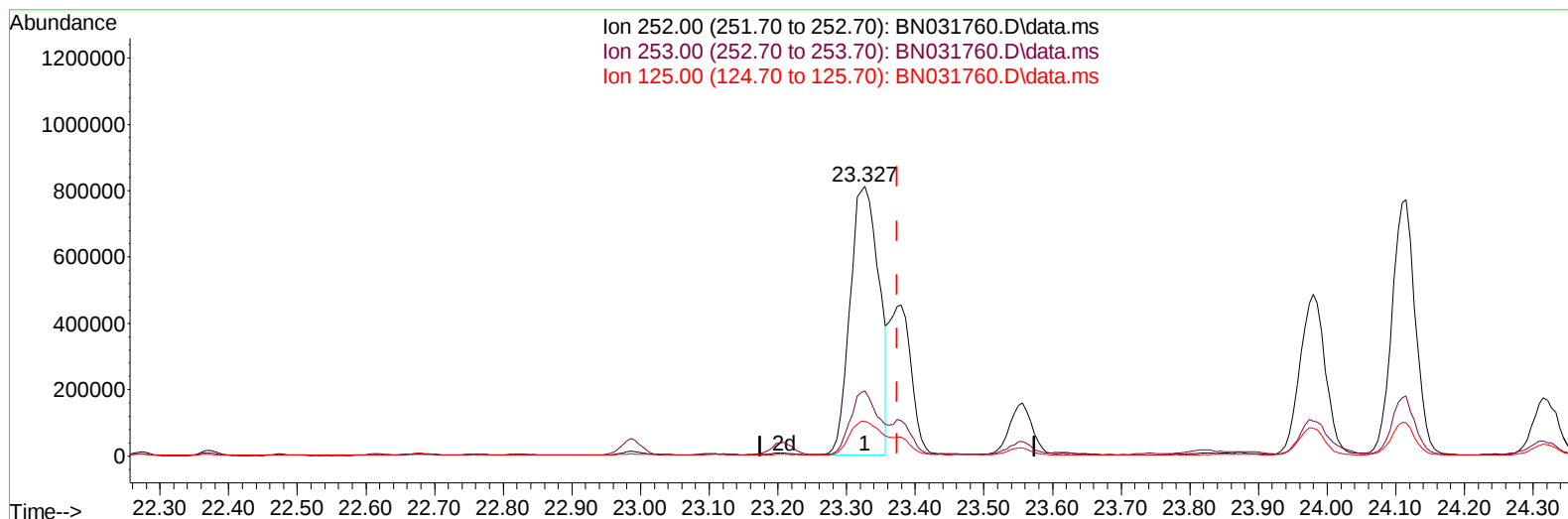
ClientSampleId :

BHBL2

Manual IntegrationsAPPROVED

Quant Time: Jun 03 15:48:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 29 15:13:07 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/04/2024
 Supervised By :mohammad ahmed 06/05/2024



TIC: BN031760.D\data.ms

(91) Benzo(k)fluoranthene**23.327min (-0.048) 19.83 ng/u1****response 2362868**

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.00	24.24
125.00	11.80	12.78
0.00	0.00	0.00

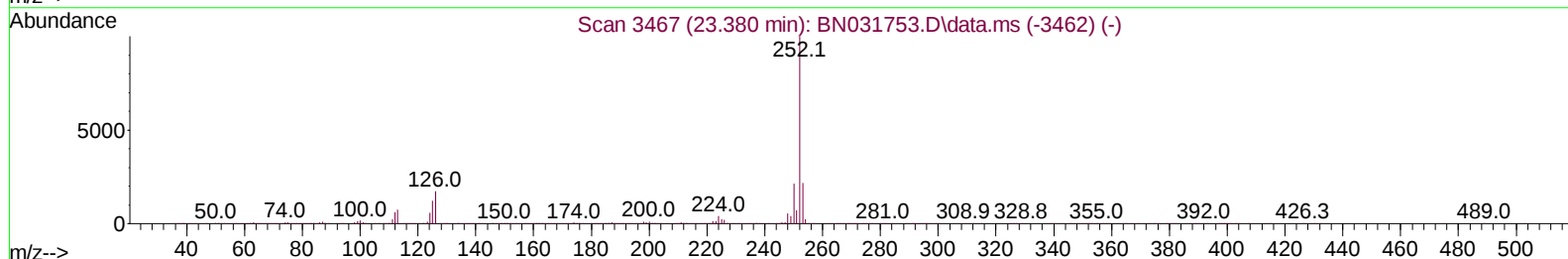
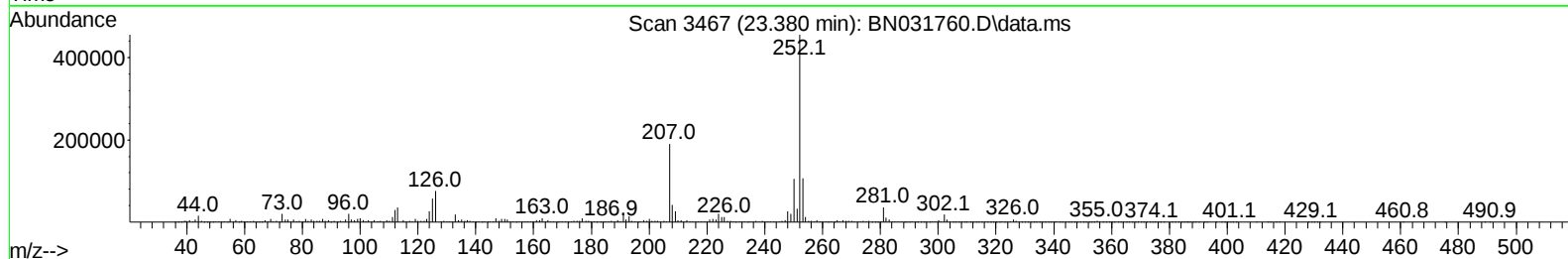
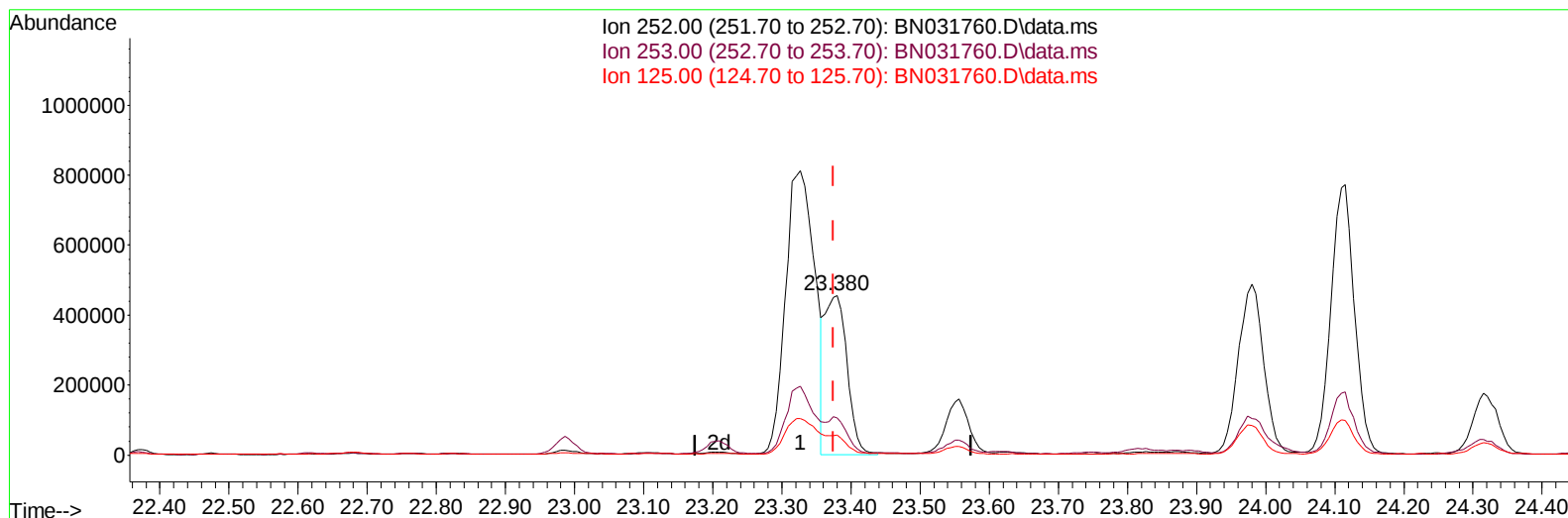
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TIC: BN031760.D\data.ms

(91) Benzo(k)fluoranthene

23.380min (+ 0.005) 8.48 ng/u1 m

response 1010050

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.00	23.47
125.00	11.80	12.52
0.00	0.00	0.00

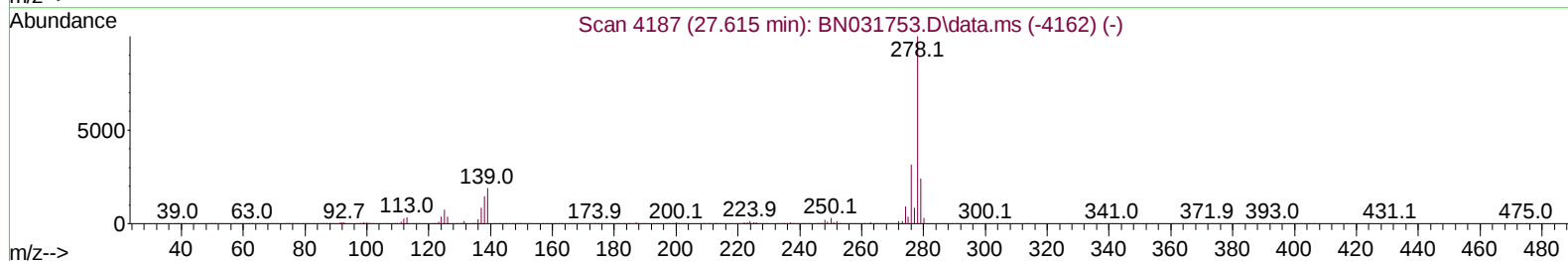
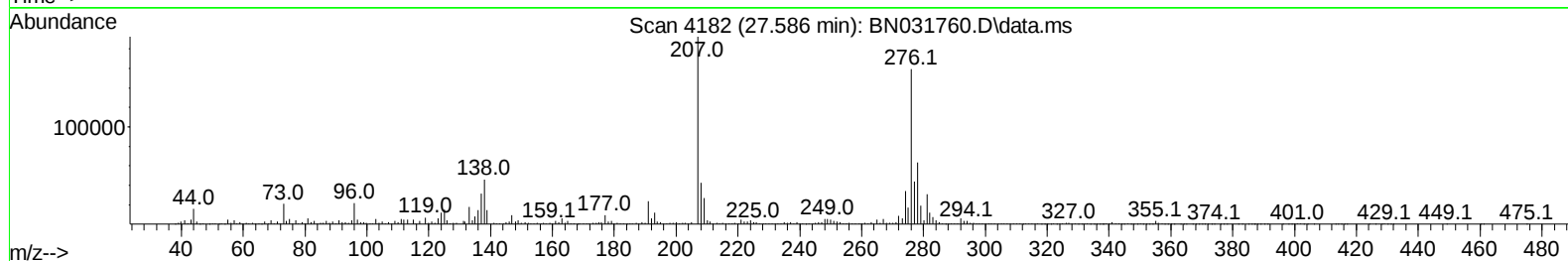
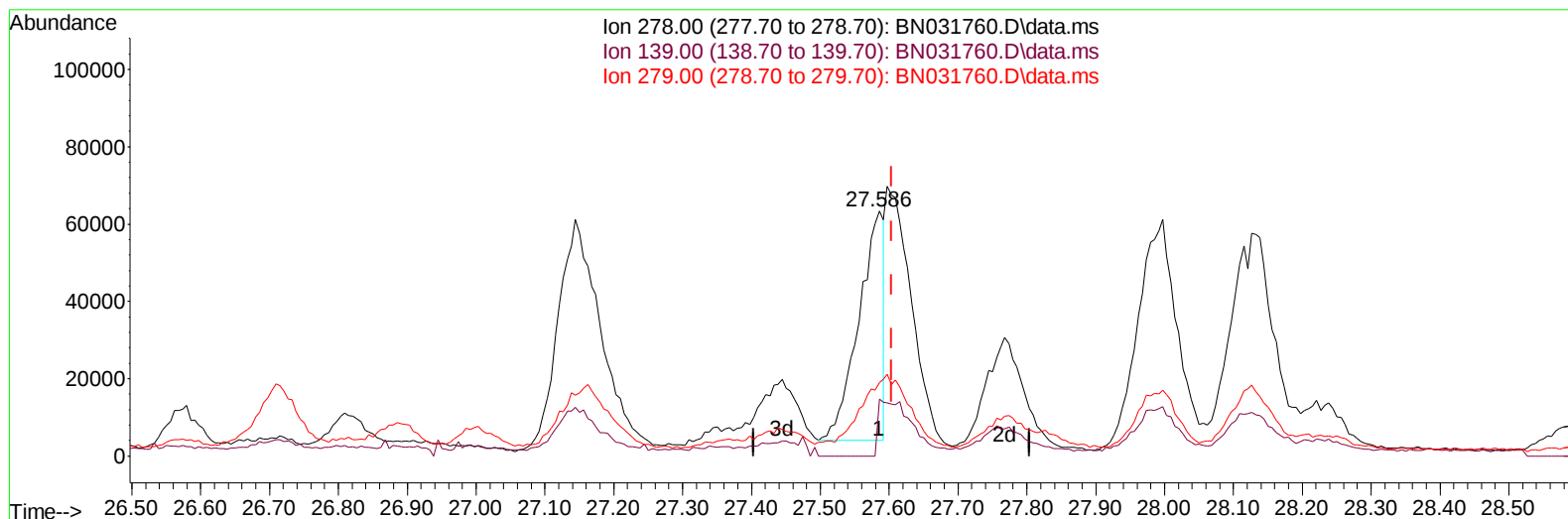
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TIC: BN031760.D\data.ms

(95) Dibenzo(a,h)anthracene

27.586min (-0.018) 1.41 ng/u1

response 150747

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	18.80	23.13#
279.00	21.70	30.00#
0.00	0.00	0.00

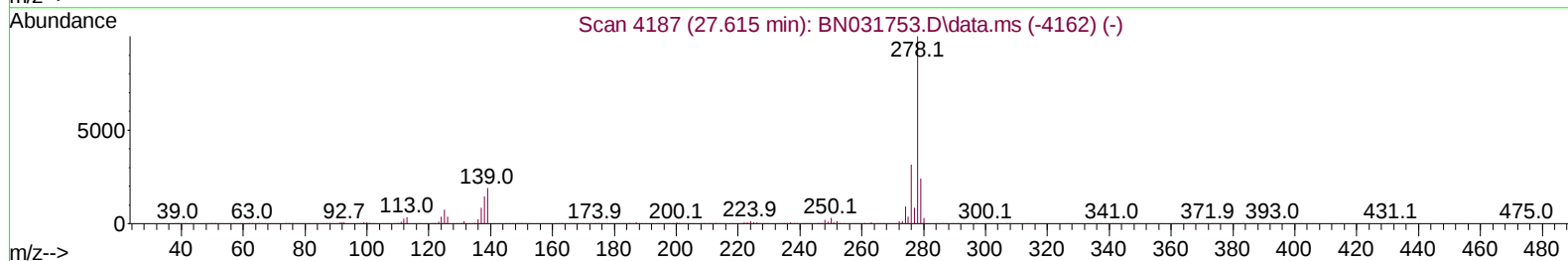
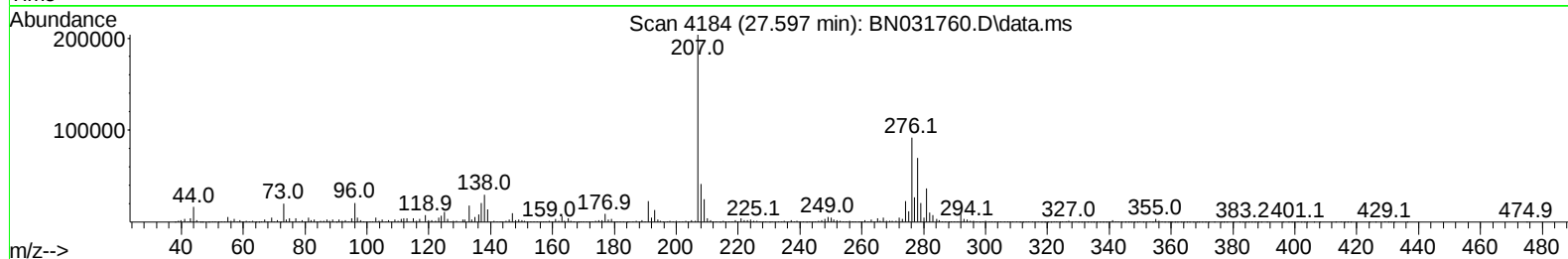
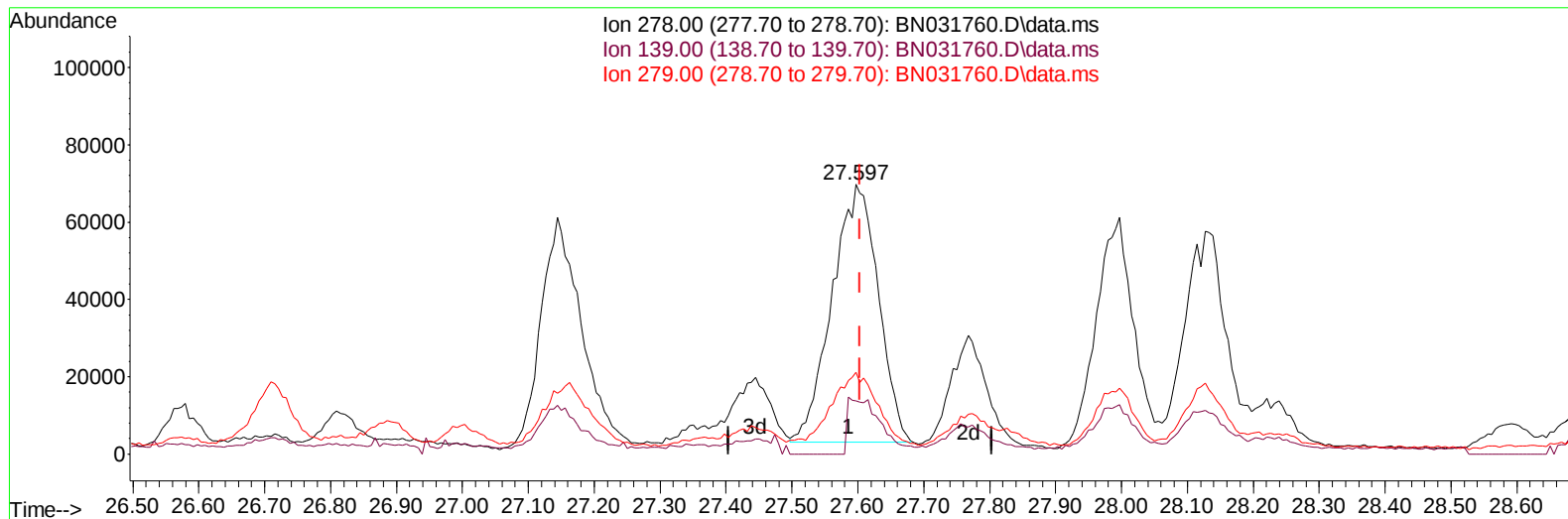
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Misc :
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TIC: BN031760.D\data.ms

(95) Dibenzo(a,h)anthracene

27.597min (-0.006) 3.04 ng/u1 m

response 324762

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	18.80	19.76
279.00	21.70	30.34#
0.00	0.00	0.00

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Instrument :
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Quant Time: Jun 03 15:52:13 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.681	152		588360	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136		2579567	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.322	164		1491787	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.068	188		2464683	20.000	ng/ul	0.00
79) Chrysene-d12	21.309	240		1703343	20.000	ng/ul	0.00
88) Perylene-d12	24.250	264		2031545	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.193	96		44427	3.124	ng/uL	0.00
4) Pyridine-d5	3.599	84		330815	8.263	ng/ul	0.00
7) Phenol-d5	6.851	99		991283	19.653	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67		624712	21.174	ng/ul	0.00
11) 2-Chlorophenol-d4	7.216	132		850380	22.096	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113		719175	17.492	ng/ul	0.00
21) Nitrobenzene-d5	8.839	128		454269	23.032	ng/ul	0.00
24) 2-Nitrophenol-d4	9.551	143		474860	23.307	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.086	165		829378	21.805	ng/ul	0.00
31) 4-Chloroaniline-d4	10.604	131		776300	13.746	ng/ul	0.00
46) Dimethylphthalate-d6	13.739	166		2480344	23.866	ng/ul	0.00
49) Acenaphthylene-d8	14.016	160		2843256	23.793	ng/ul	0.00
54) 4-Nitrophenol-d4	14.527	143		212248	10.200	ng/ul	0.00
60) Fluorene-d10	15.316	176		2024790	23.026	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200		141456	11.418	ng/ul	0.00
73) Anthracene-d10	17.168	188		2631923	24.943	ng/ul	0.00
81) Pyrene-d10	19.468	212		2570673	28.203	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.050	264		2392621	24.664	ng/ul	0.01
Target Compounds							
						Qvalue	
50) Acenaphthylene	14.045	152		295815	2.190	ng/ul	99
61) Fluorene	15.374	166		137684	1.408	ng/ul	100
72) Phenanthrene	17.110	178		2621866	20.882	ng/ul	98
74) Anthracene	17.204	178		483514	3.801	ng/ul	98
77) Carbazole	17.474	167		240609	2.218	ng/ul	99
80) Fluoranthene	19.133	202		4290895	39.390	ng/ul	100
82) Pyrene	19.498	202		3840023	34.358	ng/ul	99
85) Benzo(a)anthracene	21.292	228		1917333	16.817	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.221	149		108751	1.413	ng/ul	99
87) Chrysene	21.351	228		1895588	17.800	ng/ul	97
90) Benzo(b)fluoranthene	23.327	252		2362868	19.459	ng/ul	95
91) Benzo(k)fluoranthene	23.380	252		1010050m	8.477	ng/ul	
93) Benzo(a)pyrene	24.115	252		1824250	16.040	ng/ul	96
94) Indeno(1,2,3-cd)pyrene	27.562	276		1265046	9.653	ng/ul	98
95) Dibenzo(a,h)anthracene	27.597	278		324762m	3.035	ng/ul	
96) Benzo(g,h,i)perylene	28.597	276		1068663	10.258	ng/ul	98

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

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