

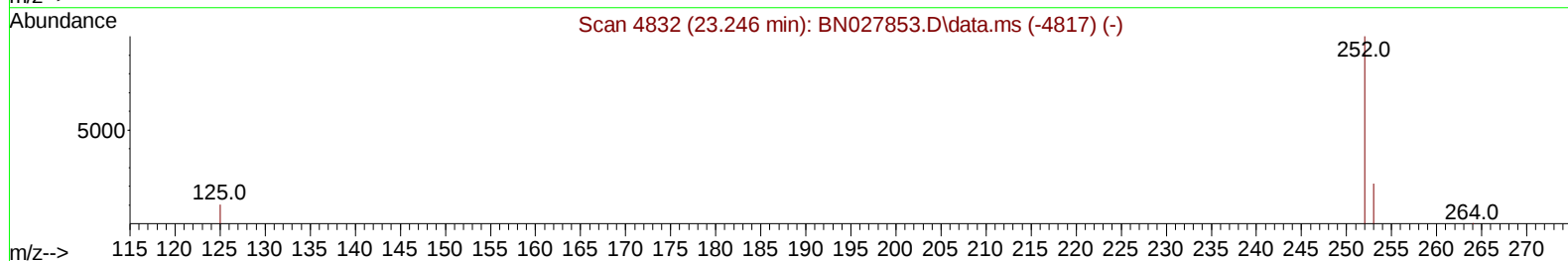
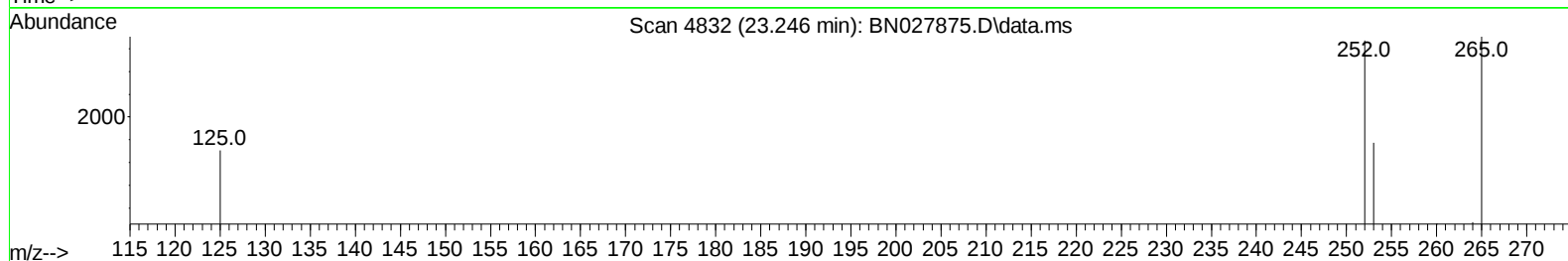
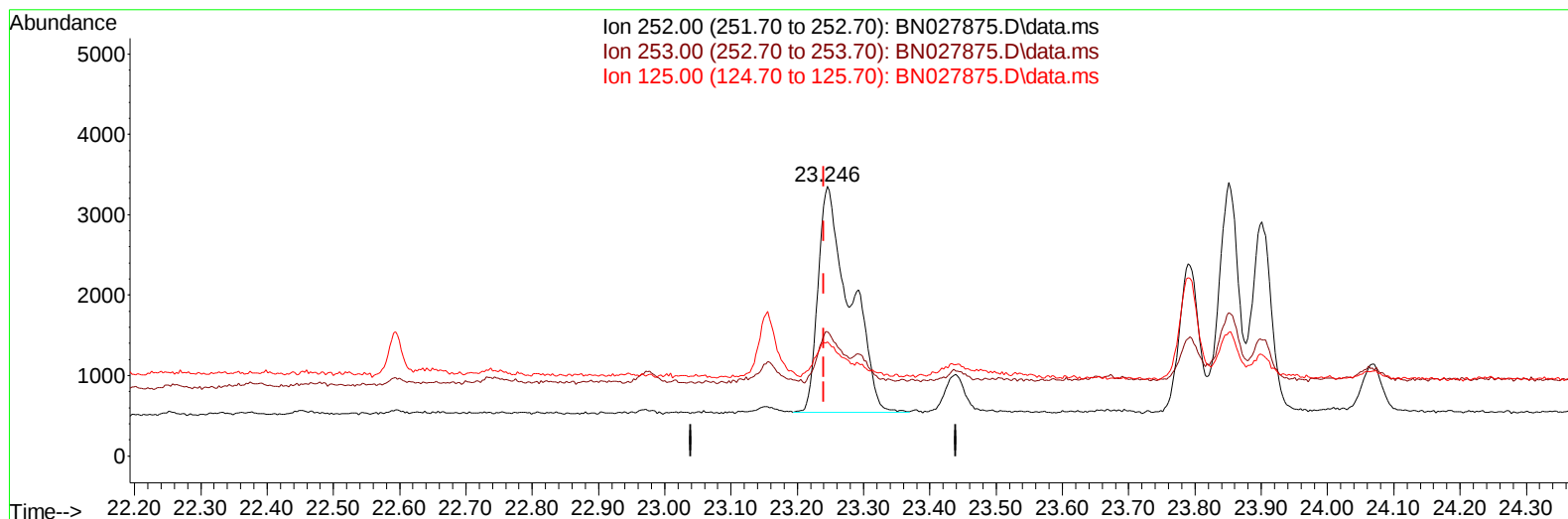
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092623\
 Data File : BN027875.D
 Acq On : 26 Sep 2023 23:21
 Operator : MA/JU
 Sample : 04410-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BBJM0

Manual IntegrationsAPPROVED

Quant Time: Sep 27 03:02:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 03:58:00 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023
 Supervised By :mohammad ahmed 09/28/2023



TIC: BN027875.D\data.ms

(24) Benzo(b)fluoranthene

23.246min (+ 0.006) 0.09 ng/u1

response 9360

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	25.10	46.23
125.00	15.70	42.29#
0.00	0.00	0.00

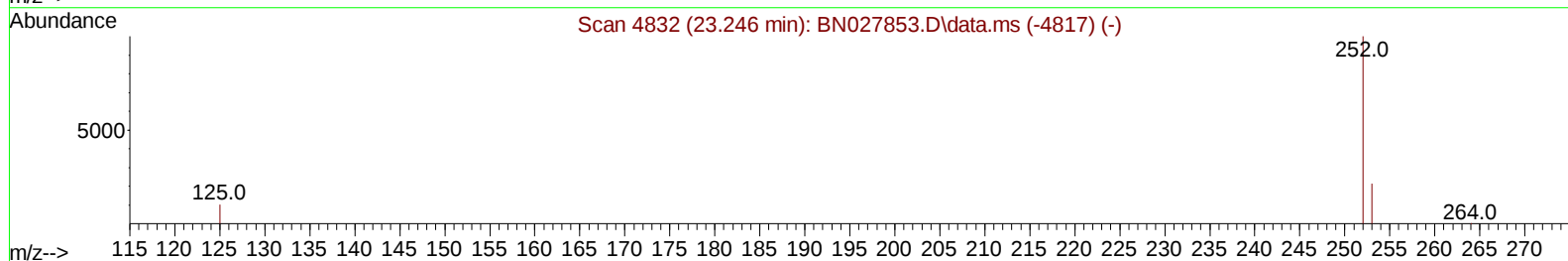
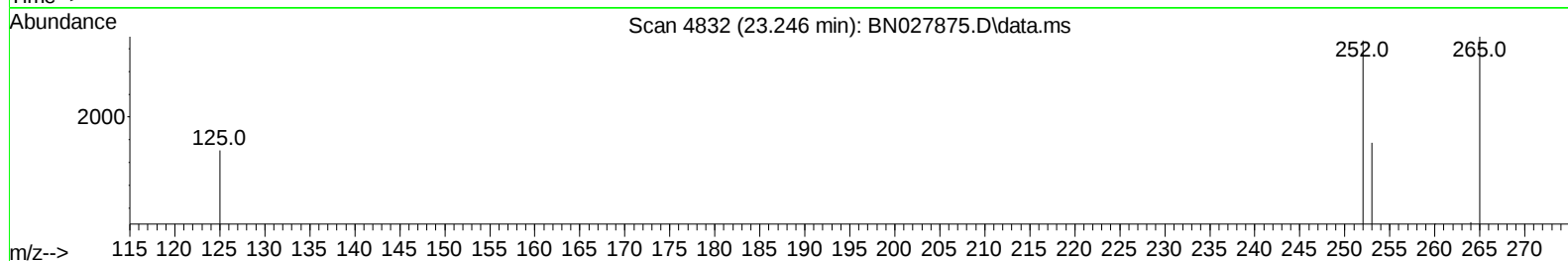
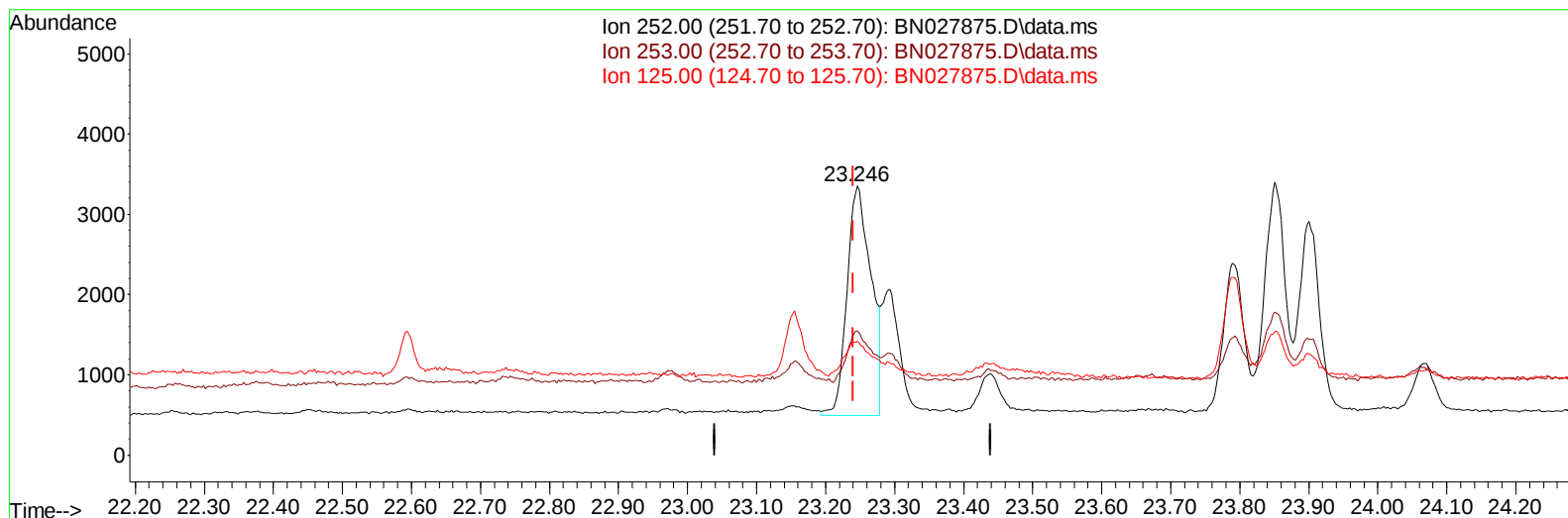
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TIC: BN027875.D\data.ms

(24) Benzo(b)fluoranthene

23.246min (+ 0.006) 0.07 ng/u1 m

response 6928

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	25.10	46.23
125.00	15.70	42.29#
0.00	0.00	0.00

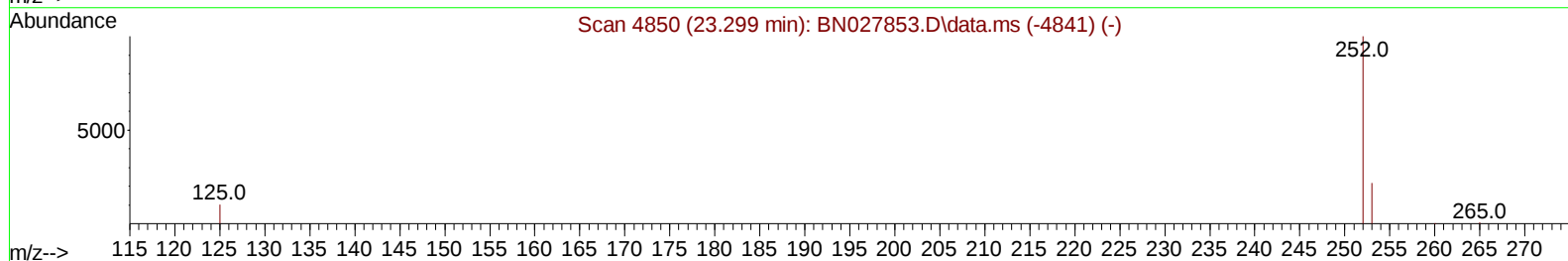
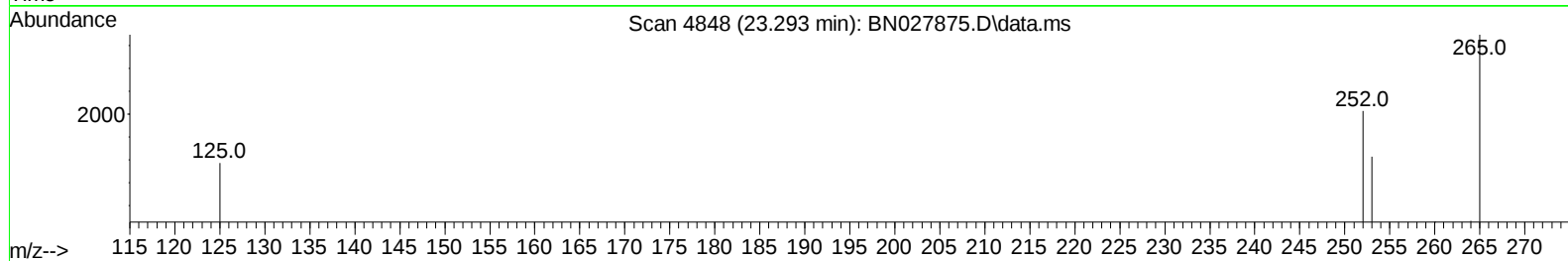
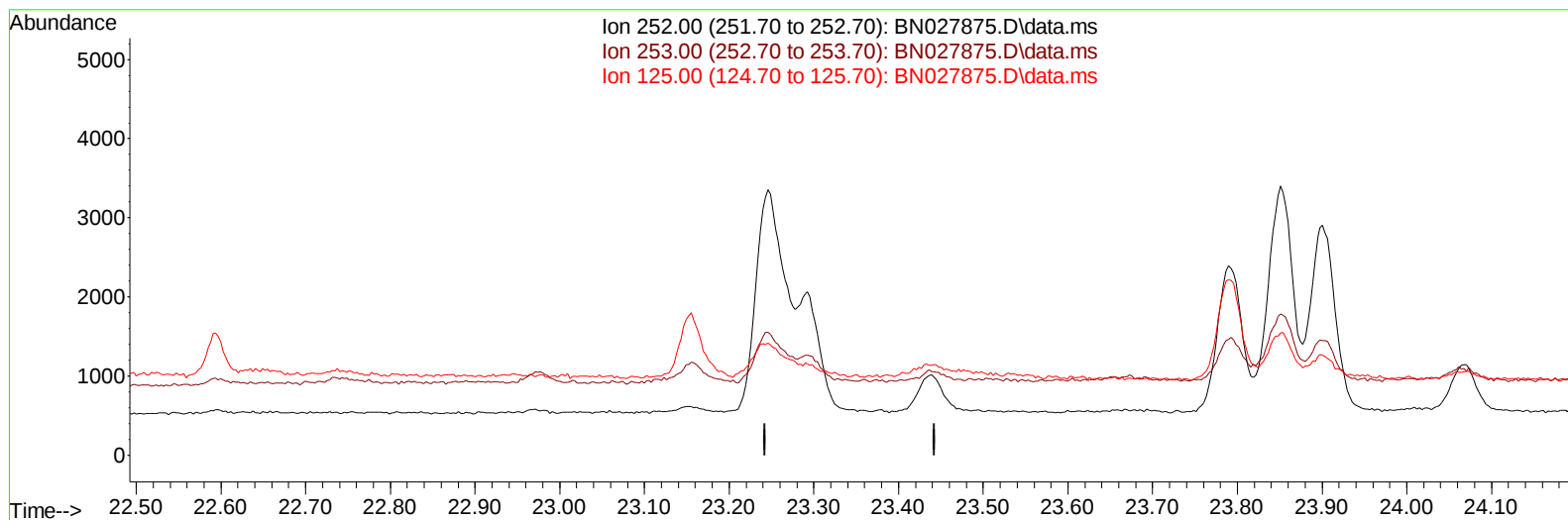
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TIC: BN027875.D\data.ms

(25) Benzo(k)fluoranthene

23.292min (-23.292) 0.00 ng/u1

response 0

Ion	Exp%	Act%
252.00	100.00	0.00
253.00	24.90	0.00#
125.00	15.80	0.00#
0.00	0.00	0.00

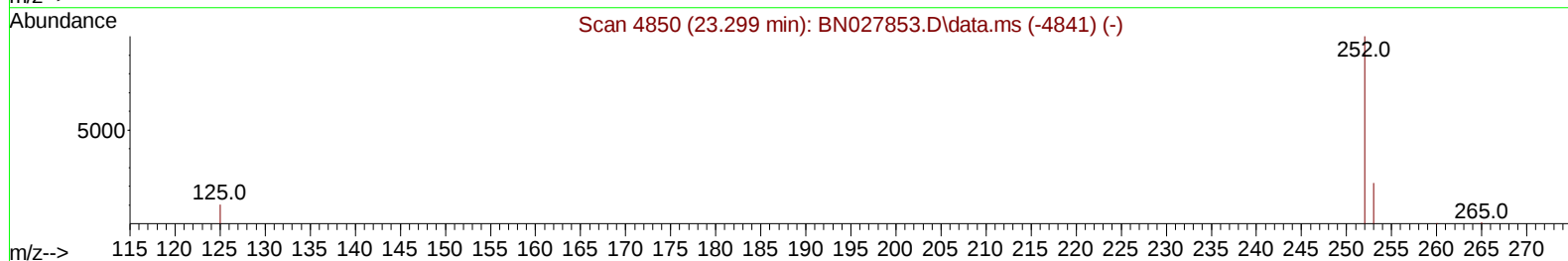
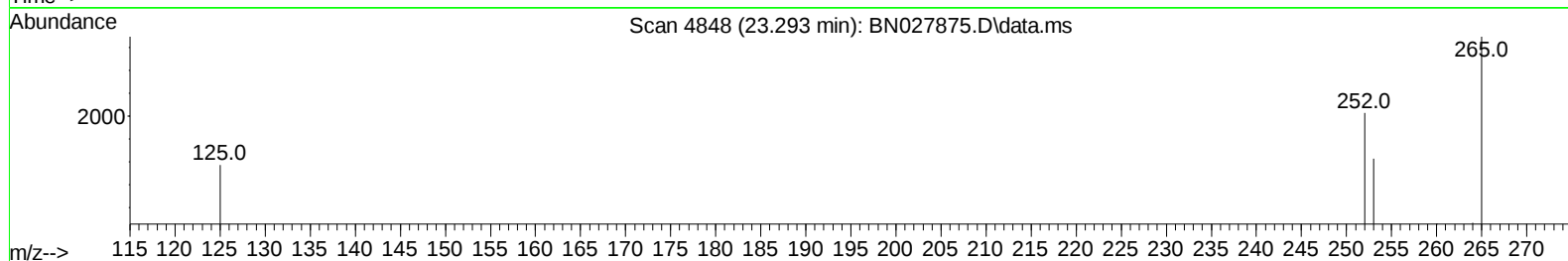
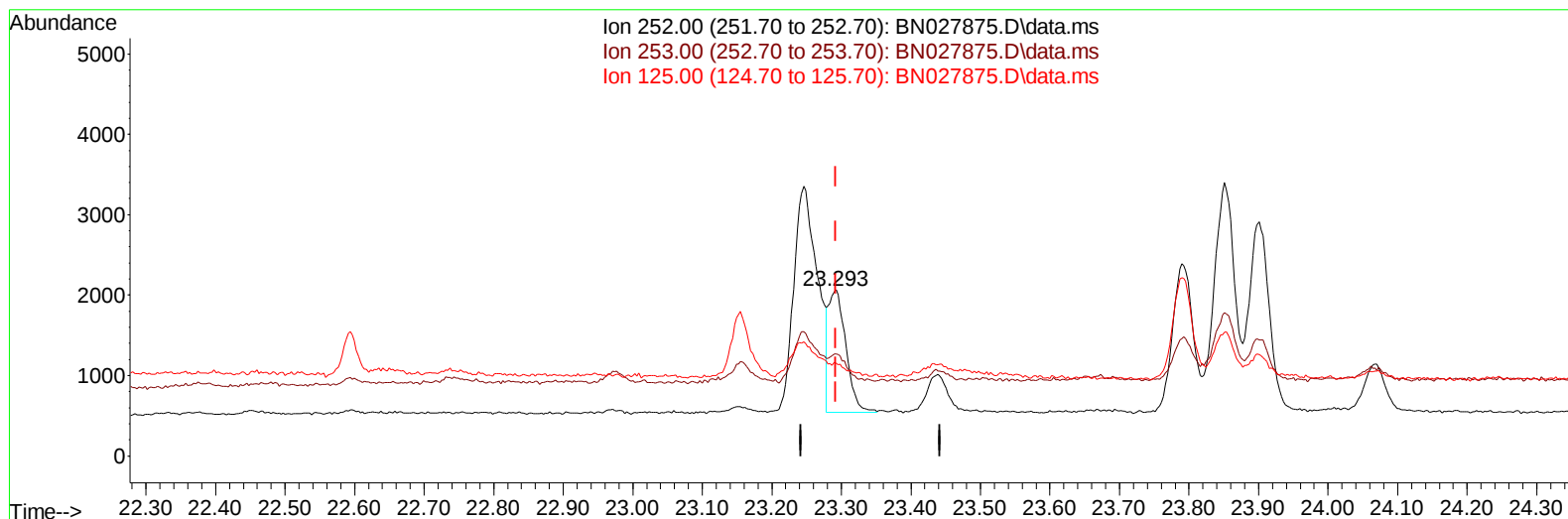
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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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(25) Benzo(k)fluoranthene

23.293min (+ 0.000) 0.03 ng/u1 m

response 2639

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	24.90	61.29#
125.00	15.80	55.91#
0.00	0.00	0.00

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 ALS Vial : 24 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.962	152		3366	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136		11816	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.606	164		7924	0.400	ng/ul	0.01
13) Phenanthrene-d10	17.358	188		18587	0.400	ng/ul	0.02
17) Chrysene-d12	21.539	240		18425	0.400	ng/ul	0.02
23) Perylene-d12	24.012	264		22515	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.330	96		13525	3.190	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152		5366	0.300	ng/ul	0.00
18) Fluoranthene-d10	19.380	212		18378	0.327	ng/ul	0.02
Target Compounds							
						Qvalue	
2) 1,4-Dioxane	3.368	88		1980	0.449	ng/ul	99
15) Phenanthrene	17.400	178		3557	0.061	ng/ul#	90
16) Anthracene	17.489	178		1958	0.040	ng/ul#	79
19) Fluoranthene	19.413	202		8379	0.112	ng/ul	96
20) Pyrene	19.780	202		7774	0.101	ng/ul	96
21) Benzo(a)anthracene	21.521	228		5047	0.071	ng/ul#	91
22) Chrysene	21.577	228		5029	0.068	ng/ul#	89
24) Benzo(b)fluoranthene	23.246	252		6928m	0.069	ng/ul	
25) Benzo(k)fluoranthene	23.293	252		2639m	0.027	ng/ul	
26) Benzo(a)pyrene	23.901	252		4701	0.058	ng/ul#	53
27) Indeno(1,2,3-cd)pyrene	26.606	276		3662	0.035	ng/ul#	98
29) Benzo(g,h,i)perylene	27.421	276		3737	0.043	ng/ul#	82

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

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