

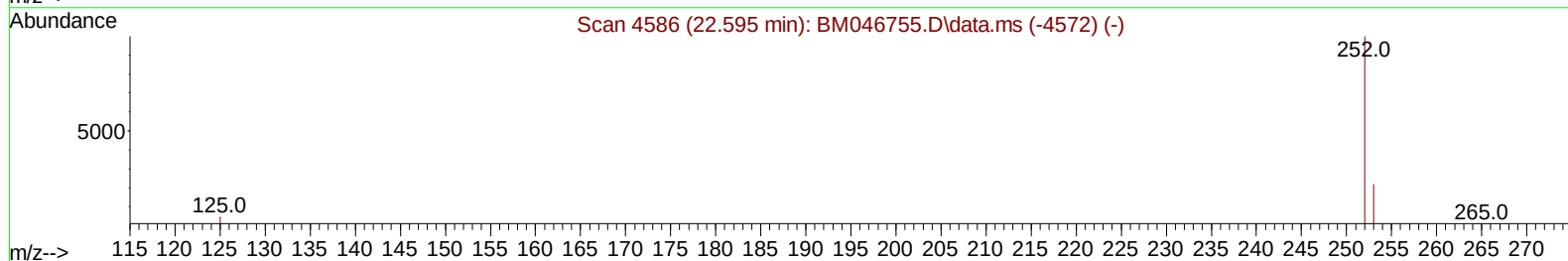
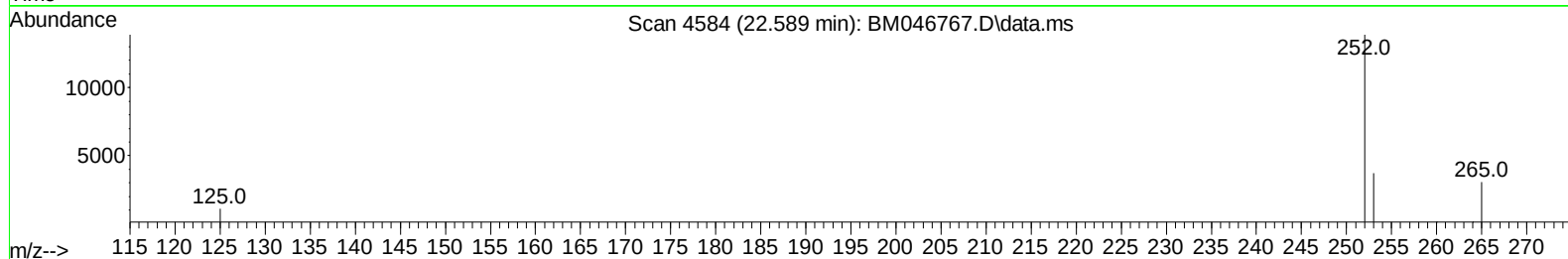
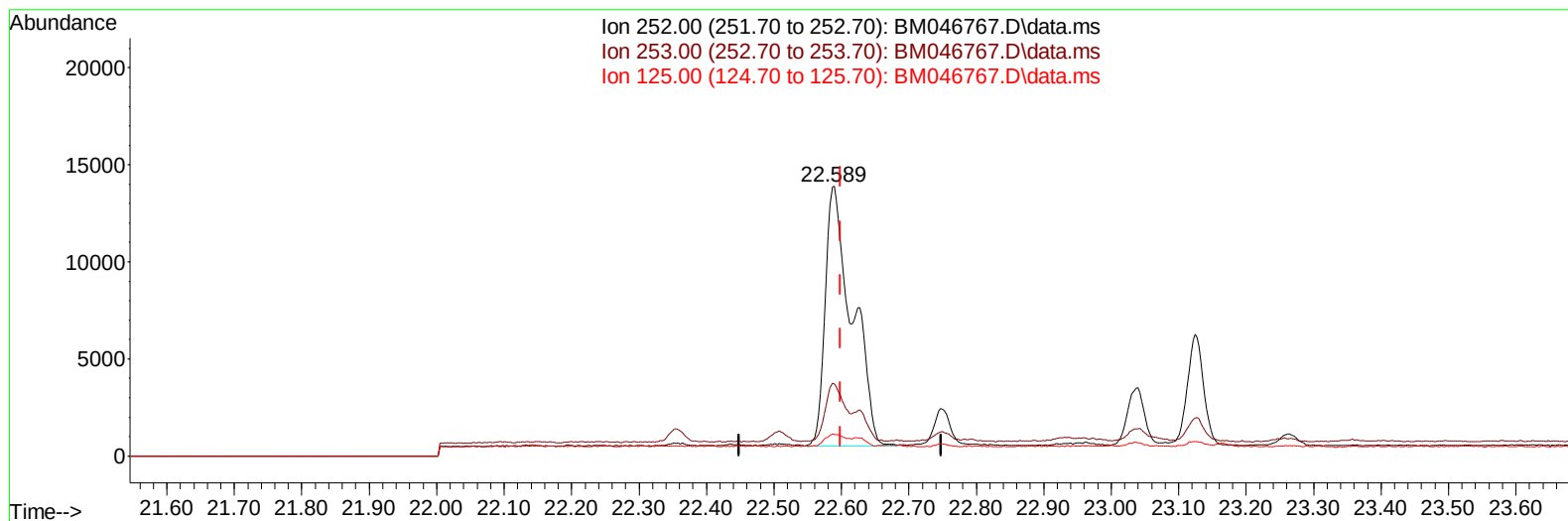
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072224\
Data File : BM046767.D
Acq On : 23 Jul 2024 03:08
Operator : MA/JU
Sample : P3201-13DL 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4CH7DL

Manual IntegrationsAPPROVED

Quant Time: Jul 23 03:38:04 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 19 11:38:04 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/23/2024
Supervised By :mohammad ahmed 07/23/2024



TIC: BM046767.D\data.ms

(24) Benzo(b)fluoranthene

22.589min (-0.009) 0.94 ng/u1

response 36986

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	28.10	26.79
125.00	8.90	7.94
0.00	0.00	0.00

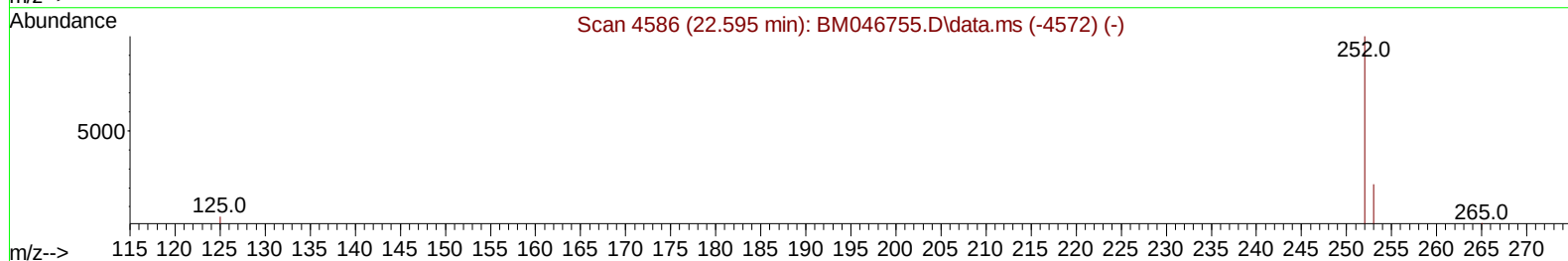
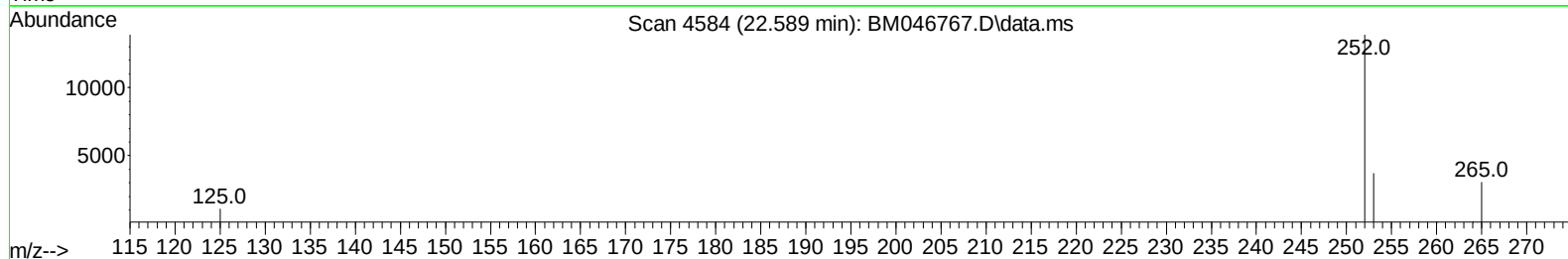
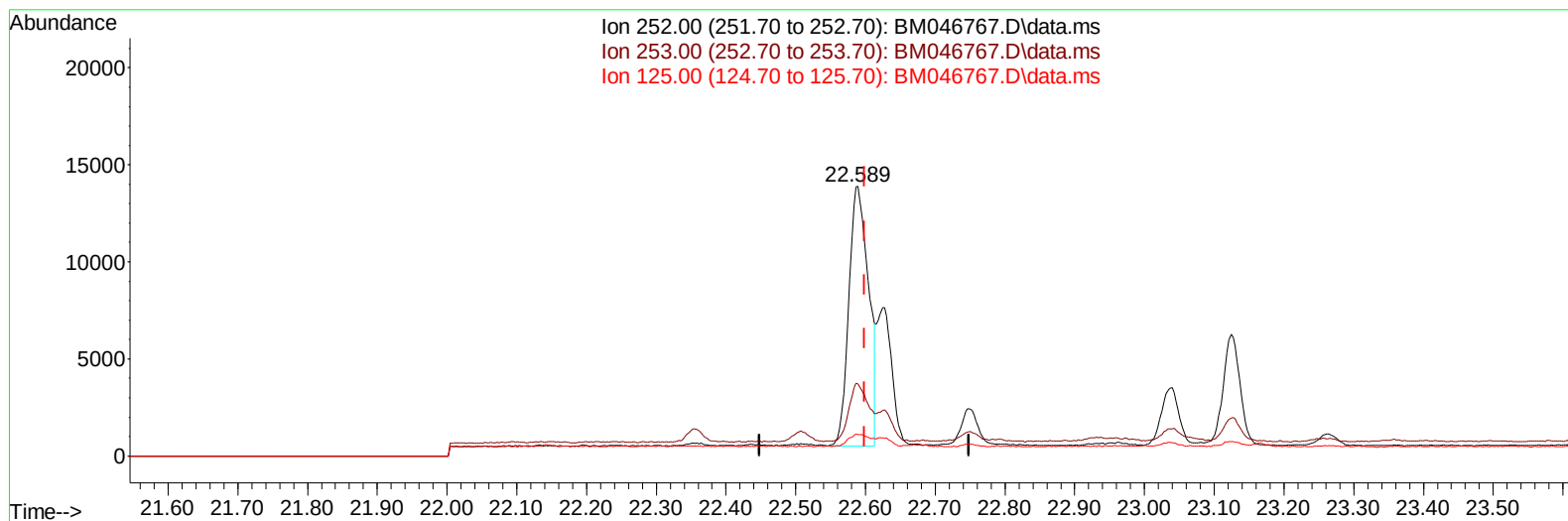
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TIC: BM046767.D\data.ms

(24) Benzo(b)fluoranthene

22.589min (-0.009) 0.66 ng/ul m

response 26197

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	28.10	26.79
125.00	8.90	7.94
0.00	0.00	0.00

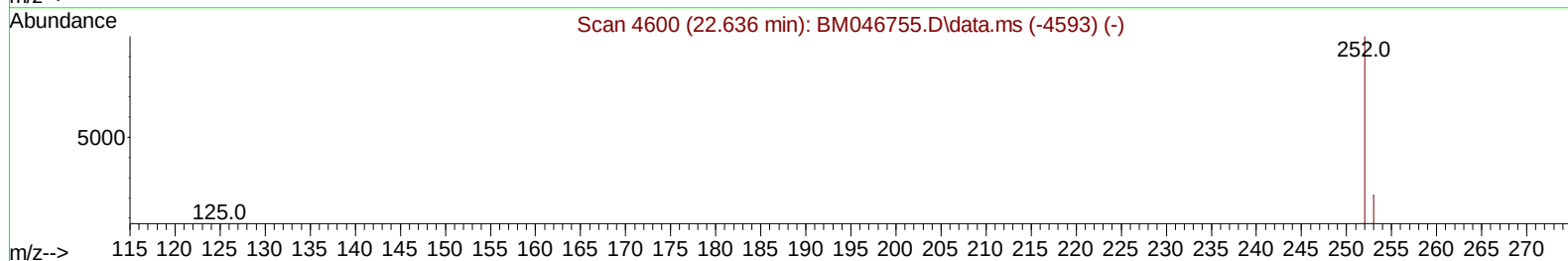
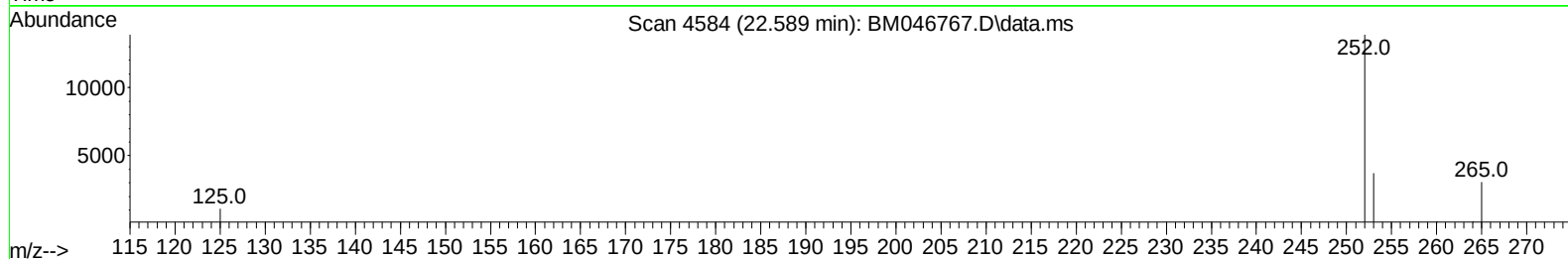
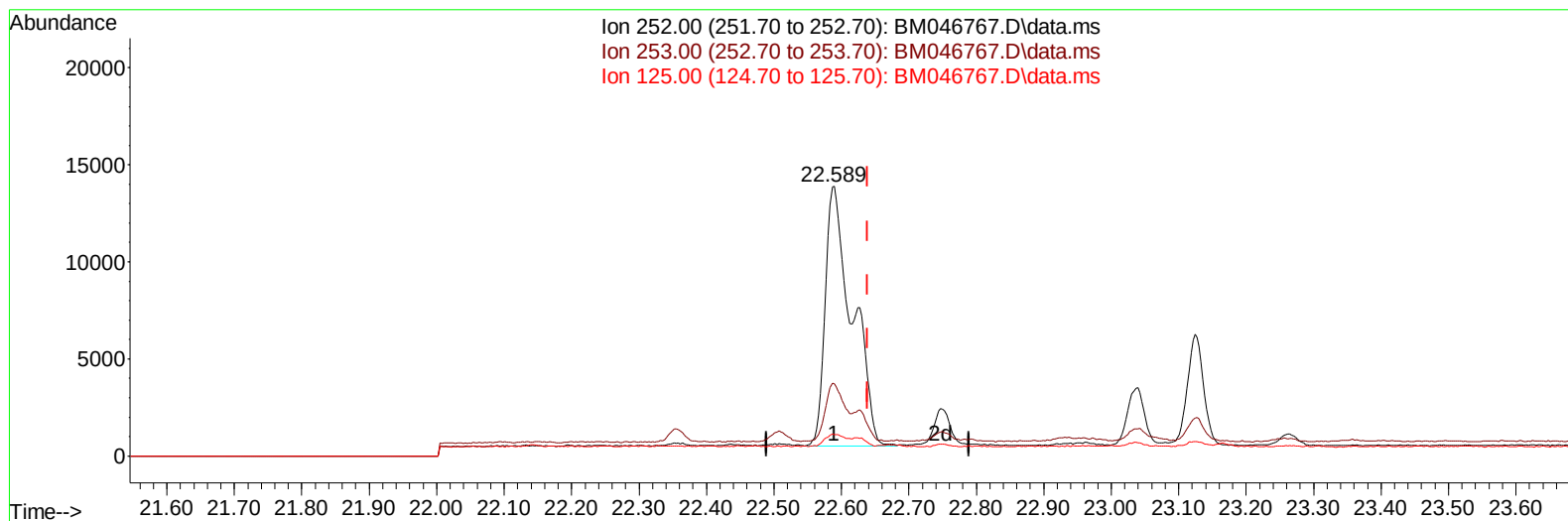
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Sample : P3201-13DL 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4CH7DL

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TIC: BM046767.D\data.ms

(25) Benzo(k)fluoranthene

22.589min (-0.050) 0.95 ng/u1

response 36986

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	28.00	26.79
125.00	8.70	7.94
0.00	0.00	0.00

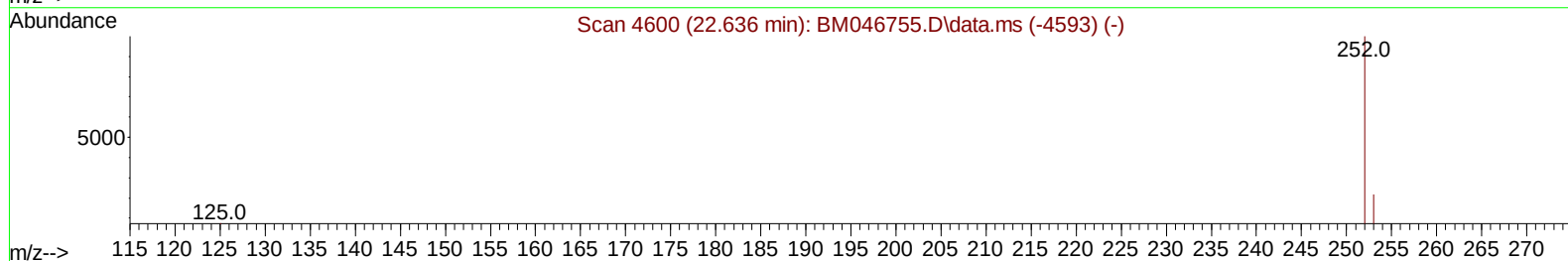
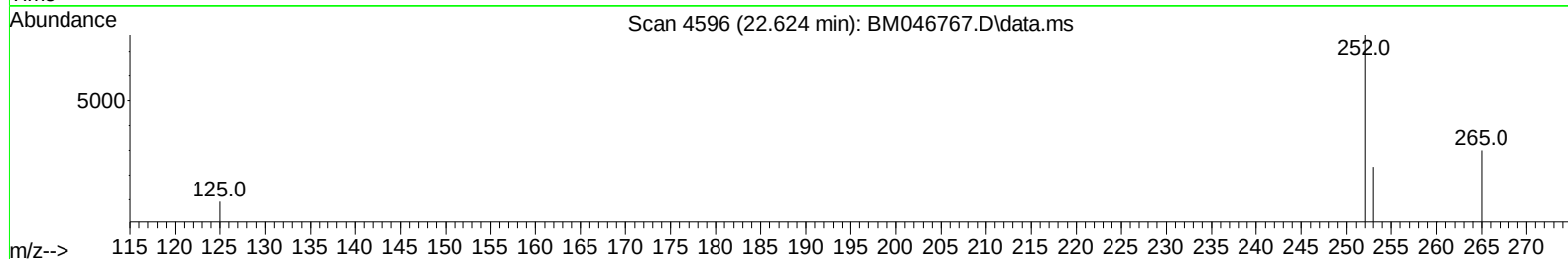
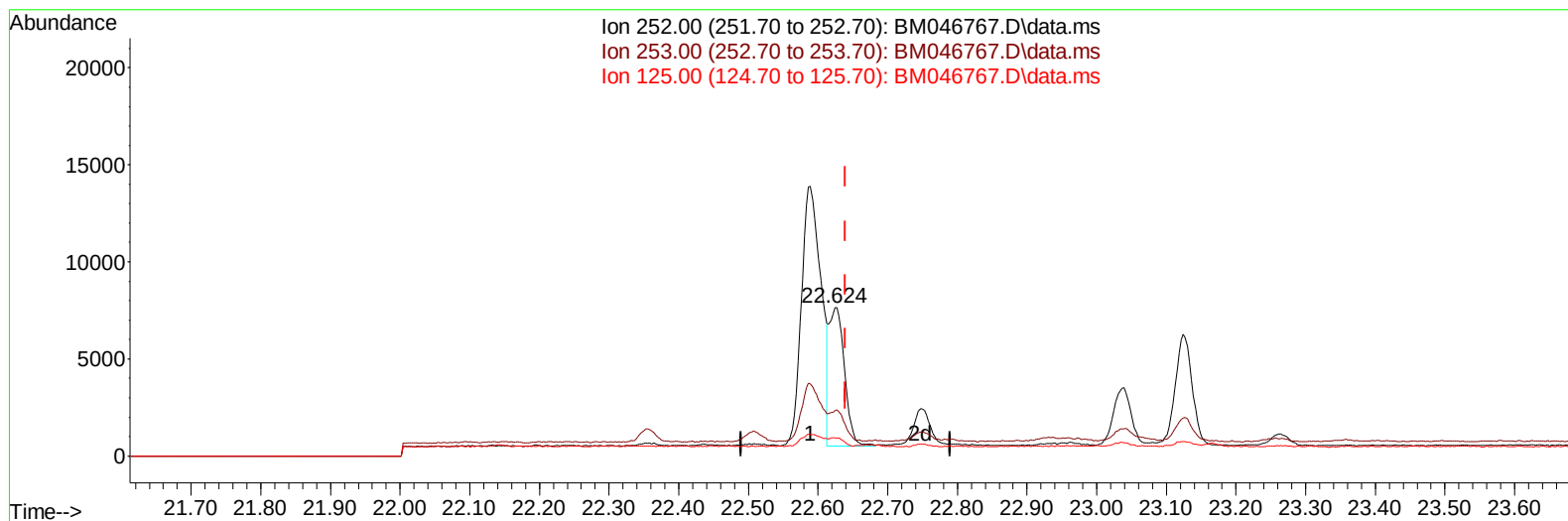
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Instrument :
BNA_M
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TIC: BM046767.D\data.ms

(25) Benzo(k)fluoranthene

22.624min (-0.015) 0.28 ng/ul m

response 10897

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	28.00	30.56
125.00	8.70	12.17#
0.00	0.00	0.00

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 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.480	152	3308	0.400	ng/ul	0.00
4) Naphthalene-d8	10.238	136	8455	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.123	164	5131	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.880	188	10511	0.400	ng/ul	0.00
17) Chrysene-d12	21.087	240	8682	0.400	ng/ul	0.00
23) Perylene-d12	23.218	264	9890	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.109	96	1118	0.411	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.833	152	388	0.028	ng/ul	-0.02
18) Fluoranthene-d10	18.918	212	814	0.031	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.288	128	518	0.024	ng/ul#	68
10) Acenaphthylene	13.841	152	1092	0.047	ng/ul#	92
11) Acenaphthene	14.188	153	937	0.060	ng/ul	97
12) Fluorene	15.182	166	1121	0.058	ng/ul#	91
15) Phenanthrene	16.922	178	21938	0.731	ng/ul	98
16) Anthracene	17.010	178	3904	0.133	ng/ul	95
19) Fluoranthene	18.951	202	42811	1.200	ng/ul	99
20) Pyrene	19.313	202	29016	0.752	ng/ul	99
21) Benzo(a)anthracene	21.070	228	16859	0.439	ng/ul	99
22) Chrysene	21.122	228	18868	0.508	ng/ul	98
24) Benzo(b)fluoranthene	22.589	252	26197m	0.663	ng/ul	
25) Benzo(k)fluoranthene	22.624	252	10897m	0.281	ng/ul	
26) Benzo(a)pyrene	23.124	252	9647	0.297	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.326	276	10520	0.213	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.336	278	3425	0.087	ng/ul#	69
29) Benzo(g,h,i)perylene	25.967	276	1663	0.042	ng/ul#	61

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

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