

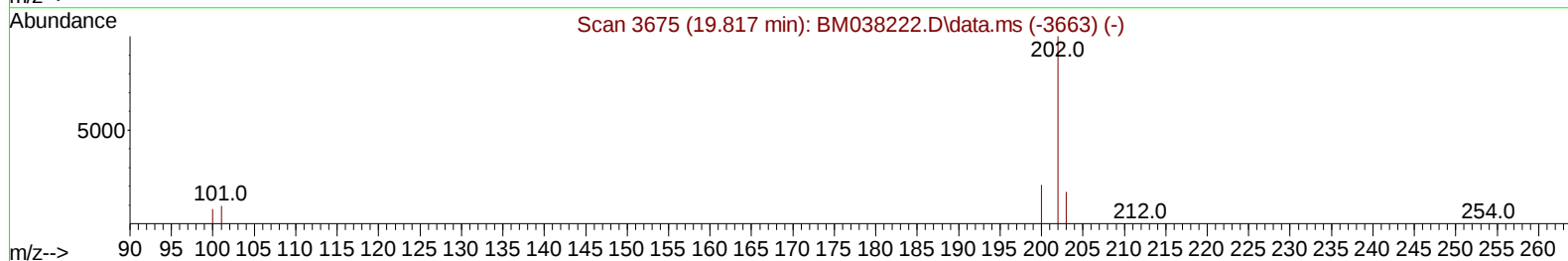
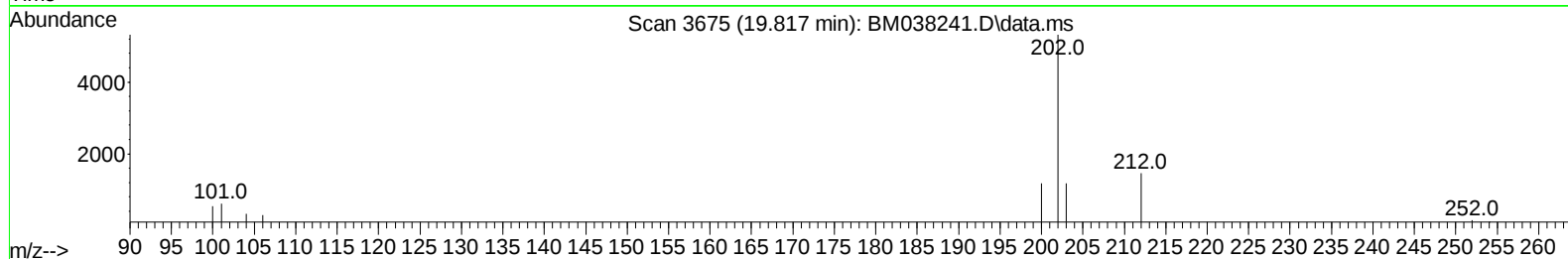
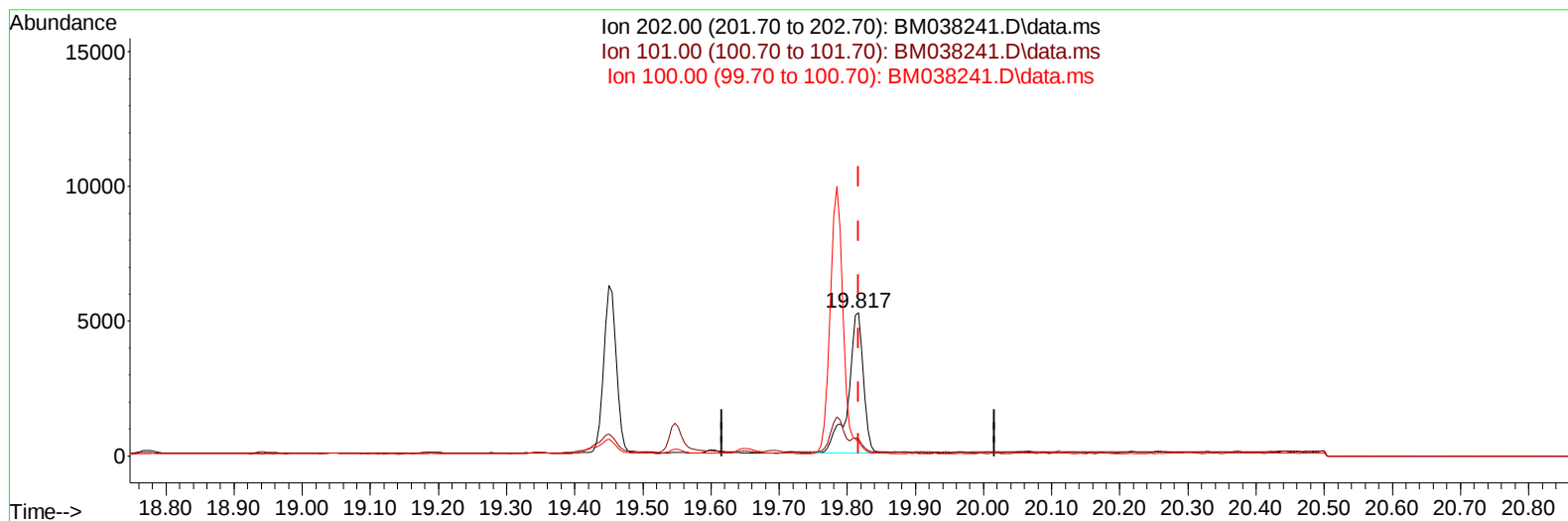
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122222\
Data File : BM038241.D
Acq On : 23 Dec 2022 20:21
Operator : CG/JU
Sample : N6014-05
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBXR6

Manual IntegrationsAPPROVED

Quant Time: Dec 23 23:24:39 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM122222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 22 23:36:45 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/24/2022
Supervised By :Yogesh Patel 12/24/2022



TIC: BM038241.D\data.ms

(20) Pyrene

19.817min (-0.000) 0.18 ng/u1

response 8330

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.10	11.58
100.00	9.20	10.08
0.00	0.00	0.00

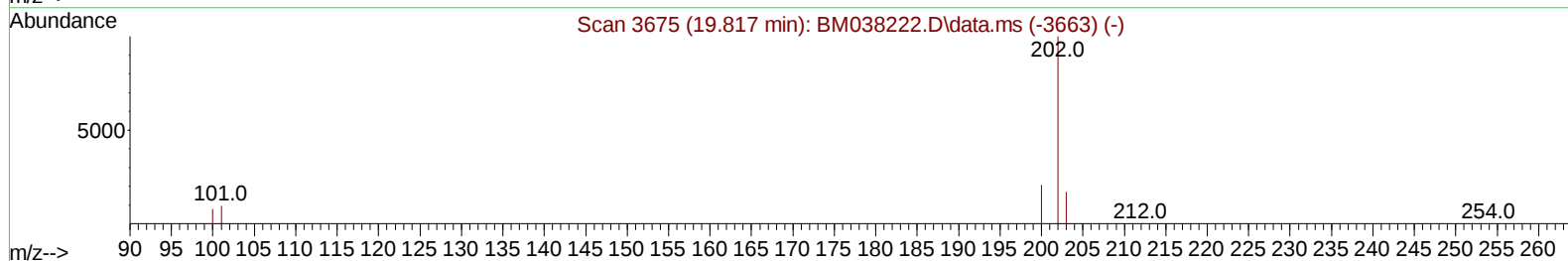
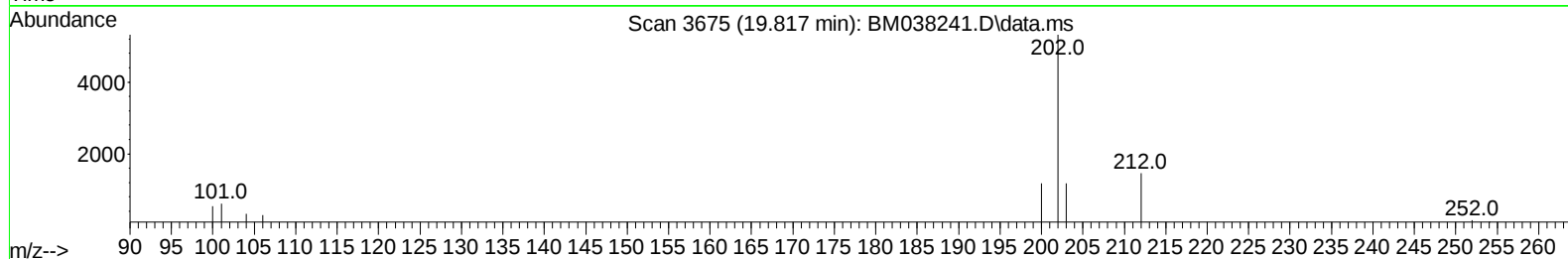
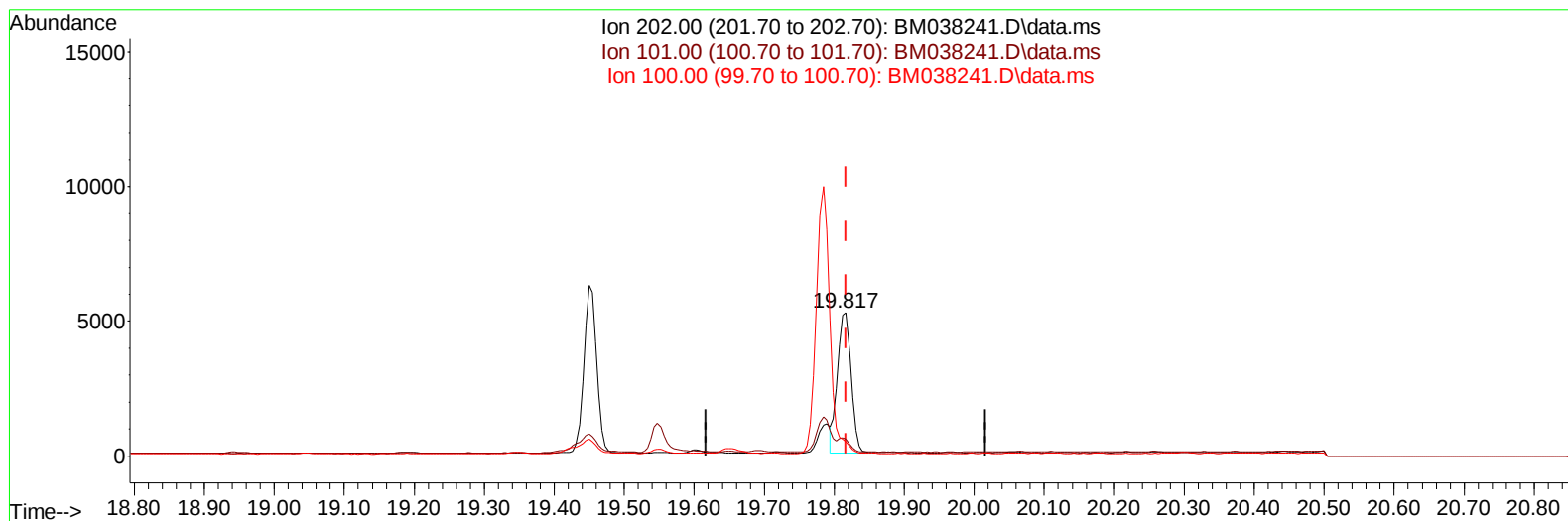
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TIC: BM038241.D\data.ms

(20) Pyrene

19.817min (-0.000) 0.16 ng/u1 m

response 6992

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.10	11.58
100.00	9.20	10.08
0.00	0.00	0.00

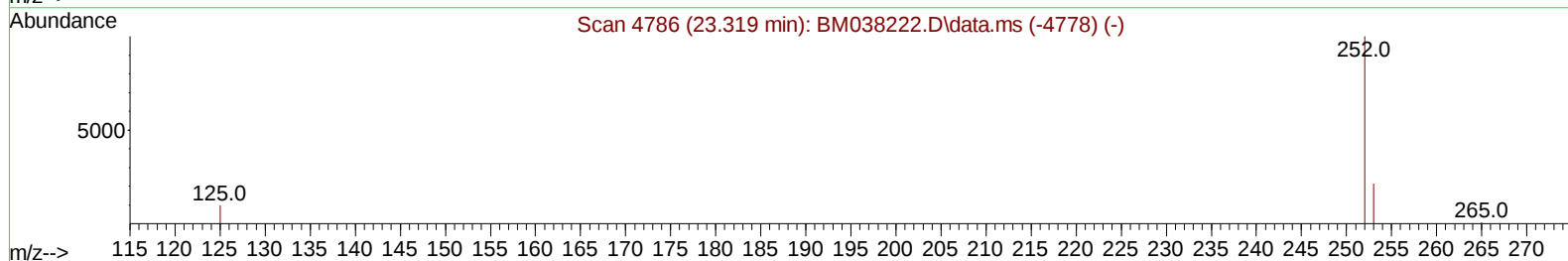
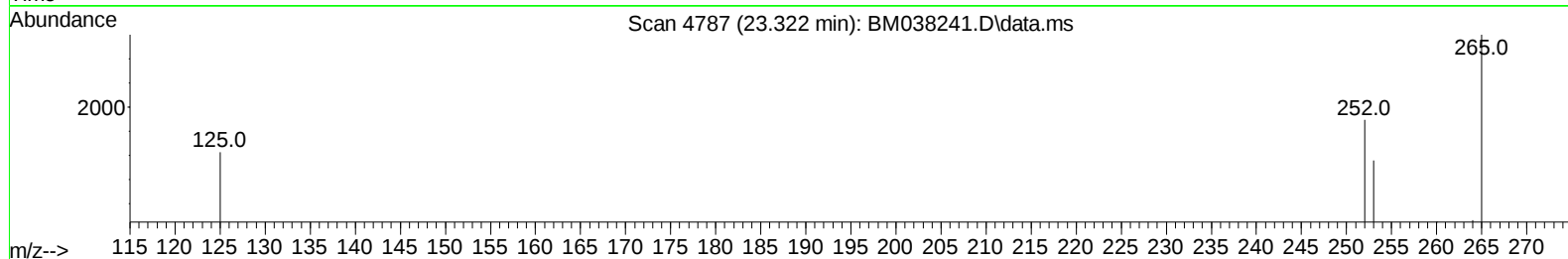
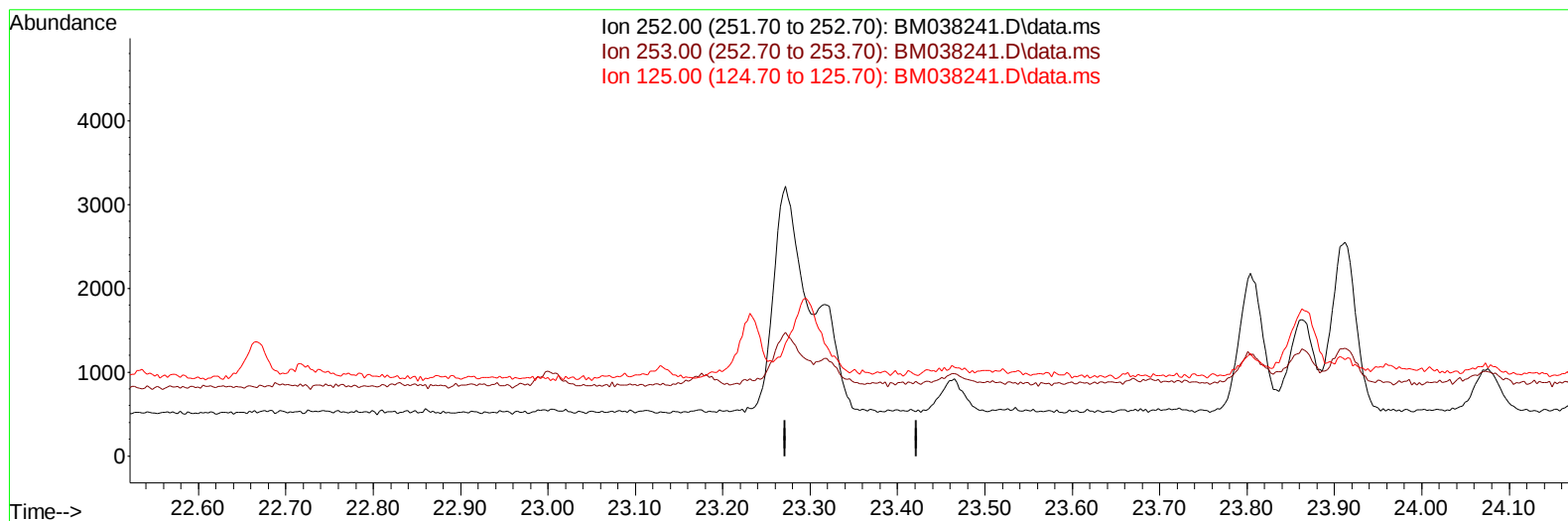
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Sample : N6014-05
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BM038241.D\data.ms

(25) Benzo(k)fluoranthene

23.322min (-23.322) 0.00 ng/u1

response 0

Ion	Exp%	Act%
252.00	100.00	0.00
253.00	28.10	0.00#
125.00	18.00	0.00#
0.00	0.00	0.00

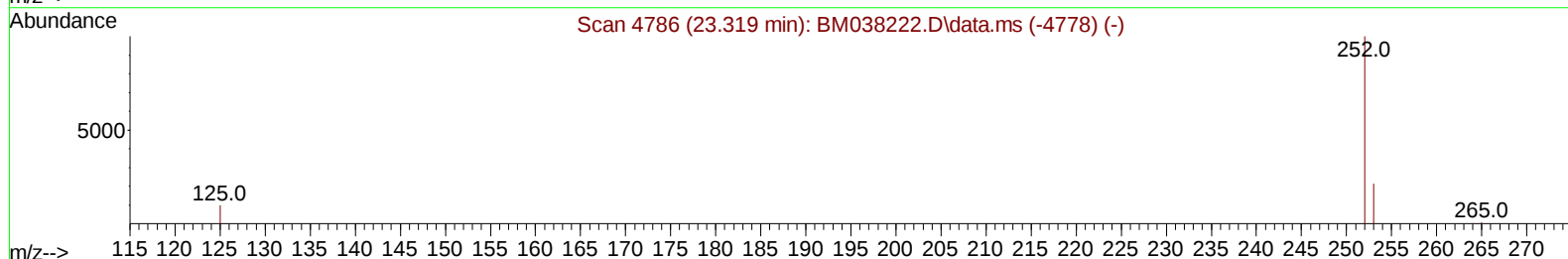
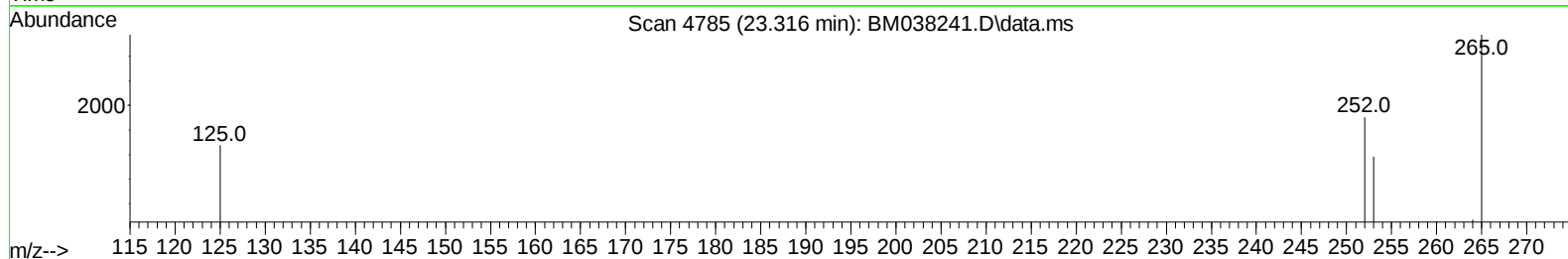
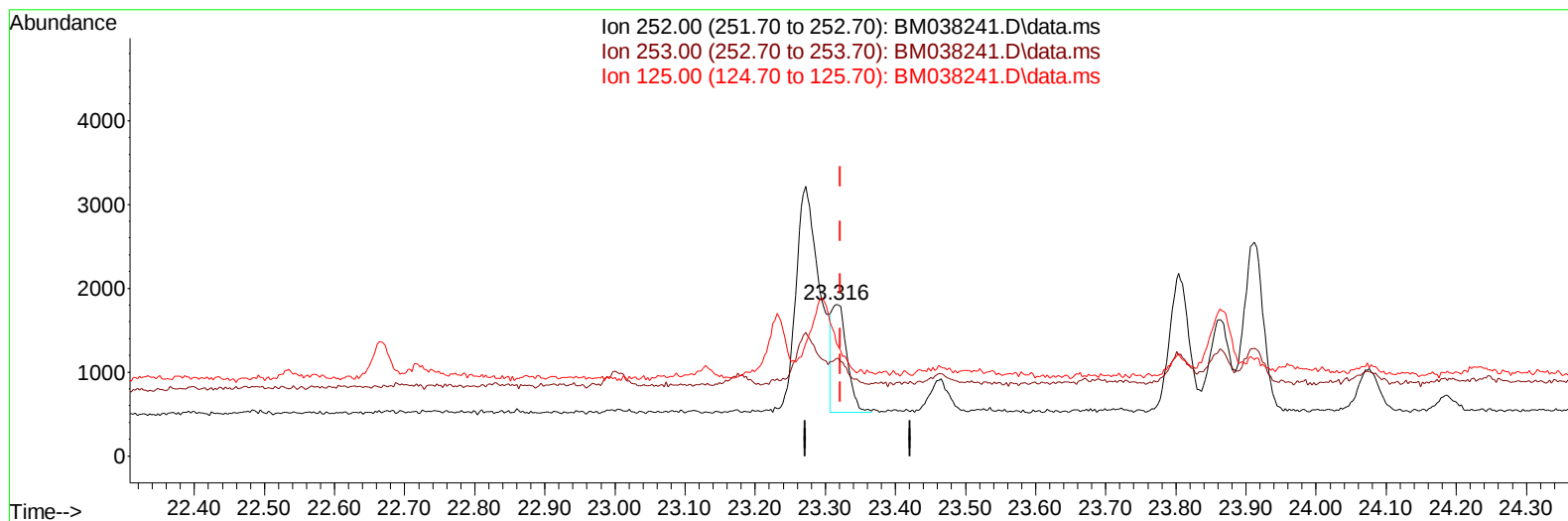
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 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
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TIC: BM038241.D\data.ms

(25) Benzo(k)fluoranthene

23.316min (-0.006) 0.05 ng/u1 m

response 1918

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	28.10	64.48#
125.00	18.00	74.48#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBXR6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/24/2022
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.034	152		4026	0.400	ng/ul	0.00
4) Naphthalene-d8	10.845	136		12075	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.661	164		6701	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.400	188		13115	0.400	ng/ul	0.00
17) Chrysene-d12	21.571	240		8147	0.400	ng/ul	0.00
23) Perylene-d12	24.023	264		8348	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.393	96		14625	3.571	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.429	152		4177	0.235	ng/ul	0.00
18) Fluoranthene-d10	19.422	212		7376	0.241	ng/ul	0.00
Target Compounds							
						Qvalue	
15) Phenanthrene	17.442	178		3183	0.073	ng/ul#	91
19) Fluoranthene	19.450	202		8029	0.180	ng/ul#	91
20) Pyrene	19.817	202		6992m	0.155	ng/ul	
21) Benzo(a)anthracene	21.553	228		3302	0.092	ng/ul#	88
22) Chrysene	21.609	228		4224	0.121	ng/ul#	91
24) Benzo(b)fluoranthene	23.272	252		6155	0.167	ng/ul#	59
25) Benzo(k)fluoranthene	23.316	252		1918m	0.054	ng/ul	
26) Benzo(a)pyrene	23.912	252		3858	0.120	ng/ul#	58
27) Indeno(1,2,3-cd)pyrene	26.582	276		3343	0.075	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.592	278		875	0.025	ng/ul#	1
29) Benzo(g,h,i)perylene	27.370	276		3451	0.092	ng/ul#	75

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

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