

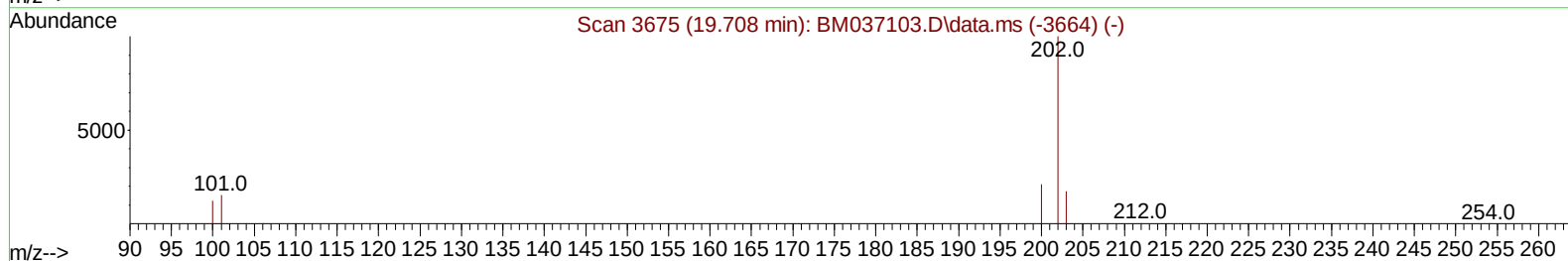
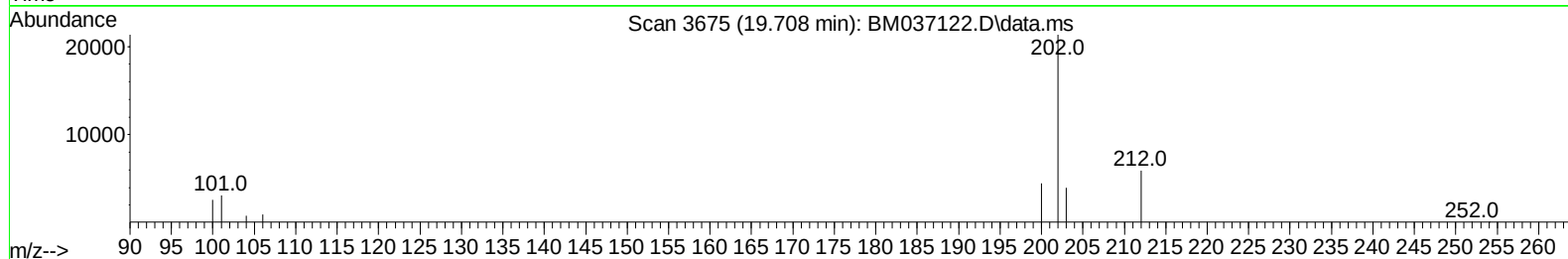
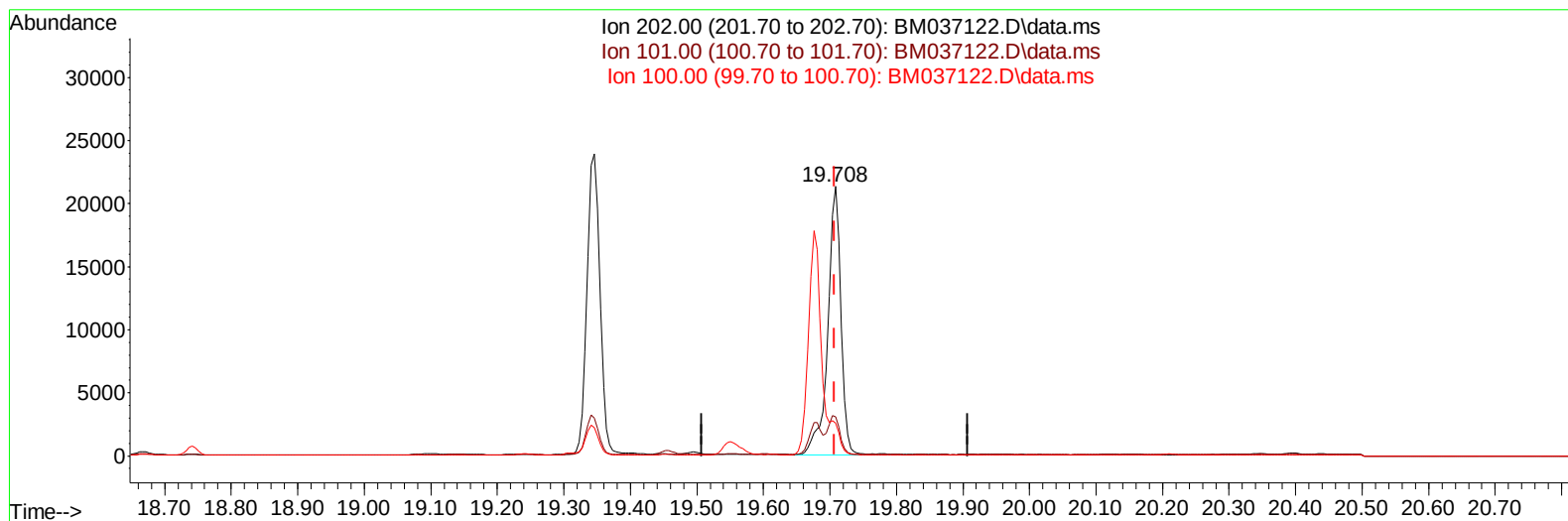
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101722\
Data File : BM037122.D
Acq On : 17 Oct 2022 21:12
Operator : CG/JU
Sample : N4984-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGNB6

Manual IntegrationsAPPROVED

Quant Time: Oct 17 23:56:12 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 17 02:26:13 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/18/2022
Supervised By :mohammad ahmed 10/19/2022



TIC: BM037122.D\data.ms

(20) Pyrene

19.708min (+ 0.001) 0.64 ng/u1

response 29406

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	15.00	14.62
100.00	12.20	12.23
0.00	0.00	0.00

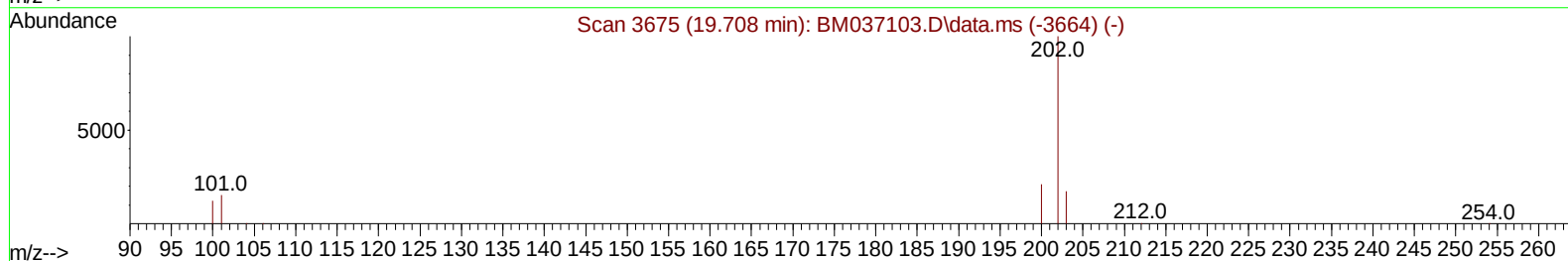
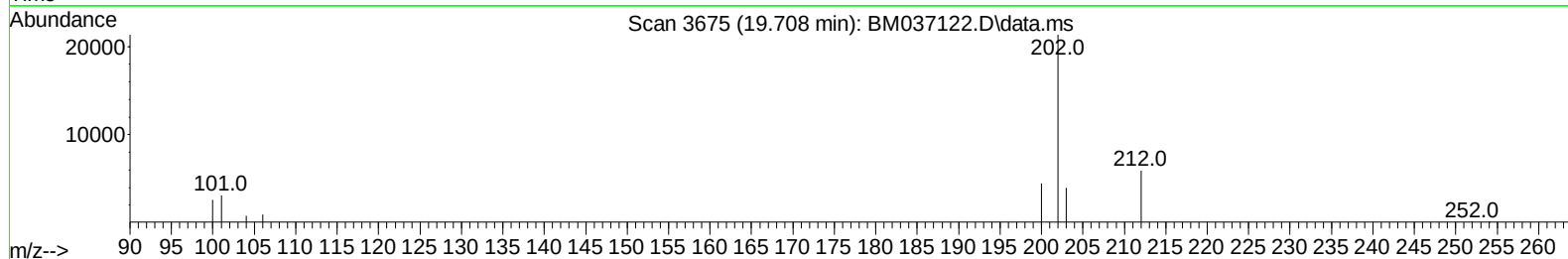
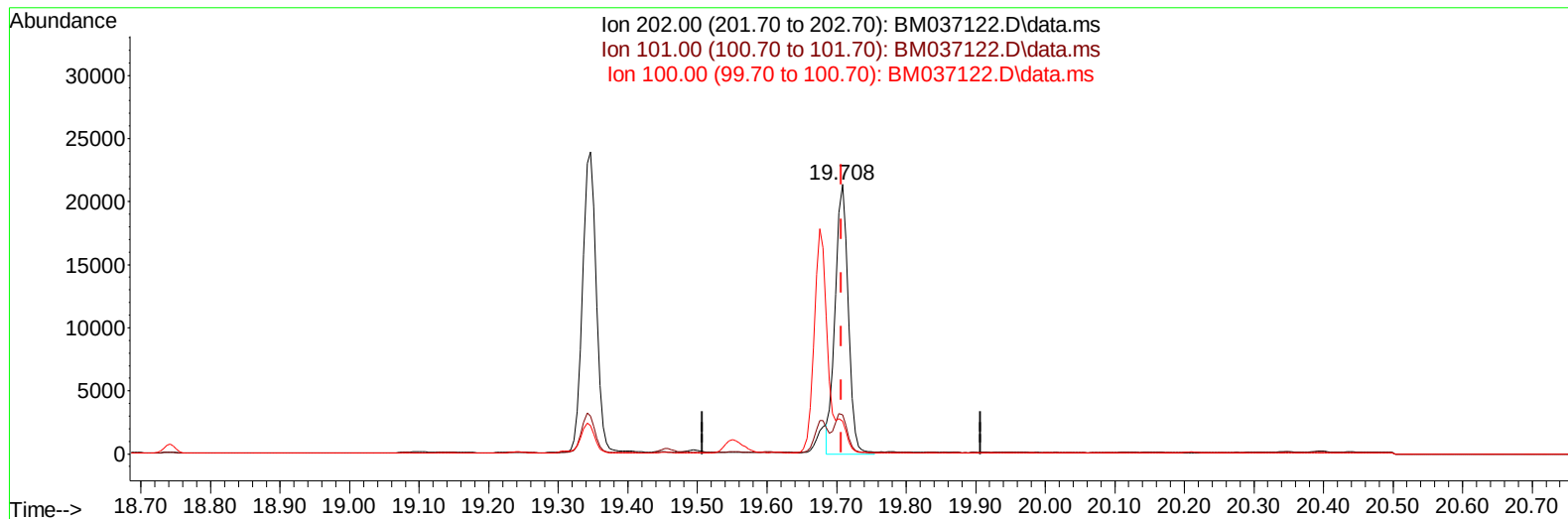
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TIC: BM037122.D\data.ms

(20) Pyrene

19.708min (+ 0.001) 0.60 ng/ul m

response 27467

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	15.00	14.62
100.00	12.20	12.23
0.00	0.00	0.00

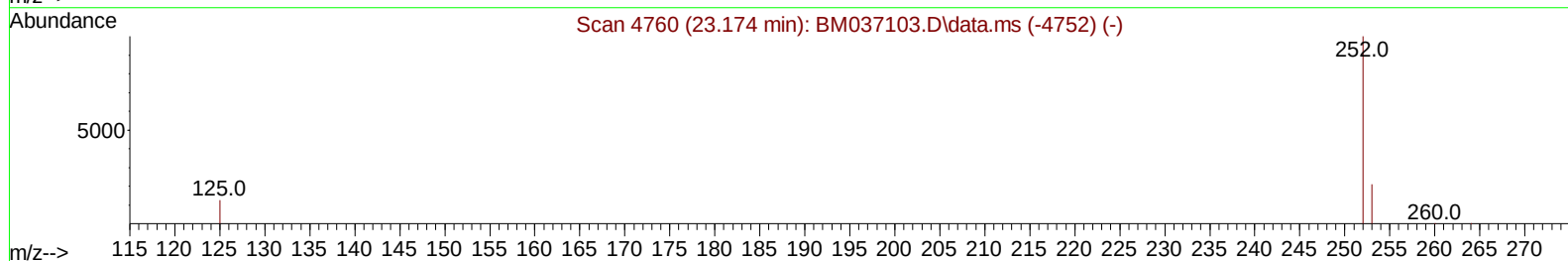
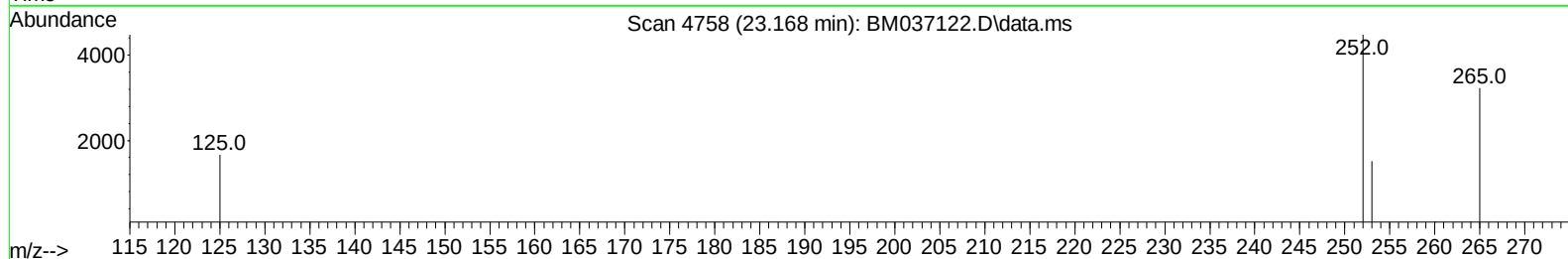
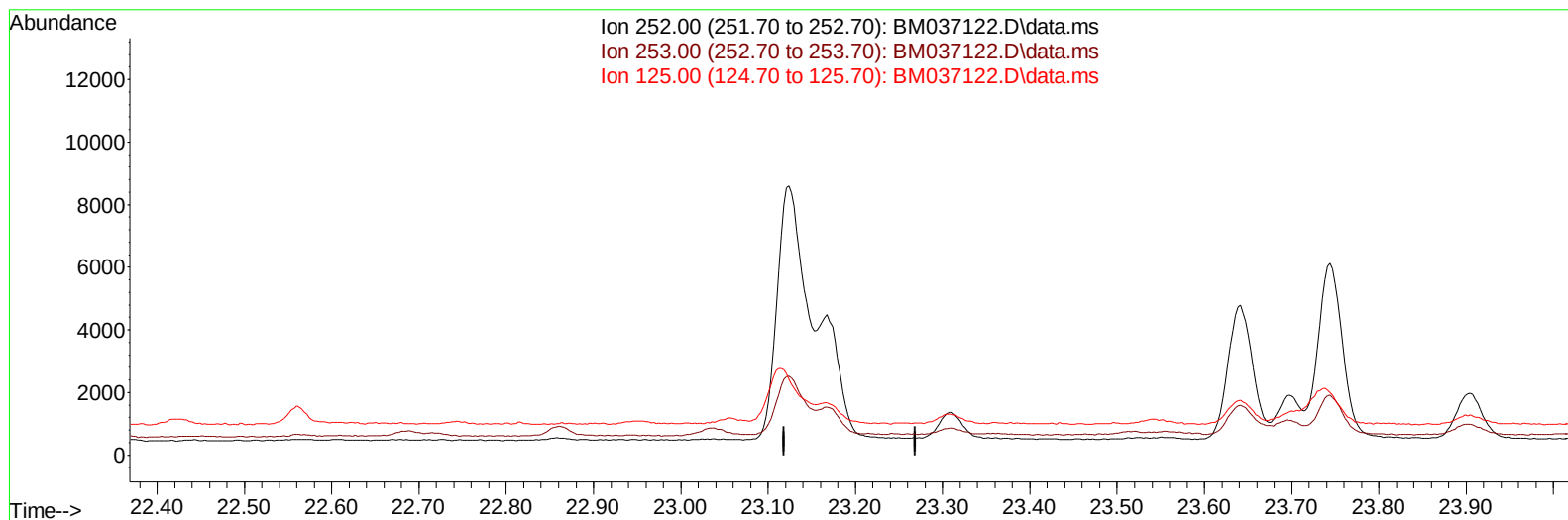
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TIC: BM037122.D\data.ms

(25) Benzo(k)fluoranthene

23.169min (-23.169) 0.00 ng/u1

response 0

Ion	Exp%	Act%
252.00	100.00	0.00
253.00	31.50	0.00#
125.00	22.90	0.00#
0.00	0.00	0.00

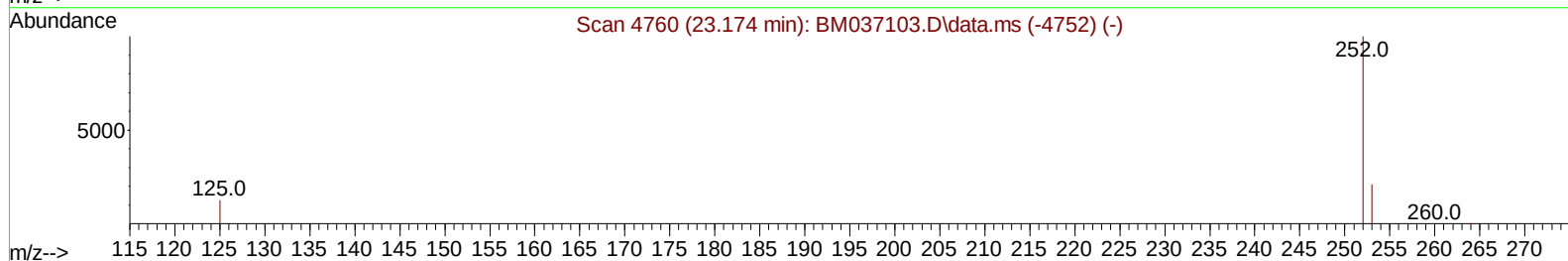
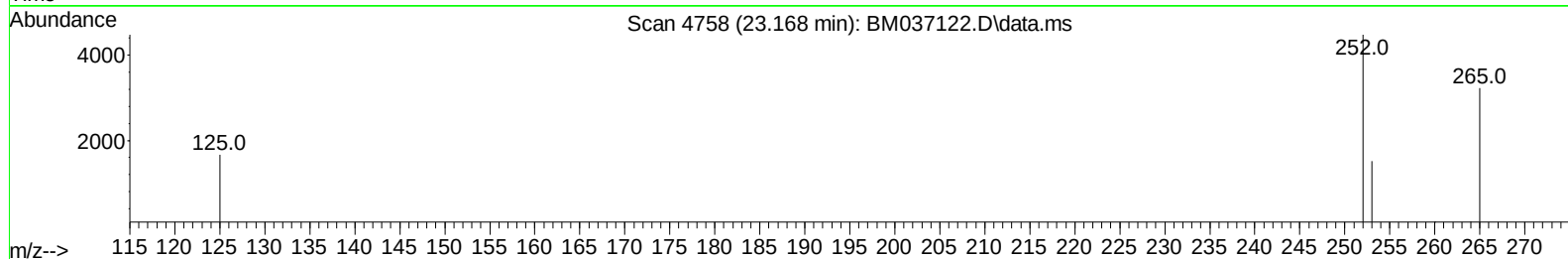
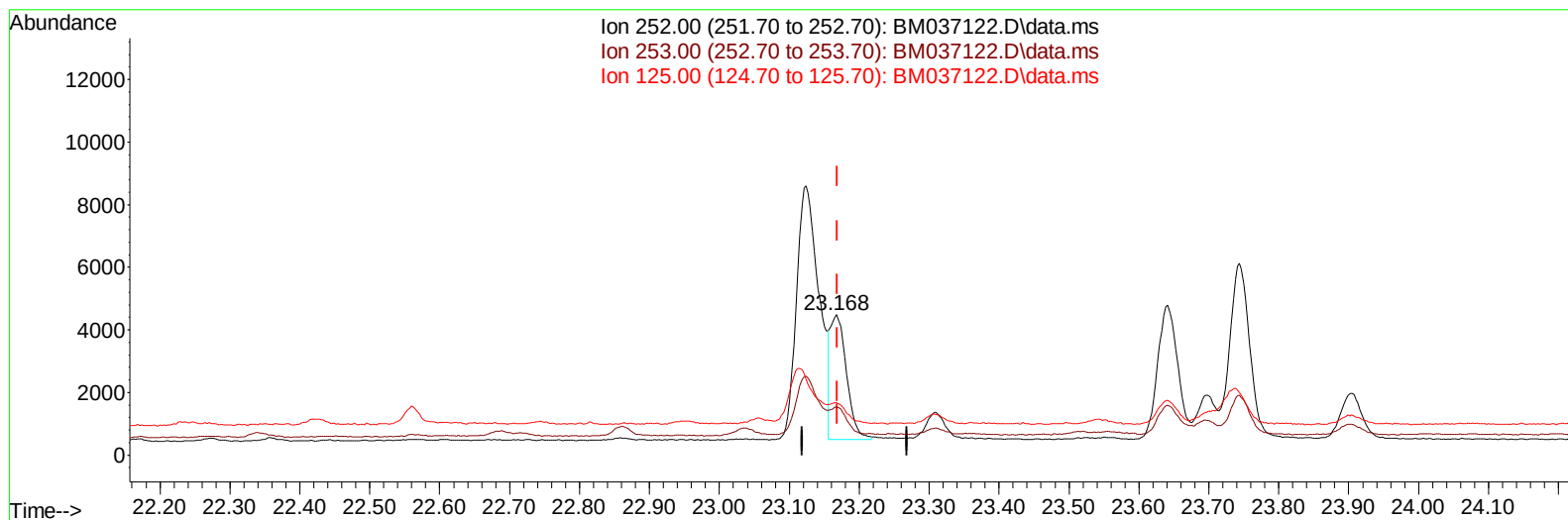
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TIC: BM037122.D\data.ms

(25) Benzo(k)fluoranthene

23.168min (-0.001) 0.15 ng/ul m

response 6303

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	31.50	34.02
125.00	22.90	37.30#
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.926	152		4566	0.400	ng/ul	0.00
4) Naphthalene-d8	10.726	136		16878	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.553	164		8756	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.293	188		17150	0.400	ng/ul	0.00
17) Chrysene-d12	21.467	240		11025	0.400	ng/ul	0.00
23) Perylene-d12	23.849	264		9855	0.400	ng/ul	# 0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.335	96		30259	4.699	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.316	152		6251	0.245	ng/ul	0.00
18) Fluoranthene-d10	19.313	212		11403	0.346	ng/ul	0.00
Target Compounds							
							Qvalue
15) Phenanthrene	17.335	178		12840	0.235	ng/ul	99
16) Anthracene	17.424	178		2504	0.055	ng/ul#	90
19) Fluoranthene	19.346	202		32305	0.737	ng/ul	99
20) Pyrene	19.708	202		27467m	0.601	ng/ul	
21) Benzo(a)anthracene	21.449	228		10647	0.267	ng/ul	98
22) Chrysene	21.502	228		13935	0.326	ng/ul	99
24) Benzo(b)fluoranthene	23.124	252		18337	0.417	ng/ul	93
25) Benzo(k)fluoranthene	23.168	252		6303m	0.147	ng/ul	
26) Benzo(a)pyrene	23.744	252		10904	0.305	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	26.325	276		8621	0.170	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.332	278		2108	0.052	ng/ul#	66
29) Benzo(g,h,i)perylene	27.087	276		7384	0.166	ng/ul	97

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

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