

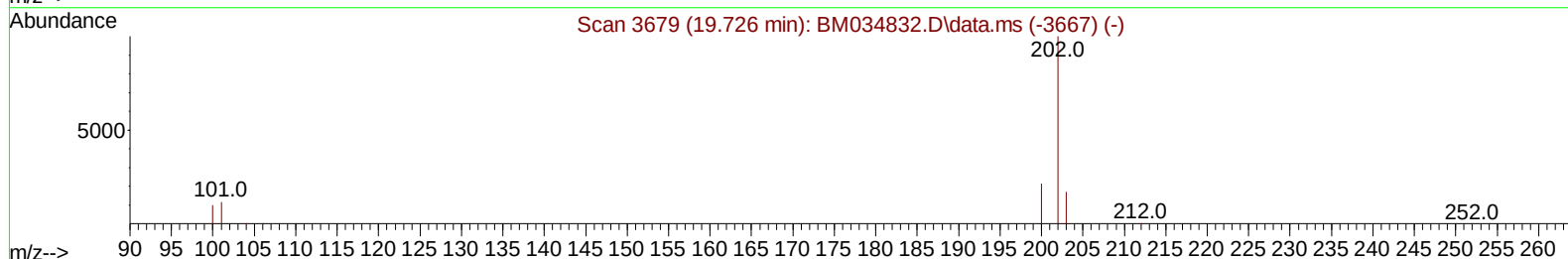
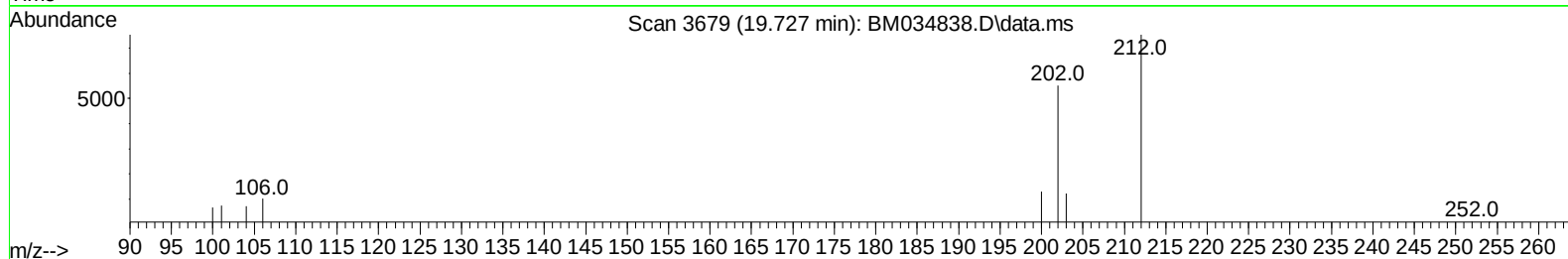
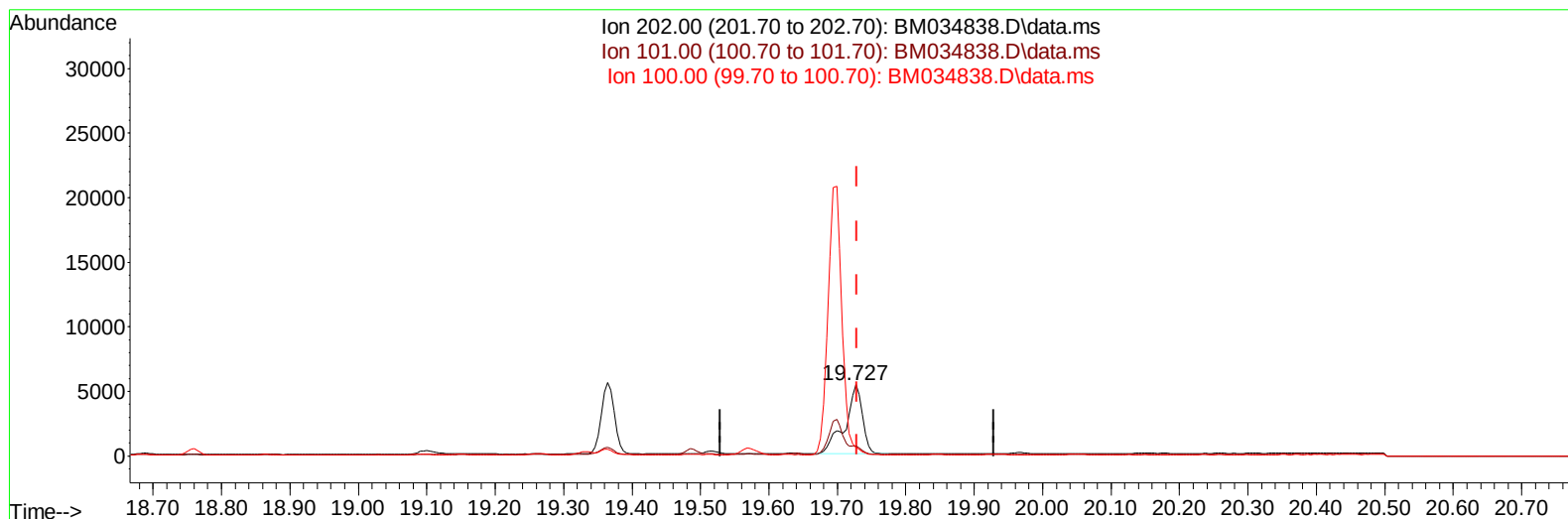
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042622\
Data File : BM034838.D
Acq On : 26 Apr 2022 16:04
Operator : CG/JU
Sample : N2475-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A0BN4

Manual IntegrationsAPPROVED

Quant Time: Apr 26 17:25:11 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 22 16:54:34 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/27/2022
Supervised By :Yogesh Patel 05/03/2022



TIC: BM034838.D\data.ms

(20) Pyrene

19.727min (-0.002) 0.17 ng/u1

response 9270

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.60	13.63
100.00	9.70	12.27#
0.00	0.00	0.00

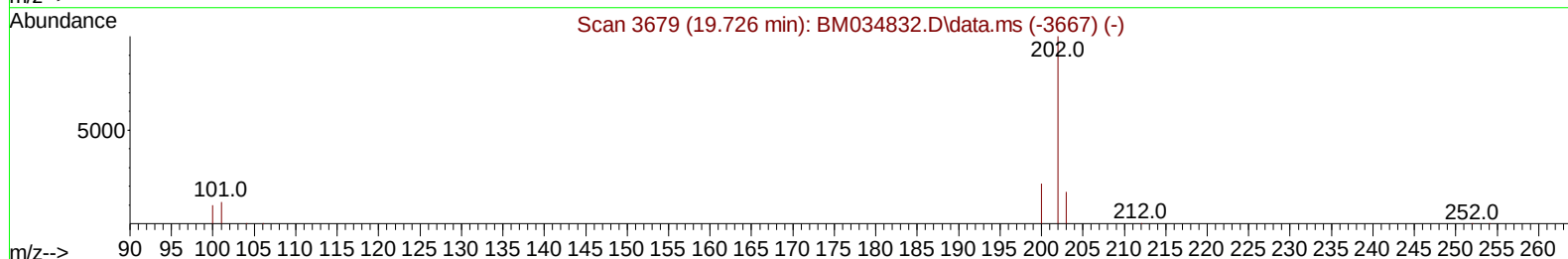
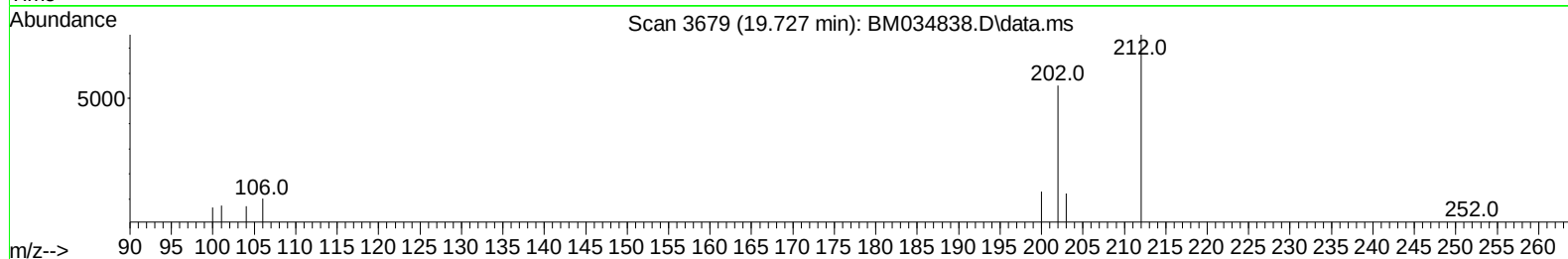
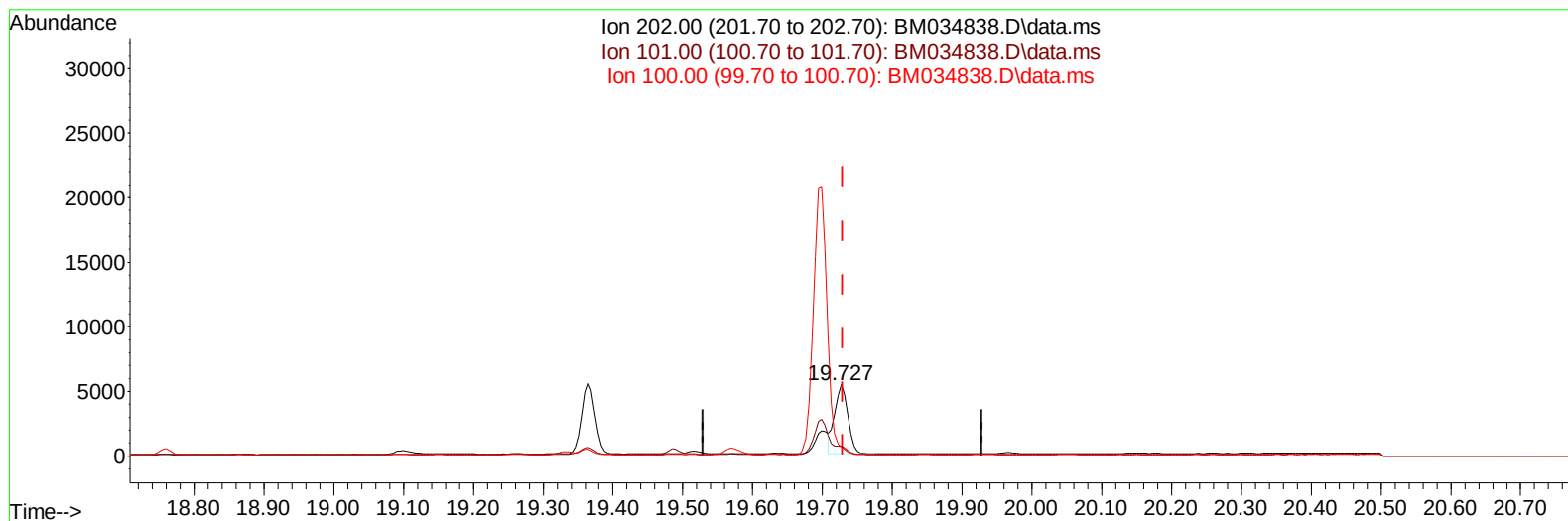
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(20) Pyrene

19.727min (-0.002) 0.13 ng/ul m

response 6941

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.60	13.63
100.00	9.70	12.27#
0.00	0.00	0.00

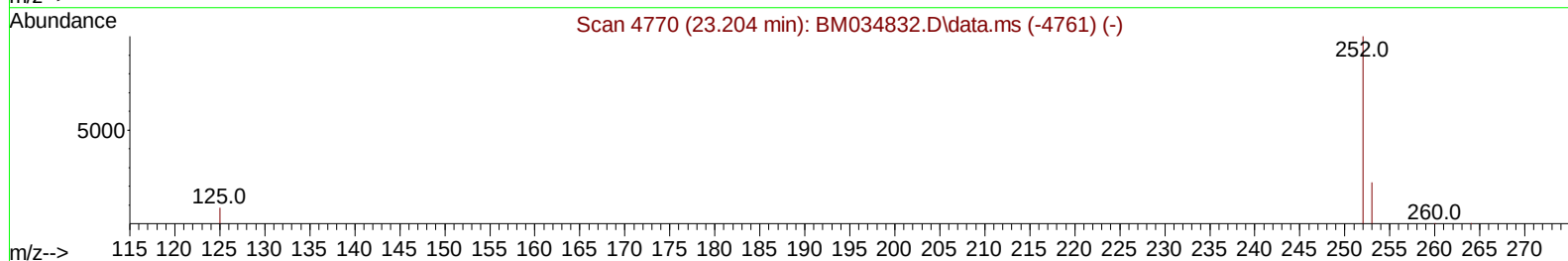
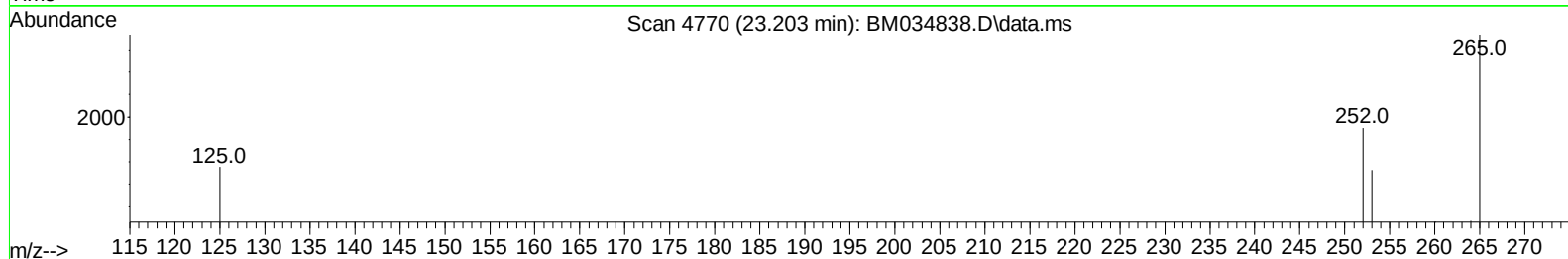
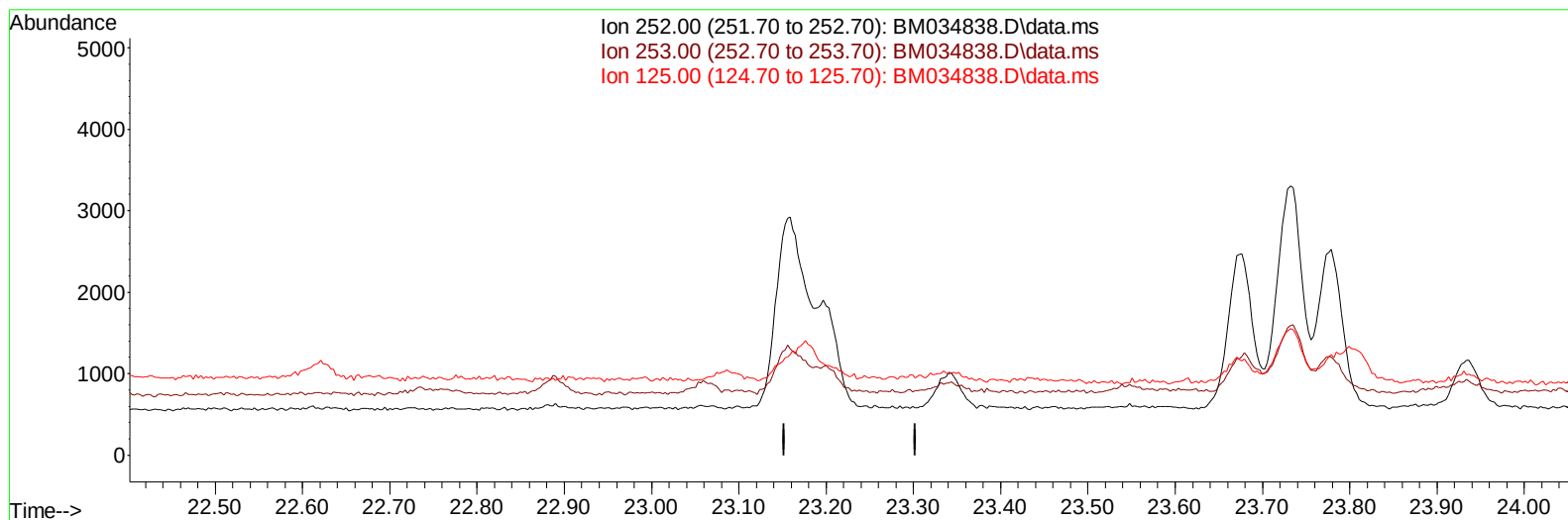
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(25) Benzo(k)fluoranthene

23.202min (-23.202) 0.00 ng/u1

response 0

Ion	Exp%	Act%
252.00	100.00	0.00
253.00	33.50	0.00#
125.00	15.00	0.00#
0.00	0.00	0.00

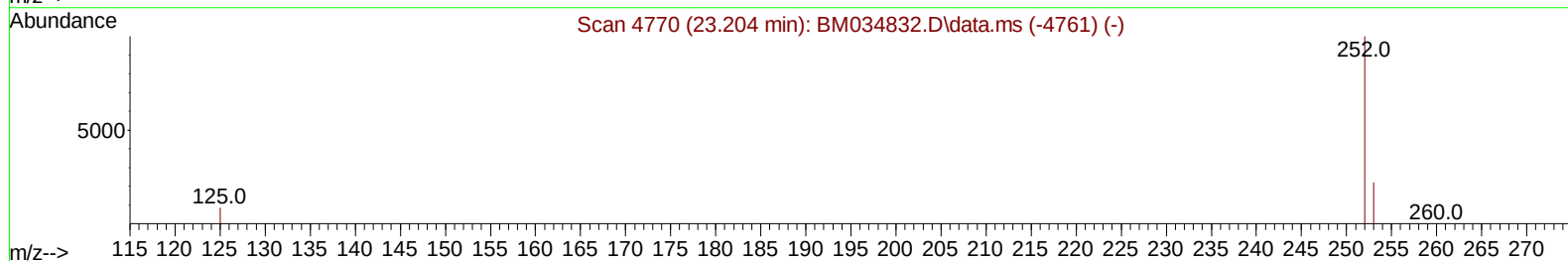
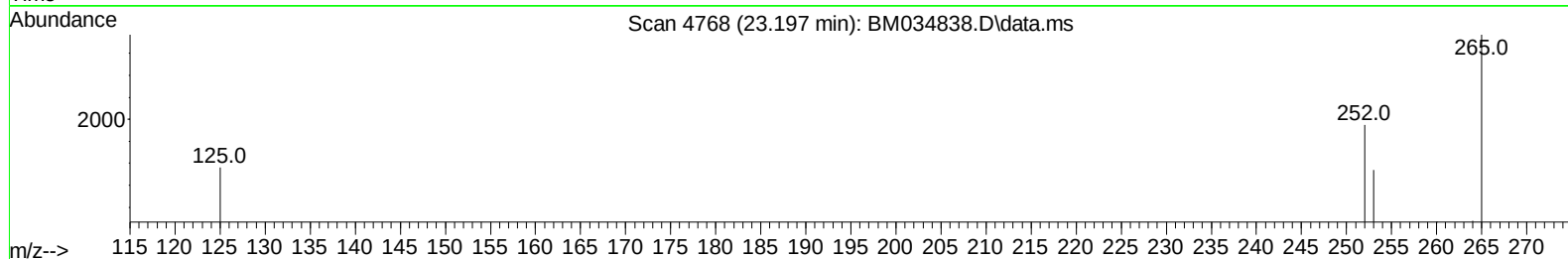
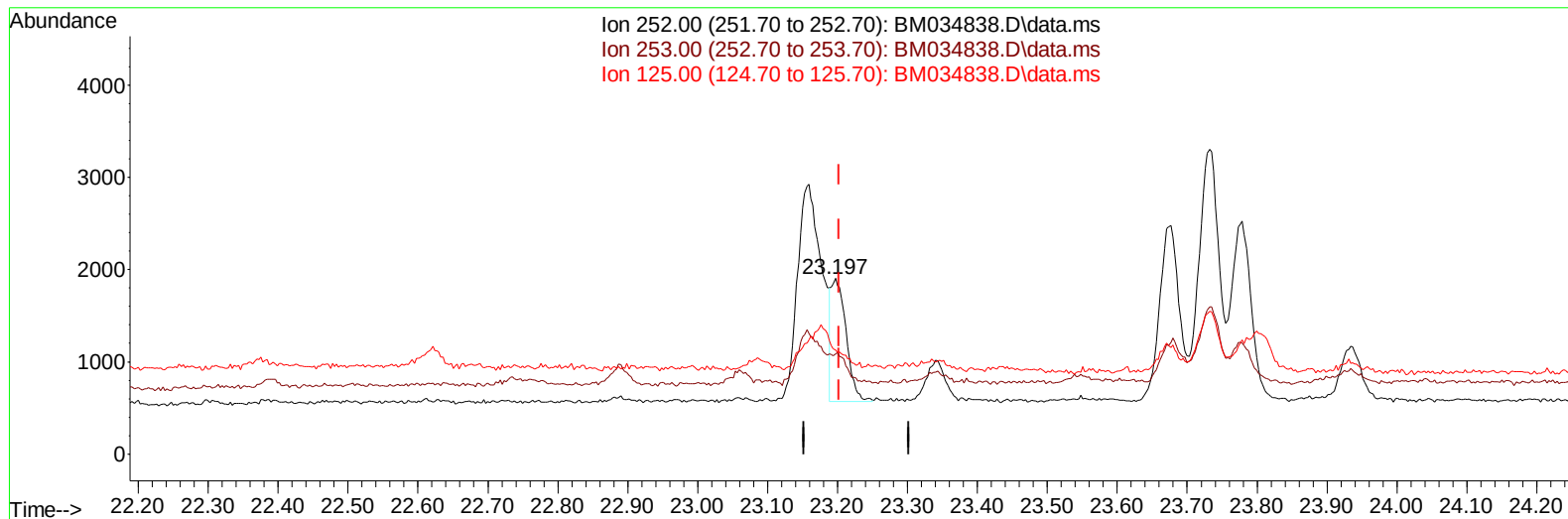
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(25) Benzo(k)fluoranthene

23.197min (-0.005) 0.03 ng/ul m

response 1978

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	33.50	56.91#
125.00	15.00	59.17#
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.942	152		5690	0.400	ng/ul	0.00
4) Naphthalene-d8	10.743	136		14820	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.571	164		8176	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.310	188		15099	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.487	240		11500	0.400	ng/ul	# 0.00
23) Perylene-d12	23.884	264		12548	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.382	96		32103	5.101	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152		8358	0.401	ng/ul	0.00
18) Fluoranthene-d10	19.336	212		16490	0.455	ng/ul	0.00
Target Compounds							
							Qvalue
15) Phenanthrene	17.352	178		3583	0.072	ng/ul	97
16) Anthracene	17.445	178		2245	0.050	ng/ul	97
19) Fluoranthene	19.364	202		7372	0.134	ng/ul	99
20) Pyrene	19.727	202		6941m	0.125	ng/ul	
21) Benzo(a)anthracene	21.470	228		3561	0.080	ng/ul#	89
22) Chrysene	21.525	228		3672	0.077	ng/ul#	88
24) Benzo(b)fluoranthene	23.159	252		5642	0.098	ng/ul#	63
25) Benzo(k)fluoranthene	23.197	252		1978m	0.035	ng/ul	
26) Benzo(a)pyrene	23.779	252		3649	0.078	ng/ul#	70
27) Indeno(1,2,3-cd)pyrene	26.355	276		4103	0.063	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.375	278		1110	0.021	ng/ul#	1
29) Benzo(g,h,i)perylene	27.127	276		4871	0.086	ng/ul#	86

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

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