

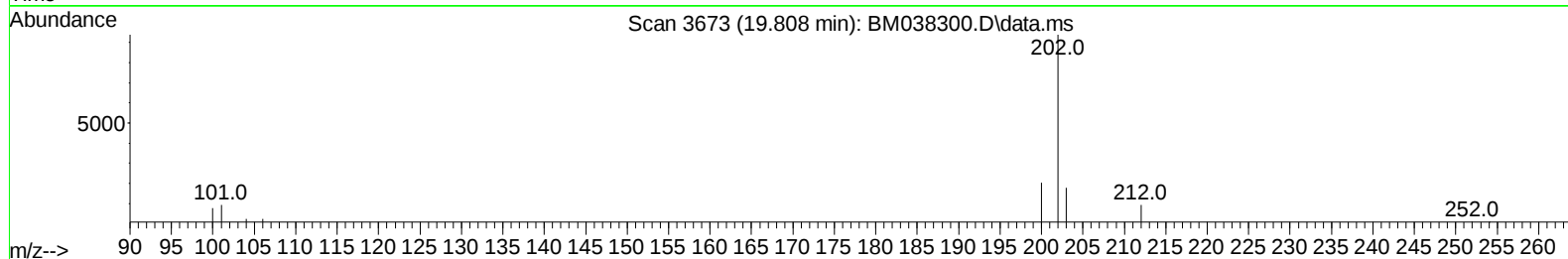
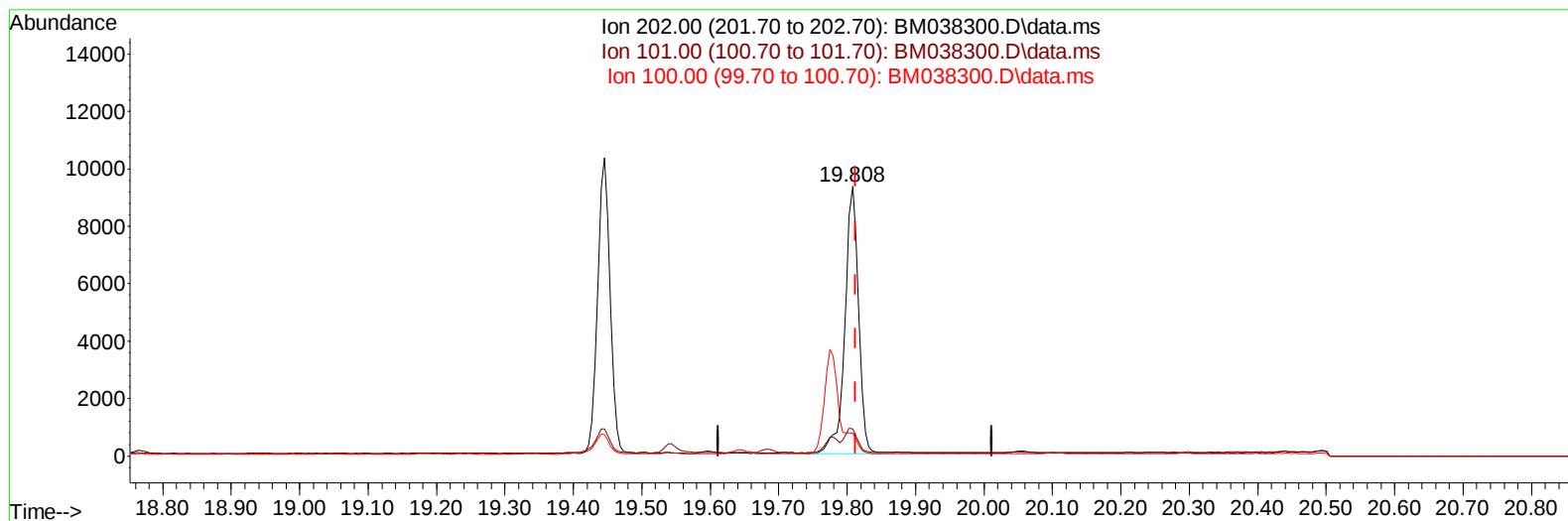
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122222\
Data File : BM038300.D
Acq On : 25 Dec 2022 09:28
Operator : CG/JU
Sample : N6018-16DL 5X
Misc :
ALS Vial : 109 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBYK1DL

Manual IntegrationsAPPROVED

Quant Time: Dec 25 23:54:53 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 : Sat Dec 24 05:23:31 2022
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/26/2022
Supervised By :Sohil Jodhani 12/26/2022



TIC: BM038300.D\data.ms

(20) Pyrene

19.808min (-0.005) 0.24 ng/u1

response 12552

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.10	10.03
100.00	9.20	8.30
0.00	0.00	0.00

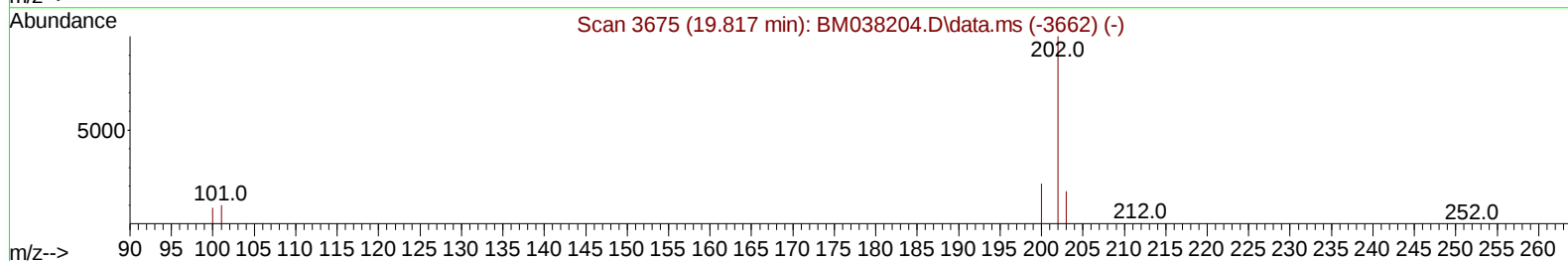
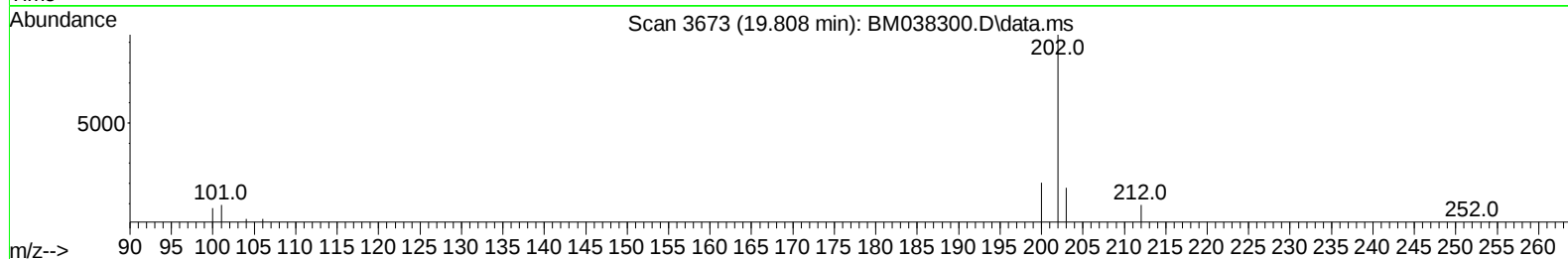
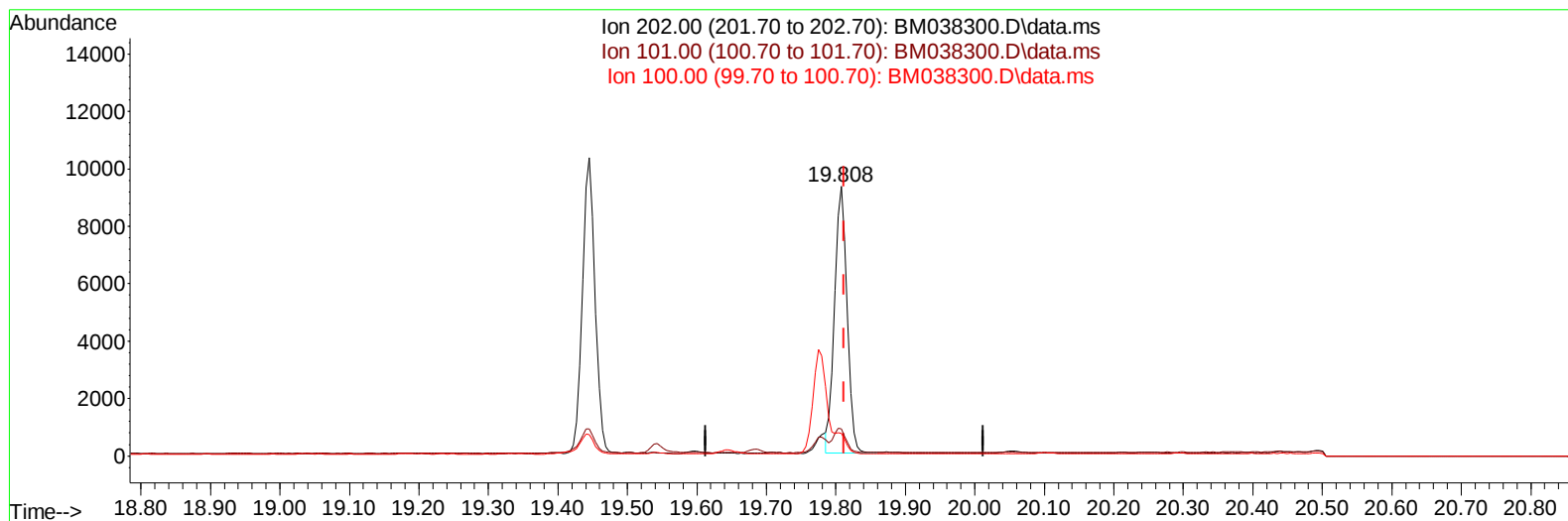
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TIC: BM038300.D\data.ms

(20) Pyrene

19.808min (-0.005) 0.22 ng/ul m

response 11843

Ion	Exp%	Act%
202.00	100.00	100.00
101.00	11.10	10.03
100.00	9.20	8.30
0.00	0.00	0.00

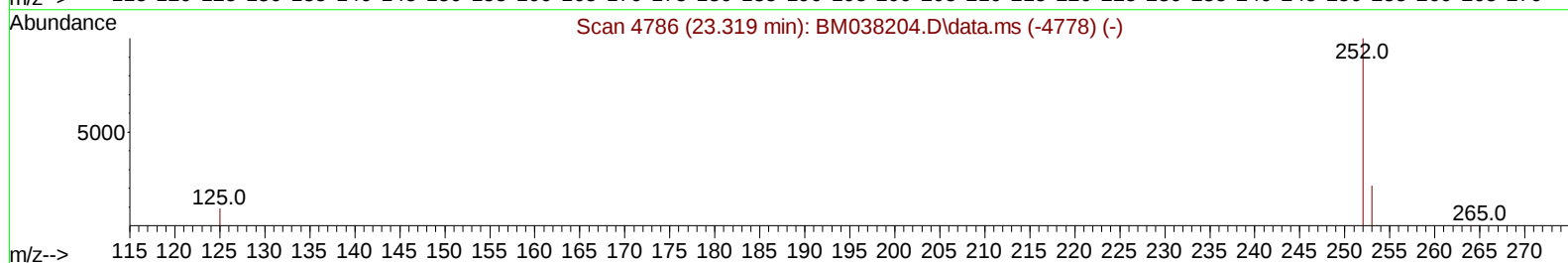
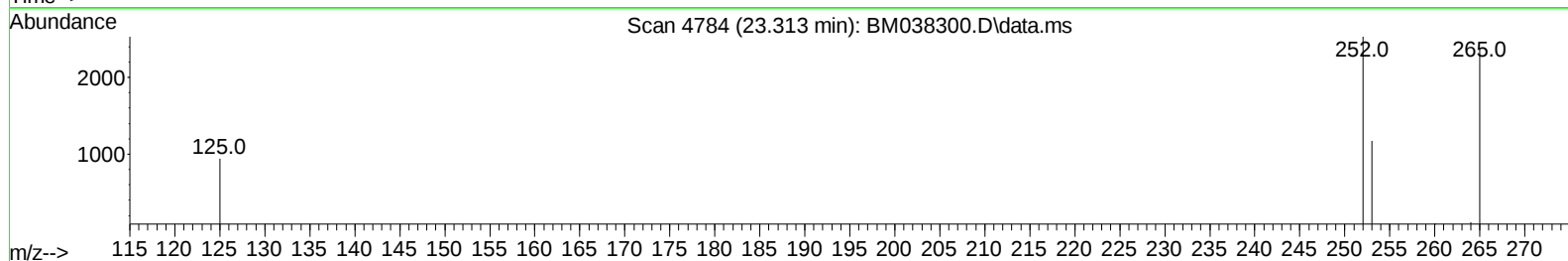
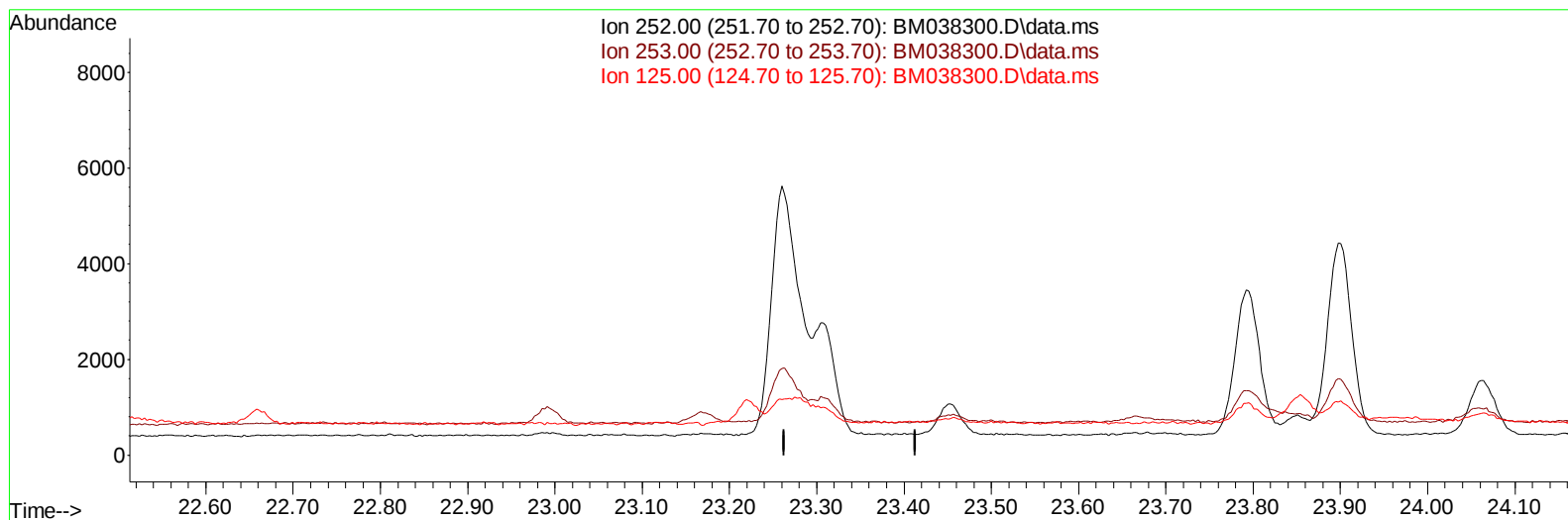
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TIC: BM038300.D\data.ms

(25) Benzo(k)fluoranthene

23.313min (-23.313) 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

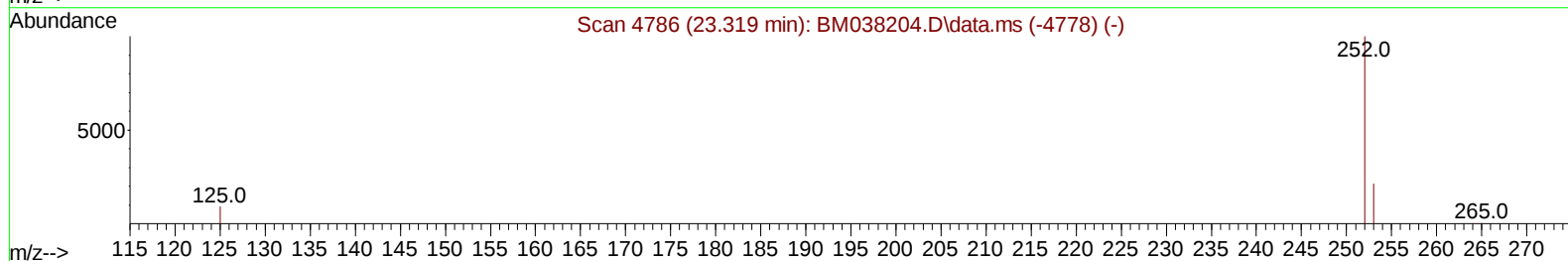
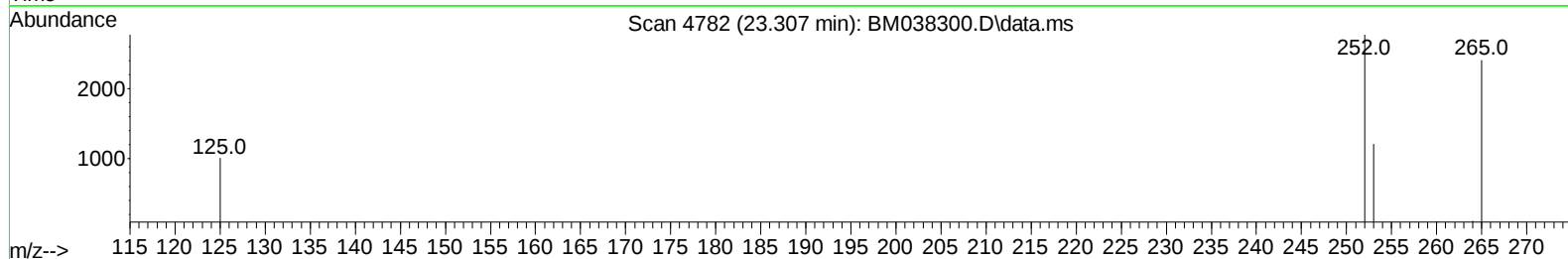
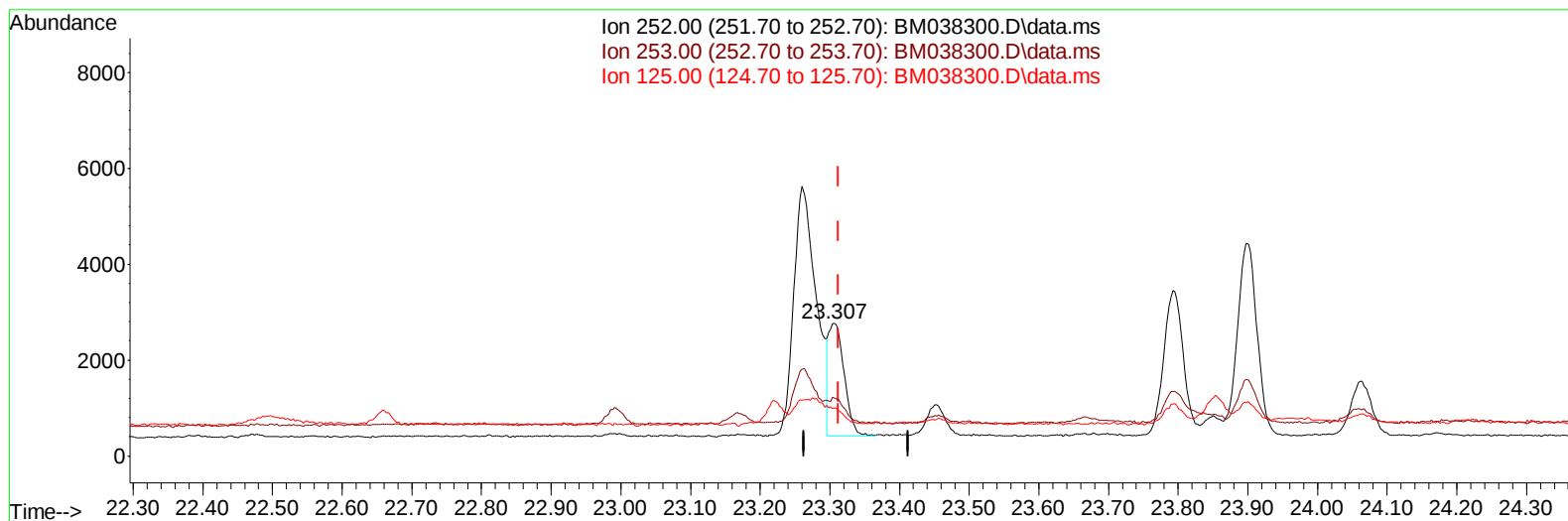
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 Misc :
 ALS Vial : 109 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBYK1DL

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TIC: BM038300.D\data.ms

(25) Benzo(k)fluoranthene

23.307min (-0.006) 0.10 ng/u1 m

response 3648

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	28.10	43.43#
125.00	18.00	36.17#
0.00	0.00	0.00

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 ALS Vial : 109 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 DBYK1DL

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/26/2022
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.026	152		3510	0.400	ng/ul	0.00
4) Naphthalene-d8	10.840	136		10534	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.652	164		6409	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.392	188		12710	0.400	ng/ul	0.00
17) Chrysene-d12	21.565	240		9527	0.400	ng/ul	0.00
23) Perylene-d12	24.009	264		8938	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.389	96		3676	1.030	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.418	152		1144	0.074	ng/ul	0.00
18) Fluoranthene-d10	19.413	212		2391	0.067	ng/ul	0.00
Target Compounds							
						Qvalue	
5) Naphthalene	10.889	128		650	0.021	ng/ul#	71
7) 2-Methylnaphthalene	12.489	142		385	0.020	ng/ul	98
15) Phenanthrene	17.434	178		5940	0.140	ng/ul	97
16) Anthracene	17.527	178		1014	0.025	ng/ul#	74
19) Fluoranthene	19.445	202		13025	0.250	ng/ul	99
20) Pyrene	19.808	202		11843m	0.225	ng/ul	
21) Benzo(a)anthracene	21.548	228		6065	0.145	ng/ul	94
22) Chrysene	21.603	228		7213	0.177	ng/ul	95
24) Benzo(b)fluoranthene	23.260	252		11639	0.294	ng/ul	93
25) Benzo(k)fluoranthene	23.307	252		3648m	0.096	ng/ul	
26) Benzo(a)pyrene	23.898	252		7819	0.227	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	26.559	276		6314	0.132	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.562	278		1461	0.039	ng/ul#	18
29) Benzo(g,h,i)perylene	27.350	276		6533	0.162	ng/ul	94

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

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