

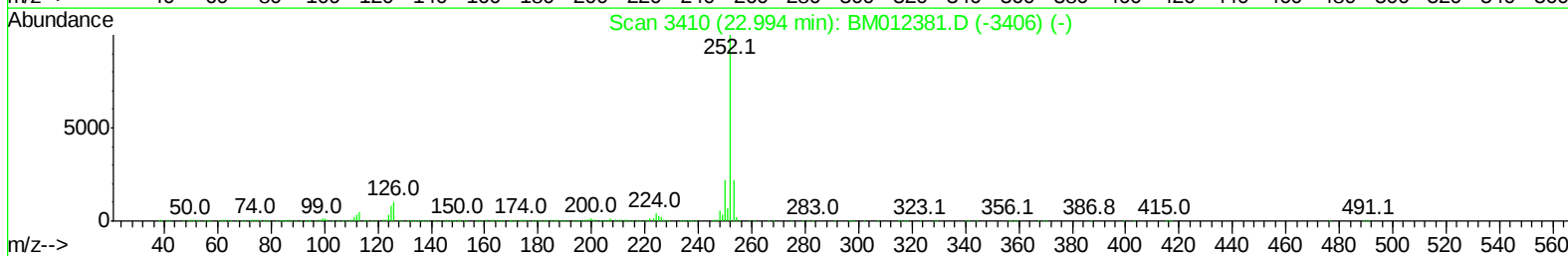
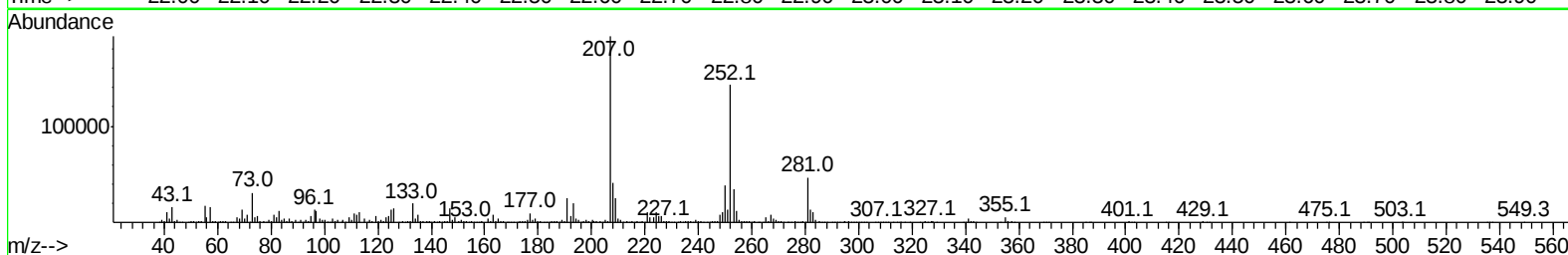
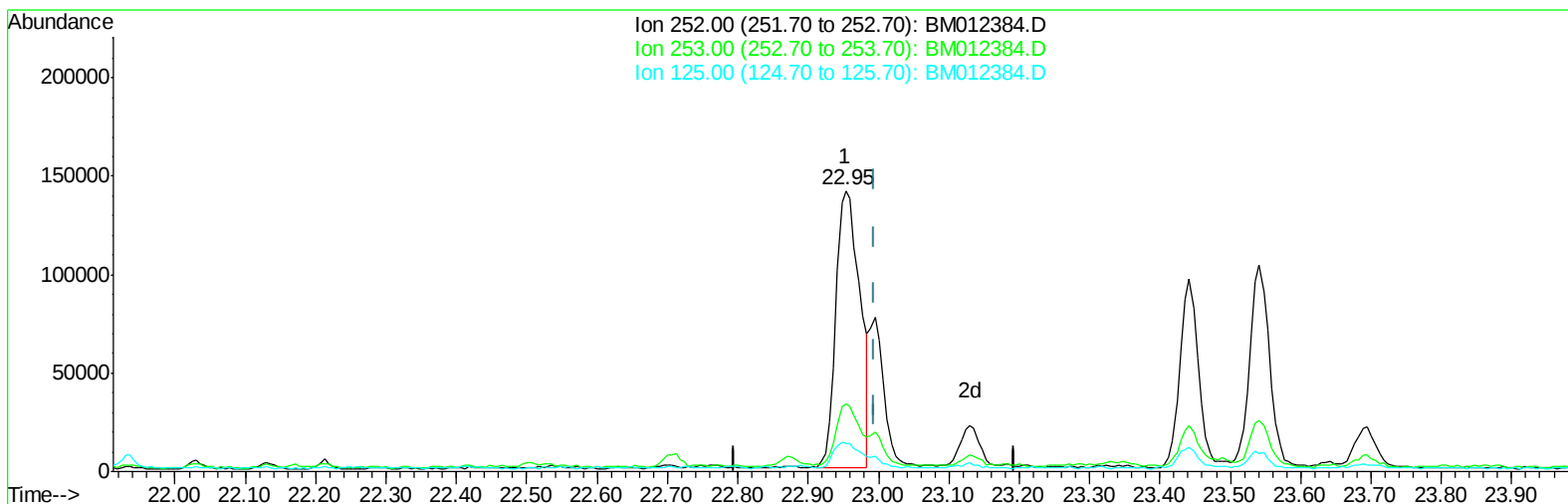
Data Path : Z:\HPCHEM1\BNA M\DATA\BM102117\
Data File : BM012384.D
Acq On : 22 Oct 2017 06:55
Operator : SJ/JU
Sample : I5771-03
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B0CQ0

Manual Integrations
APPROVED

Sohil
10/23/2017 4:00:31 PM

Quant Time: Oct 23 08:34:39 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 23 08:22:49 2017
Response via : Initial Calibration



TIC: BM012384.D

(86) Benzo(k)fluoranthene

22.953min (-0.041) 4.85ng/ul

response 334339

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	24.02
125.00	6.10	9.81#
0.00	0.00	0.00

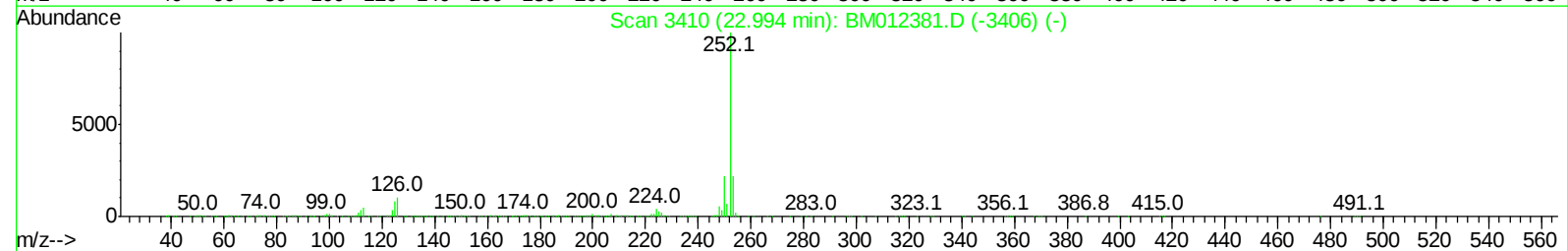
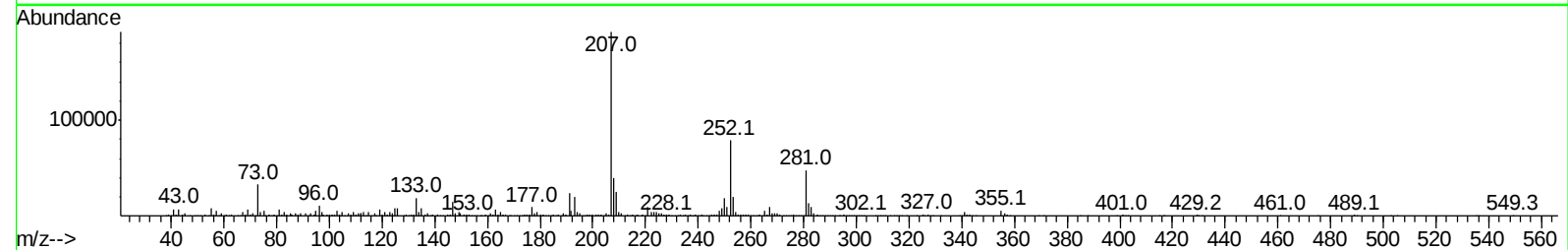
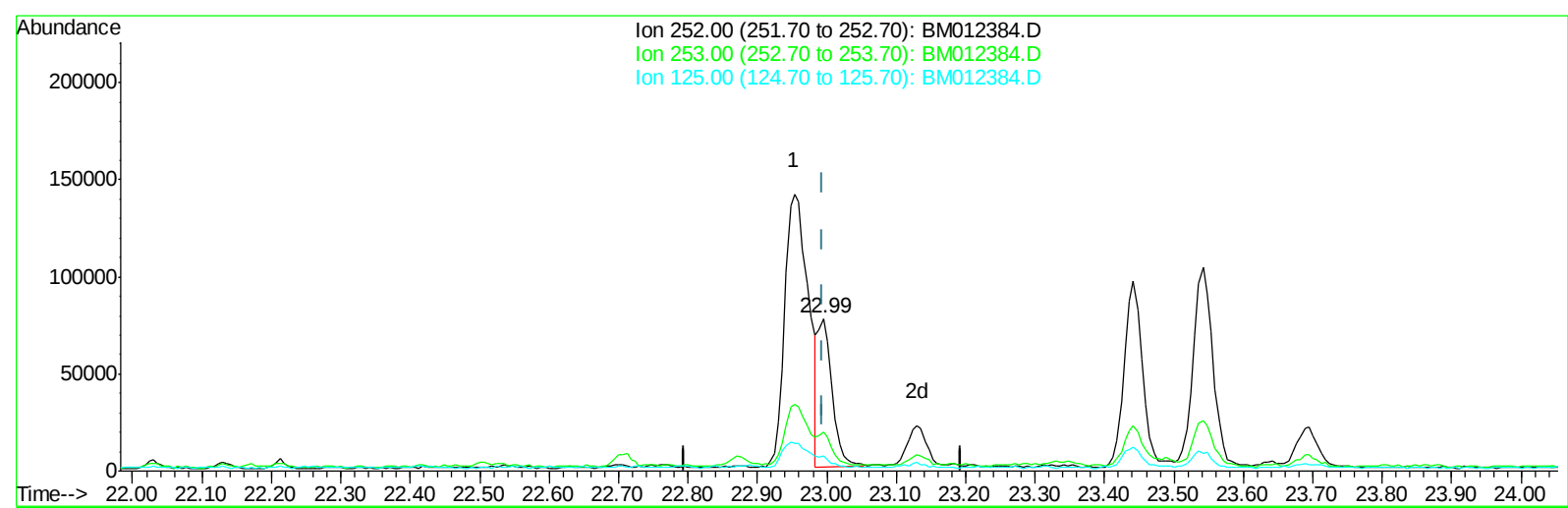
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(86) Benzo(k)fluoranthene

22.995min (+0.000) 1.58ng/ul m

response 108772

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	25.77
125.00	6.10	10.26#
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	242531	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	1078619	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	675757	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	1550559	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	1353974	20.00	ng/ul	0.00
83) Perylene-d12	23.64	264	1095695	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	18654	3.66	ng/uL	0.00
5) Phenol-d5	6.94	99	412559	20.96	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	264716	22.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	370294	22.18	ng/ul	0.00
13) 4-Methylphenol-d8	8.48	113	371753	21.94	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	176885	21.49	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	211117	21.95	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	399414	21.75	ng/ul	0.00
29) 4-Chloroaniline-d4	10.71	131	426867	24.84	ng/ul	0.00
43) Dimethylphthalate-d6	13.82	166	1418394	23.75	ng/ul	0.00
46) Acenaphthylene-d8	14.11	160	1670199	23.10	ng/ul	0.00
51) 4-Nitrophenol-d4	14.68	143	108929	12.13	ng/ul	0.02
57) Fluorene-d10	15.41	176	1196427	24.41	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.57	200	170290	16.00	ng/ul	0.00
70) Anthracene-d10	17.27	188	1803871	23.53	ng/ul	0.00
76) Pyrene-d10	19.56	212	1908230	27.76	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.49	264	1222545	23.14	ng/ul	0.00

Target Compounds

						Ovalue
6) Phenol	6.97	94	102054	5.017	ng/ul	92
44) Dimethylphthalate	13.87	163	723177	12.086	ng/ul	99
69) Phenanthrene	17.21	178	547157	6.203	ng/ul	98
74) Fluoranthene	19.23	202	796976	7.487	ng/ul	99
77) Pyrene	19.59	202	830854	9.292	ng/ul	96
80) Benzo(a)anthracene	21.34	228	289648	3.288	ng/ul	95
81) Bis(2-ethylhexyl)phthalate	21.24	149	63949	1.071	ng/ul#	96
82) Chrysene	21.39	228	378416	4.575	ng/ul	98
85) Benzo(b)fluoranthene	22.95	252	334339	4.676	ng/ul#	94
86) Benzo(k)fluoranthene	22.99	252	108772m	1.579	ng/ul	
88) Benzo(a)pyrene	23.54	252	200560	2.976	ng/ul#	94
89) Indeno(1,2,3-cd)pyrene	25.98	276	155961	2.179	ng/ul	97
91) Benzo(g,h,i)perylene	26.70	276	140734	2.401	ng/ul#	97

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

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