

Quantitation Report (Qedit)

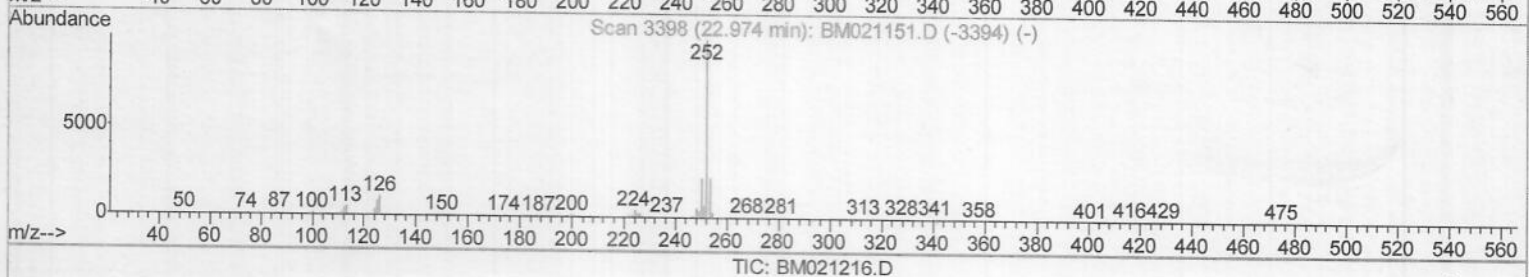
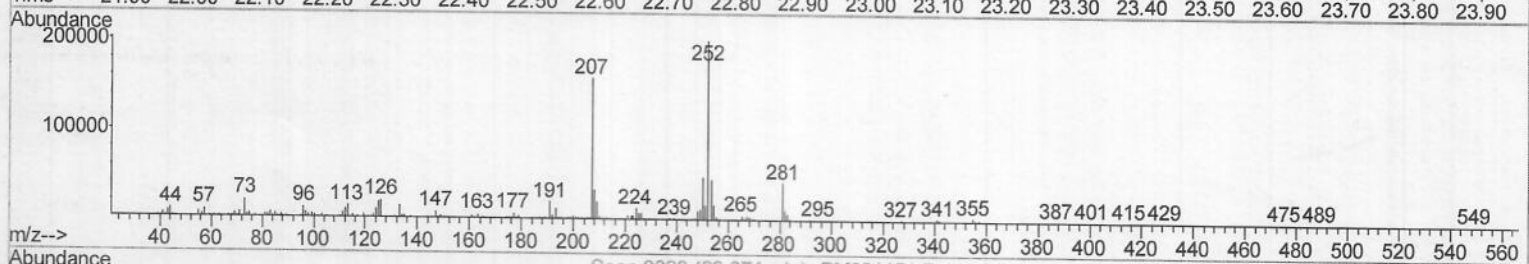
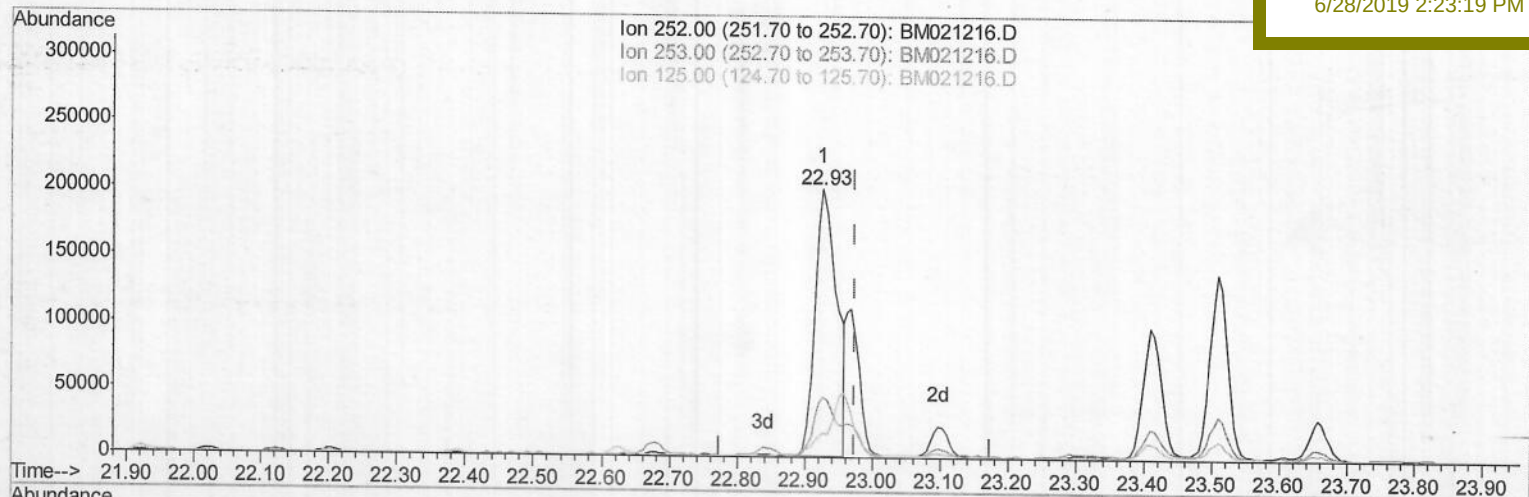
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062719\
 Data File : BM021216.D
 Acq On : 27 Jun 2019 14:44
 Operator : HP/JU
 Sample : K3502-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
ClientSampleId :
 C5AD2

Quant Time: Jun 27 15:22:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 27 15:09:50 2019
 Response via : Initial Calibration

Manual Integrations
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(88) Benzo(k)fluoranthene
 22.927min (-0.047) 5.85ng/ul
 response 445841

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	22.52
125.00	8.20	8.96
0.00	0.00	0.00

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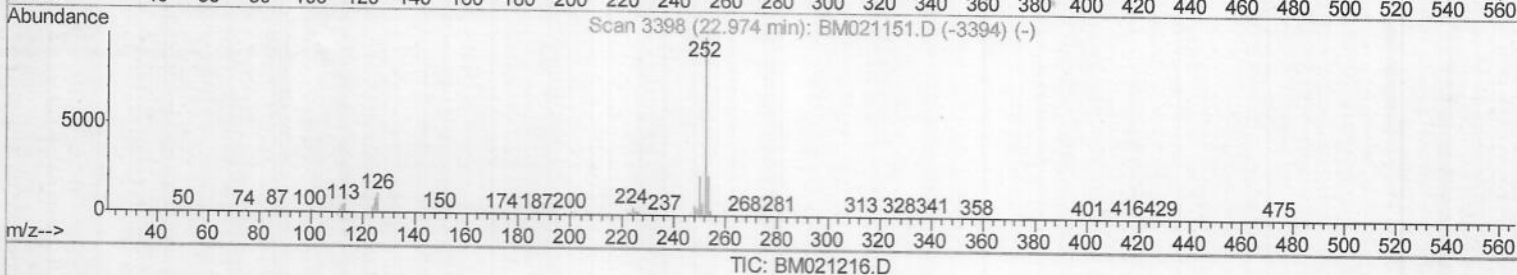
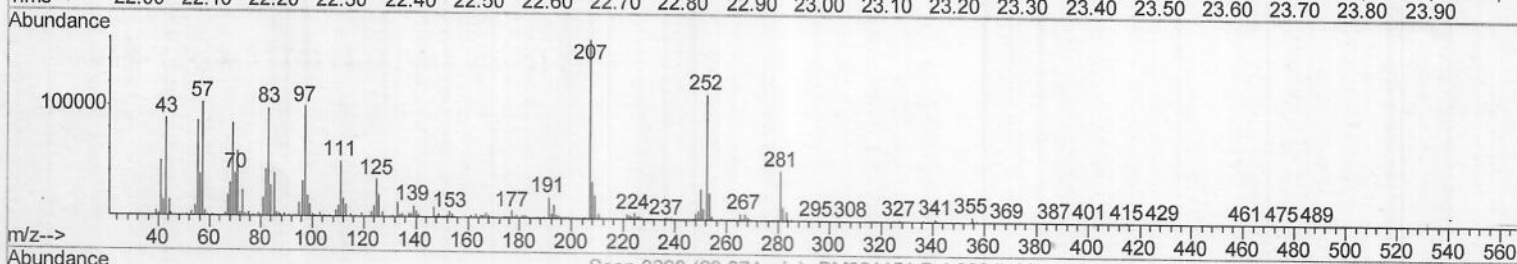
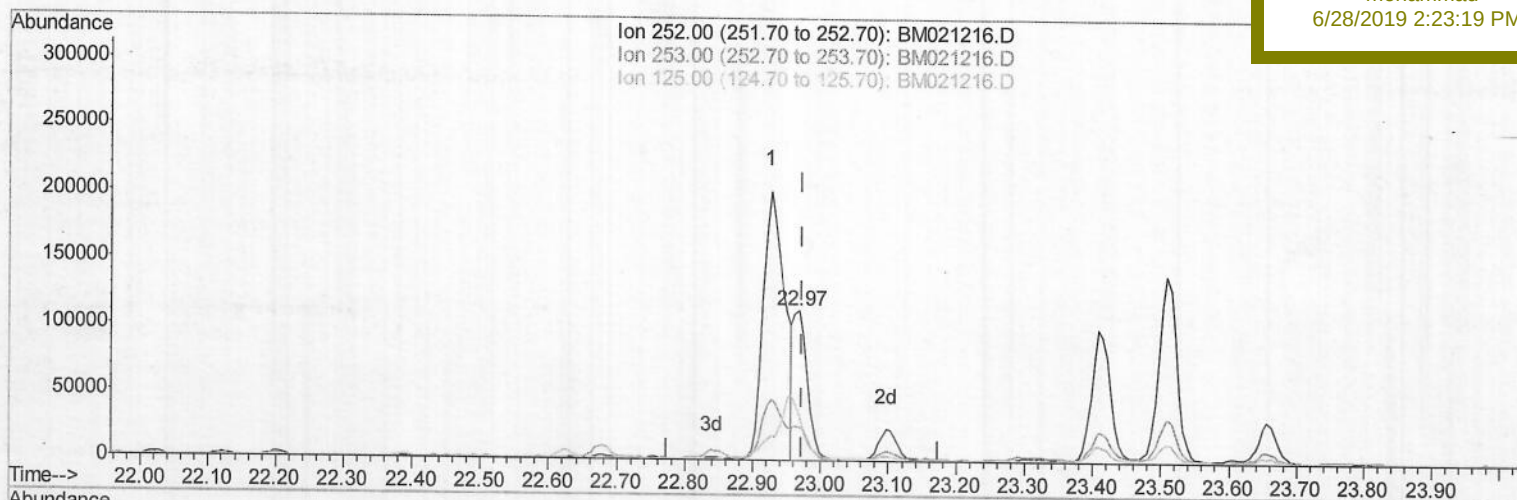
C5AD2

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(88) Benzo(k)fluoranthene

22.968min (-0.006) 2.12ng/ul m

> JU 07/01/19

response 161310

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	21.78
125.00	8.20	30.33#
0.00	0.00	0.00

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Quant Time: Jun 27 15:24:15 2019
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	244362	20.00	ng/ul	0.00
18) Naphthalene-d8	10.55	136	1100687	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.40	164	707936	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.15	188	1618975	20.00	ng/ul	0.00
77) Chrysene-d12	21.34	240	1189703	20.00	ng/ul	0.00
85) Perylene-d12	23.61	264	1250886	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	16820	3.27	ng/uL	0.00
5) Phenol-d5	6.93	99	363689	17.10	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.10	67	234079	18.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	314964	18.22	ng/ul	0.00
13) 4-Methylphenol-d8	8.46	113	294029	16.56	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	170860	19.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.64	143	195006	19.66	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	381176	18.89	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	381366	18.12	ng/ul	0.00
43) Dimethylphthalate-d6	13.82	166	1358350	21.24	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	1528019	19.51	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	138382	13.30	ng/ul	0.00
57) Fluorene-d10	15.40	176	1159823	20.84	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	144577	14.53	ng/ul	0.00
70) Anthracene-d10	17.25	188	1749621	20.80	ng/ul	0.00
78) Pyrene-d10	19.54	212	1703383	23.86	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.47	264	1491304	20.21	ng/ul	0.00

Target Compounds

					Qvalue
44) Dimethylphthalate	13.86	163	272621	4.682	ng/ul 100
69) Phenanthrene	17.19	178	589472	6.584	ng/ul 99
71) Anthracene	17.29	178	109377	1.199	ng/ul 99
76) Fluoranthene	19.21	202	1011570	10.456	ng/ul 99
79) Pyrene	19.57	202	780291	9.277	ng/ul 99
82) Benzo(a)anthracene	21.32	228	336849	4.153	ng/ul 98
84) Chrysene	21.37	228	366024	4.818	ng/ul 98
87) Benzo(b)fluoranthene	22.93	252	445841	5.627	ng/ul 98
88) Benzo(k)fluoranthene	22.97	252	161310m	2.116	ng/ul 98
90) Benzo(a)pyrene	23.51	252	255741	3.429	ng/ul 98
91) Indeno(1,2,3-cd)pyrene	25.90	276	183964	2.095	ng/ul# 95
93) Benzo(g,h,i)perylene	26.60	276	145380	2.010	ng/ul 98

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

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