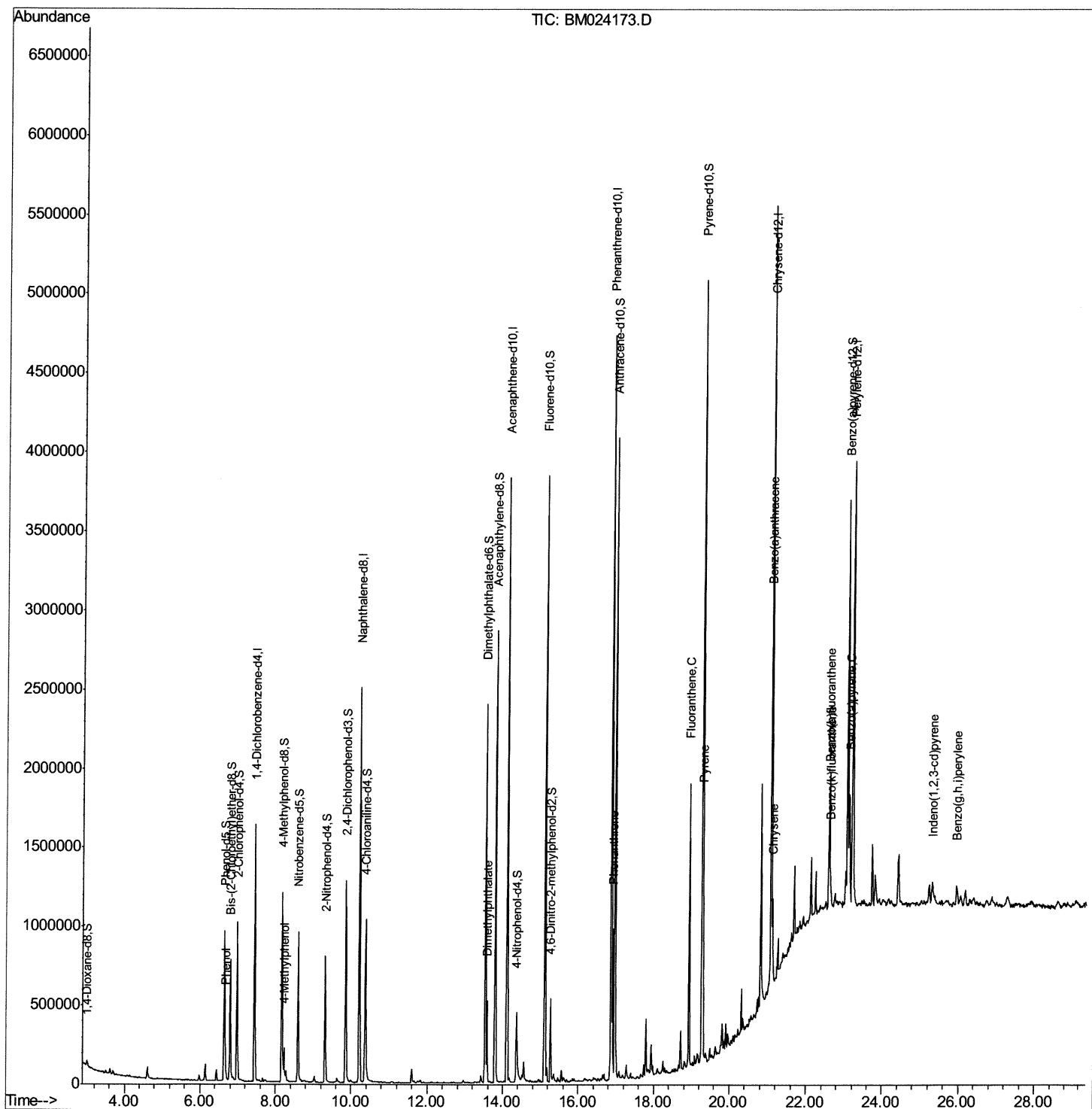


Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024173.D
Acq On : 19 Dec 2019 03:33
Operator : JU
Sample : K6171-16
Misc :
ALS Vial : 47 Sample Multiplier: 1

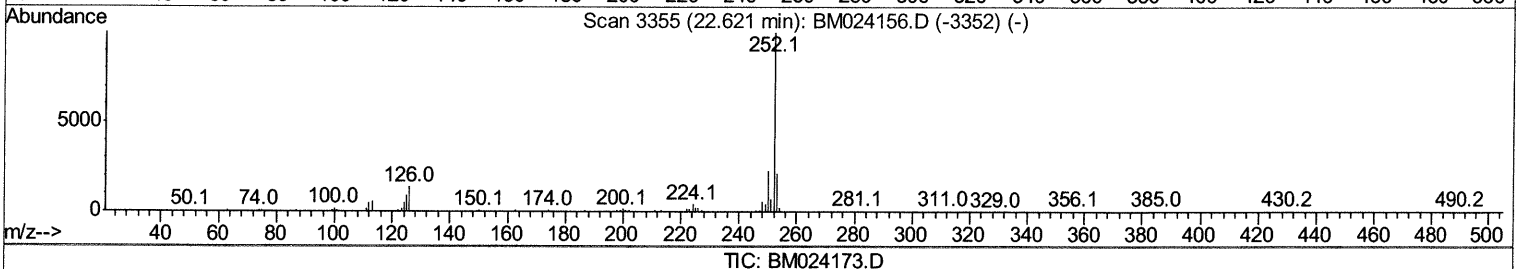
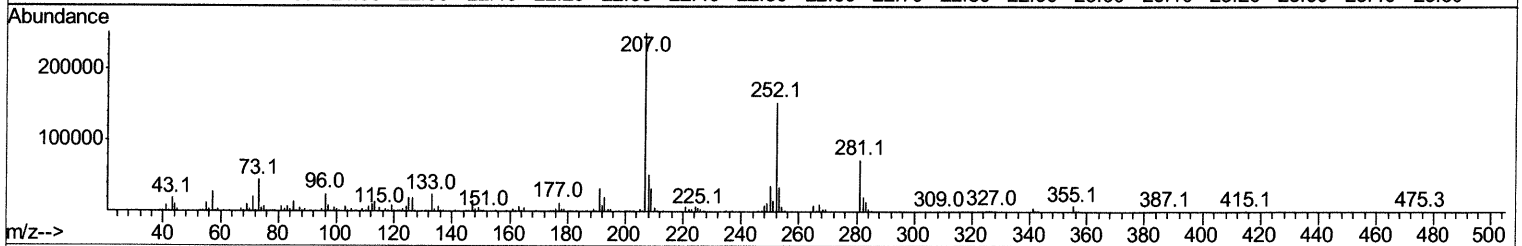
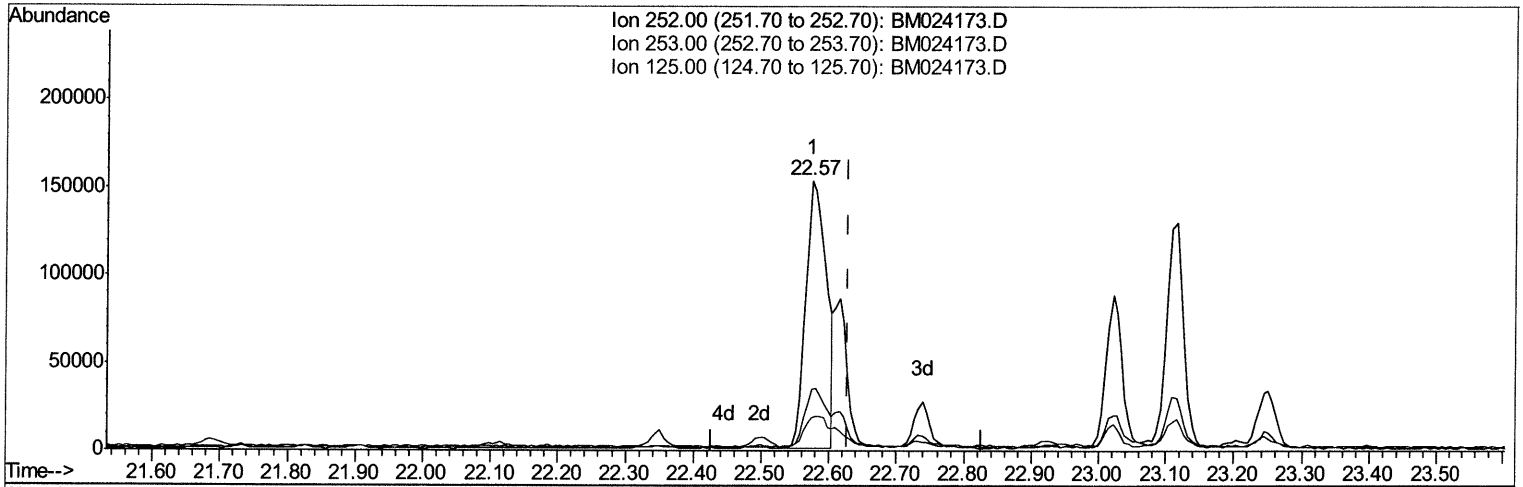
Quant Time: Dec 19 04:16:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 19 02:44:00 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
 Data File : BM024173.D
 Acq On : 19 Dec 2019 03:33
 Operator : JU
 Sample : K6171-16
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Quant Time: Dec 19 04:06:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 19 02:44:00 2019
 Response via : Initial Calibration



TIC: BM024173.D

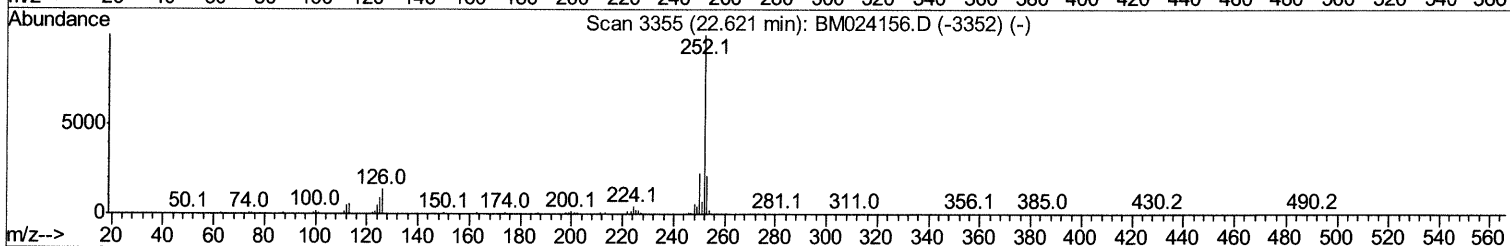
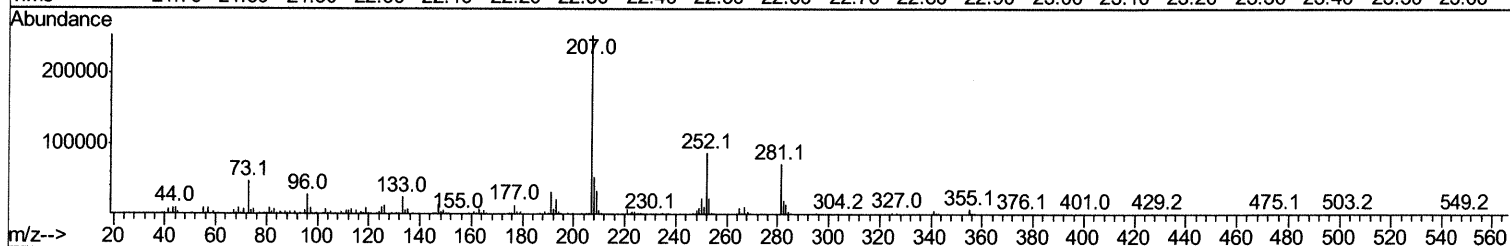
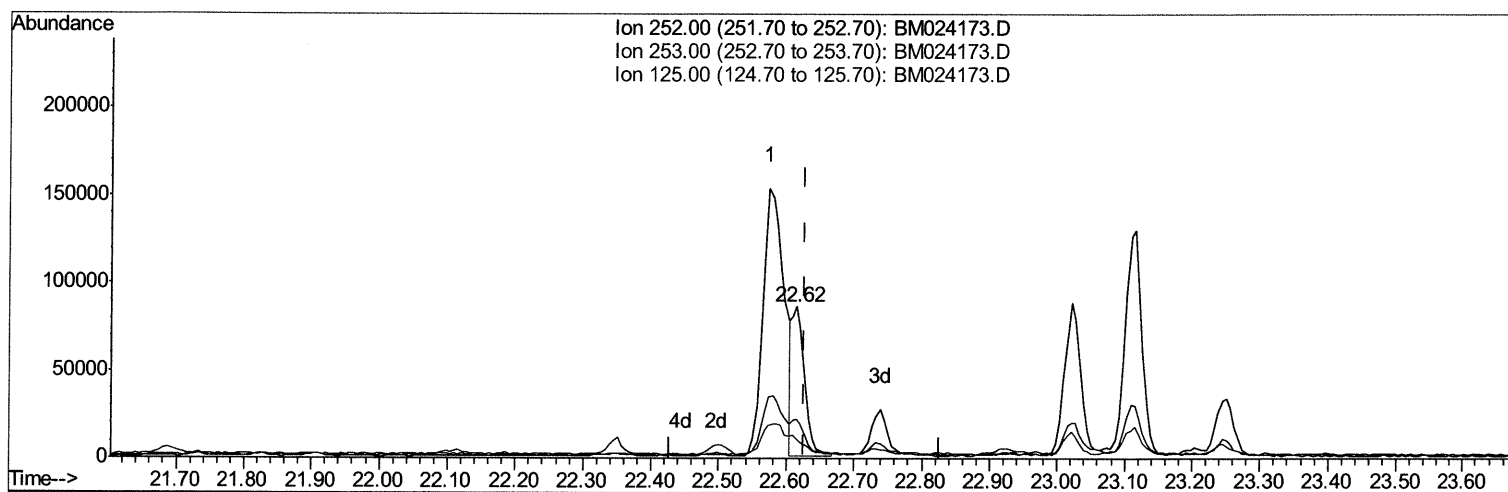
(89) Benzo(k)fluoranthene
 22.574min (-0.053) 3.01ng/ul
 response 327749

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	22.76
125.00	9.90	12.21#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
Data File : BM024173.D
Acq On : 19 Dec 2019 03:33
Operator : JU
Sample : K6171-16
Misc :
ALS Vial : 47 Sample Multiplier: 1

Quant Time: Dec 19 04:06:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 19 02:44:00 2019
Response via : Initial Calibration



TIC: BM024173.D

(89) Benzo(k)fluoranthene

22.615min (-0.012) 1.06ng/ul m JU 12/20/19

response 114878

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	25.94
125.00	9.90	13.02#
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121719\
 Data File : BM024173.D
 Acq On : 19 Dec 2019 03:33
 Operator : JU
 Sample : K6171-16
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Quant Time: Dec 19 04:16:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 19 02:44:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	442161	20.00	ng/ul	0.00
18) Naphthalene-d8	10.22	136	2048101	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	1274495	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.86	188	2579200	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	1966102	20.00	ng/ul	0.00
86) Perylene-d12	23.20	264	1854227	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.01	96	23743	2.37	ng/uL	0.00
5) Phenol-d5	6.65	99	602775	14.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.80	67	375838	14.97	ng/ul	0.00
9) 2-Chlorophenol-d4	6.99	132	468491	15.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.17	113	474290	13.86	ng/ul	0.00
19) Nitrobenzene-d5	8.60	128	236794	16.09	ng/ul	0.00
22) 2-Nitrophenol-d4	9.32	143	263179	16.26	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	490060	15.58	ng/ul	0.00
29) 4-Chloroaniline-d4	10.37	131	605787	15.86	ng/ul	0.00
44) Dimethylphthalate-d6	13.53	166	1625947	17.25	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	1943419	17.21	ng/ul	0.00
52) 4-Nitrophenol-d4	14.36	143	152697	9.02	ng/ul	0.00
58) Fluorene-d10	15.11	176	1416574	17.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	111379	7.08	ng/ul	0.00
71) Anthracene-d10	16.96	188	2109364	17.97	ng/ul	0.00
79) Pyrene-d10	19.27	212	2313888	24.71	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.07	264	1808202	19.63	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.67	94	49933	1.155	ng/ul	93
16) 4-Methylphenol	8.23	108	78553	2.201	ng/ul	86
45) Dimethylphthalate	13.58	163	317070	3.237	ng/ul	99
70) Phenanthrene	16.90	178	513311	3.555	ng/ul	99
77) Fluoranthene	18.93	202	1013589	6.440	ng/ul	99
80) Pyrene	19.30	202	786289	6.126	ng/ul	99
83) Benzo(a)anthracene	21.06	228	317448	2.534	ng/ul	99
85) Chrysene	21.11	228	285194	2.323	ng/ul	97
88) Benzo(b)fluoranthene	22.57	252	327749	2.954	ng/ul#	96
89) Benzo(k)fluoranthene	22.62	252	114878m	1.055	ng/ul	> JU 12/20/19
91) Benzo(a)pyrene	23.12	252	219724	2.189	ng/ul#	95
92) Indeno(1,2,3-cd)pyrene	25.31	276	141870	1.177	ng/ul	95
94) Benzo(g,h,i)perylene	25.95	276	137865	1.401	ng/ul	98

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