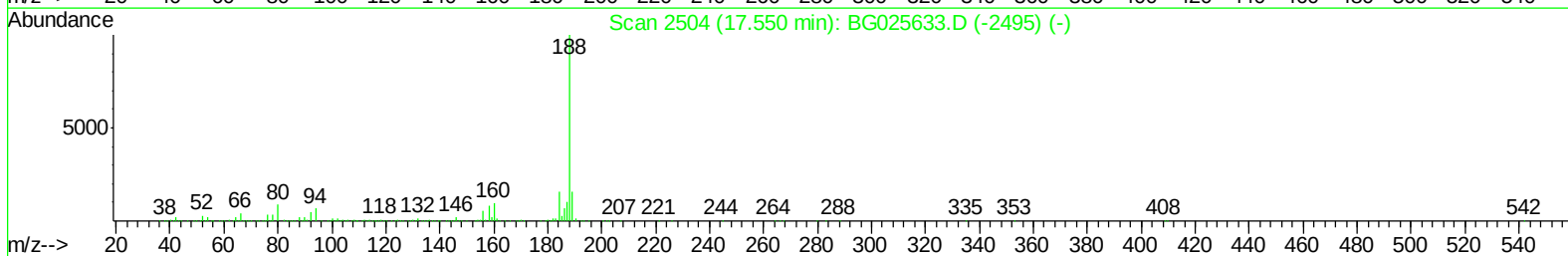
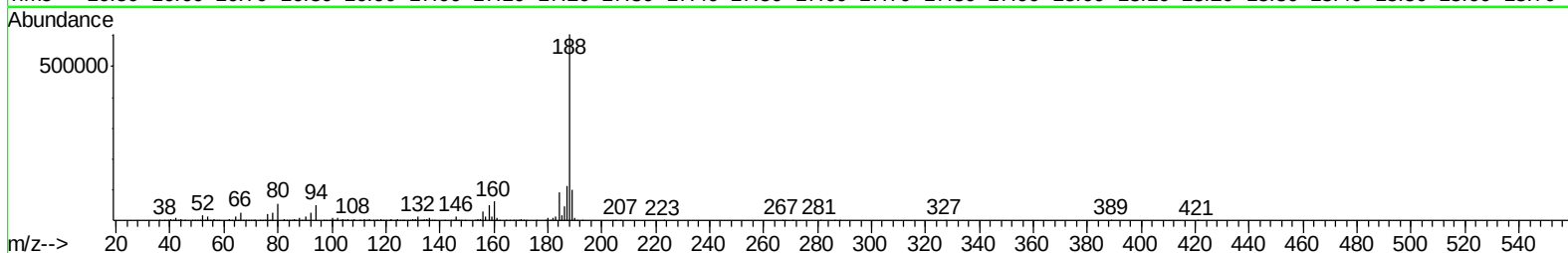
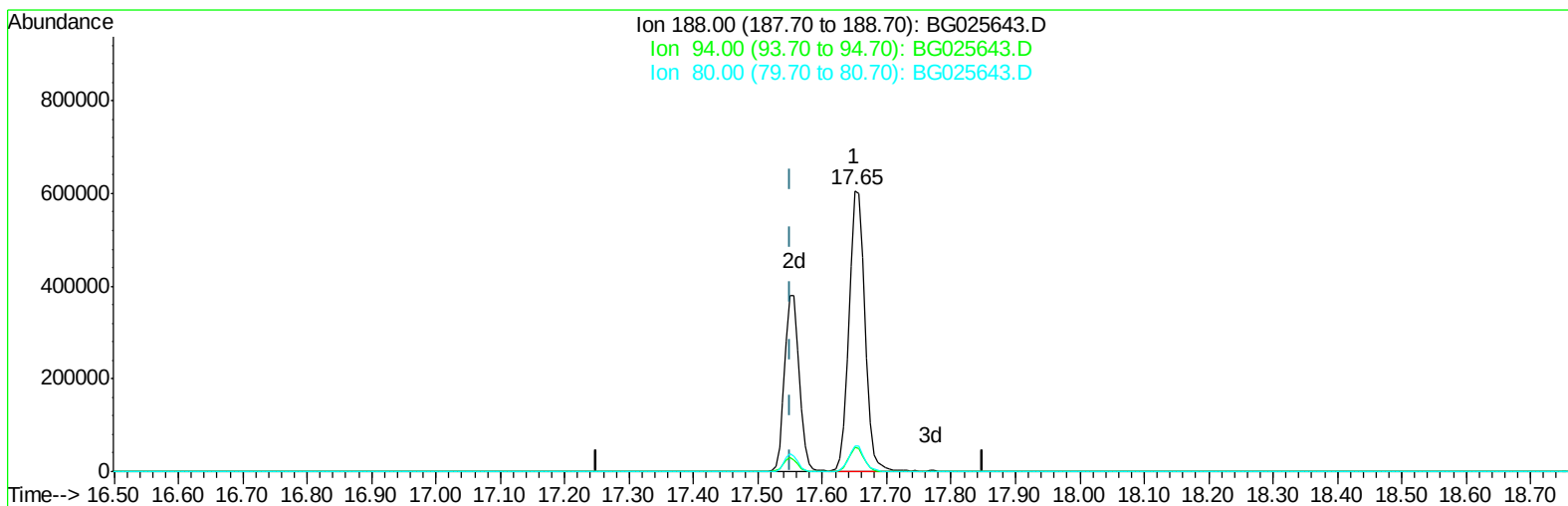


Data Path : Z:\HPCHEM1\BNA_G\Data\BG012517\
Data File : BG025643.D
Acq On : 25 Jan 2017 20:59
Operator : SJ/MA
Sample : PB96323BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jan 26 09:33:14 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 25 18:08:33 2017
Response via : Initial Calibration



TIC: BG025643.D

(61) Phenanthrene-d10 (I)

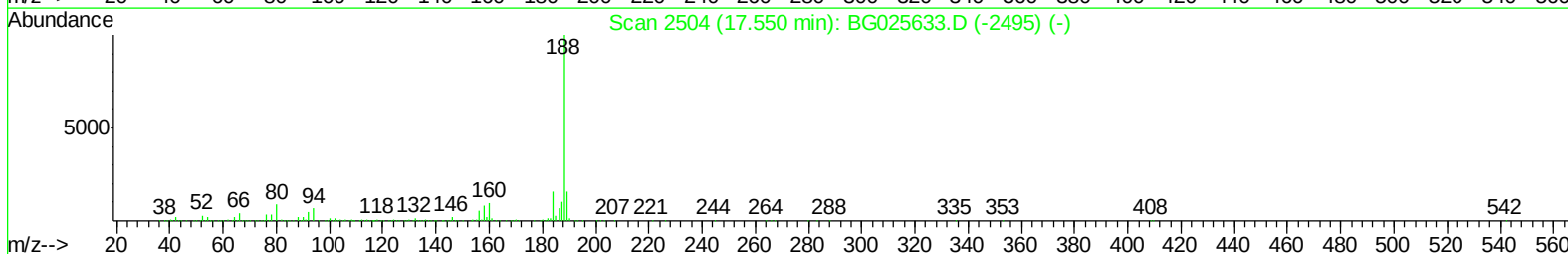
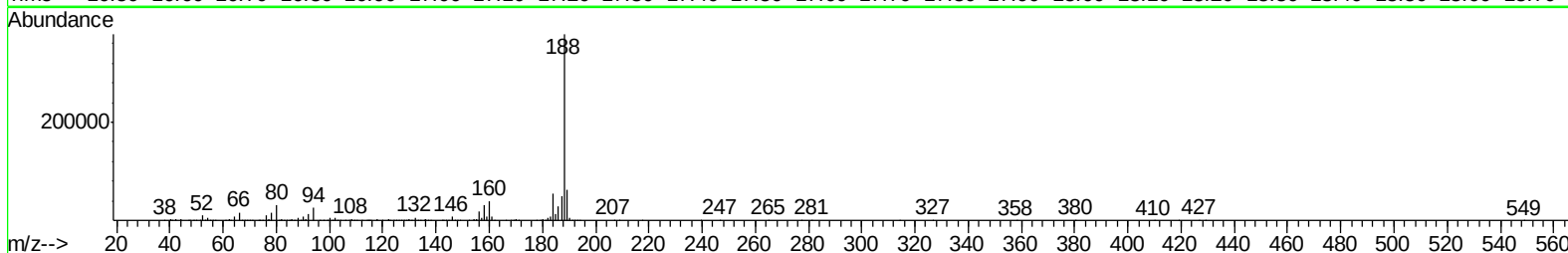
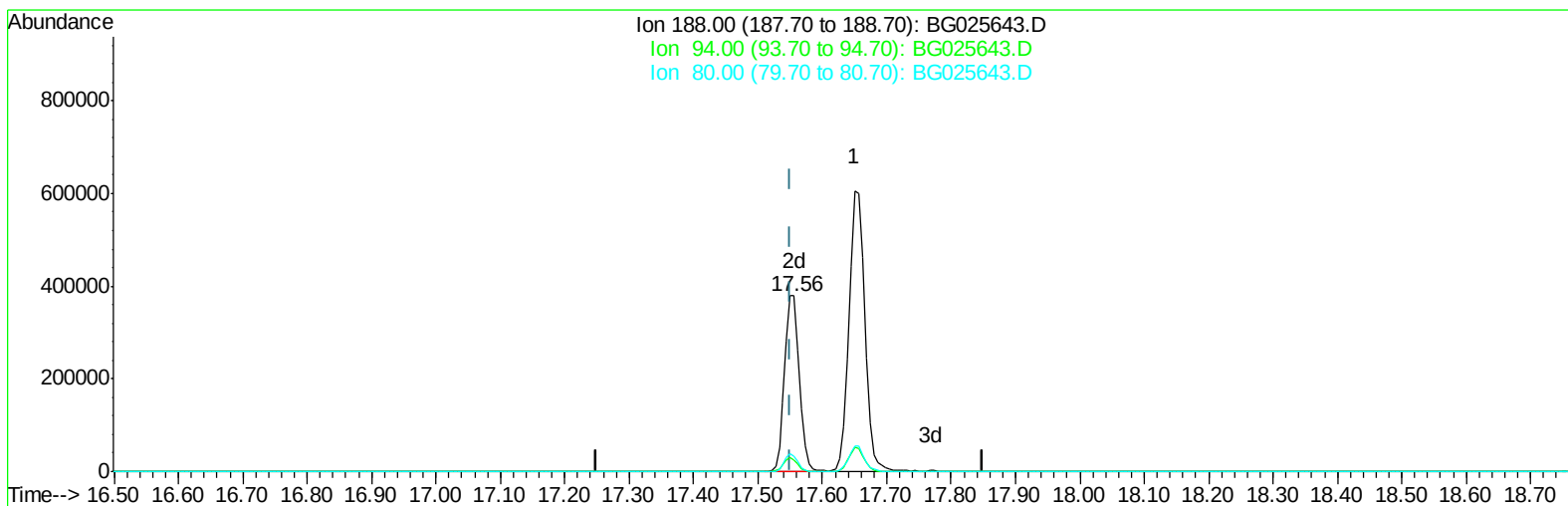
17.650min (+0.100) 20.00ng/ul

response 1030853

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	8.60
80.00	9.50	9.22
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG012517\
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Sample : PB96323BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jan 26 09:33:14 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
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TIC: BG025643.D

(61) Phenanthrene-d10 (I)

17.556min (+0.006) 20.00ng/ul m

response 610770

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	6.87#
80.00	9.50	8.56
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG012517\
 Data File : BG025643.D
 Acq On : 25 Jan 2017 20:59
 Operator : SJ/MA
 Sample : PB96323BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jan 26 09:34:34 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 25 18:08:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	62874	20.00	ng/ul	0.00
18) Naphthalene-d8	11.01	136	290658	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.81	164	253553	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.56	188	610770m	20.00	ng/ul	0.00
75) Chrysene-d12	21.85	240	736733	20.00	ng/ul	0.00
83) Perylene-d12	25.23	264	742147	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	8987	7.15	ng/uL	0.00
5) Phenol-d5	7.34	99	186629	34.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.49	67	138178	39.70	ng/ul	0.00
9) 2-Chlorophenol-d4	7.71	132	140371	36.93	ng/ul	0.00
13) 4-Methylphenol-d8	8.89	113	166006	36.24	ng/ul	0.00
19) Nitrobenzene-d5	9.36	128	75974	36.88	ng/ul	0.00
22) 2-Nitrophenol-d4	10.09	143	99138	40.33	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.64	165	159970	32.98	ng/ul	0.00
29) 4-Chloroaniline-d4	11.15	131	221440	40.85	ng/ul	0.00
43) Dimethylphthalate-d6	14.20	166	711273	39.72	ng/ul	0.00
46) Acenaphthylene-d8	14.51	160	769601	39.79	ng/ul	0.00
51) 4-Nitrophenol-d4	15.04	143	83483	34.03	ng/ul	0.00
57) Fluorene-d10	15.79	176	631280	37.40	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.94	200	124680	33.09	ng/ul	0.00
70) Anthracene-d10	17.65	188	1032375	39.70	ng/ul	0.00
76) Pyrene-d10	19.93	212	1247329	40.69	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.99	264	1268856	40.68	ng/ul	0.00

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

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