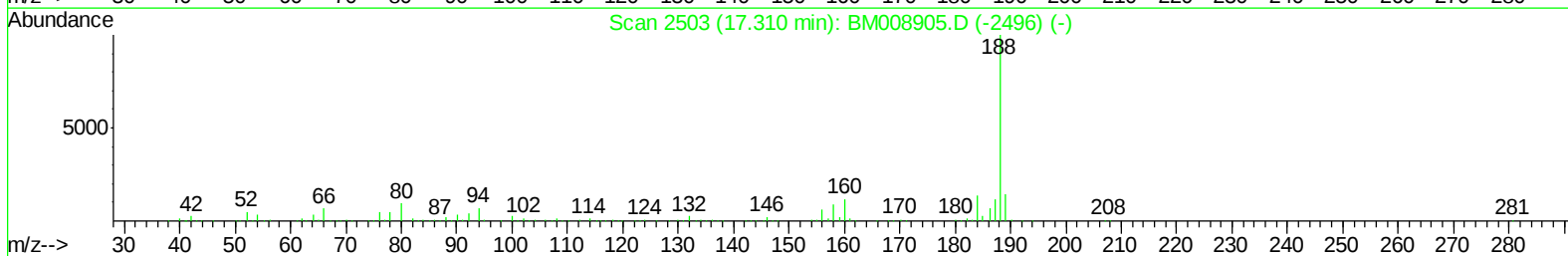
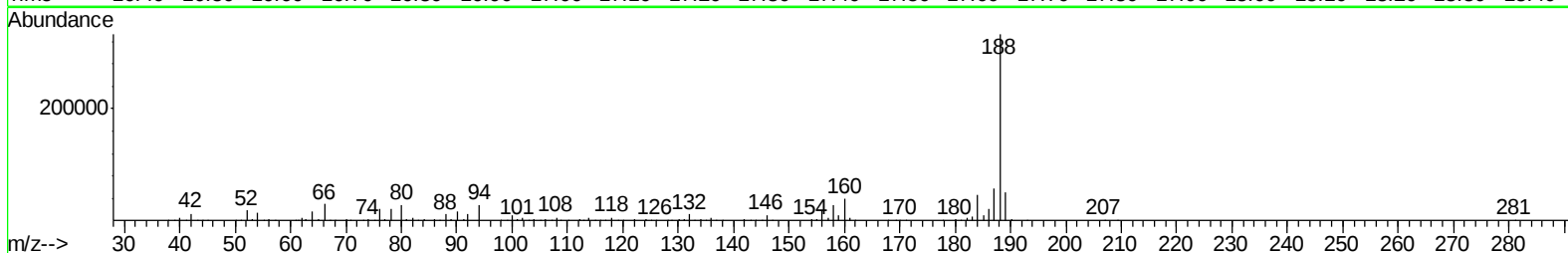
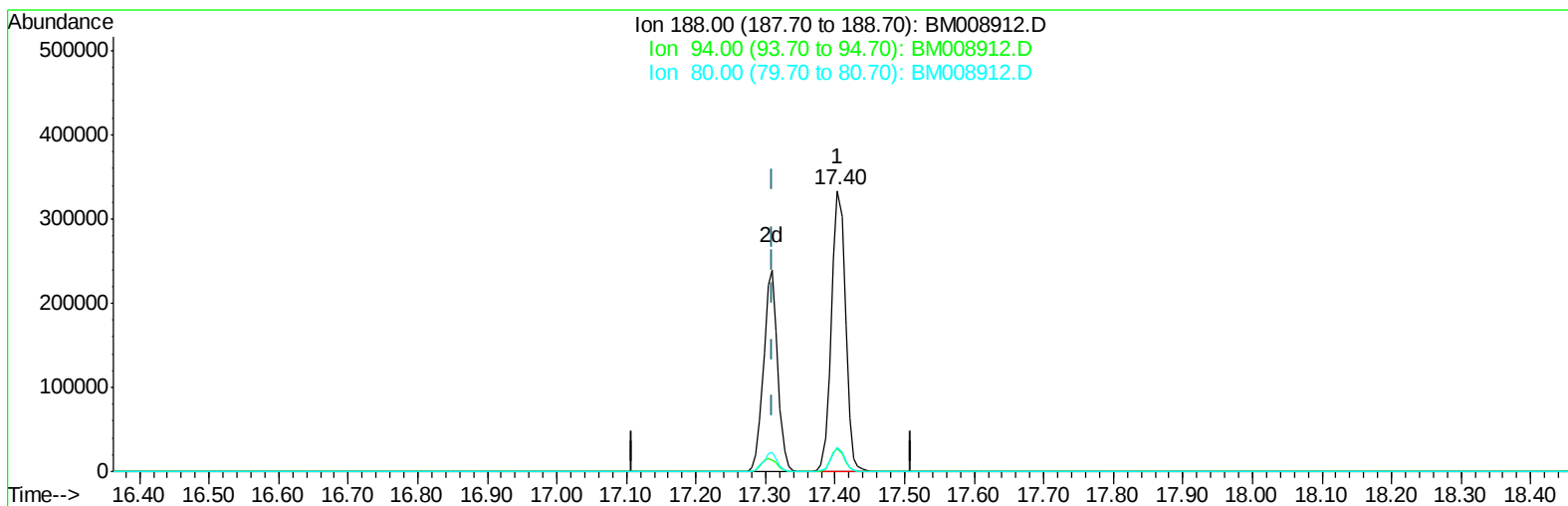


Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012717\
Data File : BM008912.D
Acq On : 27 Jan 2017 21:38
Operator : SJ/MA
Sample : I1506-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 27 22:39:30 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 27 22:34:22 2017
Response via : Initial Calibration



TIC: BM008912.D

(61) Phenanthrene-d10 (I)

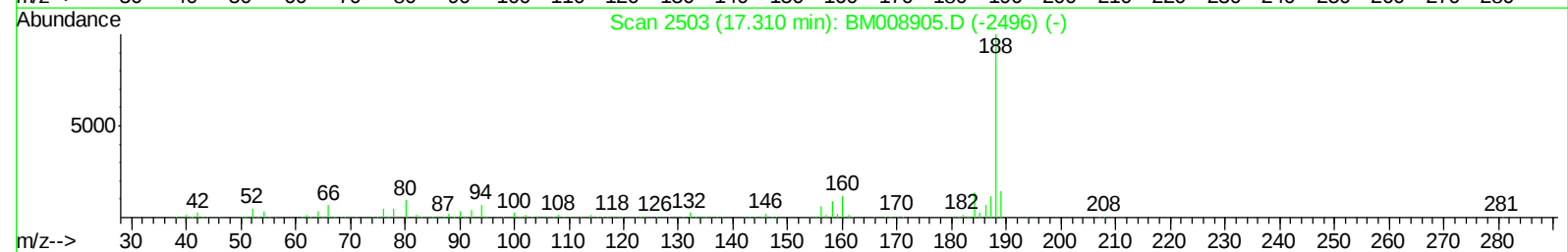
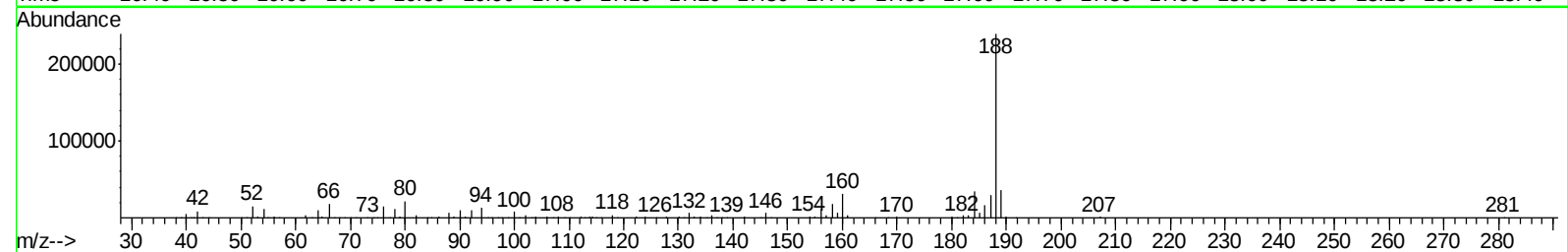
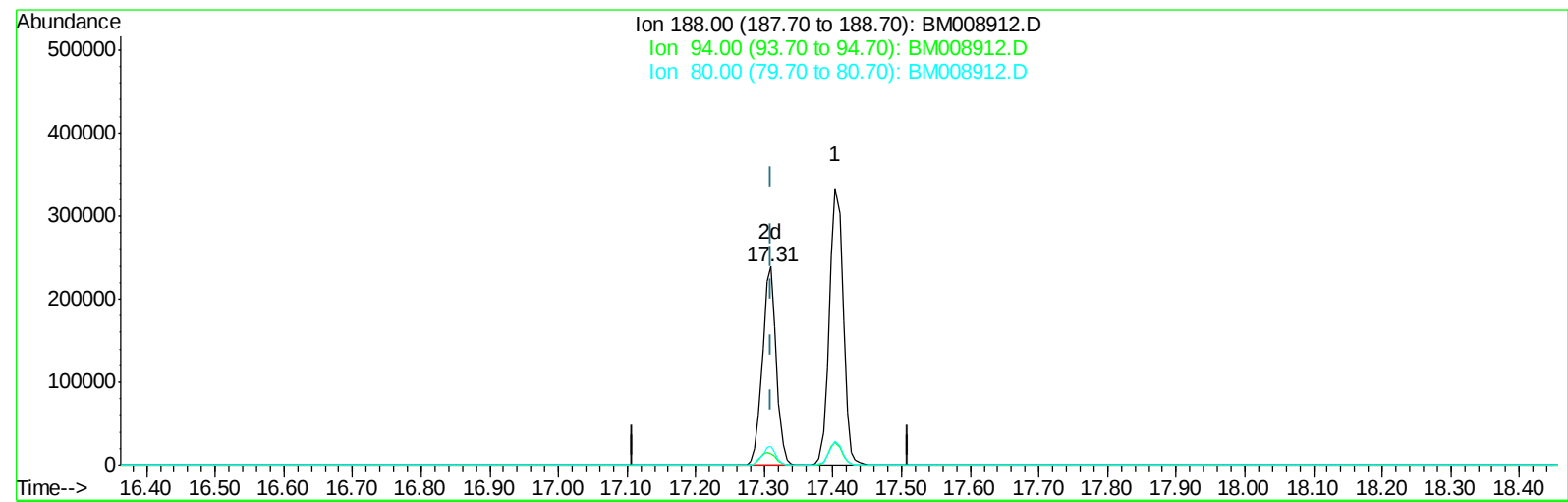
17.403min (+0.094) 20.00ng/ul

response 464541

Ion	Exp%	Act%
188.00	100	100
94.00	8.00	8.39
80.00	9.50	8.58
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012717\
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Acq On : 27 Jan 2017 21:38
Operator : SJ/MA
Sample : I1506-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 27 22:39:30 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 27 22:34:22 2017
Response via : Initial Calibration



TIC: BM008912.D

(61) Phenanthrene-d10 (I)

17.309min (-0.000) 20.00ng/ul m

response 338005

Ion	Exp%	Act%
188.00	100	100
94.00	8.00	5.94#
80.00	9.50	9.25
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012717\
 Data File : BM008912.D
 Acq On : 27 Jan 2017 21:38
 Operator : SJ/MA
 Sample : I1506-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 27 23:47:40 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 27 22:34:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	61444	20.00	ng/ul	0.00
18) Naphthalene-d8	10.73	136	226396	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.56	164	160811	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.31	188	338005m	20.00	ng/ul	0.00
75) Chrysene-d12	21.47	240	436591	20.00	ng/ul	0.00
83) Perylene-d12	23.83	264	434966	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.39	96	1367	1.10	ng/uL	0.00
5) Phenol-d5	7.07	99	27595	5.30	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.25	67	108147	26.14	ng/ul	0.00
9) 2-Chlorophenol-d4	7.45	132	68683	20.26	ng/ul	0.00
13) 4-Methylphenol-d8	8.62	113	45730	11.22	ng/ul	0.00
19) Nitrobenzene-d5	9.09	128	40164	27.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.80	143	47817	27.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.34	165	90431	24.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.86	131	6075	1.52	ng/ul	0.00
43) Dimethylphthalate-d6	13.97	166	322120	30.57	ng/ul	0.00
46) Acenaphthylene-d8	14.26	160	358307	28.74	ng/ul	0.00
51) 4-Nitrophenol-d4	14.74	143	9032	5.12	ng/ul	0.00
57) Fluorene-d10	15.55	176	285869	28.56	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.66	200	49660	24.67	ng/ul	0.00
70) Anthracene-d10	17.40	188	464541	31.43	ng/ul	0.00
76) Pyrene-d10	19.69	212	572603	29.90	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.68	264	555123	30.14	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Phenol	7.10	94	7879	1.39	ng/ul#	51
44) Dimethylphthalate	14.01	163	18384	1.75	ng/ul	99

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

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