

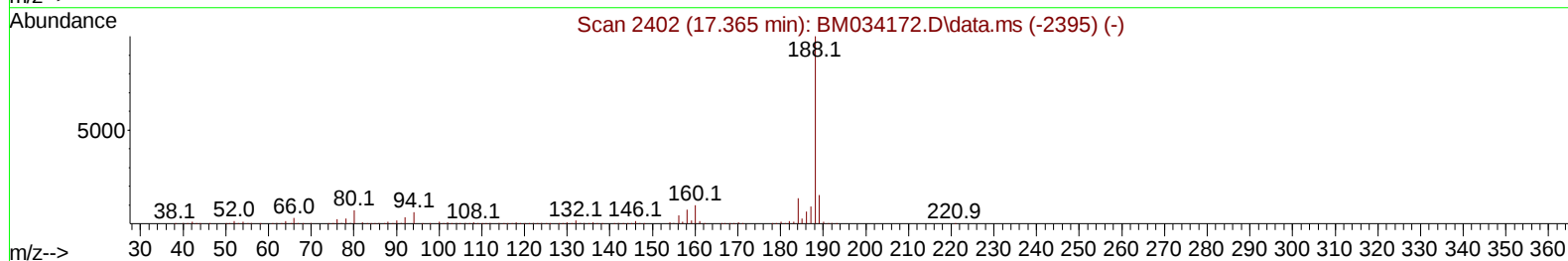
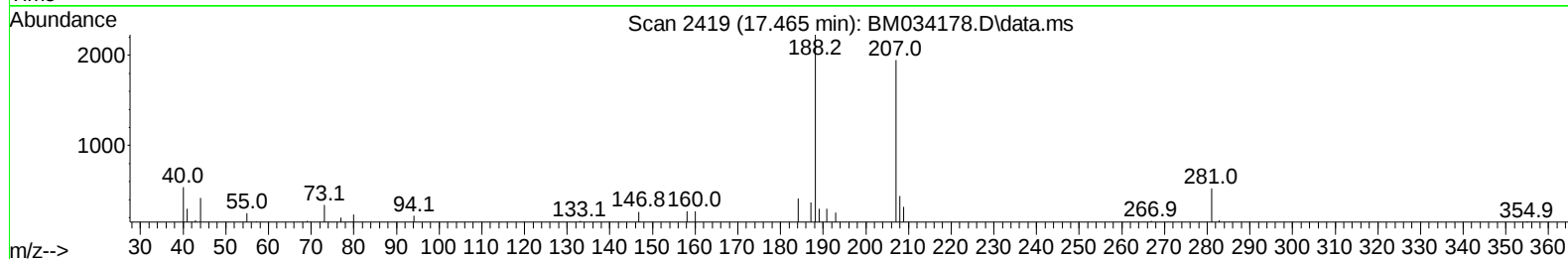
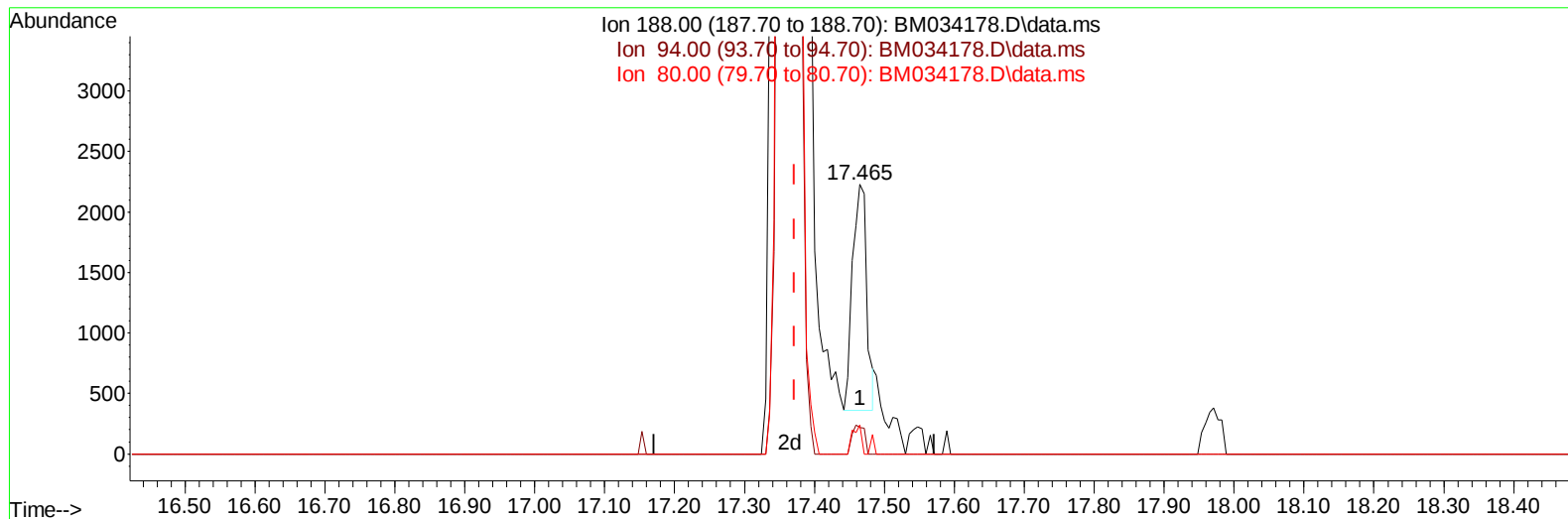
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034178.D
 Acq On : 09 Mar 2022 09:42
 Operator : CG/JU
 Sample : N1669-03
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGZR1

Manual IntegrationsAPPROVED

Quant Time: Mar 09 23:55:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 08 13:09:16 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/10/2022
 Supervised By :Yogesh Patel 03/11/2022



TIC: BM034178.D\data.ms

(64) Phenanthrene-d10 (I)

17.465min (+ 0.094) 20.00 ng/u1

response 2660

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	8.40	9.88
80.00	9.70	10.69
0.00	0.00	0.00

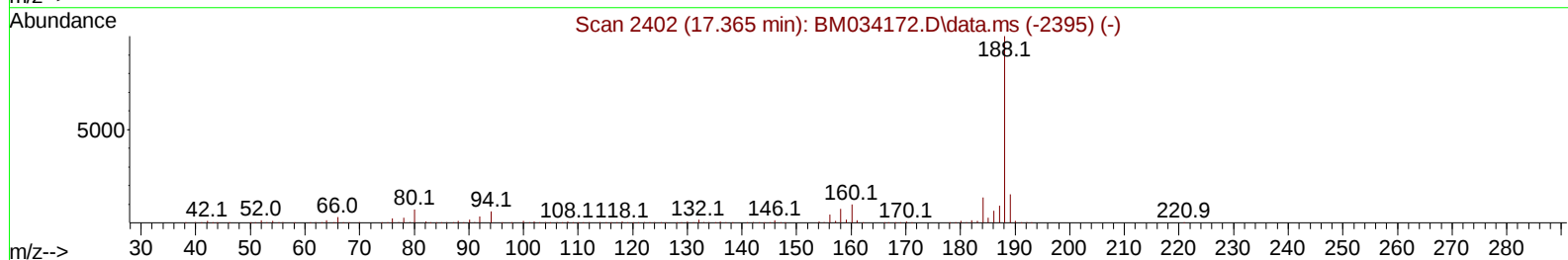
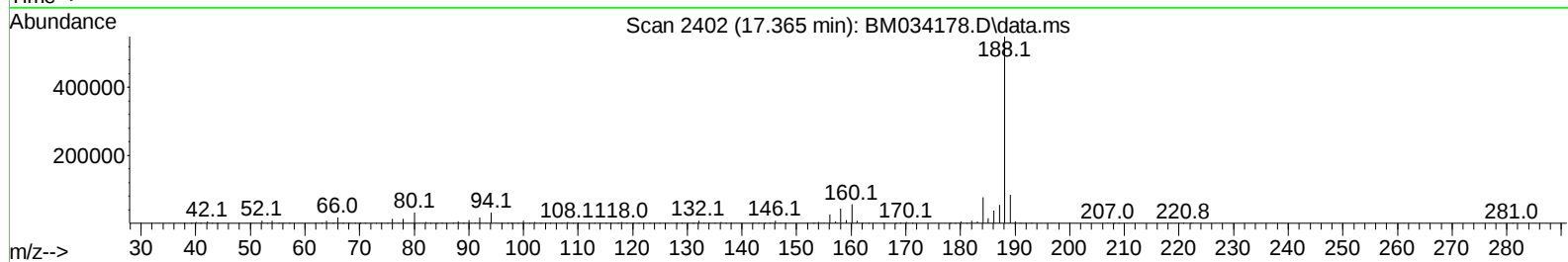
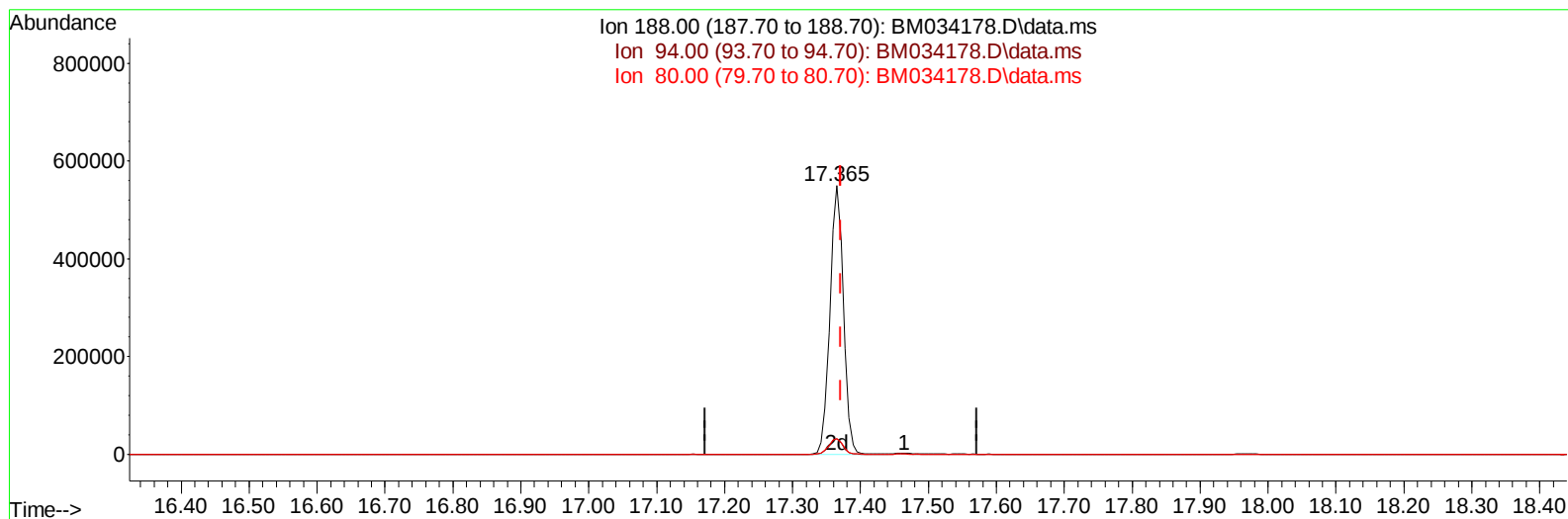
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(64) Phenanthrene-d10 (I)

17.365min (-0.006) 20.00 ng/ul m

response 754628

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	8.40	5.70#
80.00	9.70	5.86#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.989	152	177196	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	627327	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	393788	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	754628m	20.000	ng/ul	0.00
79) Chrysene-d12	21.536	240	589485	20.000	ng/ul	# 0.00
88) Perylene-d12	23.983	264	607765	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0	0.000	ng/ul	
7) Phenol-d5	0.000	99	0	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0	0.000	ng/ul	
11) 2-Chlorophenol-d4	0.000	132	0	0.000	ng/ul	
15) 4-Methylphenol-d8	0.000	113	0	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128	0	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0	0.000	ng/ul	
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	0.000	166	0	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160	0	0.000	ng/ul	
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	0.000	176	0d	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	0.000	200	0	0.000	ng/ul	
73) Anthracene-d10	0.000	188	0d	0.000	ng/ul	
81) Pyrene-d10	0.000	212	0d	0.000	ng/ul	
92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/ul	

Target Compounds Qvalue

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

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