

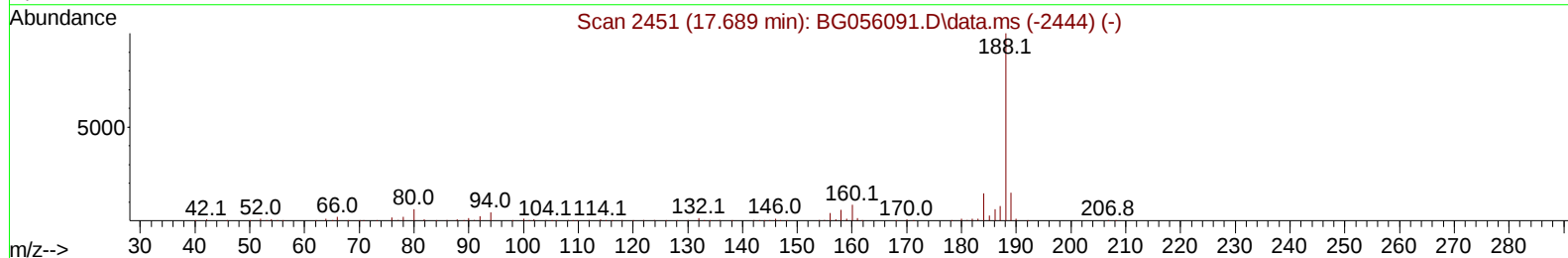
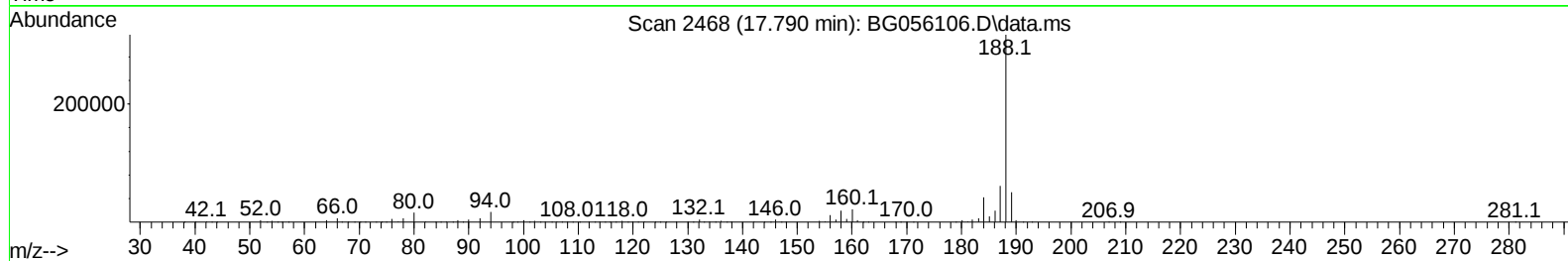
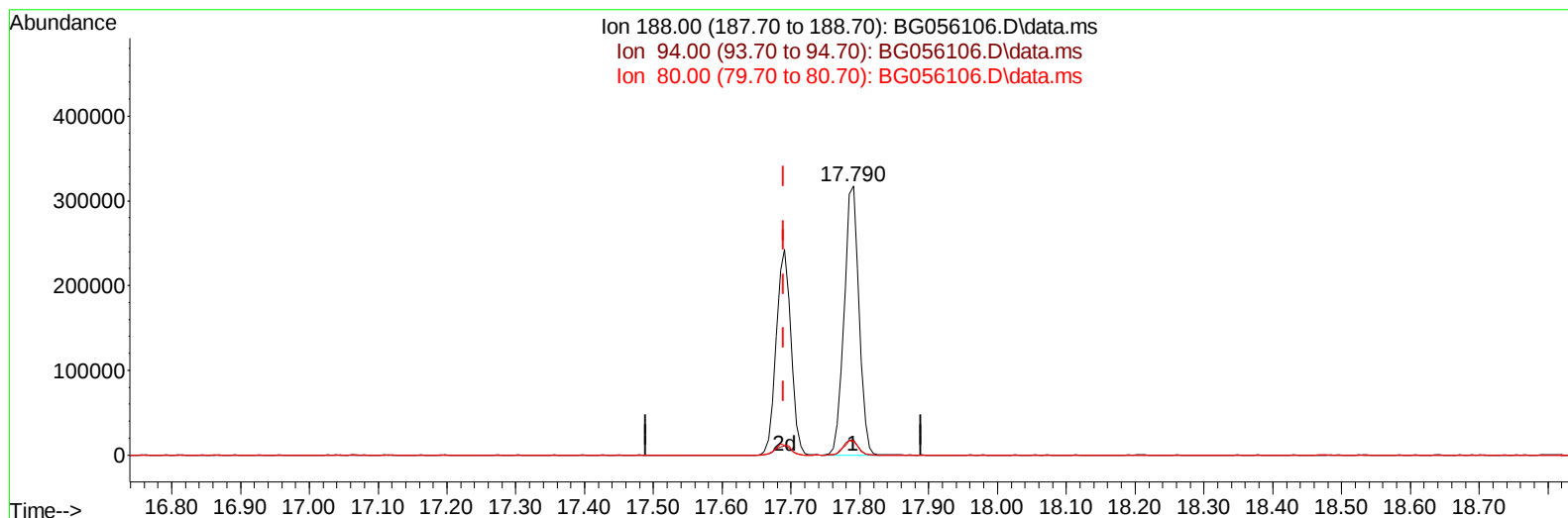
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056106.D
 Acq On : 25 Dec 2022 2:36
 Operator : CG/JU
 Sample : N6017-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBYF5

Manual IntegrationsAPPROVED

Quant Time: Dec 25 05:42:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 22:54:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/25/2022
 Supervised By :Yogesh Patel 12/25/2022



TIC: BG056106.D\data.ms

(64) Phenanthrene-d10 (I)

17.790min (+ 0.101) 20.00 ng/u1

response 474167

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	4.60	5.46
80.00	6.20	5.06
0.00	0.00	0.00

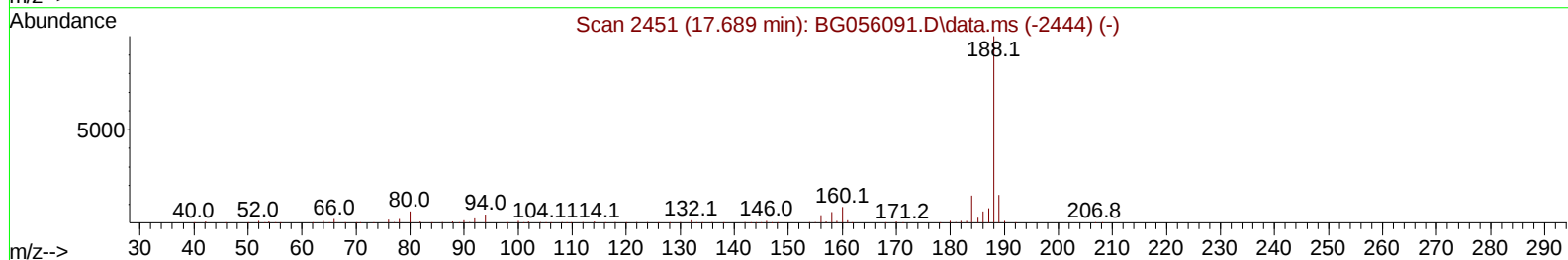
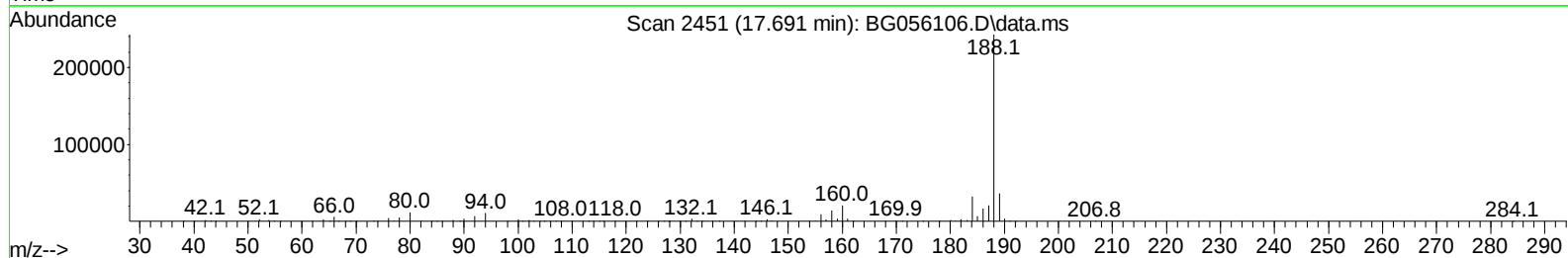
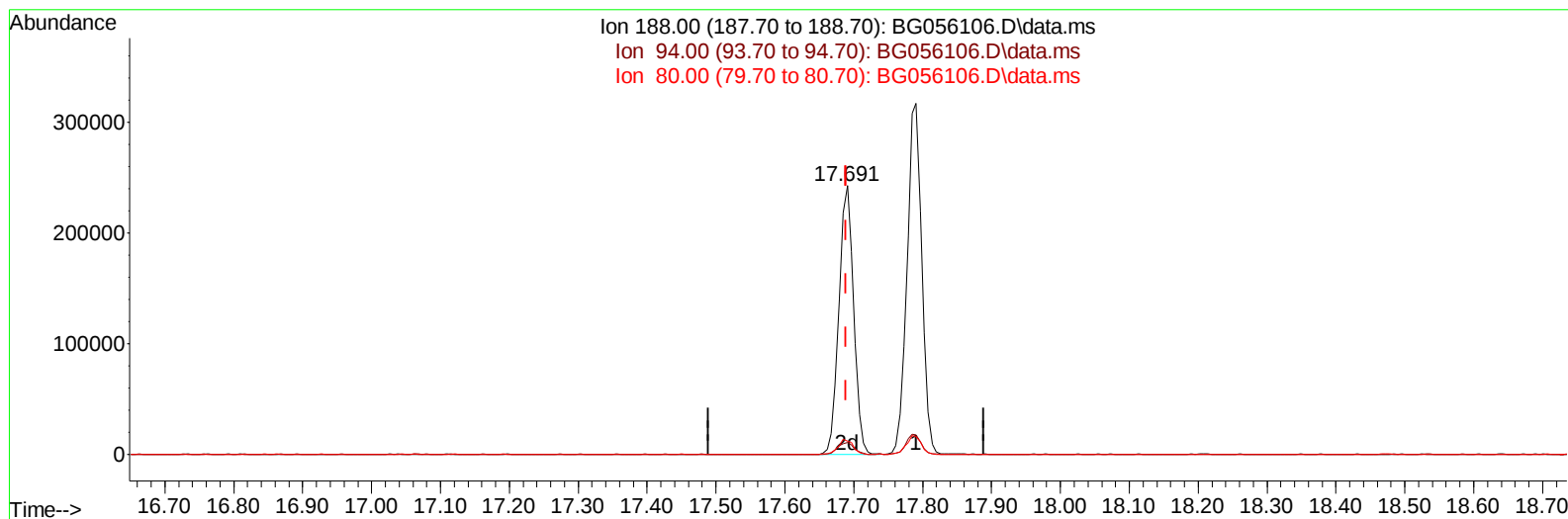
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TIC: BG056106.D\data.ms

(64) Phenanthrene-d10 (I)

17.691min (+ 0.001) 20.00 ng/ul m

response 359397

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	4.60	4.38
80.00	6.20	4.95#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.343	152	39065	20.000	ng/ul	0.00
20) Naphthalene-d8	11.181	136	160216	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.953	164	143853	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.691	188	359397m	20.000	ng/ul	0.00
79) Chrysene-d12	21.986	240	348633	20.000	ng/ul	0.00
88) Perylene-d12	25.458	264	374966	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.648	96	5058	5.420	ng/uL	0.00
4) Pyridine-d5	4.083	84	73056	24.009	ng/ul	0.00
7) Phenol-d5	7.473	99	100981	28.497	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.649	67	43250	30.967	ng/ul	0.00
11) 2-Chlorophenol-d4	7.867	132	66559	31.780	ng/ul	0.00
15) 4-Methylphenol-d8	9.036	113	68444	25.664	ng/ul	0.00
21) Nitrobenzene-d5	9.518	128	36105	30.940	ng/ul	0.00
24) 2-Nitrophenol-d4	10.246	143	46723	30.646	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.781	165	97906	28.898	ng/ul	0.00
31) 4-Chloroaniline-d4	11.304	131	81764	21.337	ng/ul	0.00
46) Dimethylphthalate-d6	14.342	166	329679	29.611	ng/ul	0.00
49) Acenaphthylene-d8	14.653	160	366195	29.124	ng/ul	0.00
54) 4-Nitrophenol-d4	15.117	143	42152	25.151	ng/ul	0.00
60) Fluorene-d10	15.940	176	315483	28.707	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.040	200	60677	26.263	ng/ul	0.00
73) Anthracene-d10	17.790	188	474167	28.550	ng/ul	0.00
81) Pyrene-d10	20.052	212	587882	30.533	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.217	264	567856	29.795	ng/ul	0.00

Target Compounds Qvalue

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

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