

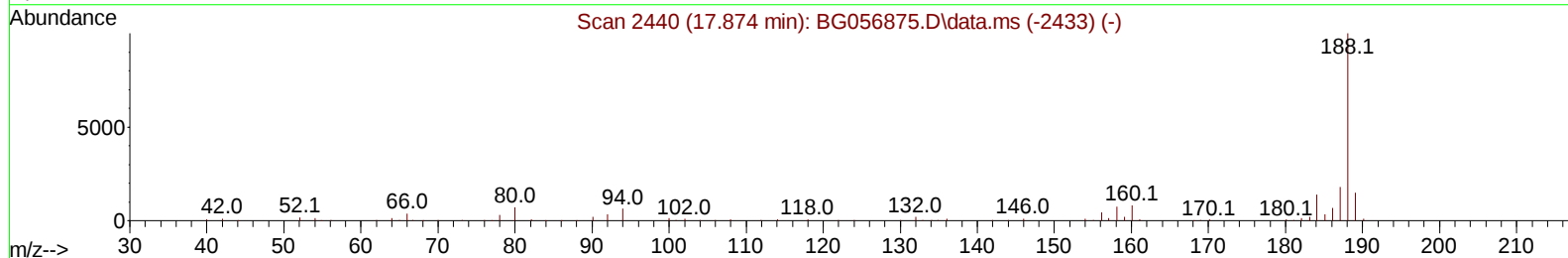
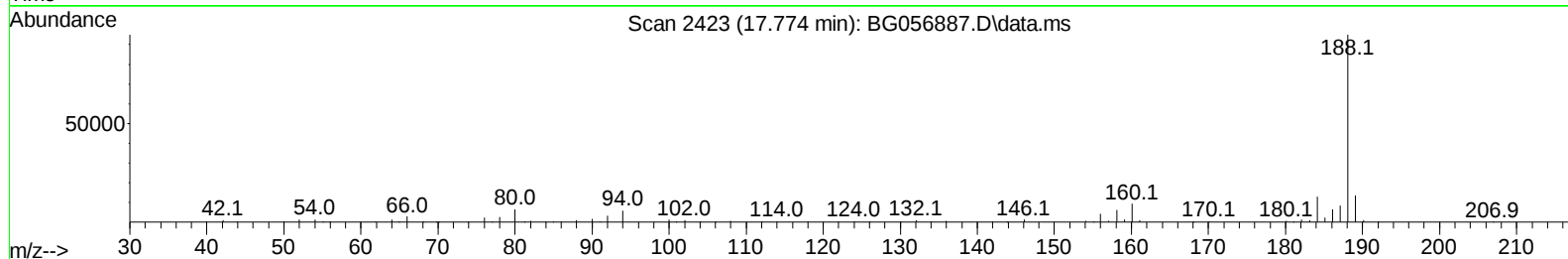
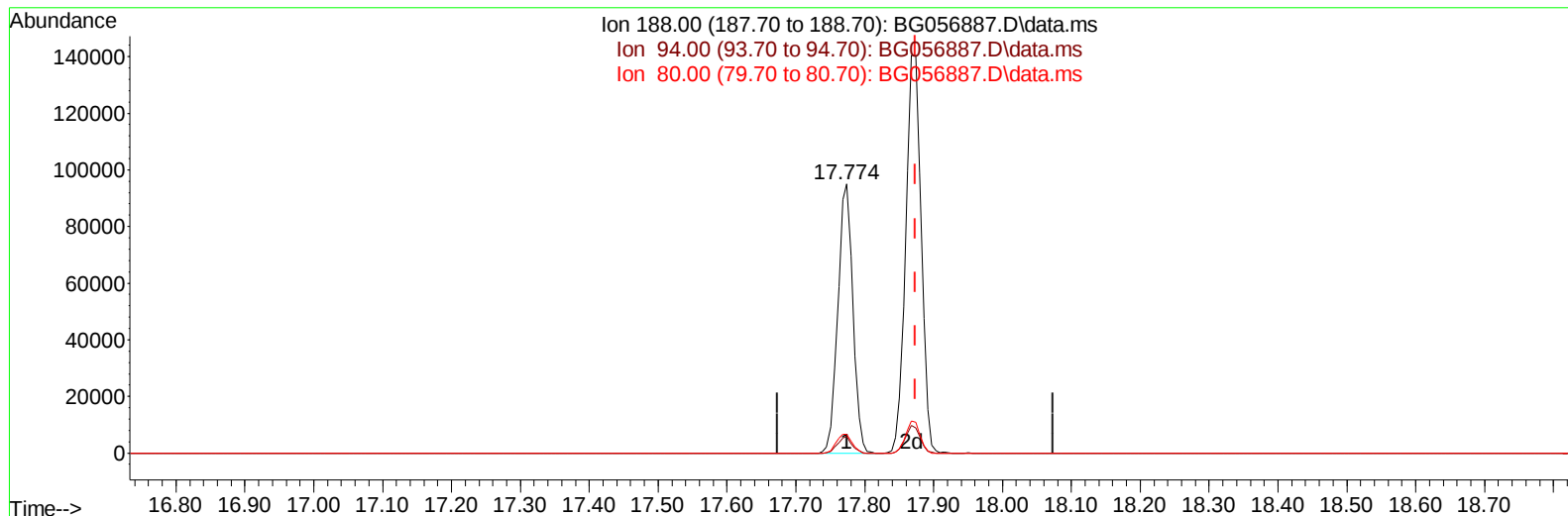
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030823\
 Data File : BG056887.D
 Acq On : 8 Mar 2023 19:12
 Operator : CG/JU
 Sample : PB151258BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK258

Manual IntegrationsAPPROVED

Quant Time: Mar 09 00:00:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 08 23:58:51 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/09/2023
 Supervised By :Jagrut Upadhyay 03/09/2023



TIC: BG056887.D\data.ms

(73) Anthracene-d10 (S)

17.774min (-0.100) 20.69 ng/u1

response 142942

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	6.70	6.35
80.00	6.10	7.00
0.00	0.00	0.00

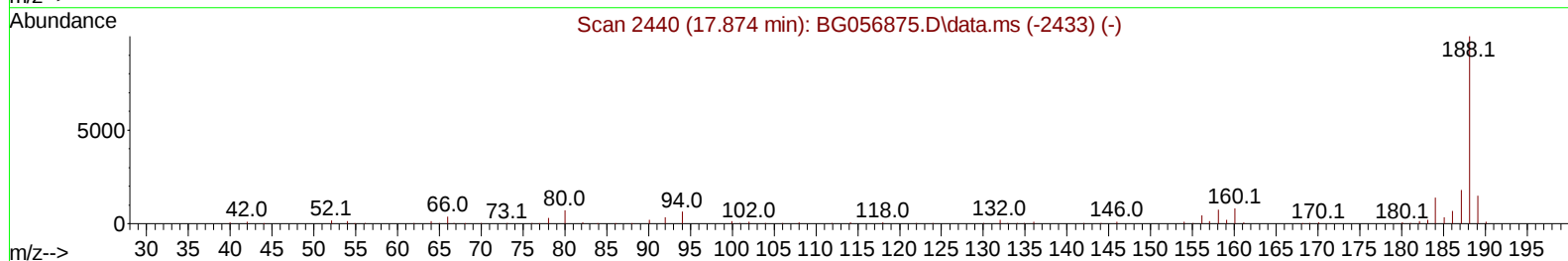
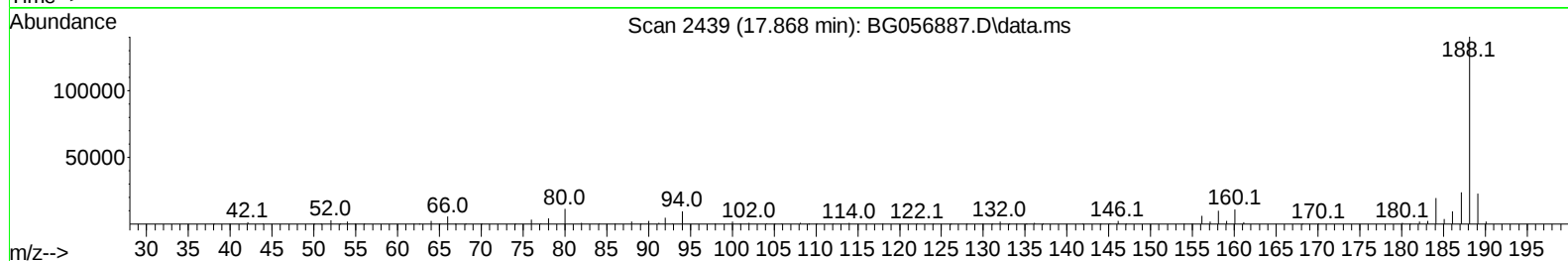
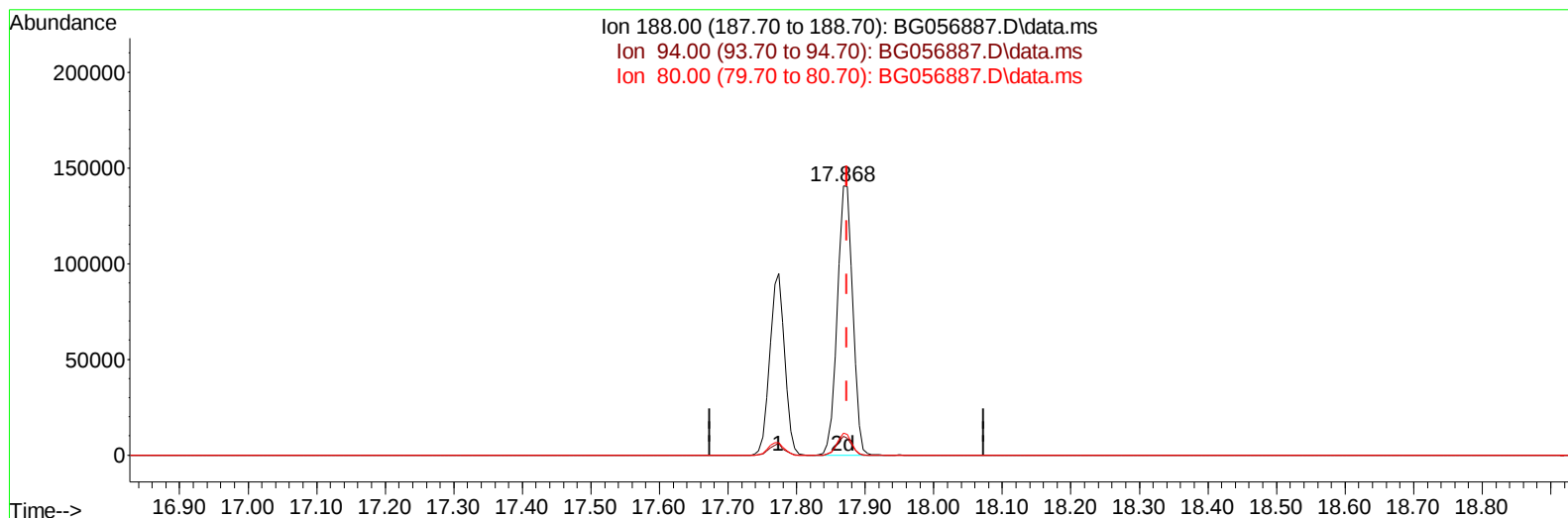
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(73) Anthracene-d10 (S)

17.868min (-0.006) 31.85 ng/ul m

response 220062

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	6.70	6.98
80.00	6.10	8.08#
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	8.420	152	13796	20.000	ng/ul	0.00
20) Naphthalene-d8	11.258	136	58671	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.030	164	49602	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.774	188	142942	20.000	ng/ul	0.00
79) Chrysene-d12	22.092	240	146394	20.000	ng/ul	0.00
88) Perylene-d12	25.647	264	152913	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.690	96	2340	6.578	ng/uL	0.00
4) Pyridine-d5	4.131	84	34928	30.181	ng/ul	0.00
7) Phenol-d5	7.539	99	42145	29.806	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.721	67	28077	31.613	ng/ul	0.00
11) 2-Chlorophenol-d4	7.938	132	27712	31.315	ng/ul	0.00
15) 4-Methylphenol-d8	9.102	113	30673	28.749	ng/ul	0.00
21) Nitrobenzene-d5	9.589	128	14707	30.223	ng/ul	0.00
24) 2-Nitrophenol-d4	10.318	143	18937	32.895	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.864	165	31697	28.426	ng/ul	0.00
31) 4-Chloroaniline-d4	11.381	131	42640	29.020	ng/ul	0.00
46) Dimethylphthalate-d6	14.419	166	136487	32.169	ng/ul	0.00
49) Acenaphthylene-d8	14.730	160	138061	29.636	ng/ul	0.00
54) 4-Nitrophenol-d4	15.189	143	24278	33.769	ng/ul	0.00
60) Fluorene-d10	16.017	176	120601	30.654	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.117	200	29853	28.898	ng/ul	0.00
73) Anthracene-d10	17.868	188	220062m	31.854	ng/ul	0.00
81) Pyrene-d10	20.136	212	284283	31.881	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.400	264	270068	32.255	ng/ul	0.00

Target Compounds Qvalue

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

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