

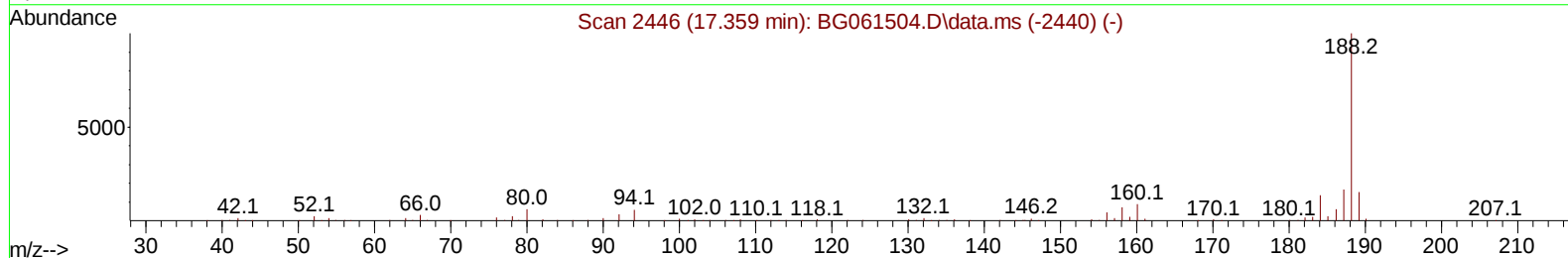
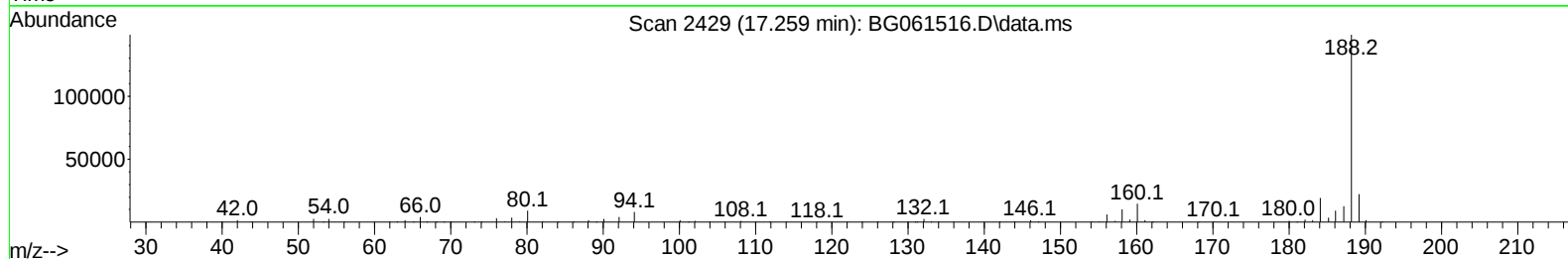
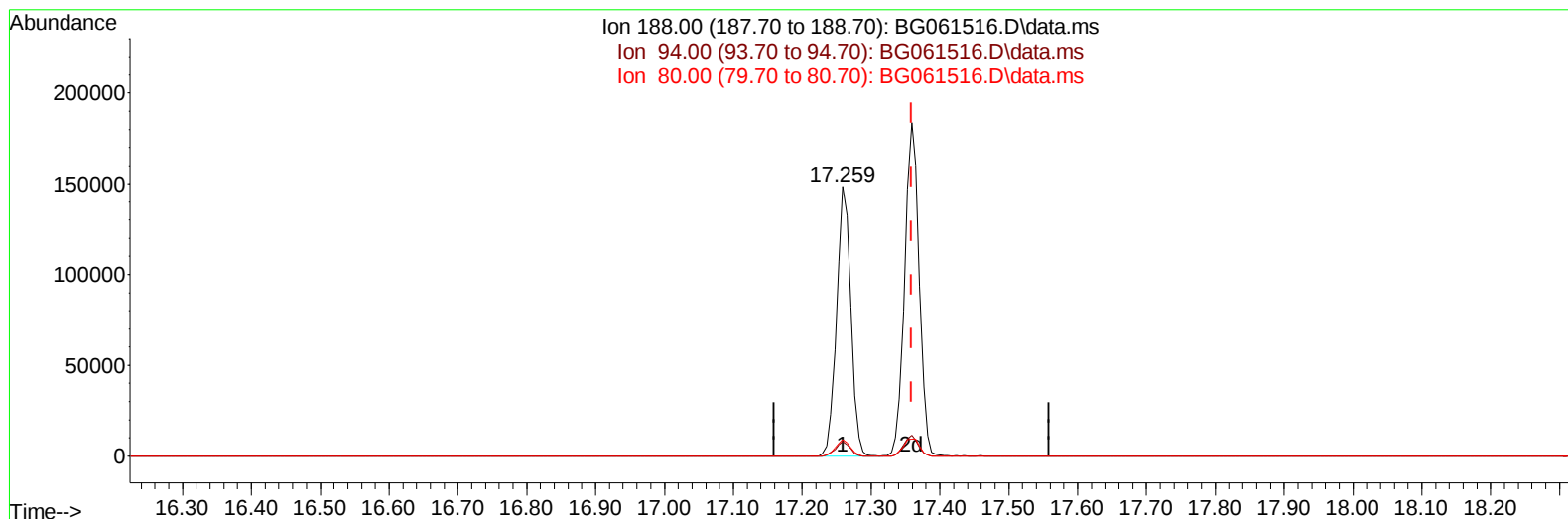
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060624\
Data File : BG061516.D
Acq On : 6 Jun 2024 19:05
Operator : MA/JU
Sample : PB161346BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK346

Manual IntegrationsAPPROVED

Quant Time: Jun 06 22:44:18 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 06 12:15:45 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/07/2024
Supervised By :mohammad ahmed 06/08/2024



TIC: BG061516.D\data.ms

(73) Anthracene-d10 (S)

17.259min (-0.100) 22.68 ng/u1

response 213199

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	6.60	5.43
80.00	6.40	6.08
0.00	0.00	0.00

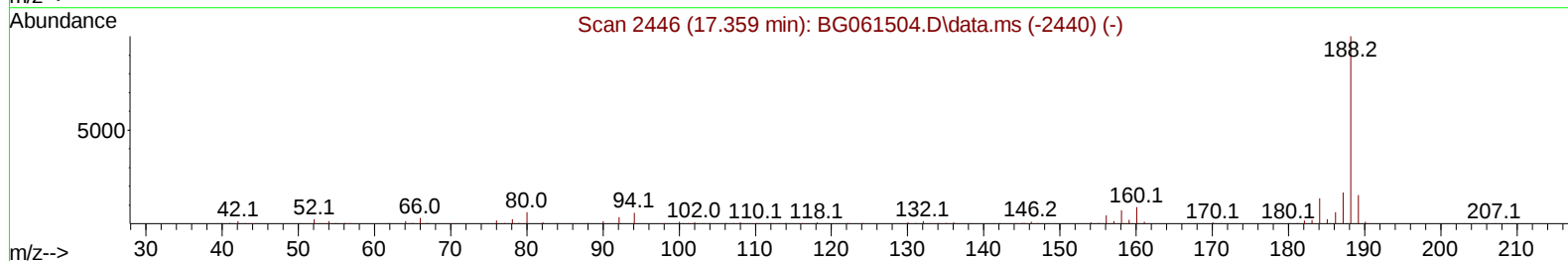
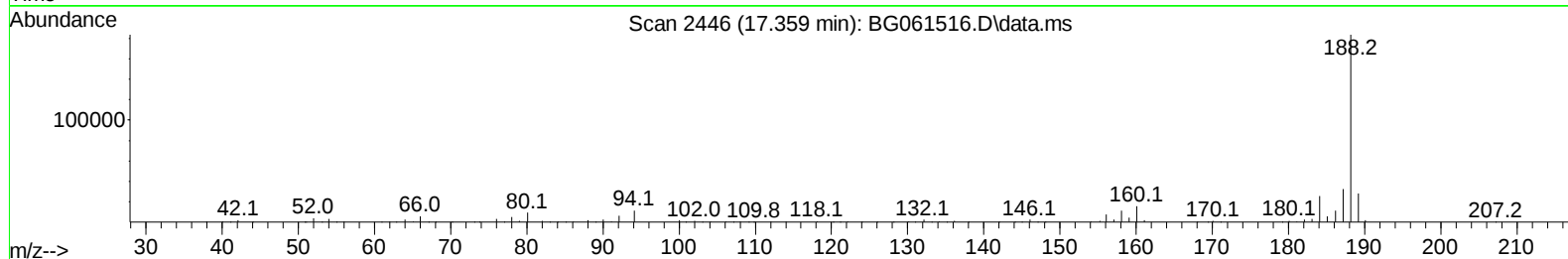
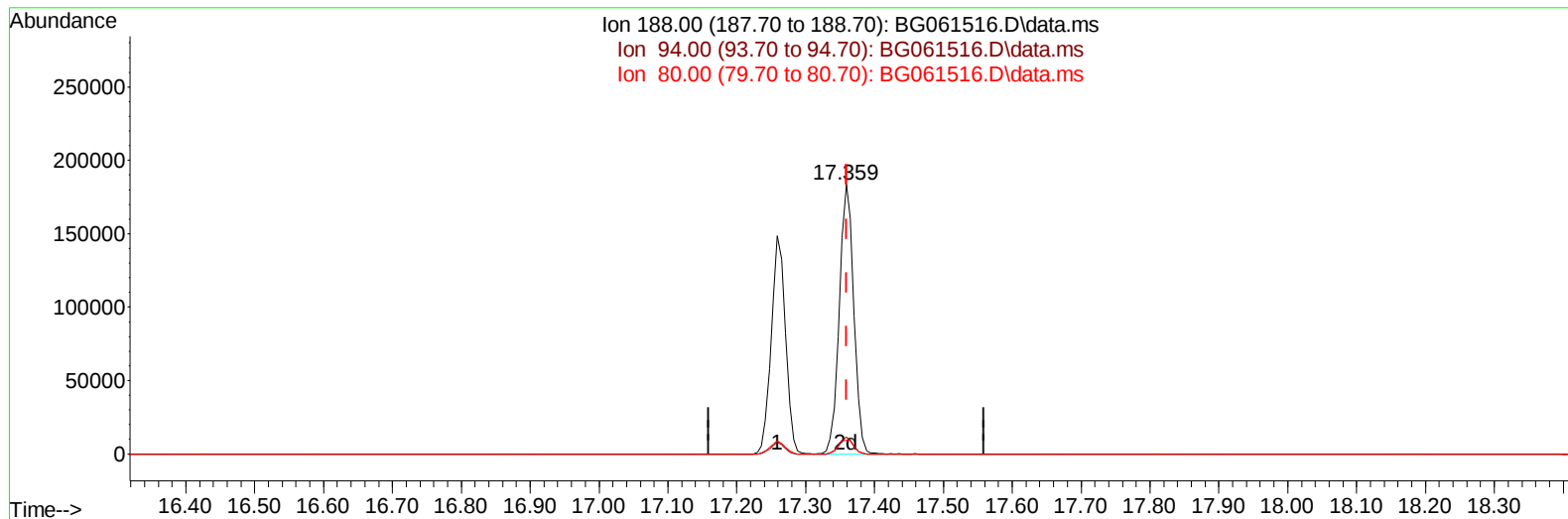
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TIC: BG061516.D\data.ms

(73) Anthracene-d10 (S)

17.359min (-0.000) 28.56 ng/u1 m

response 268409

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	6.60	6.21
80.00	6.40	5.04#
0.00	0.00	0.00

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 Operator : MA/JU
 Sample : PB161346BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK346

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/07/2024
 Supervised By :mohammad ahmed 06/08/2024

Quant Time: Jun 06 22:53:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 06 12:15:45 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.876	152		36970	20.000	ng/ul	0.00
20) Naphthalene-d8	10.667	136		130608	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.509	164		94244	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.259	188		213199	20.000	ng/ul	0.00
79) Chrysene-d12	21.519	240		235447	20.000	ng/ul	0.00
88) Perylene-d12	24.609	264		265059	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.369	96		3188	4.826	ng/uL	0.01
4) Pyridine-d5	3.769	84		48335	20.566	ng/ul	0.00
7) Phenol-d5	7.071	99		66012	21.743	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.206	67		41764	21.886	ng/ul	0.00
11) 2-Chlorophenol-d4	7.418	132		55726	24.801	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113		55633	21.490	ng/ul	0.00
21) Nitrobenzene-d5	9.039	128		28001	27.845	ng/ul	0.00
24) 2-Nitrophenol-d4	9.756	143		31789	27.011	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.308	165		59444	25.909	ng/ul	0.00
31) 4-Chloroaniline-d4	10.808	131		72259	24.073	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166		188067	25.730	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160		205279	27.014	ng/ul	0.00
54) 4-Nitrophenol-d4	14.762	143		18204	20.144	ng/ul	0.00
60) Fluorene-d10	15.502	176		159493	25.274	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.655	200		23602	18.304	ng/ul	0.00
73) Anthracene-d10	17.359	188		268409m	28.556	ng/ul	0.00
81) Pyrene-d10	19.656	212		324091	26.808	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.398	264		367760	28.860	ng/ul	0.00

Target Compounds Qvalue

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

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