

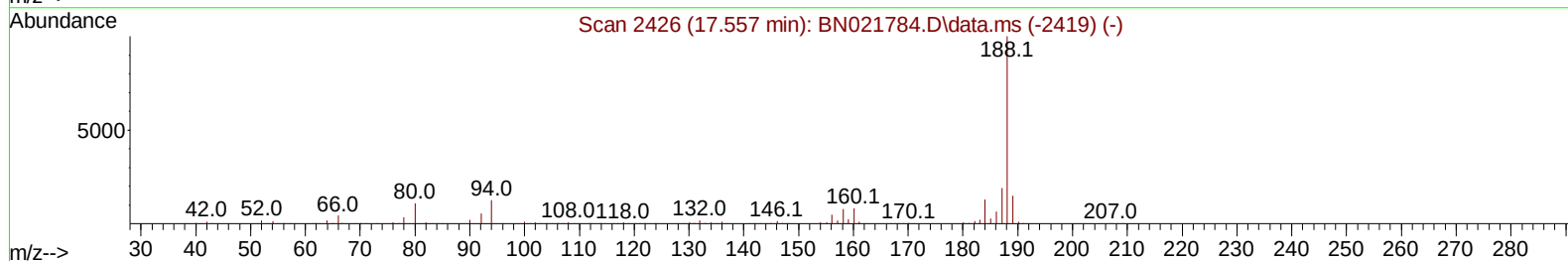
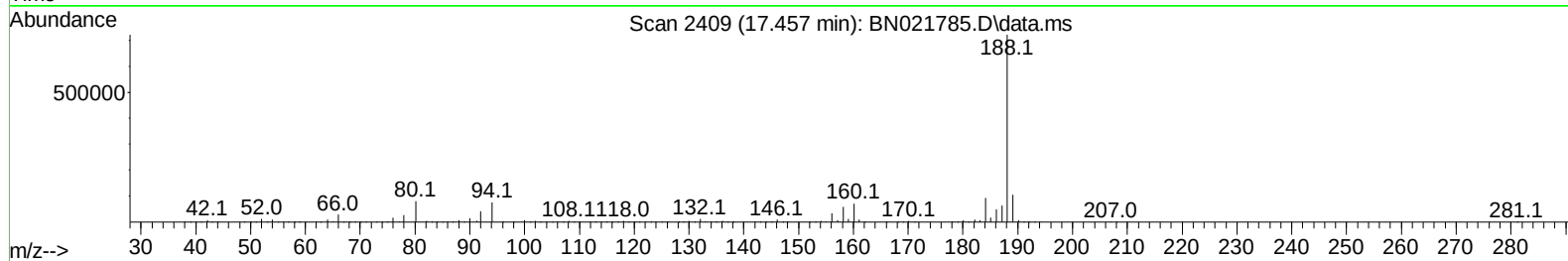
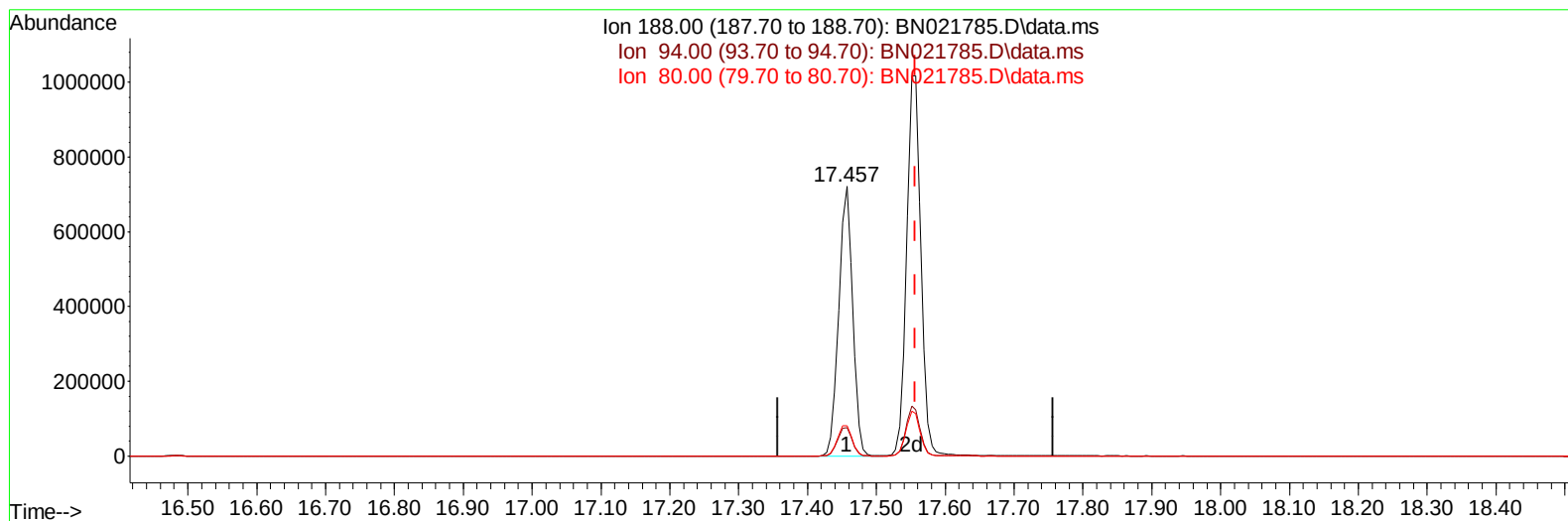
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091622\
Data File : BN021785.D
Acq On : 16 Sep 2022 12:50
Operator : CG/JU
Sample : PB147637BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK637

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/19/2022
Supervised By :mohammad ahmed 09/19/2022

Quant Time: Sep 17 00:20:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081922.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 17 00:17:38 2022
Response via : Initial Calibration



TIC: BN021785.D\data.ms

(73) Anthracene-d10 (S)

17.457min (-0.100) 22.69 ng/u1

response 1009084

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	10.90	10.59
80.00	9.70	11.27
0.00	0.00	0.00

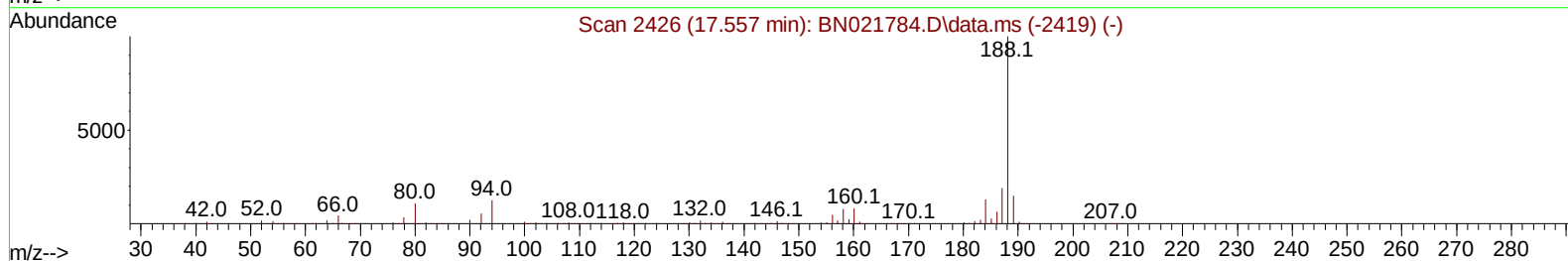
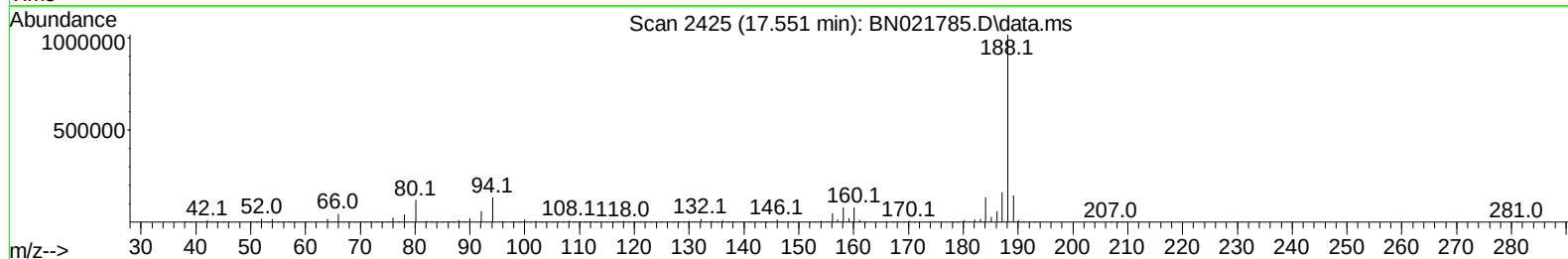
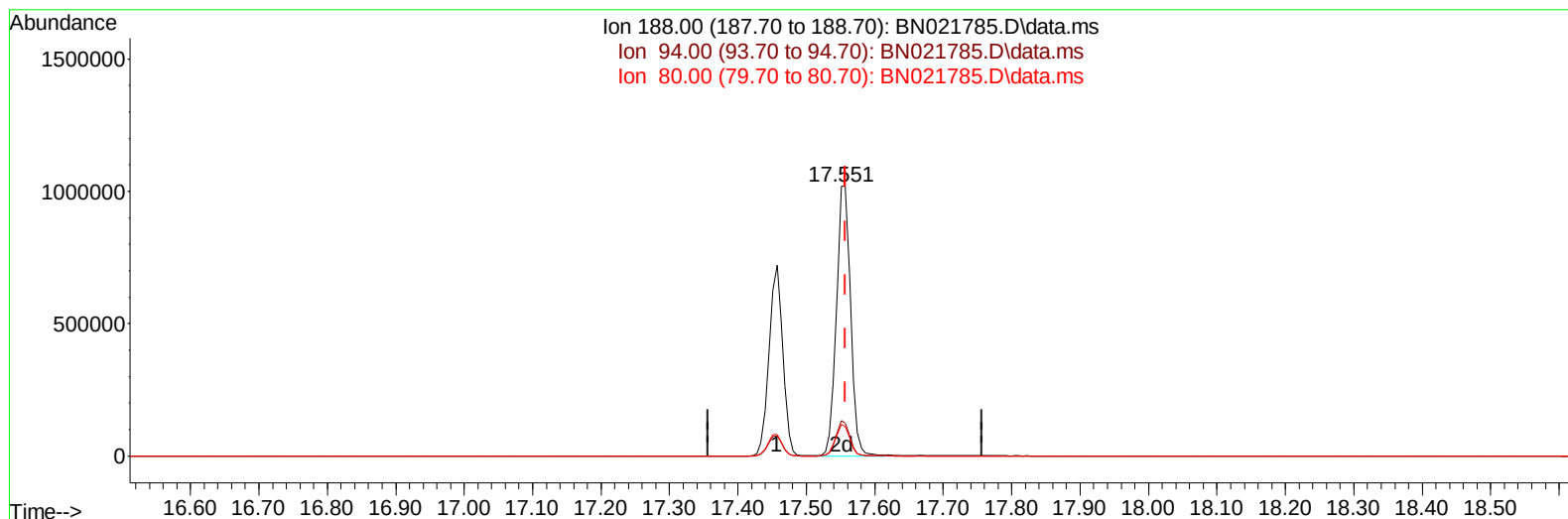
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TIC: BN021785.D\data.ms

(73) Anthracene-d10 (S)

17.551min (-0.006) 33.09 ng/ul m

response 1471429

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	10.90	13.16#
80.00	9.70	11.73#
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	8.046	152	159202	20.000	ng/ul	0.00
20) Naphthalene-d8	10.875	136	739871	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	473985	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.457	188	1009084	20.000	ng/ul	0.00
79) Chrysene-d12	21.645	240	867940	20.000	ng/ul	# 0.00
88) Perylene-d12	24.151	264	793378	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	25141	5.929	ng/uL	0.00
4) Pyridine-d5	3.787	84	317905	27.509	ng/ul	0.00
7) Phenol-d5	7.199	99	426571	29.543	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	272632	31.193	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	332029	31.606	ng/ul	0.00
15) 4-Methylphenol-d8	8.763	113	348313	30.744	ng/ul	0.00
21) Nitrobenzene-d5	9.222	128	171023	33.927	ng/ul	0.00
24) 2-Nitrophenol-d4	9.952	143	180347	40.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.493	165	318985	31.990	ng/ul	0.00
31) 4-Chloroaniline-d4	11.016	131	516252	30.485	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	1110245	34.270	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	1258557	33.323	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	180621	28.017	ng/ul	0.00
60) Fluorene-d10	15.698	176	933722	33.123	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	112013	27.752	ng/ul	0.00
73) Anthracene-d10	17.551	188	1471429m	33.089	ng/ul	0.00
81) Pyrene-d10	19.851	212	1639724	39.726	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.992	264	1382285	36.758	ng/ul	0.00

Target Compounds Qvalue

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

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