

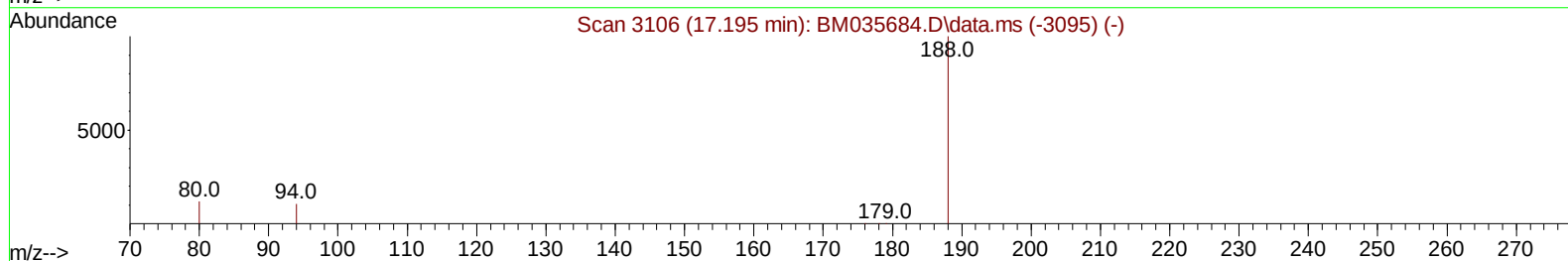
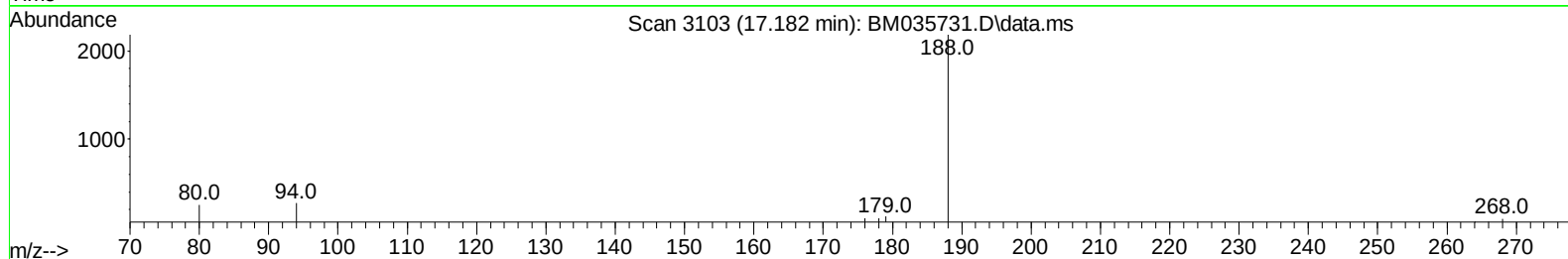
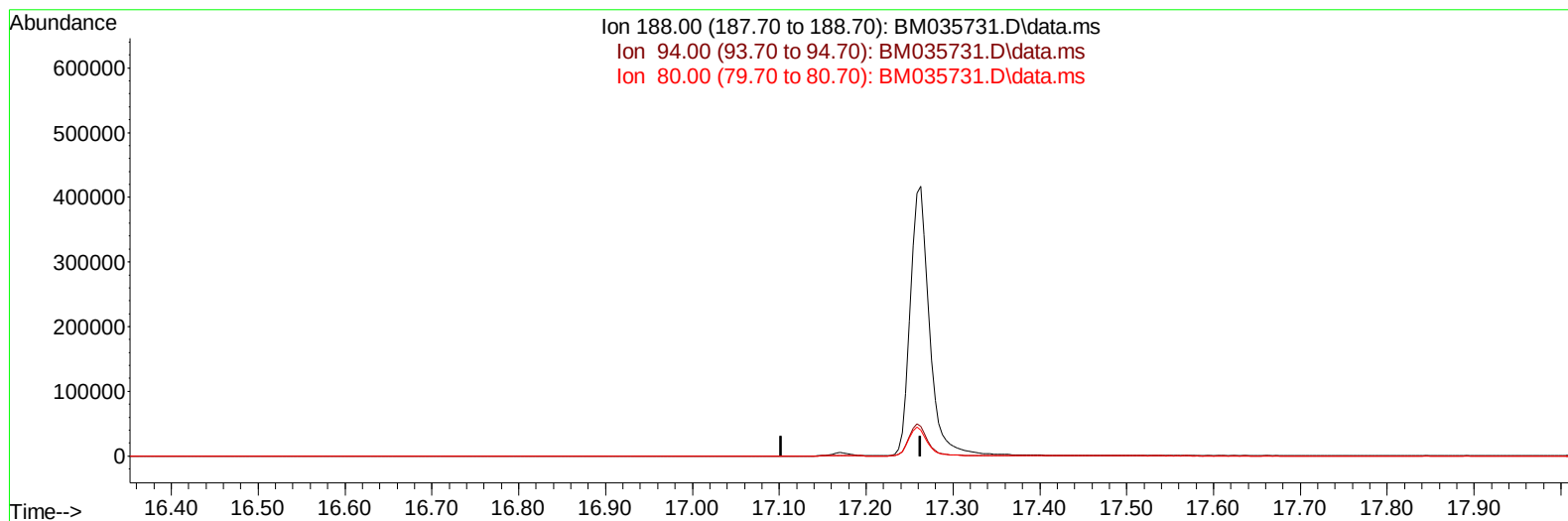
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
 Data File : BM035731.D
 Acq On : 28 Jun 2022 19:25
 Operator : CG/JU
 Sample : N3401-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AD7

Manual IntegrationsAPPROVED

Quant Time: Jun 29 04:27:40 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 29 04:24:18 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/29/2022
 Supervised By :mohammad ahmed 07/01/2022



TIC: BM035731.D\data.ms

(13) Phenanthrene-d10 (I)

17.182min (-17.182) 0.00 ng/ul

response 0

Ion	Exp%	Act%
188.00	100.00	0.00
94.00	11.40	0.00#
80.00	12.70	0.00#
0.00	0.00	0.00

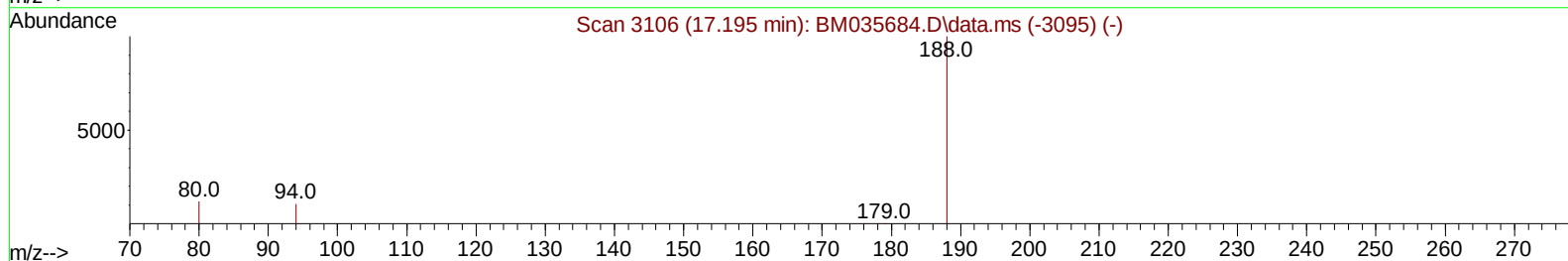
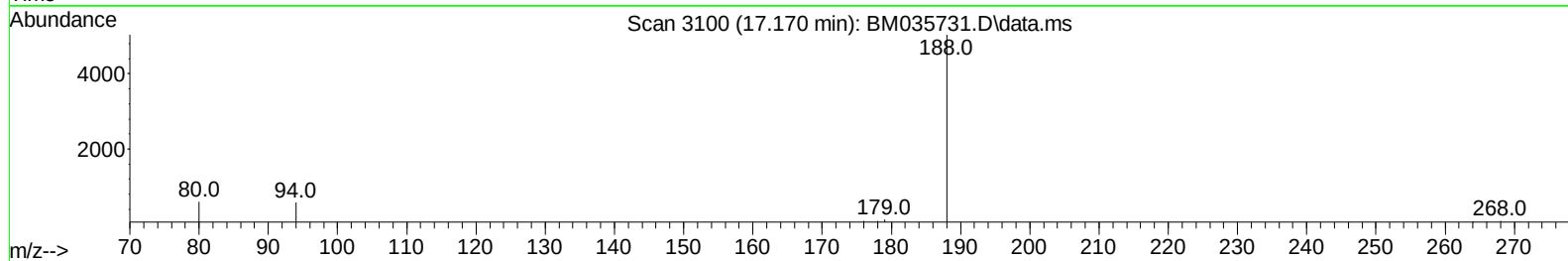
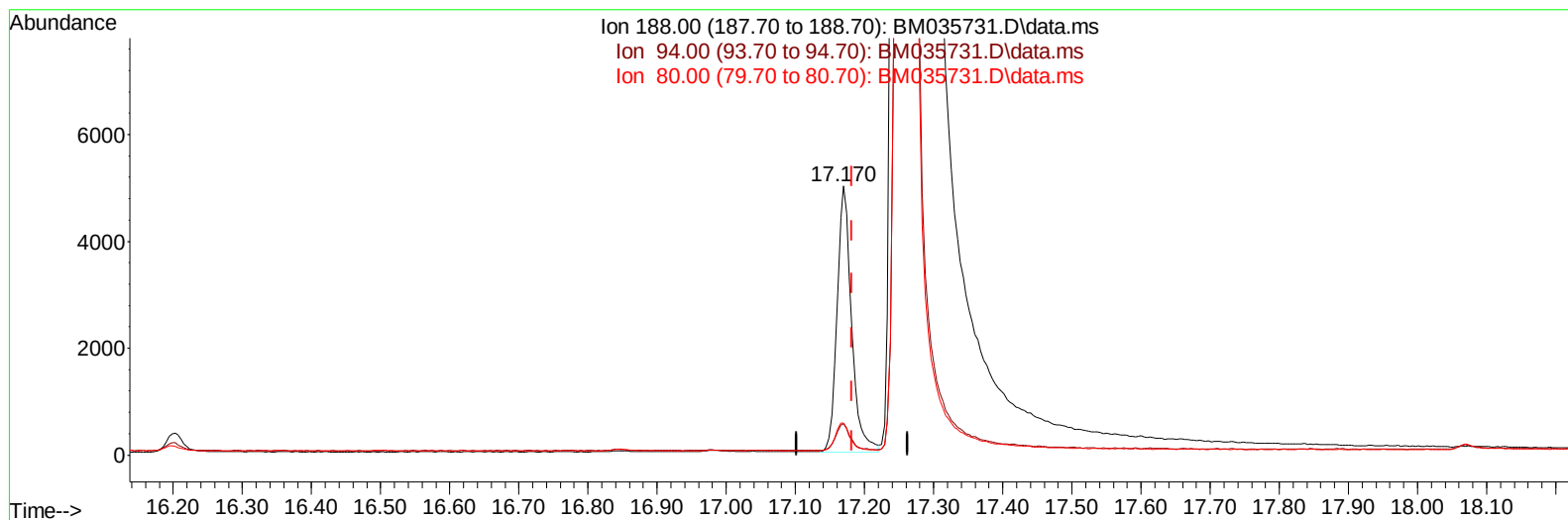
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(13) Phenanthrene-d10 (I)

17.170min (-0.013) 0.40 ng/ul m

response 7264

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	11.40	11.66
80.00	12.70	12.00
0.00	0.00	0.00

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 ALS Vial : 10 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.795	152		2142	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.578	136		6901	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.422	164		3713	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.170	188		7264m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.374	240		6514	0.400	ng/ul	# 0.00
23) Perylene-d12	23.689	264		8189	0.400	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.209	96		13041	4.486	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152		3317	0.307	ng/ul	0.01
18) Fluoranthene-d10	19.205	212		7442	0.339	ng/ul	0.00
Target Compounds							Qvalue
21) Benzo(a)anthracene	21.368	228		590	0.026	ng/ul#	31
22) Chrysene	21.418	228		924	0.032	ng/ul#	46
24) Benzo(b)fluoranthene	22.984	252		1368	0.039	ng/ul#	1
25) Benzo(k)fluoranthene	23.031	252		1192	0.033	ng/ul#	1

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

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