

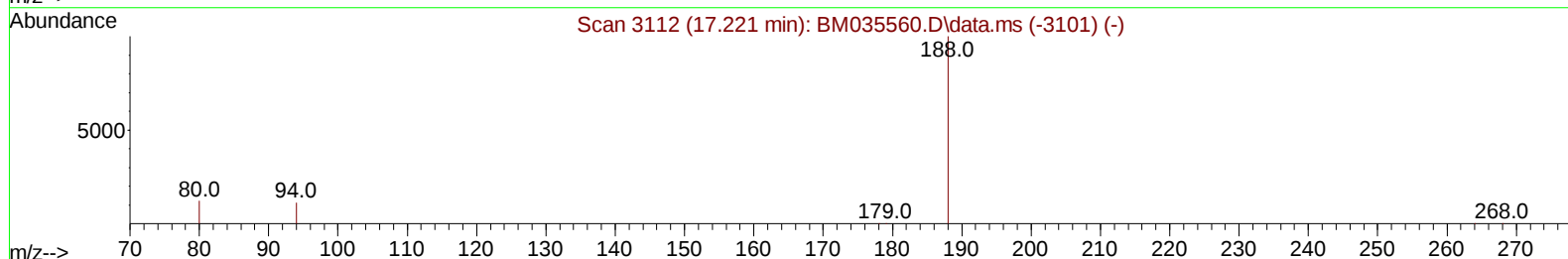
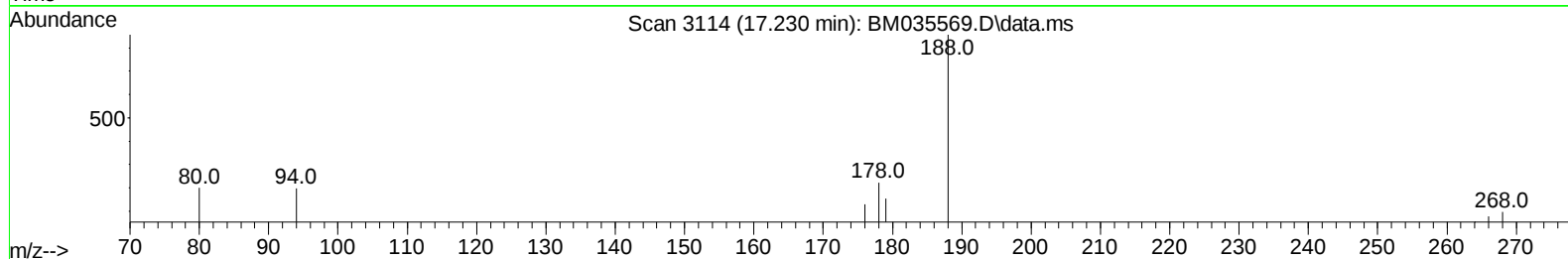
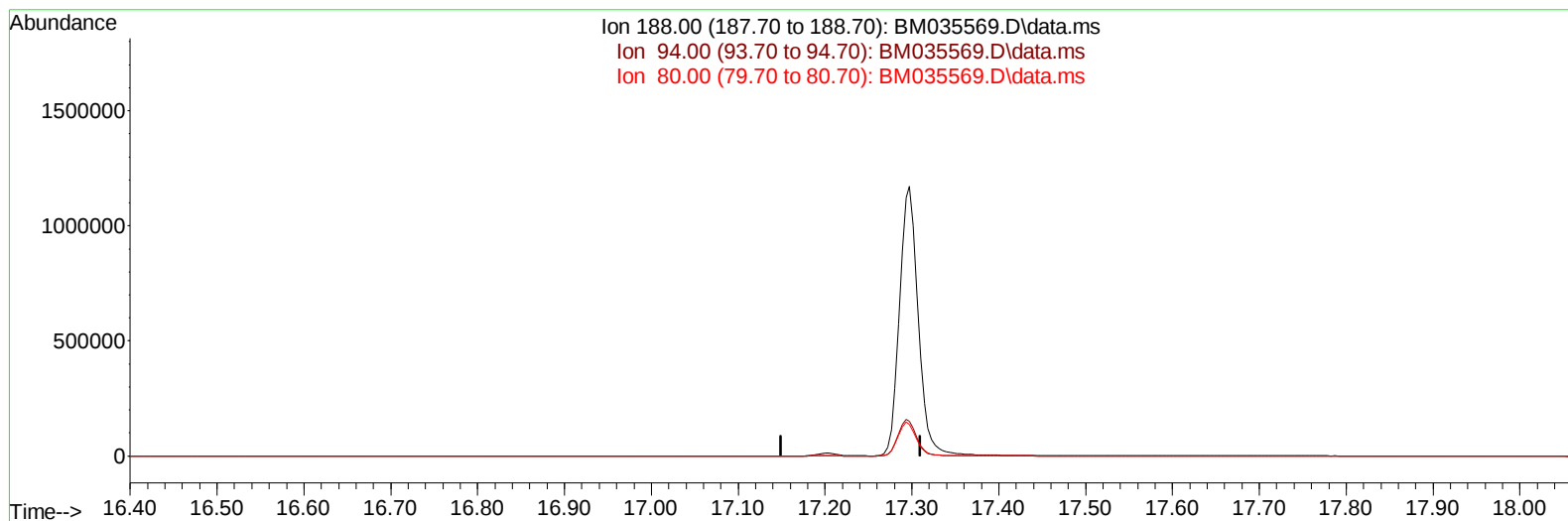
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035569.D
 Acq On : 21 Jun 2022 13:07
 Operator : CG/JU
 Sample : N3285-12
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4Q5

Manual IntegrationsAPPROVED

Quant Time: Jun 22 01:32:27 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/22/2022
 Supervised By :Yogesh Patel 06/25/2022



TIC: BM035569.D\data.ms

(13) Phenanthrene-d10 (I)

17.230min (-17.230) 0.00 ng/u1

response 0

Ion	Exp%	Act%
188.00	100.00	0.00
94.00	12.00	0.00#
80.00	13.50	0.00#
0.00	0.00	0.00

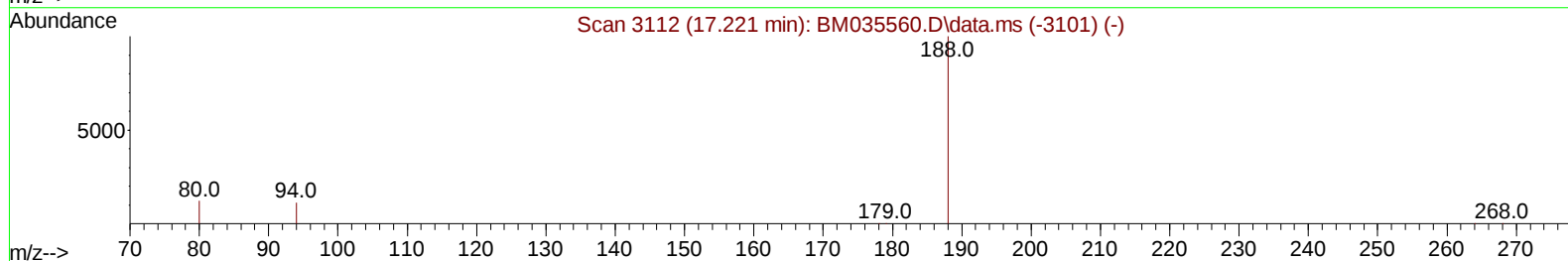
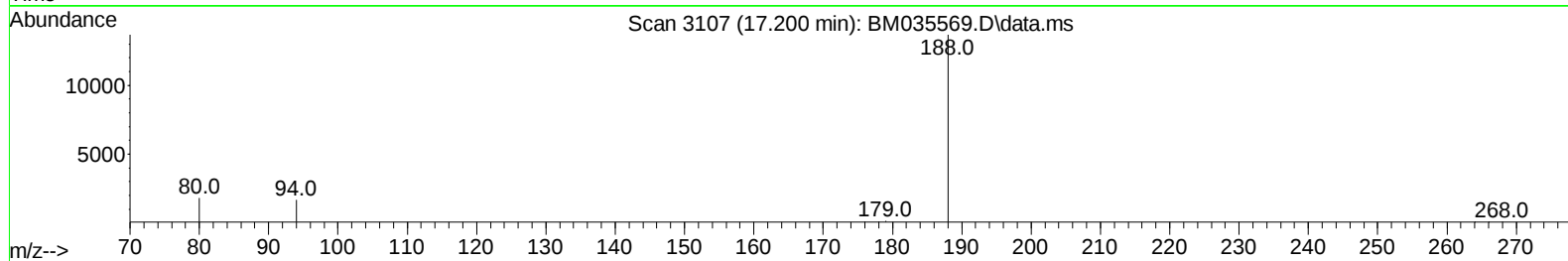
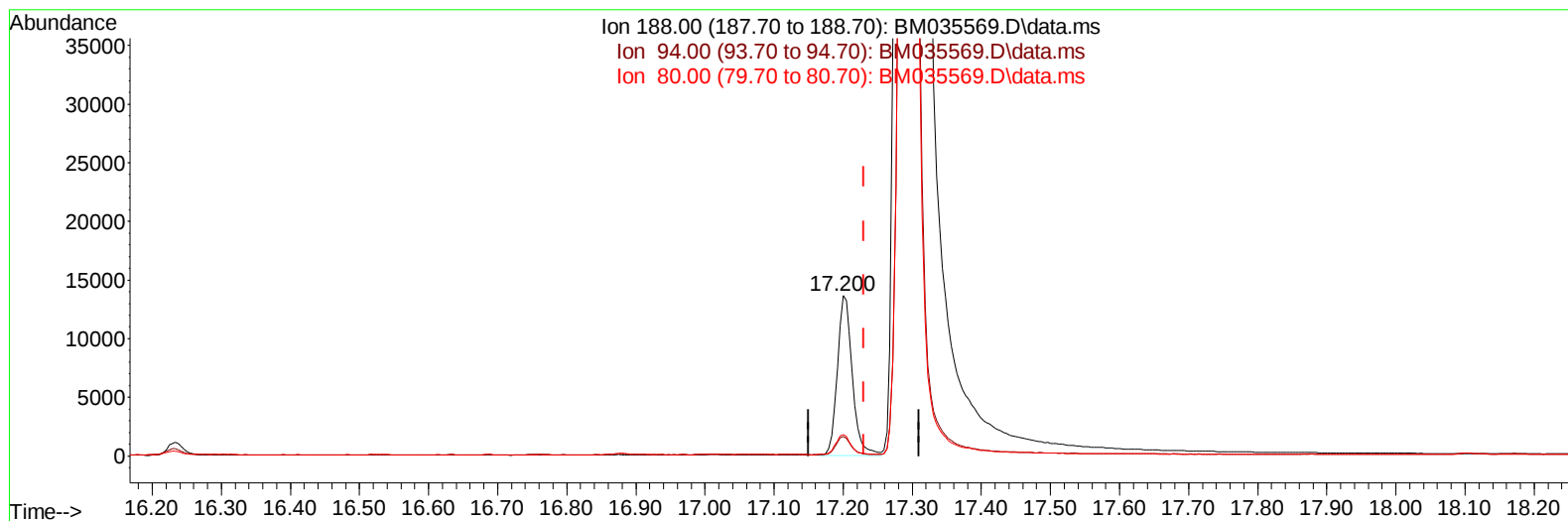
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TIC: BM035569.D\data.ms

(13) Phenanthrene-d10 (I)

17.200min (-0.030) 0.40 ng/ul m

response 20314

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	12.00	12.13
80.00	13.50	13.17
0.00	0.00	0.00

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 ALS Vial : 48 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.825	152		5253	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.611	136		19480	0.400	ng/ul	#-0.03
9) Acenaphthene-d10	14.456	164		10444	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.200	188		20314m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.406	240		16401	0.400	ng/ul	0.00
23) Perylene-d12	23.735	264		19877	0.400	ng/ul	#-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96		30473	4.123	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.233	152		9714	0.336	ng/ul	-0.01
18) Fluoranthene-d10	19.234	212		20633	0.374	ng/ul	-0.02

Target Compounds Qvalue

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

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