

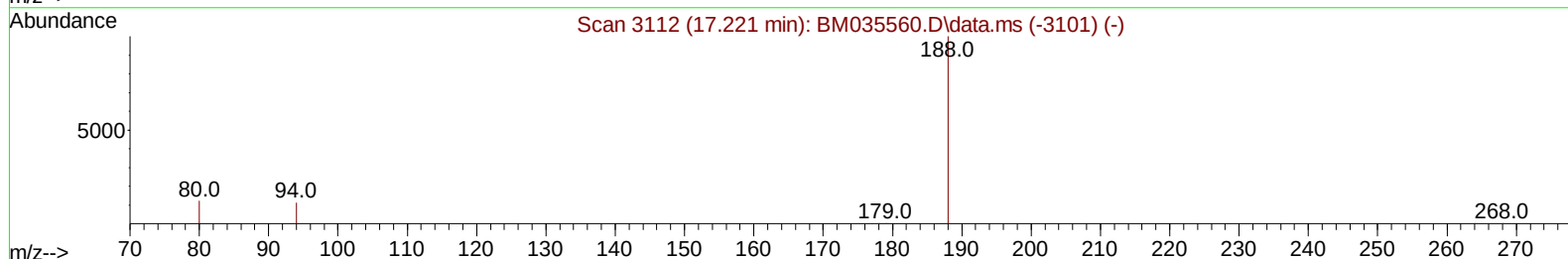
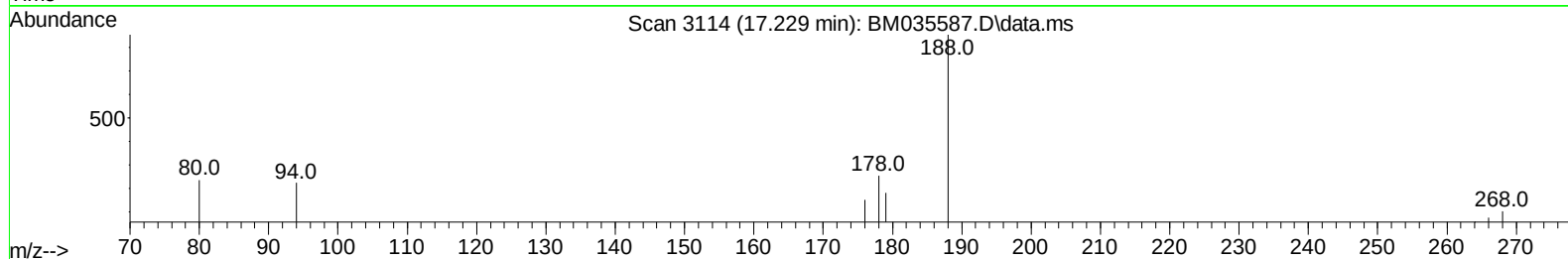
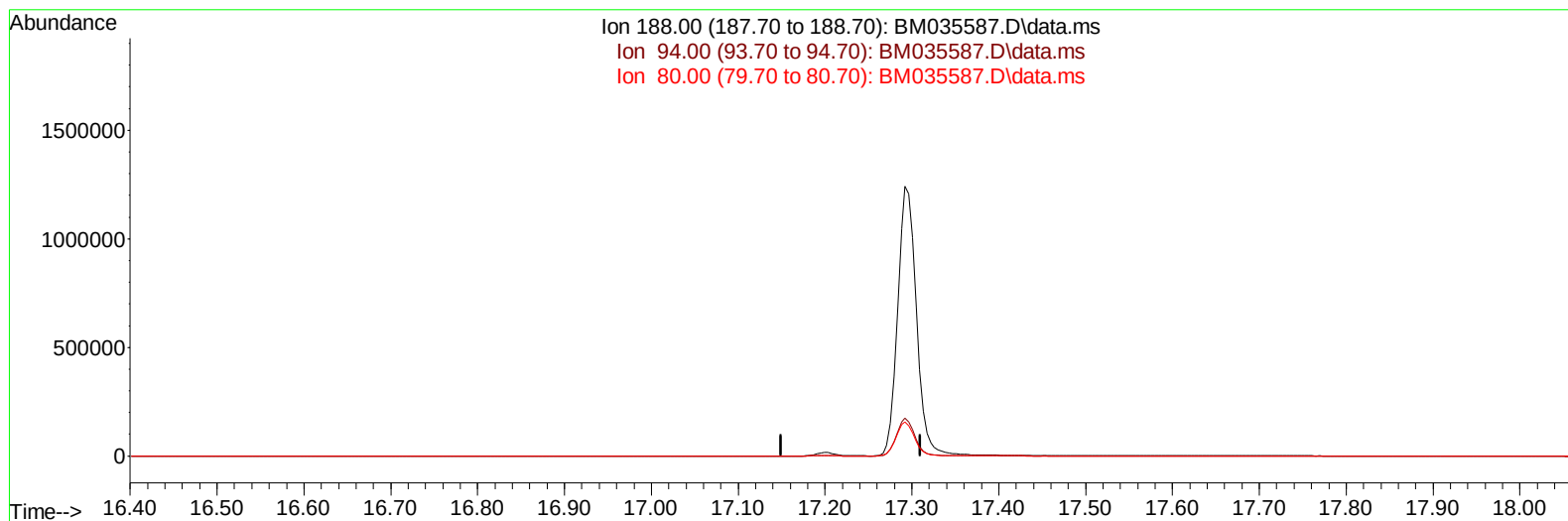
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035587.D
 Acq On : 22 Jun 2022 00:38
 Operator : CG/JU
 Sample : N3285-21
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4R3

Manual IntegrationsAPPROVED

Quant Time: Jun 22 01:33:43 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: BM035587.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

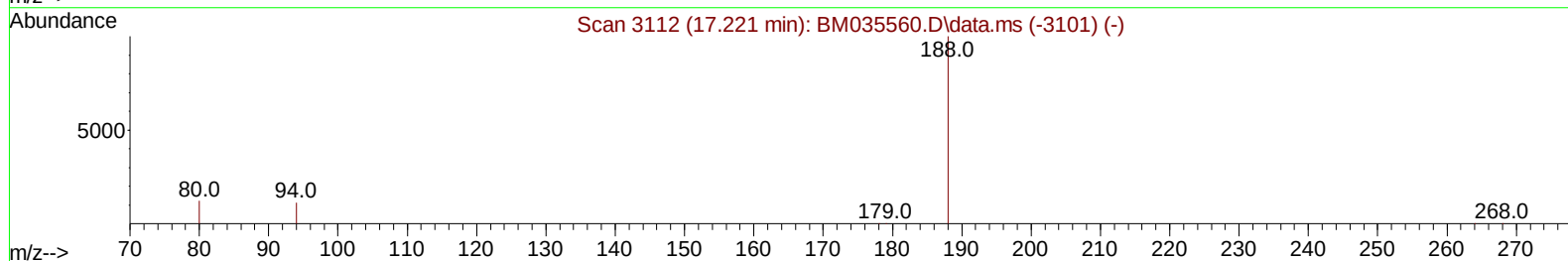
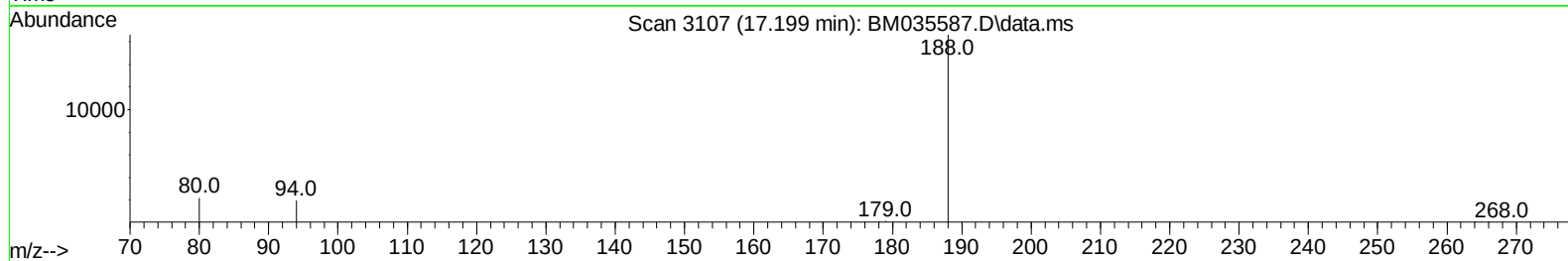
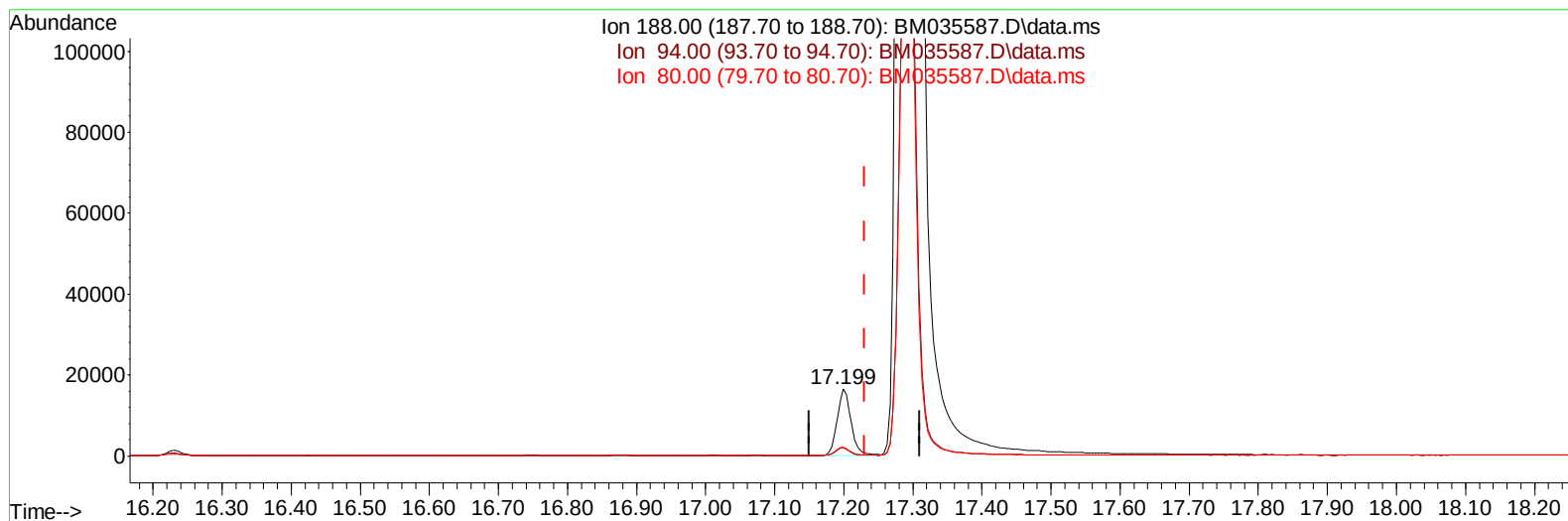
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TIC: BM035587.D\data.ms

(13) Phenanthrene-d10 (I)

17.199min (-0.031) 0.40 ng/ul m

response 23315

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	12.00	11.84
80.00	13.50	13.32
0.00	0.00	0.00

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 ALS Vial : 11 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	7.820	152		5319	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.611	136		19217	0.400	ng/ul	#-0.03
9) Acenaphthene-d10	14.450	164		12195	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.199	188		23315m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.400	240		16367	0.400	ng/ul	-0.01
23) Perylene-d12	23.730	264		18606	0.400	ng/ul	#-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96		30039	4.014	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.217	152		10315	0.361	ng/ul	-0.03
18) Fluoranthene-d10	19.233	212		21347	0.388	ng/ul	-0.02

Target Compounds Qvalue

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

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