

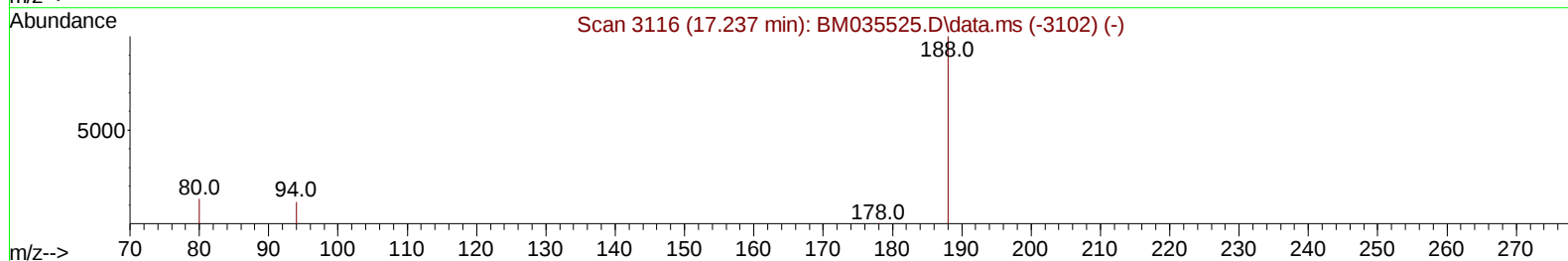
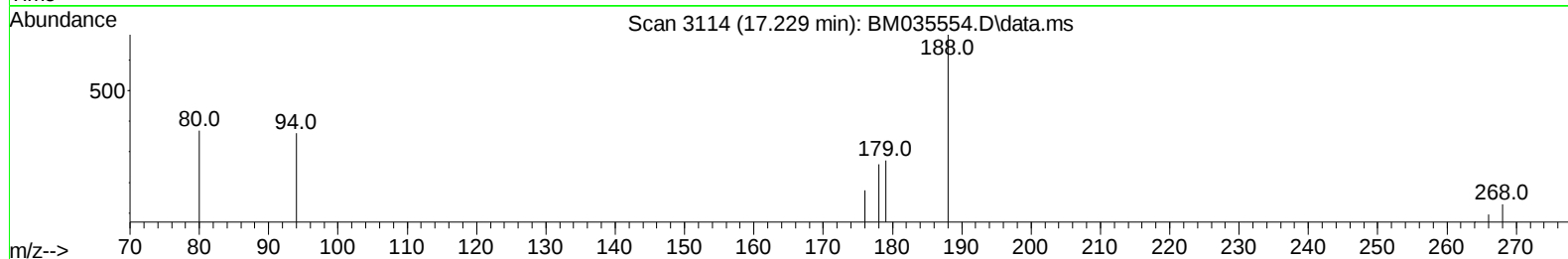
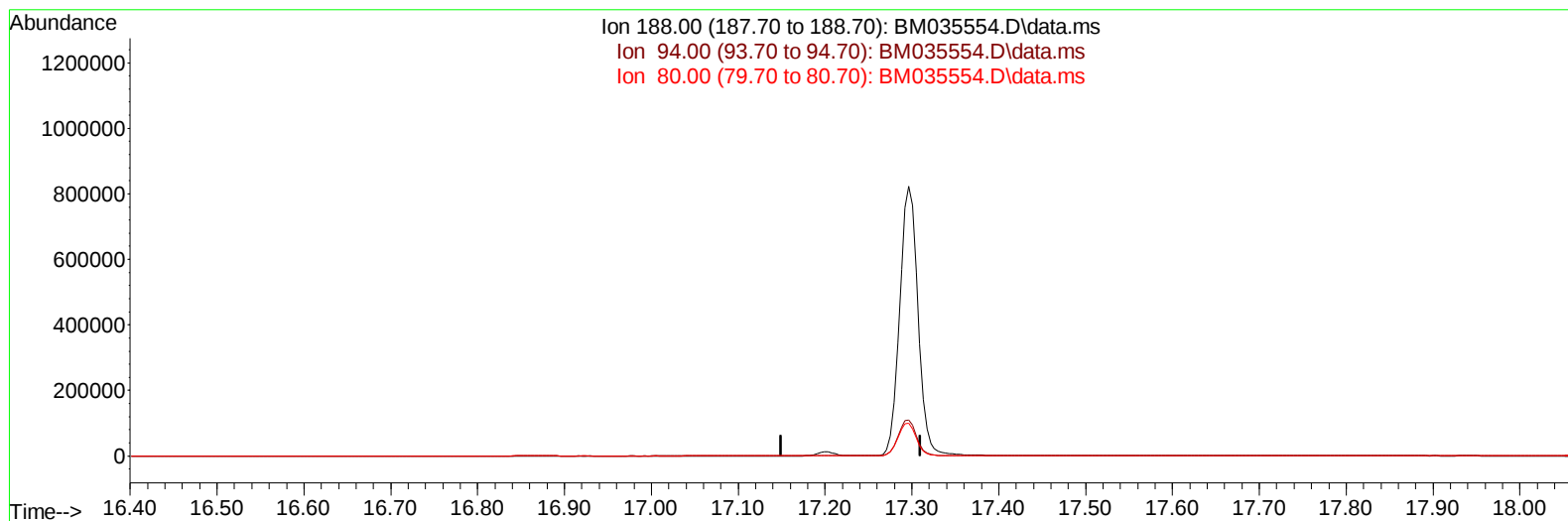
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035554.D
 Acq On : 21 Jun 2022 03:32
 Operator : CG/JU
 Sample : N3285-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4Q2

Manual IntegrationsAPPROVED

Quant Time: Jun 21 05:09:43 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 06/21/2022
 Supervised By :Yogesh Patel 06/25/2022



TIC: BM035554.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

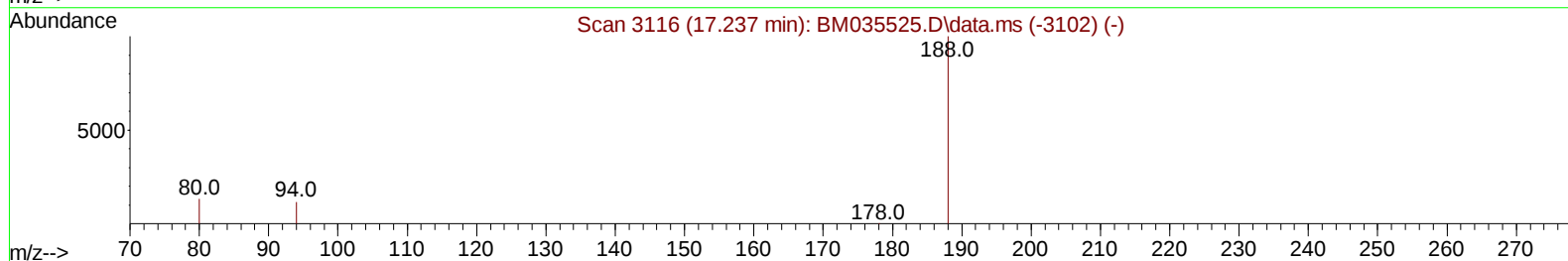
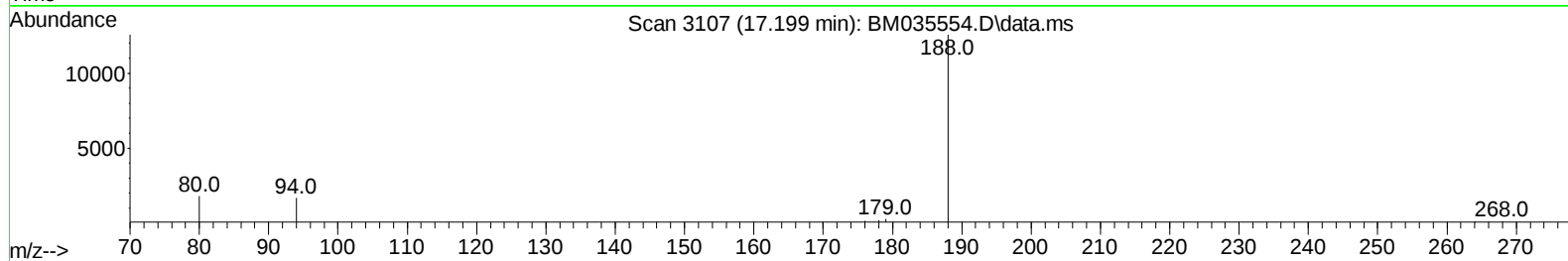
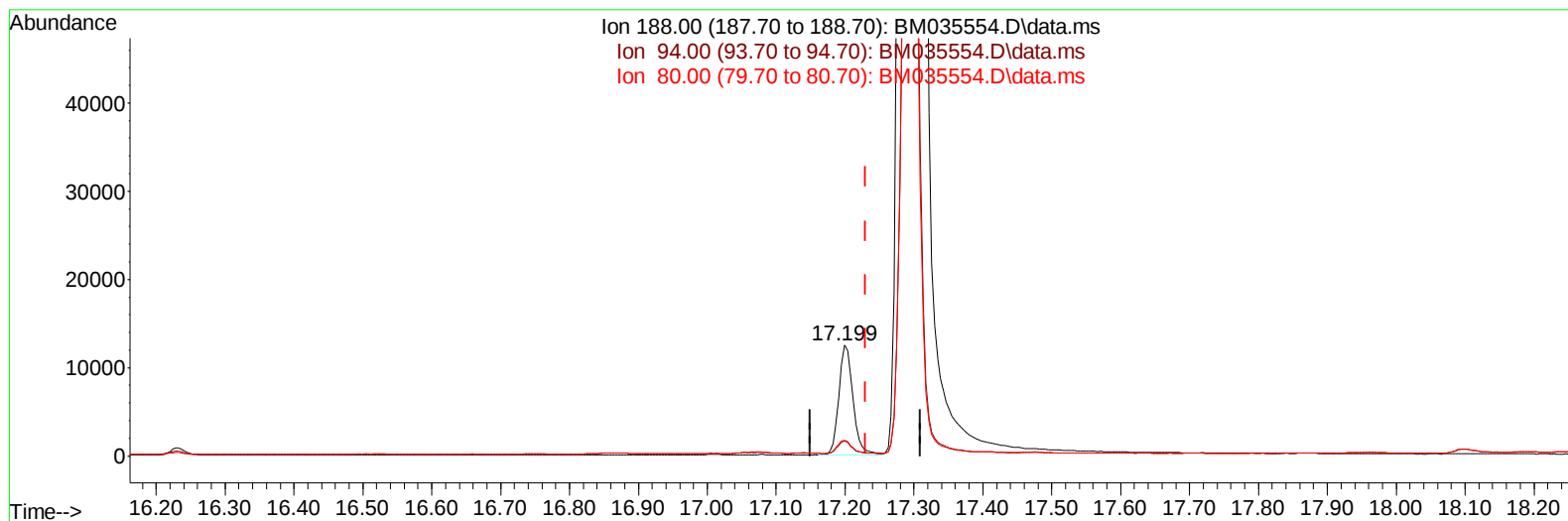
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(13) Phenanthrene-d10 (I)

17.199min (-0.030) 0.40 ng/ul m

response 17783

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	12.00	13.34
80.00	13.50	14.36
0.00	0.00	0.00

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 Misc :
 ALS Vial : 32 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		4605	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.611	136		15641	0.400	ng/ul	#-0.03
9) Acenaphthene-d10	14.455	164		8783	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.199	188		17783m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.392	240		16298	0.400	ng/ul	-0.02
23) Perylene-d12	23.730	264		17902	0.400	ng/ul	#-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.268	96		25704	3.967	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.217	152		6566	0.282	ng/ul	-0.03
18) Fluoranthene-d10	19.233	212		16507	0.302	ng/ul	-0.02

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

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