

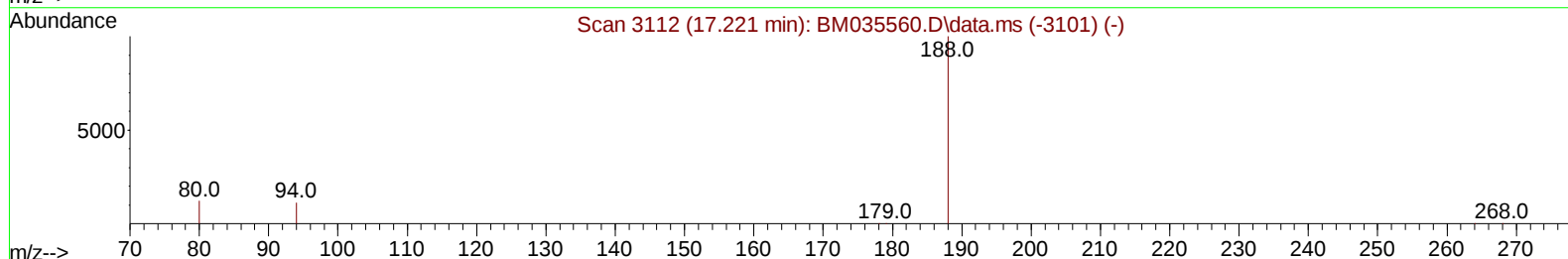
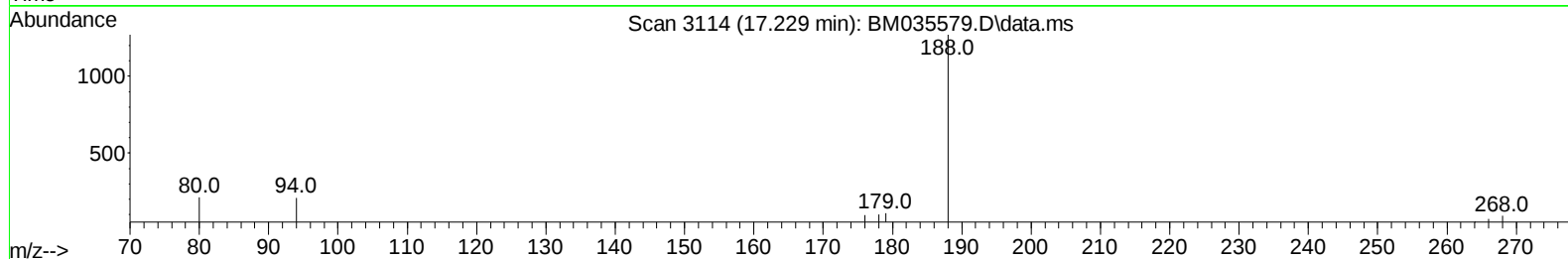
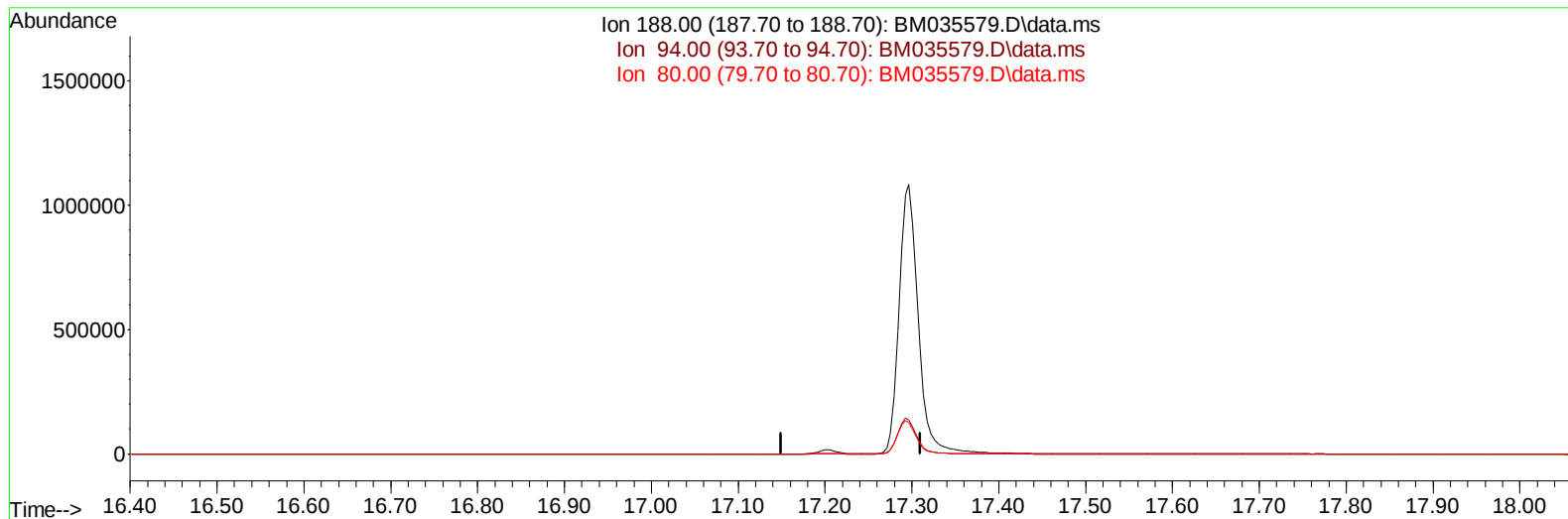
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
Data File : BM035579.D
Acq On : 21 Jun 2022 19:48
Operator : CG/JU
Sample : PB145592BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK592

Manual IntegrationsAPPROVED

Quant Time: Jun 22 01:33:12 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: BM035579.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

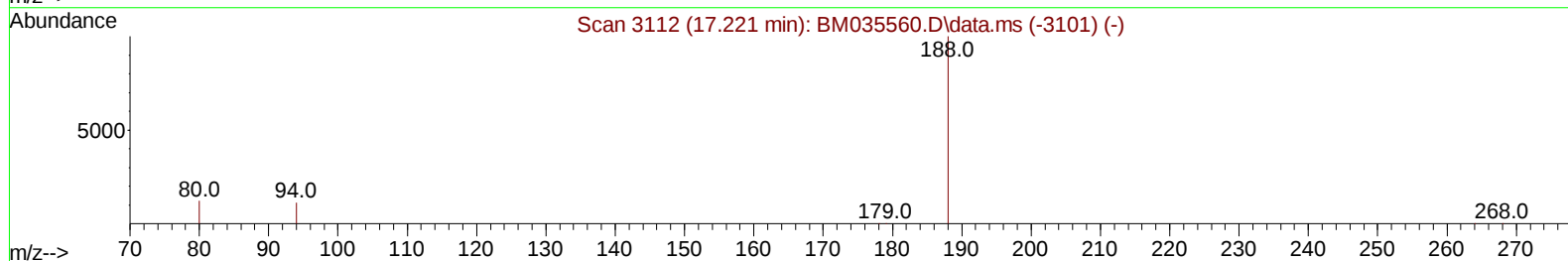
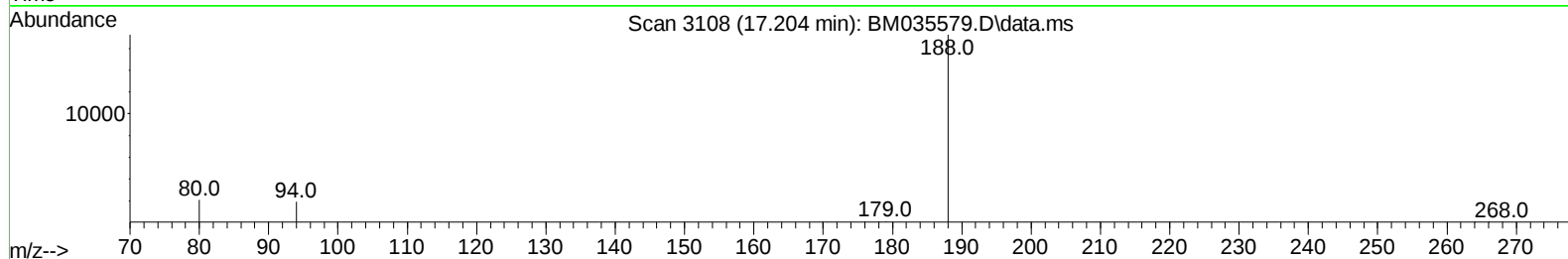
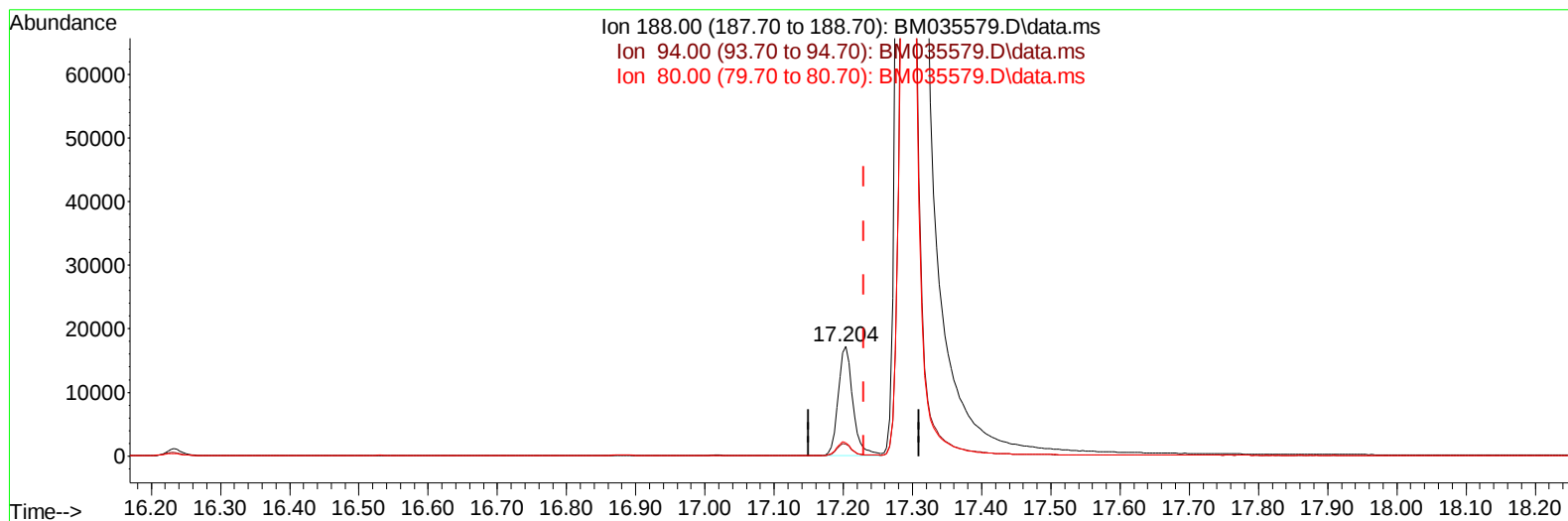
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TIC: BM035579.D\data.ms

(13) Phenanthrene-d10 (I)

17.204min (-0.026) 0.40 ng/ul m

response 25301

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	12.00	11.05
80.00	13.50	12.01
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : PB145592BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 SBLK592

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.825	152		5725	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.611	136		19741	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.455	164		11843	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.204	188		25301m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.415	240		16392	0.400	ng/ul	0.00
23) Perylene-d12	23.733	264		18162	0.400	ng/ul	#-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96		38163	4.738	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152		8181	0.279	ng/ul	0.00
18) Fluoranthene-d10	19.238	212		17589	0.319	ng/ul	-0.01

Target Compounds Qvalue

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

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