

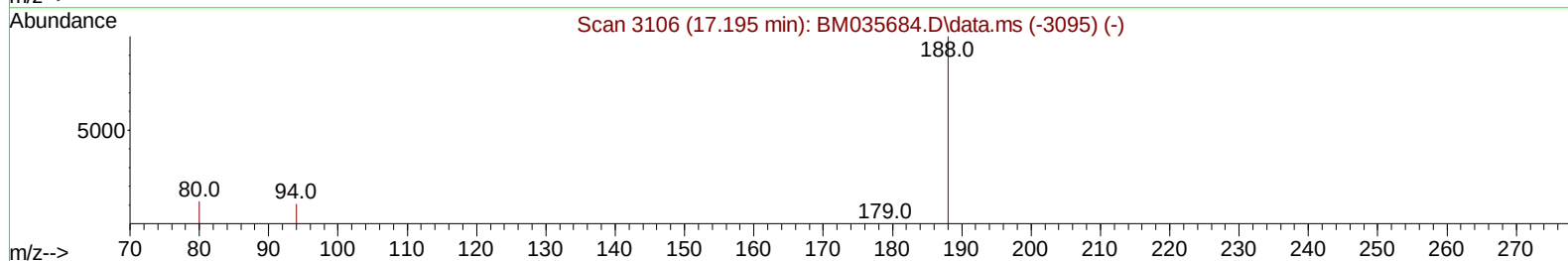
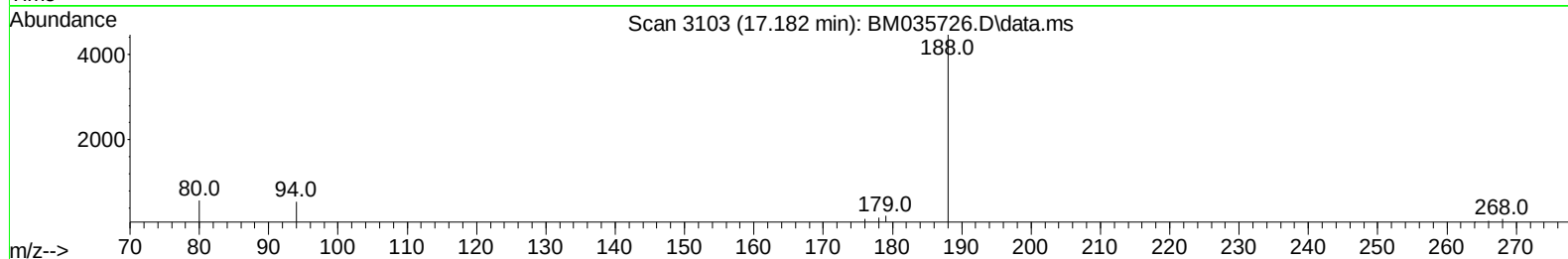
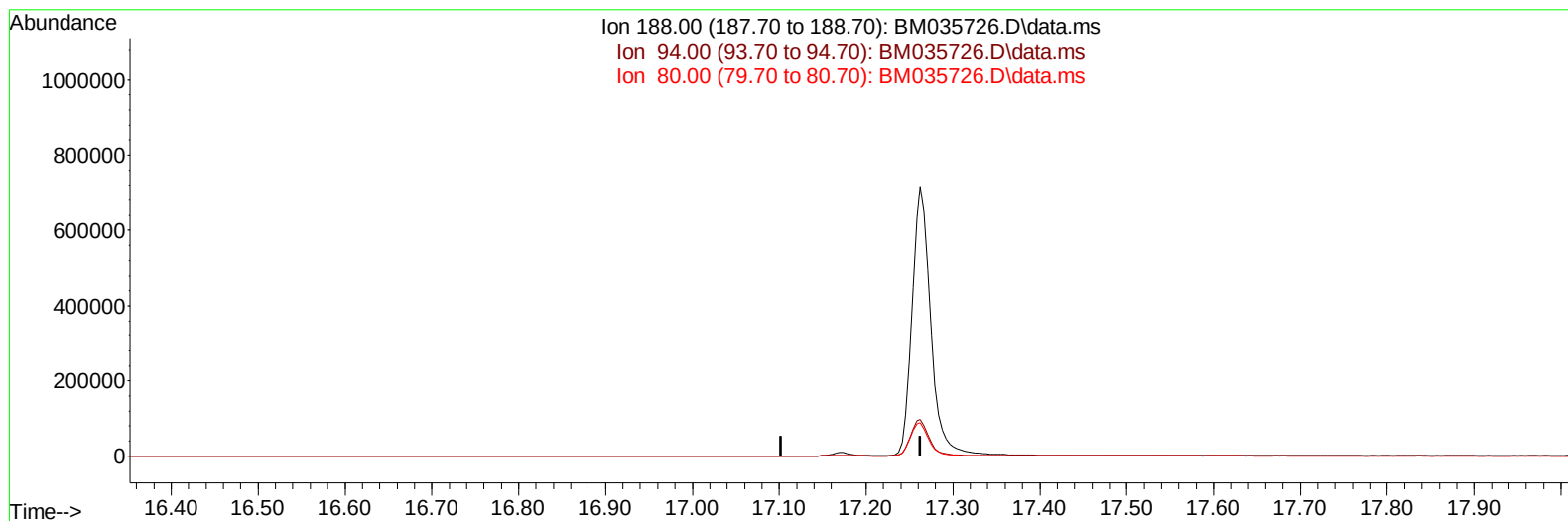
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
Data File : BM035726.D
Acq On : 28 Jun 2022 16:17
Operator : CG/JU
Sample : N3401-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AE0

Manual IntegrationsAPPROVED

Quant Time: Jun 29 04:27:11 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: BM035726.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

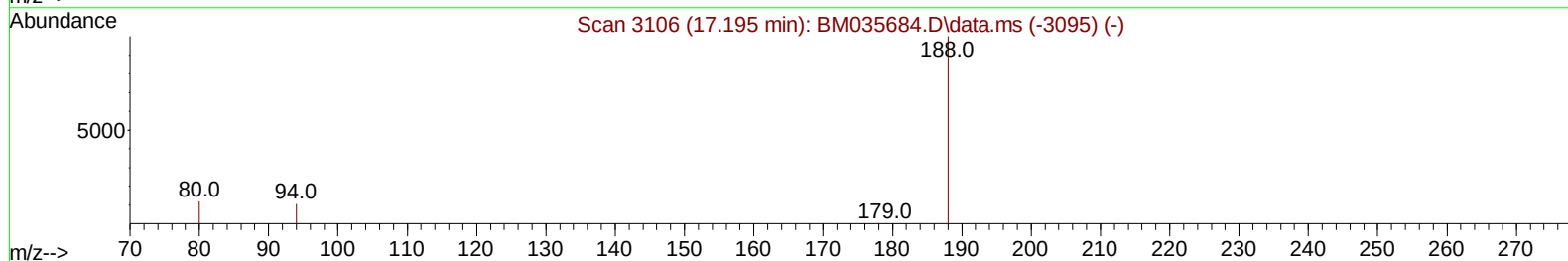
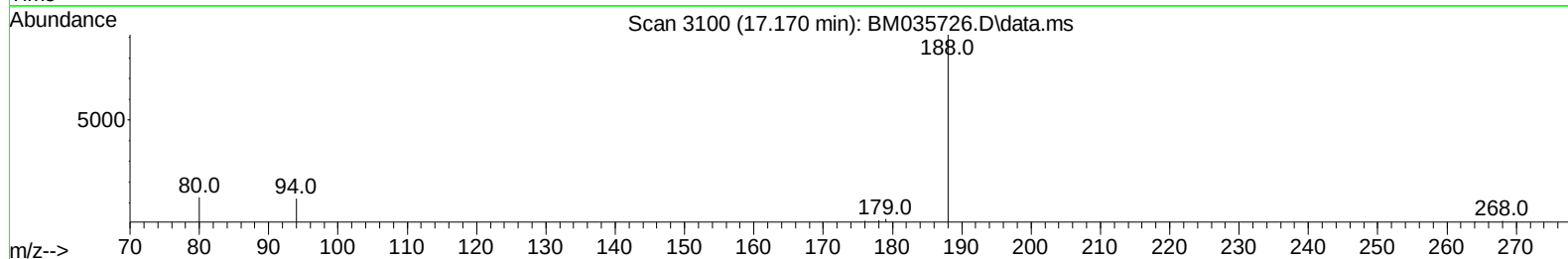
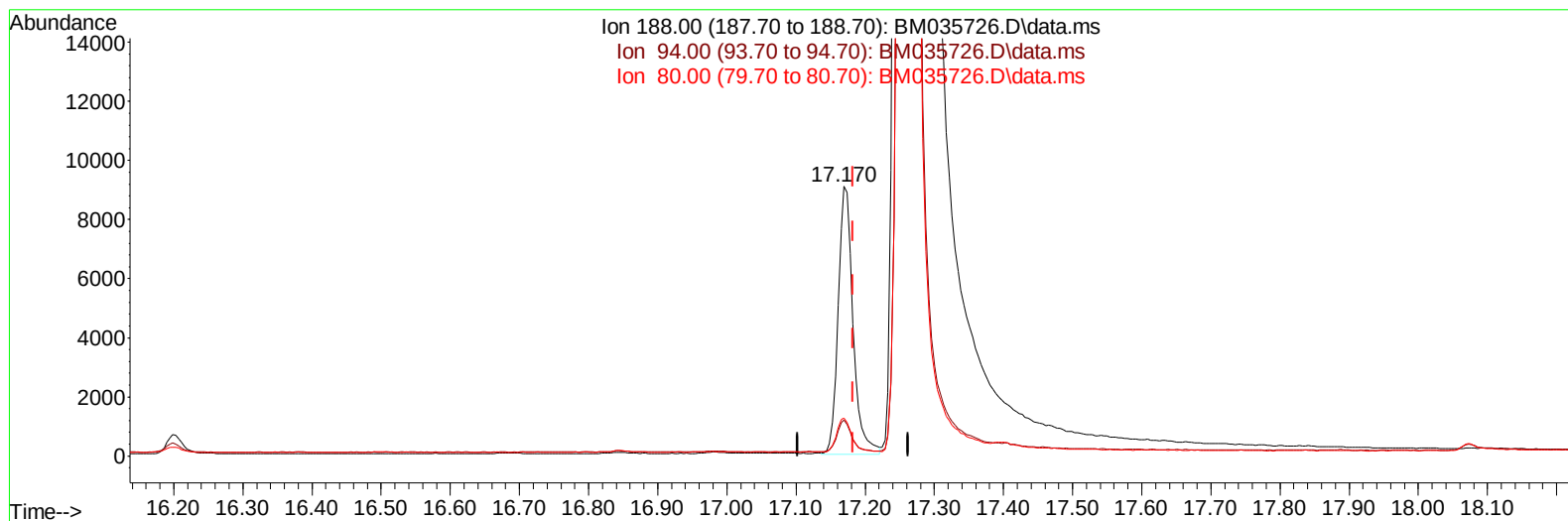
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(13) Phenanthrene-d10 (I)

17.170min (-0.013) 0.40 ng/ul m

response 13474

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	11.40	13.31
80.00	12.70	14.01
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : N3401-01
 Misc :
 ALS Vial : 5 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	7.791	152		3854	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.578	136		12128	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.422	164		6858	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.170	188		13474m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.377	240		12178	0.400	ng/ul	# 0.00
23) Perylene-d12	23.695	264		15612	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.213	96		29068	5.558	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152		5252	0.277	ng/ul	0.01
18) Fluoranthene-d10	19.205	212		12985	0.316	ng/ul	0.00
Target Compounds							Qvalue
22) Chrysene	21.421	228		2205	0.041	ng/ul#	52
24) Benzo(b)fluoranthene	22.990	252		2705	0.040	ng/ul#	1
25) Benzo(k)fluoranthene	23.034	252		2346	0.034	ng/ul#	1

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

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