

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
Data File : BM035727.D
Acq On : 28 Jun 2022 16:55
Operator : CG/JU
Sample : N3401-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

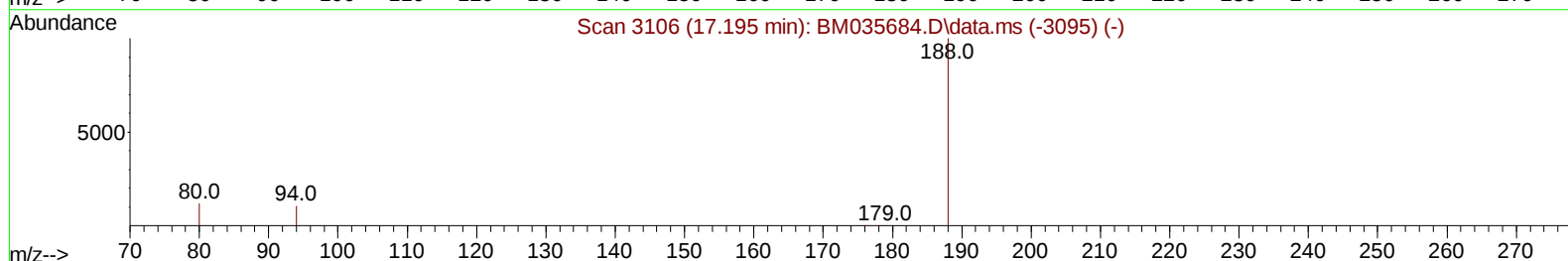
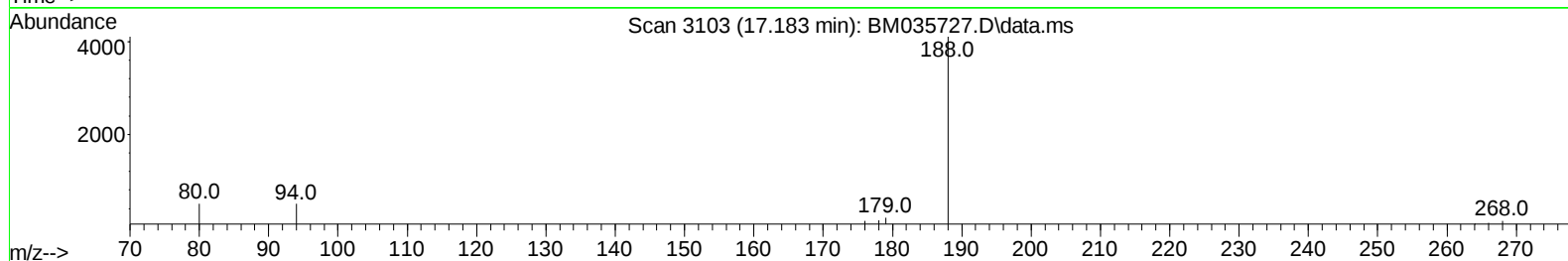
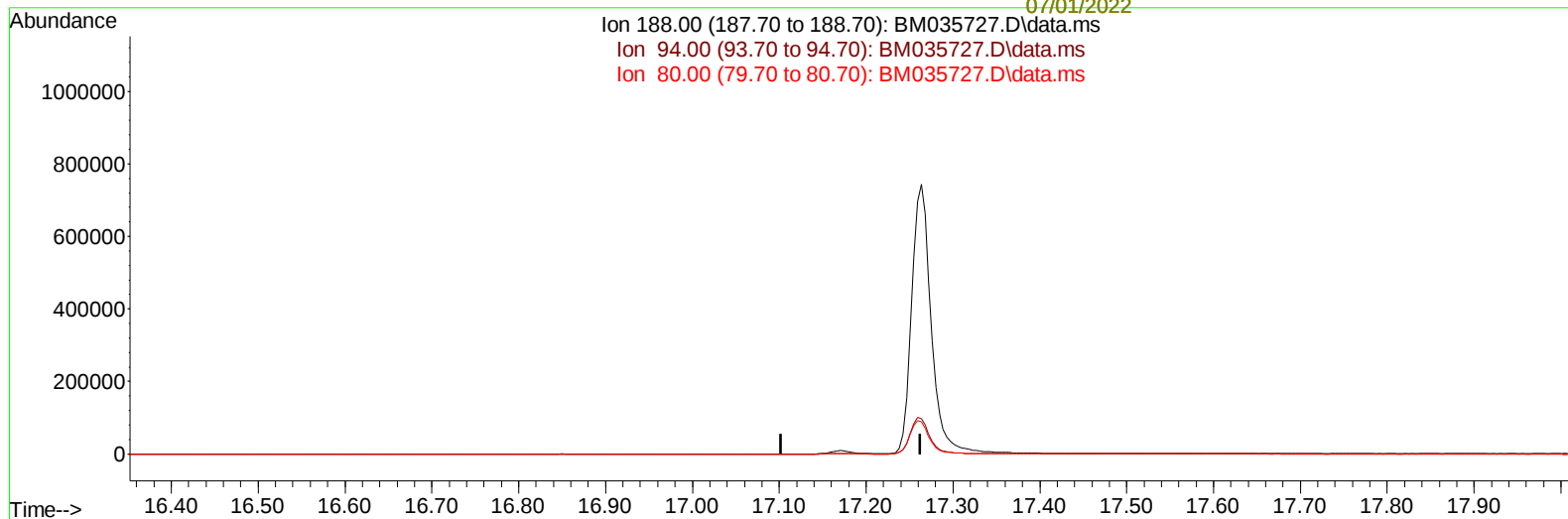
Instrument :
BNA_M
ClientSampleId :
C0AE2

Manual IntegrationsAPPROVED

Quant Time: Jun 29 04:27:17 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: BM035727.D\data.ms

(13) Phenanthrene-d10 (I)

17.182min (-17.182) 0.00 ng/u1

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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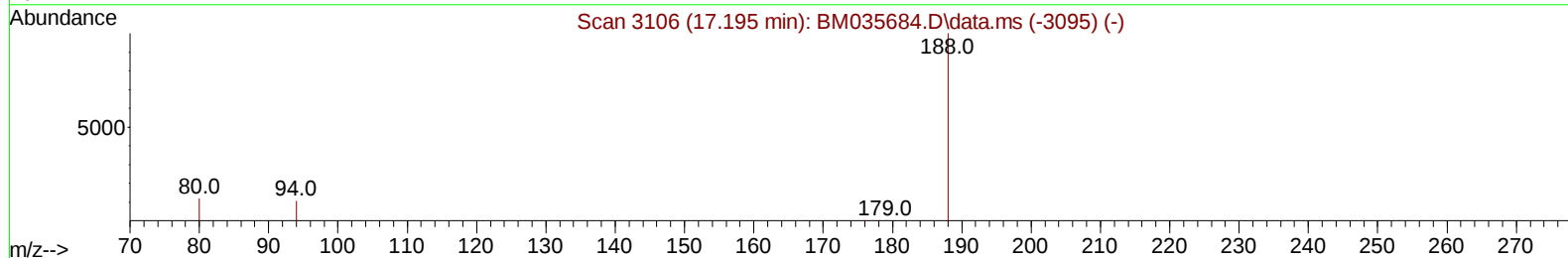
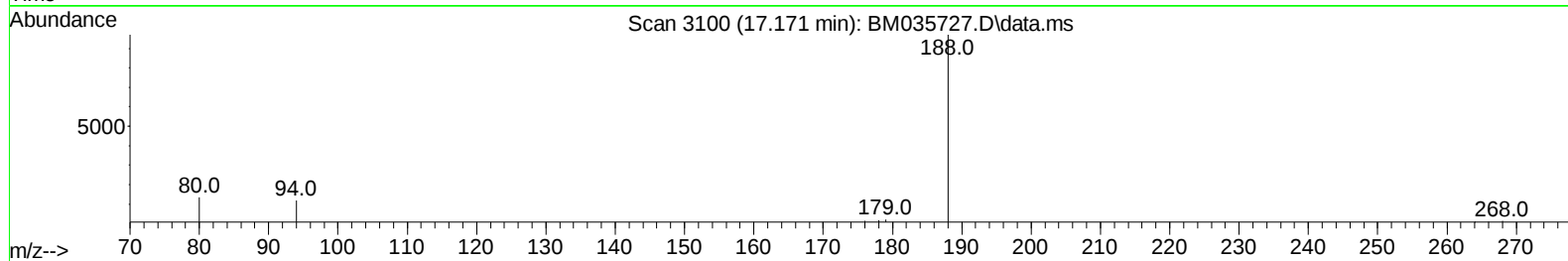
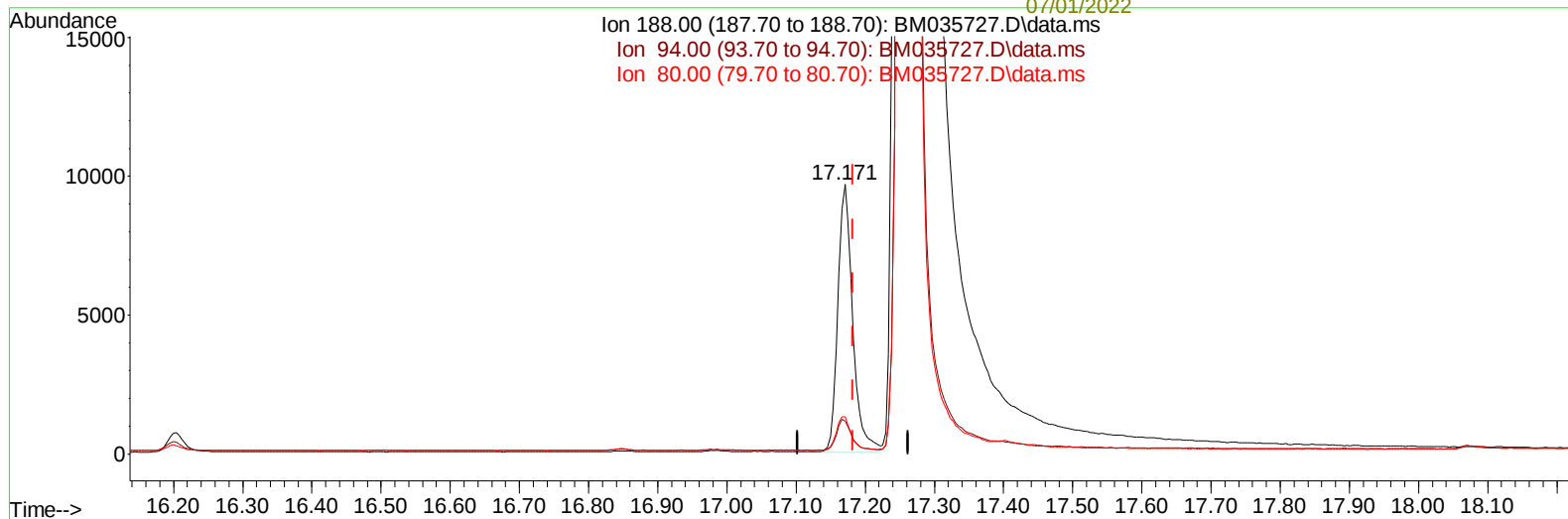
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(13) Phenanthrene-d10 (I)

17.171min (-0.012) 0.40 ng/ul m

response 14359

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	11.40	12.22
80.00	12.70	13.89
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.795	152		4167	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.578	136		13669	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.423	164		7355	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.171	188		14359m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.379	240		12583	0.400	ng/ul	# 0.00
23) Perylene-d12	23.697	264		15951	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.213	96		25640	4.534	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152		6479	0.303	ng/ul	0.01
18) Fluoranthene-d10	19.206	212		13431	0.317	ng/ul	0.00
Target Compounds							Qvalue
21) Benzo(a)anthracene	21.373	228		1419	0.033	ng/ul#	37
22) Chrysene	21.420	228		2217	0.040	ng/ul#	53
24) Benzo(b)fluoranthene	22.992	252		2774	0.041	ng/ul#	1
25) Benzo(k)fluoranthene	23.033	252		2289	0.033	ng/ul#	1

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

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