

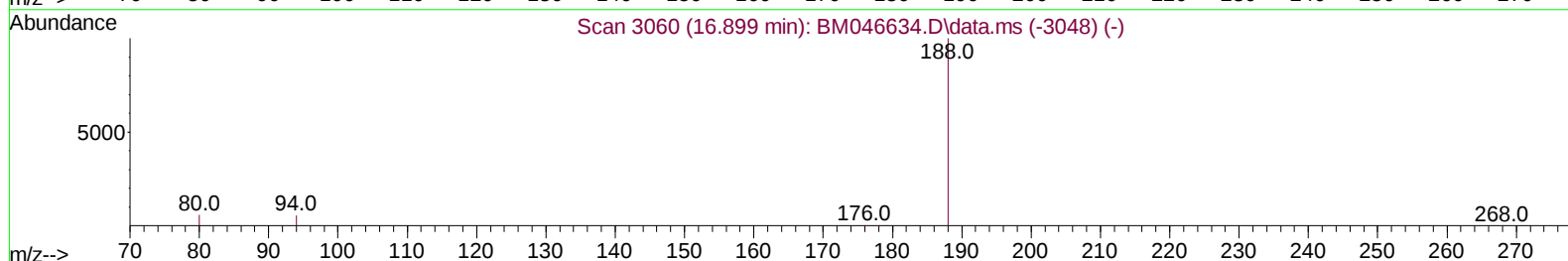
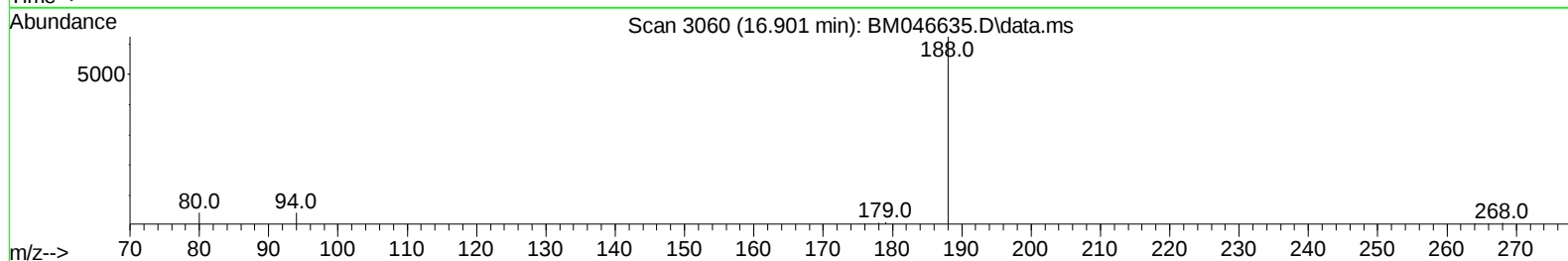
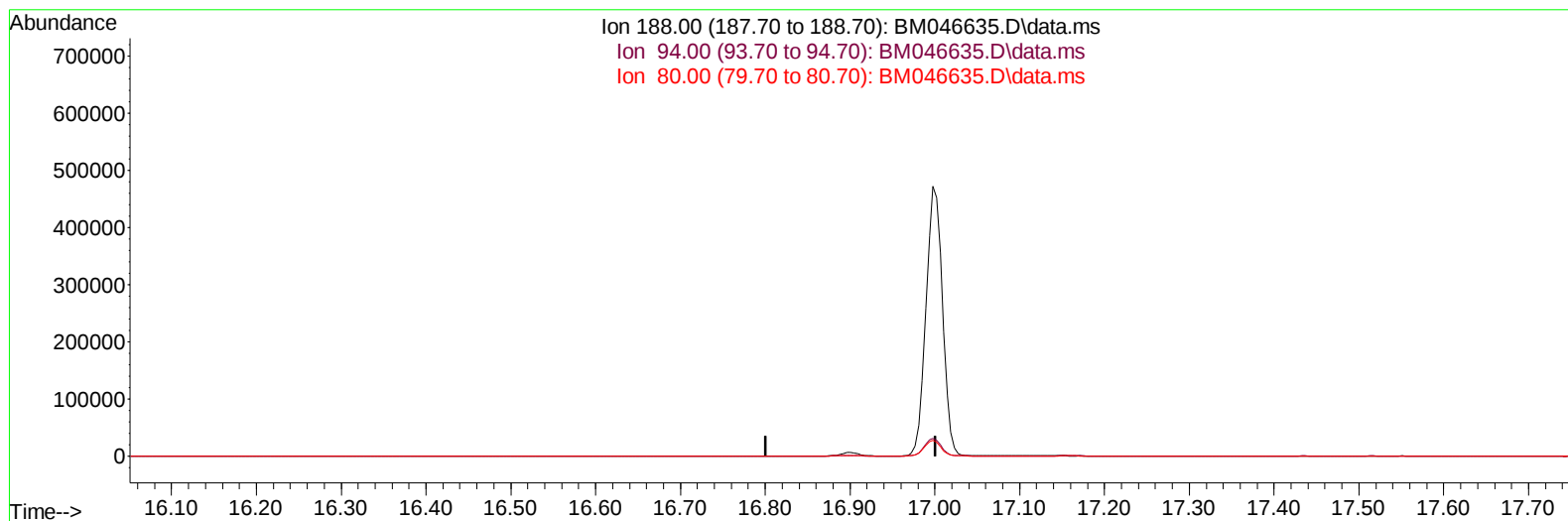
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071524\
Data File : BM046635.D
Acq On : 16 Jul 2024 16:20
Operator : MA/JU
Sample : PB162023BL
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK023

Manual IntegrationsAPPROVED

Quant Time: Jul 16 17:29:01 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 15 08:57:27 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/18/2024
Supervised By :mohammad ahmed 07/18/2024



TIC: BM046635.D\data.ms

(13) Phenanthrene-d10 (I)

16.901min (-16.901) 0.00 ng/u1

response 0

Ion	Exp%	Act%
188.00	100.00	0.00
94.00	6.40	0.00#
80.00	6.60	0.00#
0.00	0.00	0.00

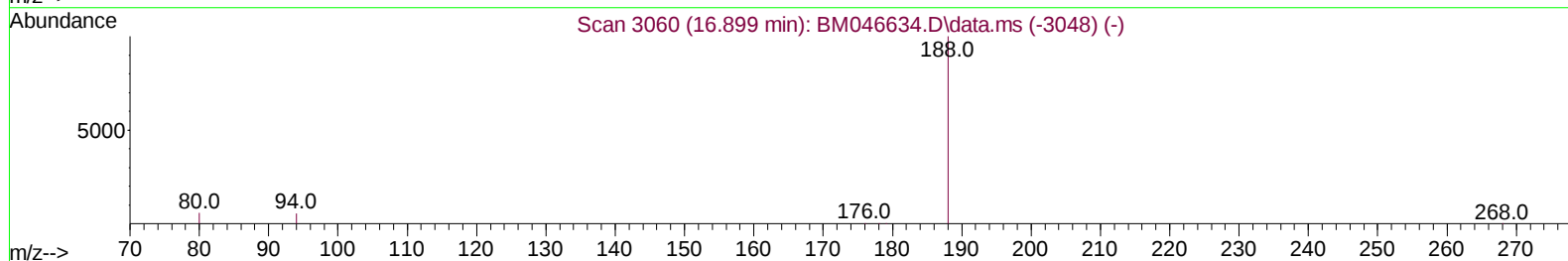
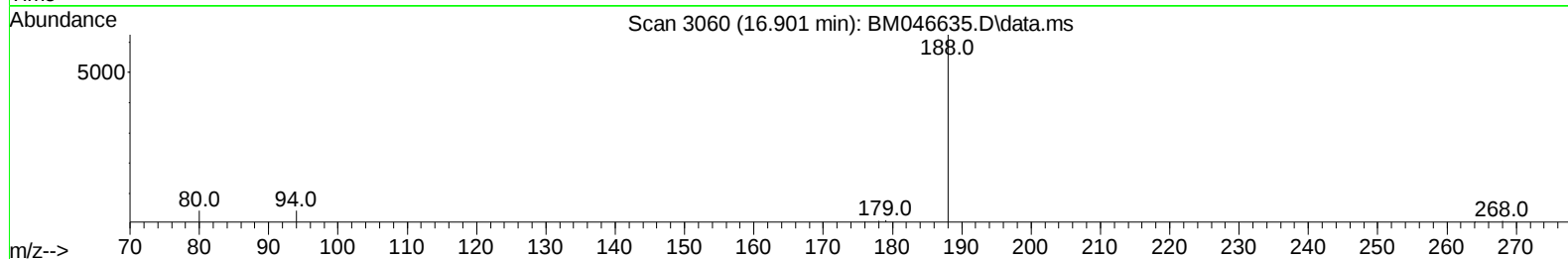
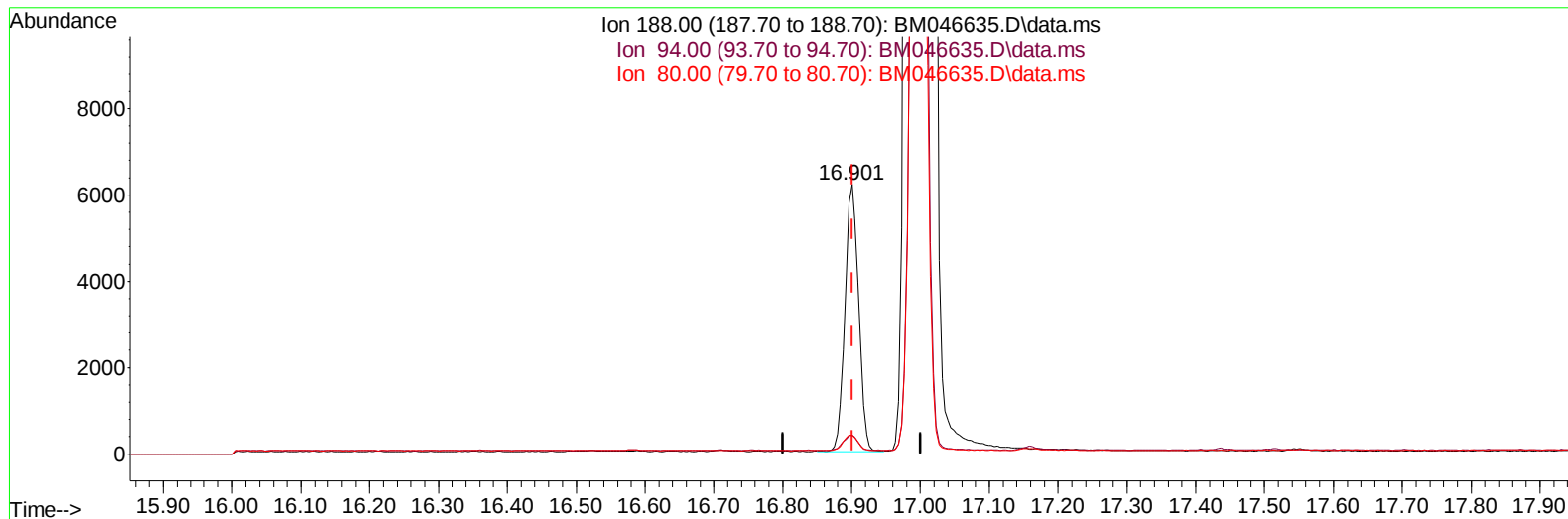
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(13) Phenanthrene-d10 (I)

16.901min (-0.000) 0.40 ng/u1 m

response 8466

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	6.40	6.89
80.00	6.60	6.87
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	7.501	152		1883	0.400	ng/ul	0.00
4) Naphthalene-d8	10.260	136		5478	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.146	164		3788	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.901	188		8466m	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240		6539	0.400	ng/ul	# 0.00
23) Perylene-d12	23.250	264		6775	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.113	96		7920	5.119	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.860	152		3086	0.349	ng/ul	0.00
18) Fluoranthene-d10	18.942	212		7555	0.380	ng/ul	0.00

Target Compounds Qvalue

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

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