

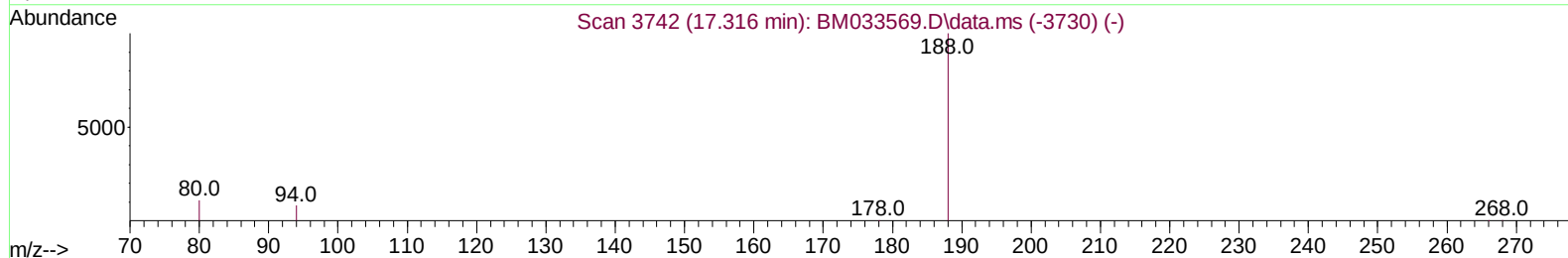
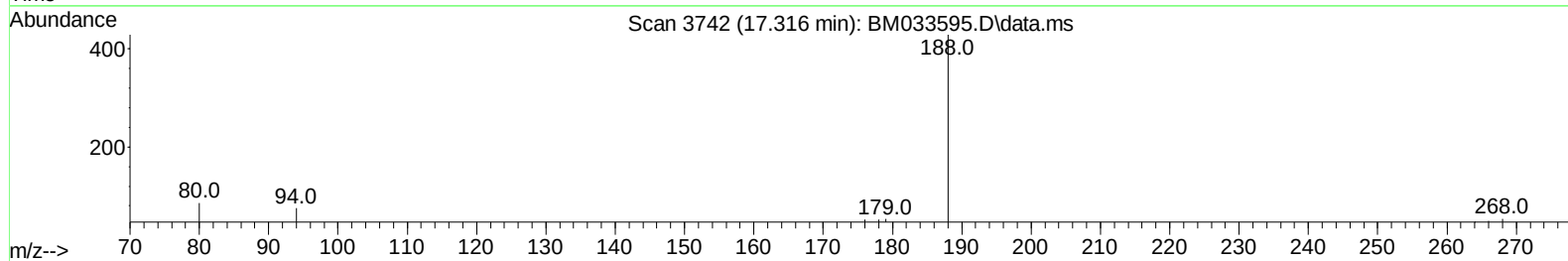
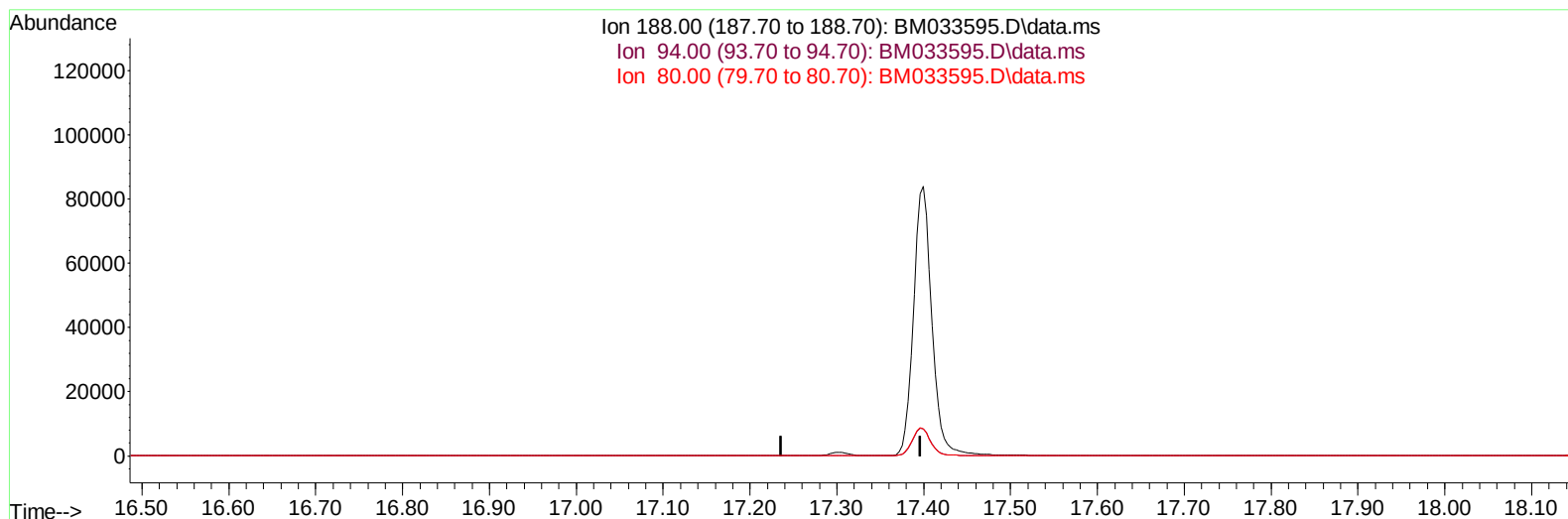
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122221\
 Data File : BM033595.D
 Acq On : 23 Dec 2021 03:39
 Operator : CG/JU
 Sample : PB141491BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK491

Manual IntegrationsAPPROVED

Quant Time: Dec 23 07:46:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM122221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 12:43:31 2021
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2021
 Supervised By :mohammad ahmed 12/29/2021



TIC: BM033595.D\data.ms

(13) Phenanthrene-d10 (I)

17.316min (-17.316) 0.00 ng/u1

response 0

Ion	Exp%	Act%
188.00	100.00	0.00
94.00	13.30	0.00#
80.00	15.50	0.00#
0.00	0.00	0.00

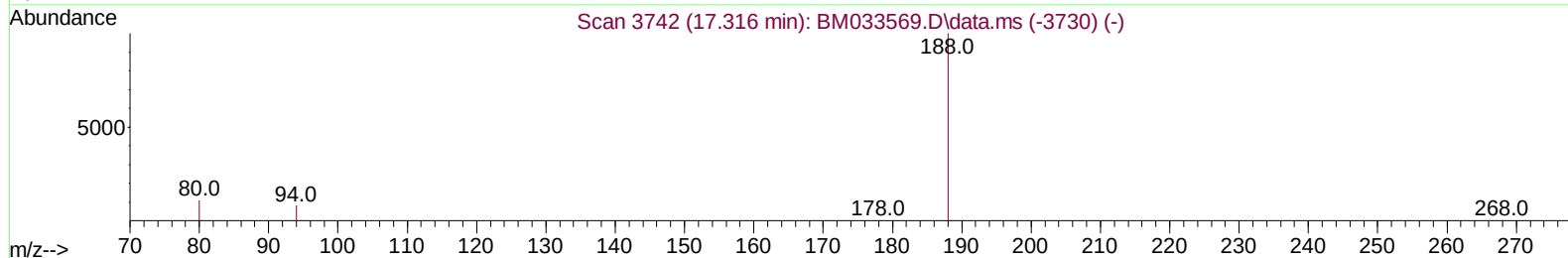
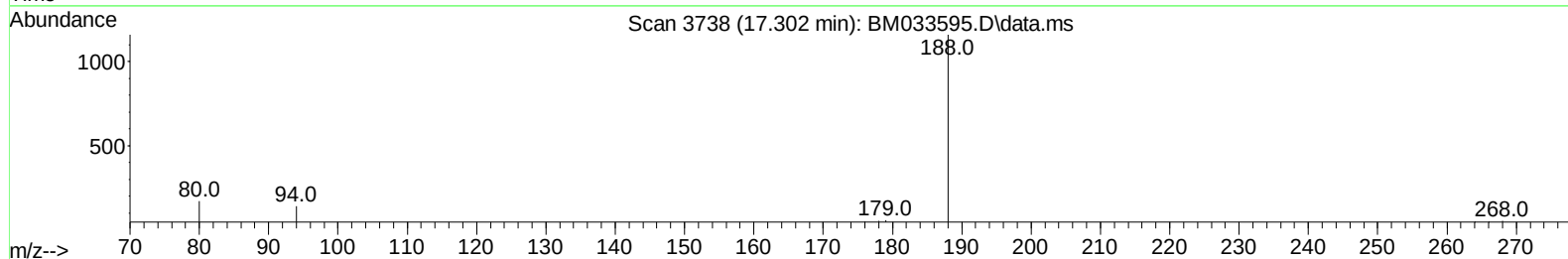
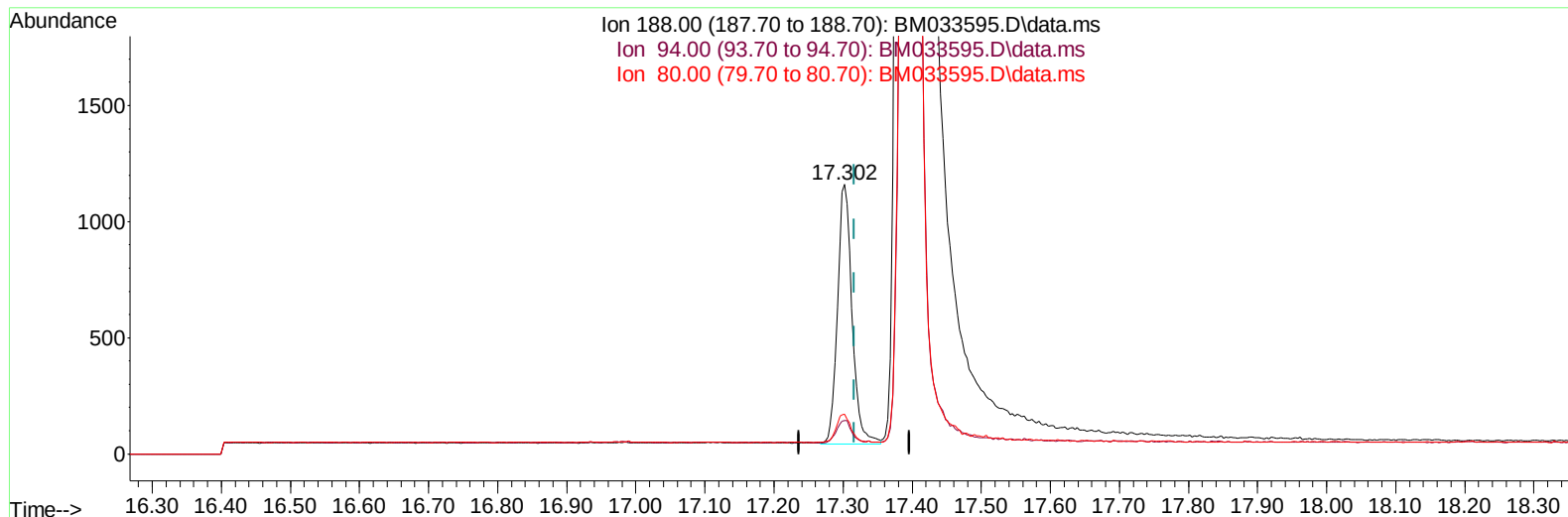
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(13) Phenanthrene-d10 (I)

17.302min (-0.014) 0.40 ng/ul m

response 1631

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	13.30	12.33
80.00	15.50	14.83
0.00	0.00	0.00

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 Acq On : 23 Dec 2021 03:39
 Operator : CG/JU
 Sample : PB141491BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.912	152		439	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.716	136		1130	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.557	164		720	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.302	188		1631m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.502	240		1734	0.400	ng/ul	0.01
23) Perylene-d12	23.875	264		2317	0.400	ng/ul	#-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.327	96		2111	5.633	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152		448	0.296	ng/ul	0.02
18) Fluoranthene-d10	19.330	212		1440	0.359	ng/ul	0.00

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

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