

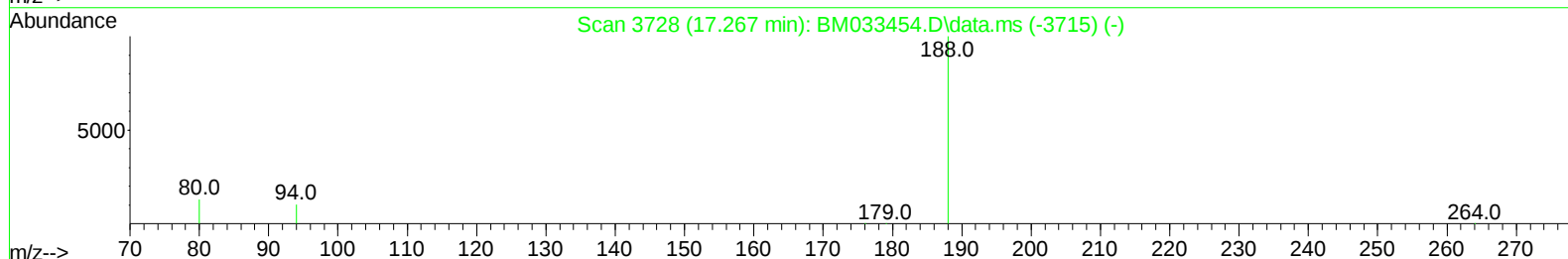
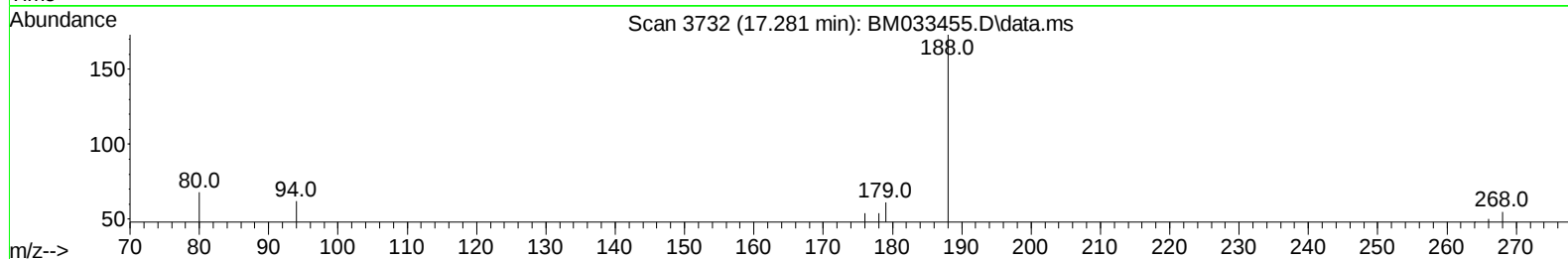
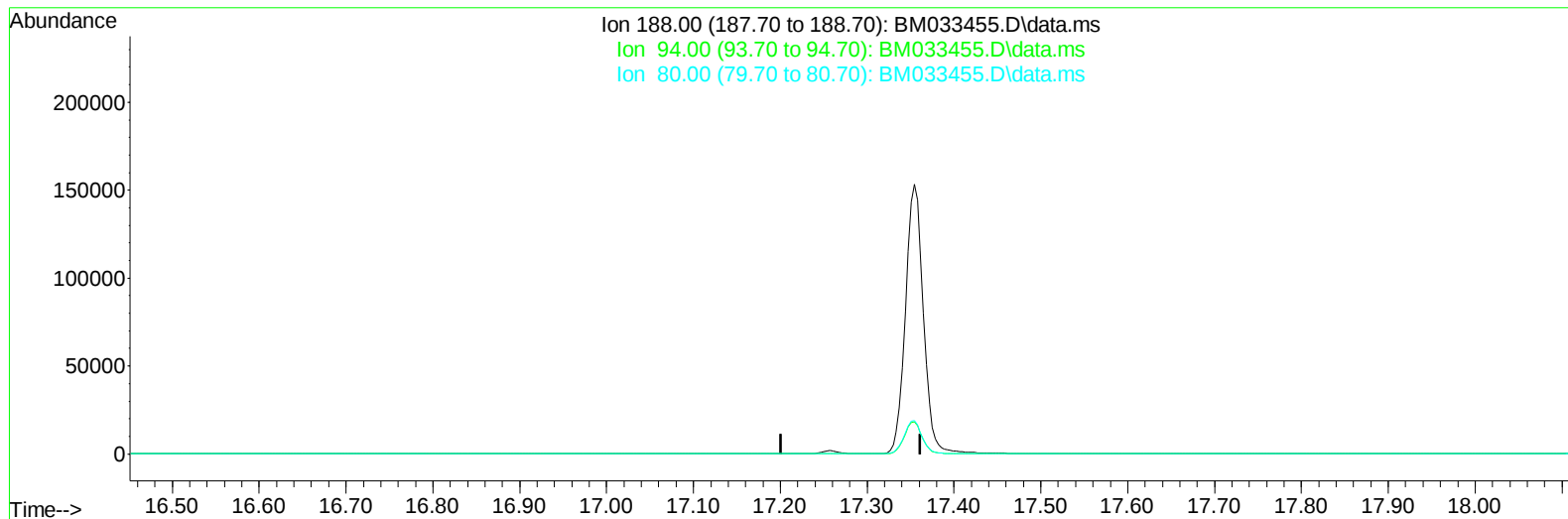
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121421\
Data File : BM033455.D
Acq On : 15 Dec 2021 00:49
Operator : CG/JU
Sample : PB141279BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK279

Manual IntegrationsAPPROVED

Quant Time: Dec 15 03:51:24 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 15 01:00:55 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/15/2021
Supervised By :mohammad ahmed 12/24/2021



TIC: BM033455.D\data.ms

(13) Phenanthrene-d10 (I)

17.281min (-17.281) 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

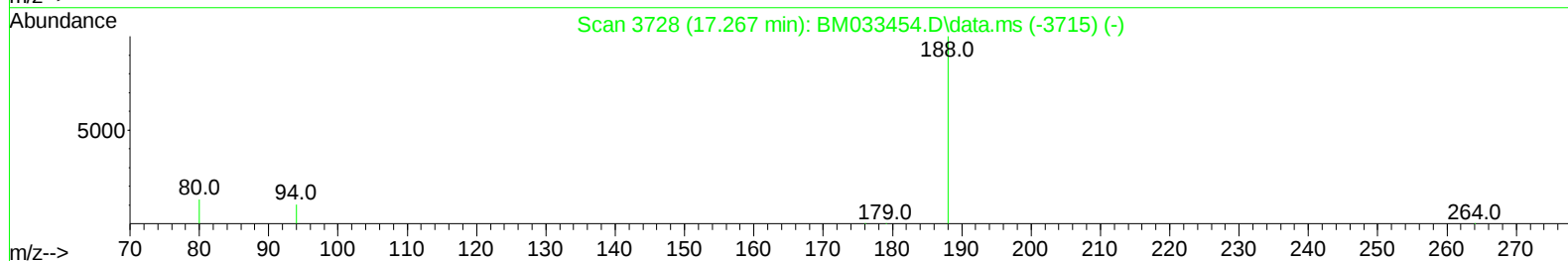
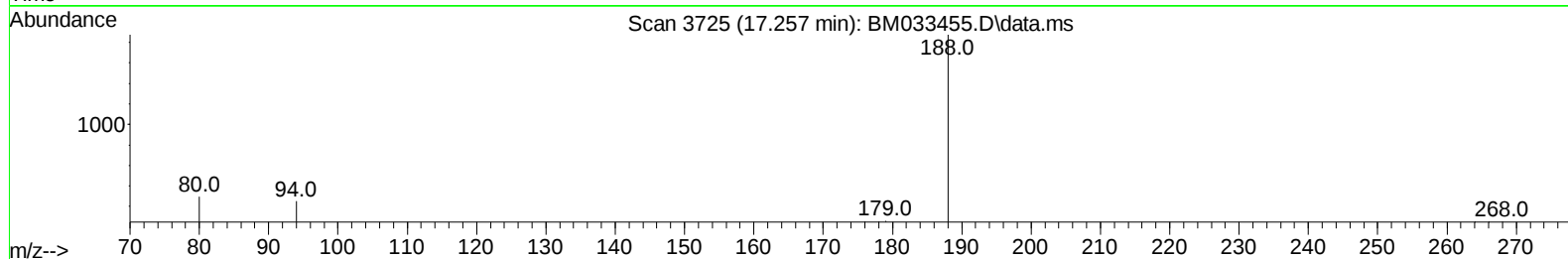
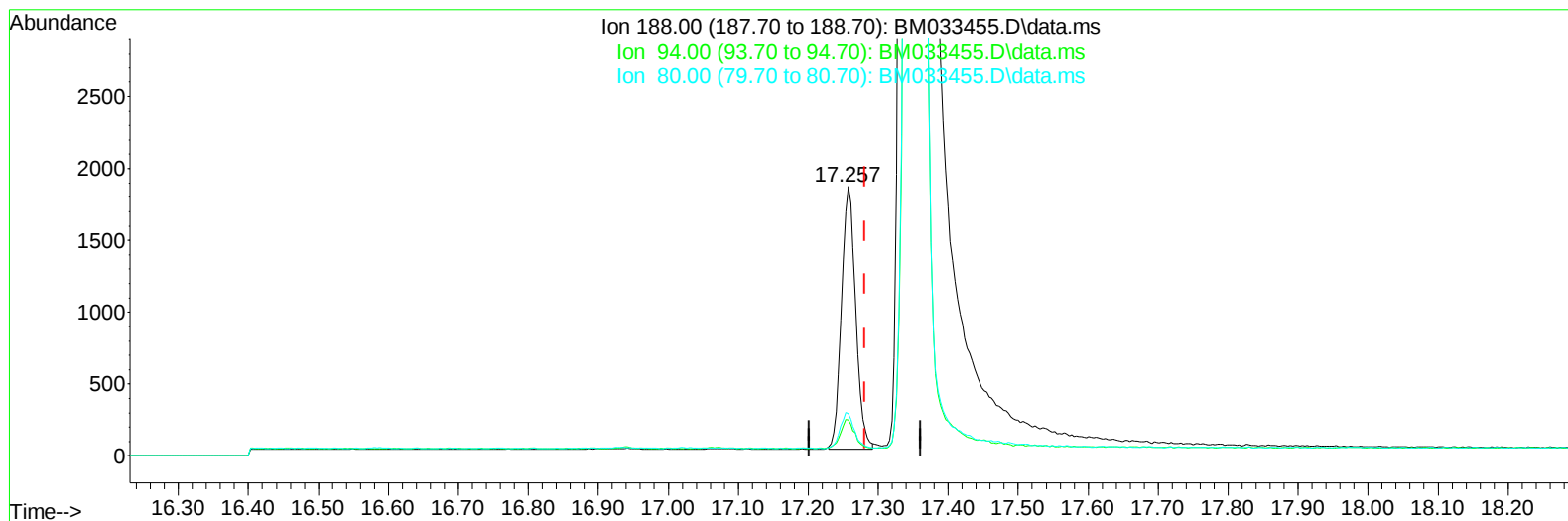
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(13) Phenanthrene-d10 (I)

17.257min (-0.024) 0.40 ng/ul m

response 2569

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	13.30	13.28
80.00	15.50	15.73
0.00	0.00	0.00

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 ALS Vial : 19 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.901	152		556	0.40	ng/ul	0.00
4) Naphthalene-d8	10.684	136		1480	0.40	ng/ul	-0.02
9) Acenaphthene-d10	14.519	164		1074	0.40	ng/ul	-0.01
13) Phenanthrene-d10	17.257	188		2569m	0.40	ng/ul	-0.02
17) Chrysene-d12	21.436	240		3285	0.40	ng/ul	0.00
23) Perylene-d12	23.735	264		3965	0.40	ng/ul	#-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.358	96		2523	5.60	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.294	152		766	0.41	ng/ul	0.00
18) Fluoranthene-d10	19.281	212		3116	0.44	ng/ul	-0.01

Target Compounds Qvalue

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

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