

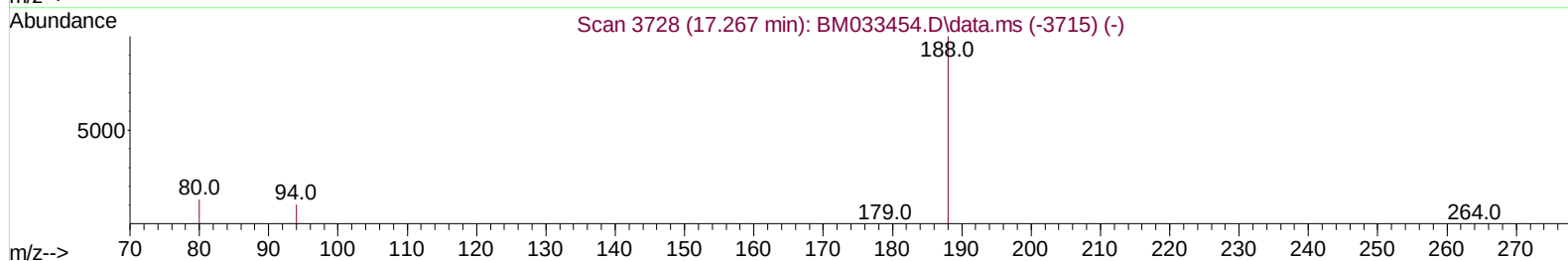
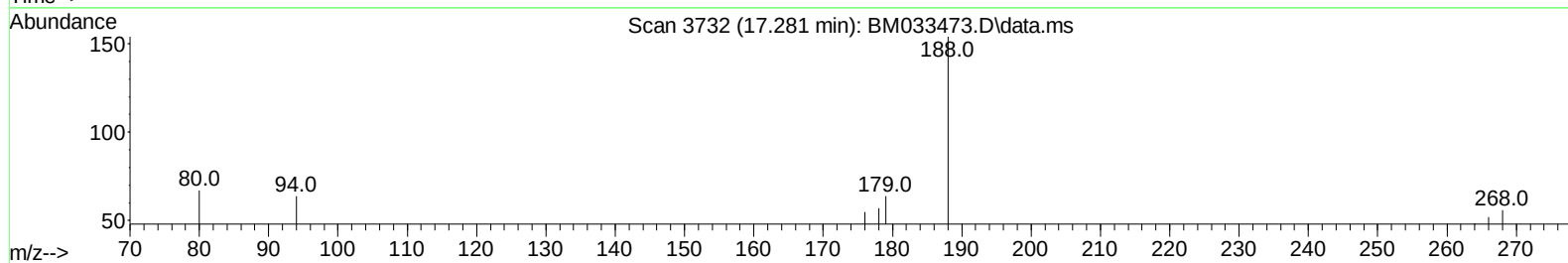
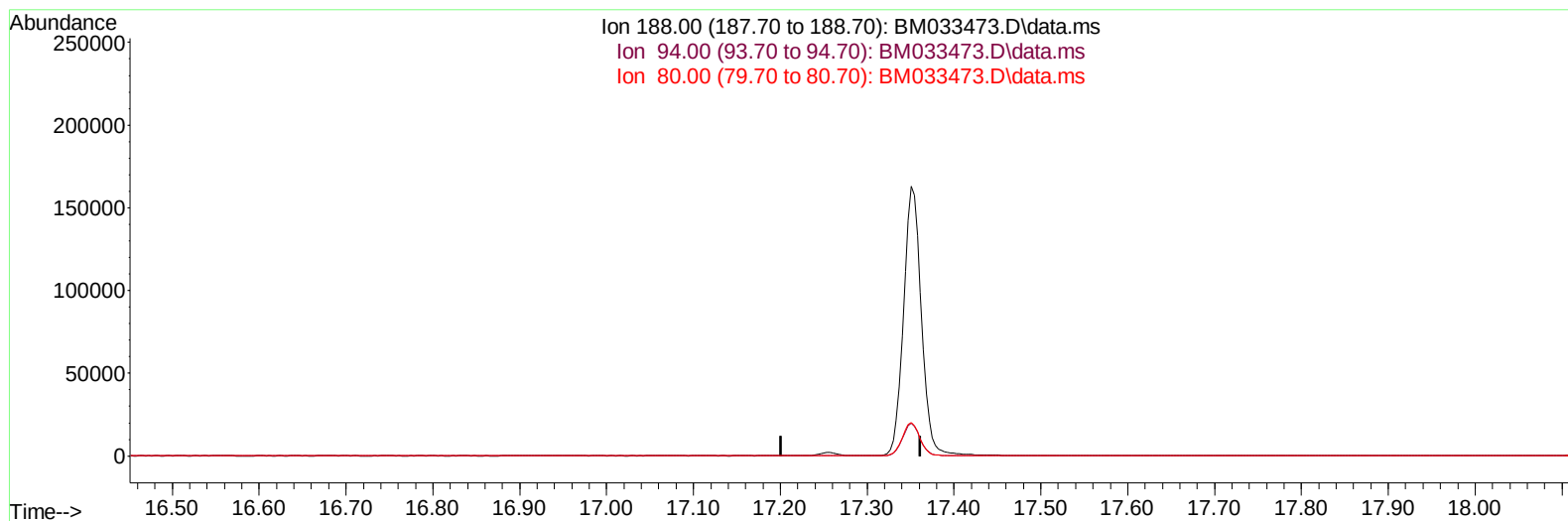
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121421\
 Data File : BM033473.D
 Acq On : 15 Dec 2021 11:33
 Operator : CG/JU
 Sample : PB141294BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK294

Manual IntegrationsAPPROVED

Quant Time: Dec 16 02:33:22 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 12/16/2021
 Supervised By :mohammad ahmed 12/24/2021



TIC: BM033473.D\data.ms

(13) Phenanthrene-d10 (I)

17.281min (-17.281) 0.00 ng/ul

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

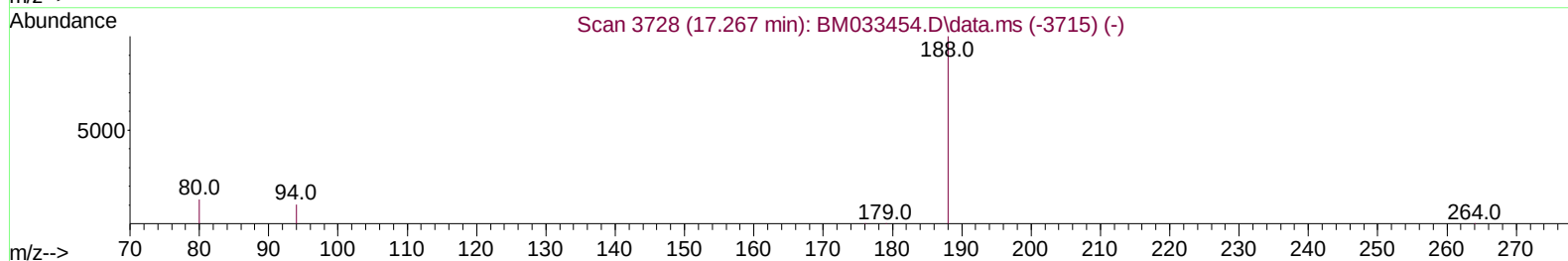
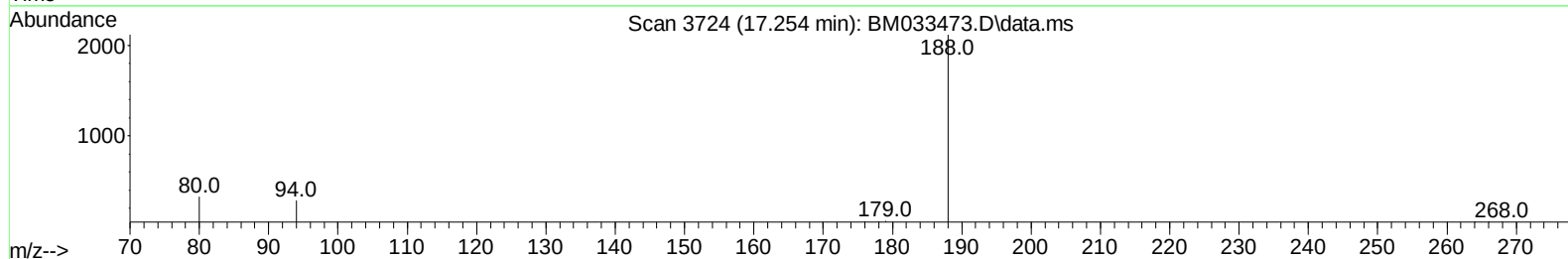
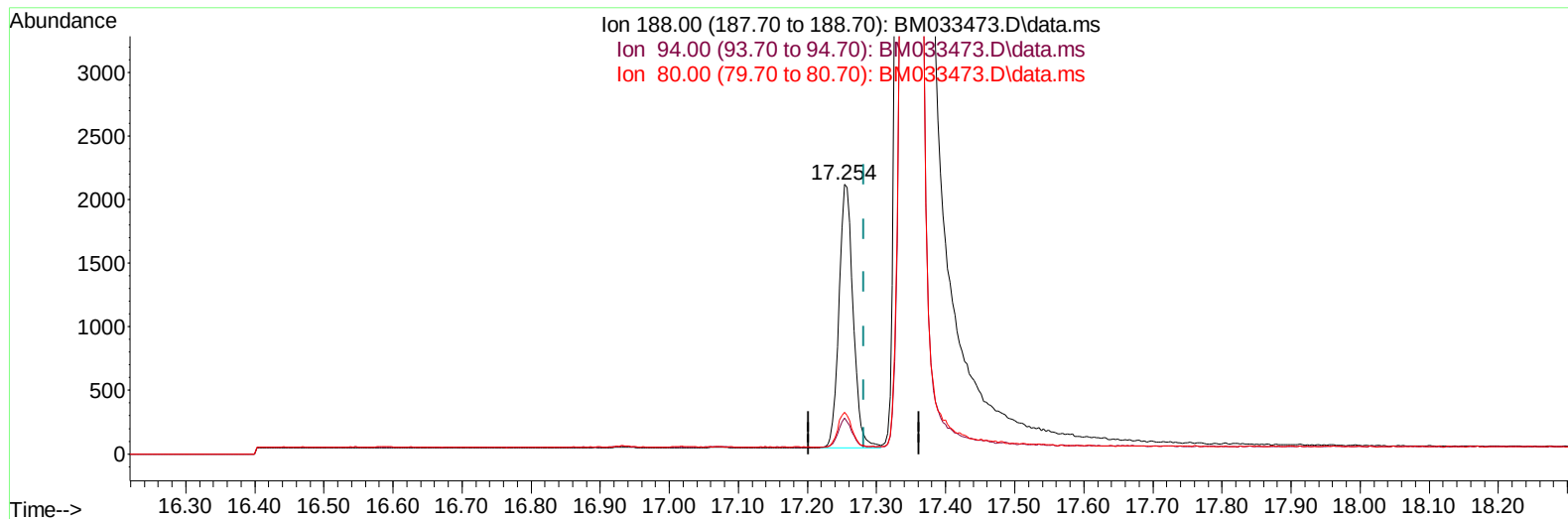
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(13) Phenanthrene-d10 (I)

17.254min (-0.028) 0.40 ng/ul m

response 2938

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	13.30	13.41
80.00	15.50	15.39
0.00	0.00	0.00

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 ALS Vial : 37 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	7.898	152		663	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.684	136		1739	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.515	164		1222	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.254	188		2938m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.434	240		3596	0.400	ng/ul	0.00
23) Perylene-d12	23.730	264		4019	0.400	ng/ul	#-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.358	96		2897	5.393	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.294	152		833	0.379	ng/ul	0.00
18) Fluoranthene-d10	19.277	212		3211	0.410	ng/ul	-0.01

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

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