

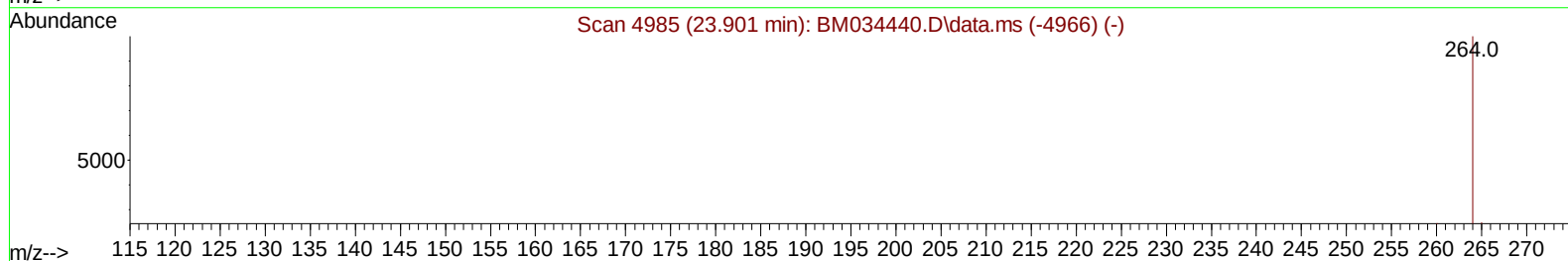
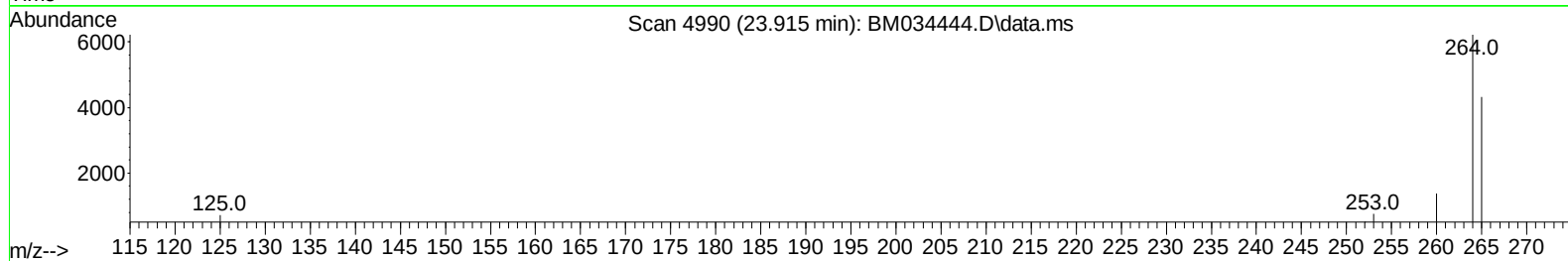
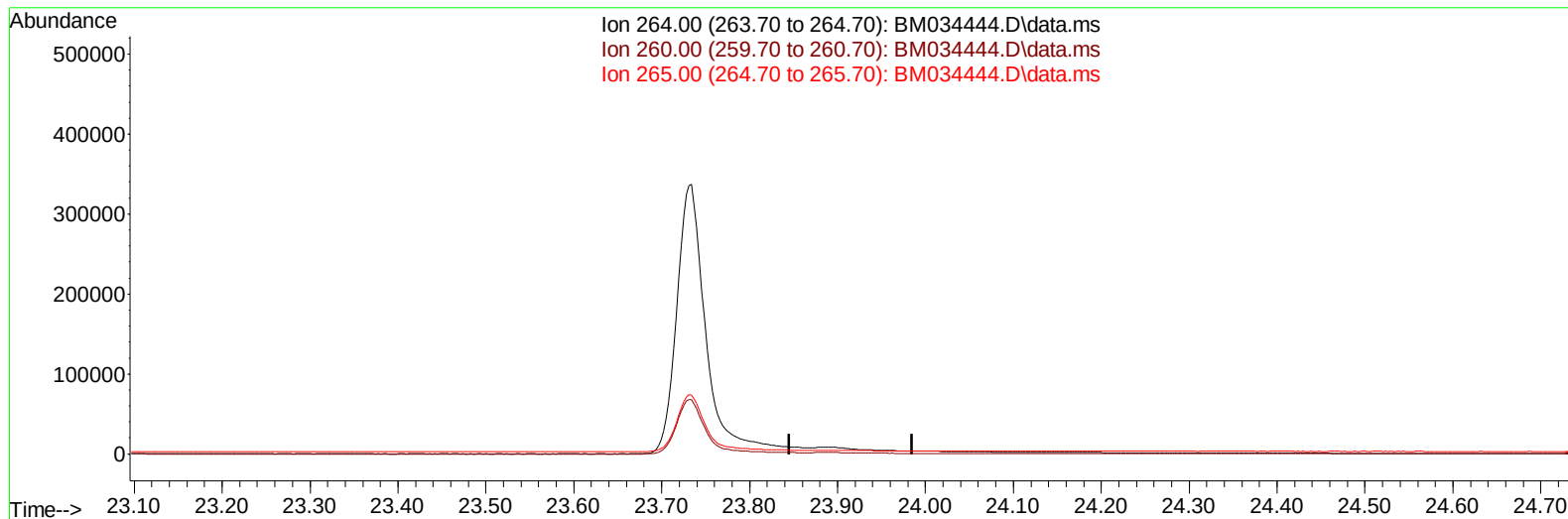
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032922\
Data File : BM034444.D
Acq On : 30 Mar 2022 17:10
Operator : CG/JU
Sample : N2097-02
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AD1

Manual IntegrationsAPPROVED

Quant Time: Mar 31 02:21:34 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM032822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 28 18:27:40 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/31/2022
Supervised By :Yogesh Patel 04/02/2022



TIC: BM034444.D\data.ms

(23) Perylene-d12 (I)

23.915min (-23.915) 0.00 ng/ul

response 0

Ion	Exp%	Act%
264.00	100.00	0.00
260.00	27.80	0.00#
265.00	95.30	0.00#
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	7.929	152		2694	0.400	ng/ul	0.00
4) Naphthalene-d8	10.735	136		7494	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.569	164		4319	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.308	188		8320	0.400	ng/ul	#-0.02
17) Chrysene-d12	21.501	240		5731	0.400	ng/ul	# 0.00
23) Perylene-d12	23.889	264		20260m	0.400	ng/ul	-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.301	96		19000	4.741	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152		3997	0.396	ng/ul	0.00
18) Fluoranthene-d10	19.334	212		9638	0.539	ng/ul	-0.01
Target Compounds							Qvalue
2) 1,4-Dioxane	3.339	88		1826	0.413	ng/ul#	73
15) Phenanthrene	17.350	178		1726	0.065	ng/ul#	15

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

```
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