

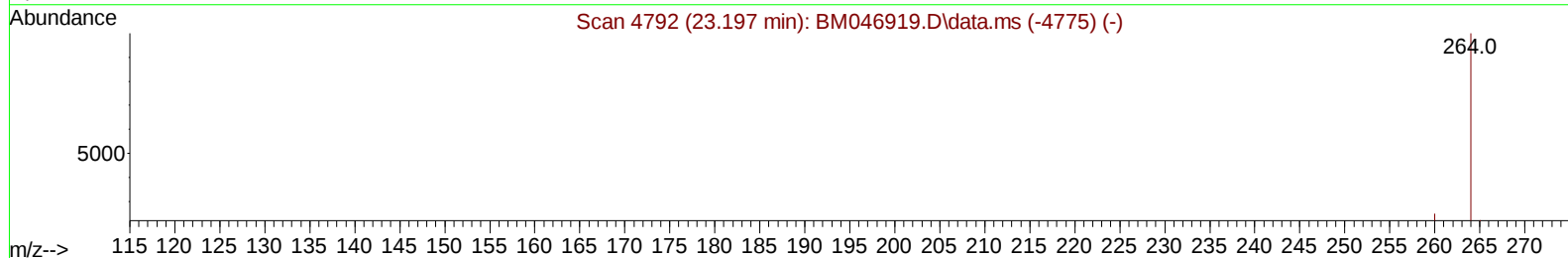
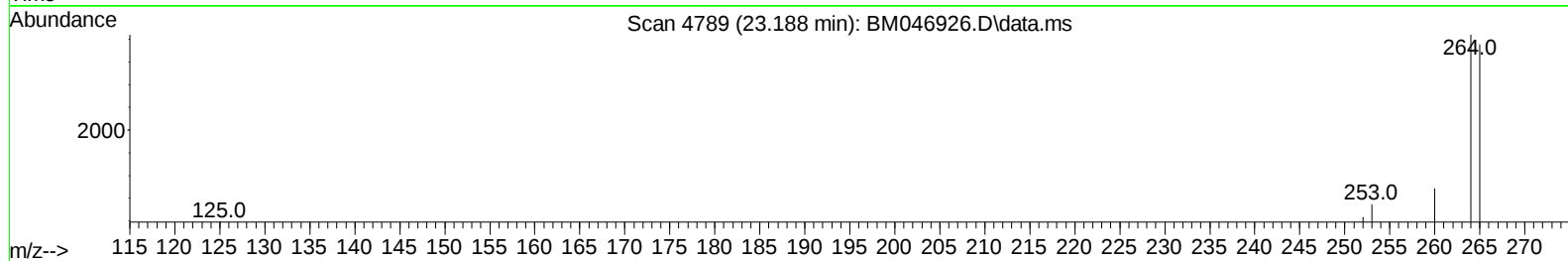
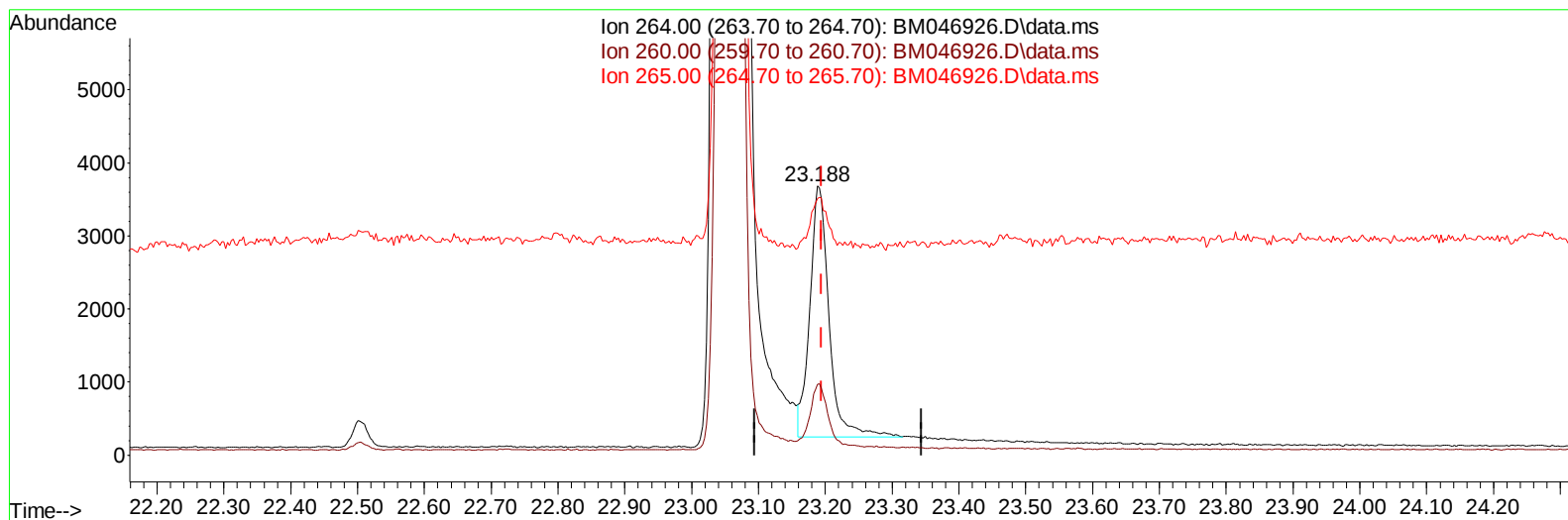
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072924\
Data File : BM046926.D
Acq On : 30 Jul 2024 19:24
Operator : MA/JU
Sample : P3336-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F7E45

Manual IntegrationsAPPROVED

Quant Time: Jul 31 00:00:56 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM072424.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 30 04:18:24 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/31/2024
Supervised By :mohammad ahmed 08/03/2024



TIC: BM046926.D\data.ms

(23) Perylene-d12 (I)

23.188min (-0.006) 0.40 ng/u1

response 6832

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	27.10	26.30
265.00	89.90	95.46
0.00	0.00	0.00

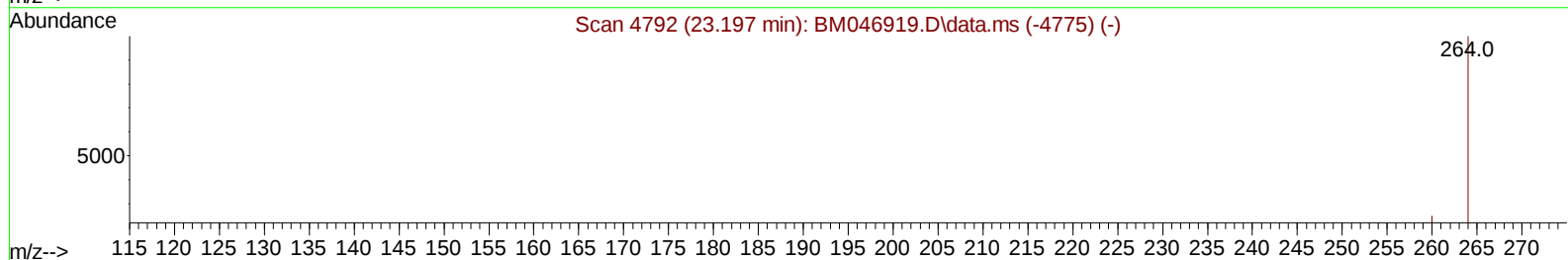
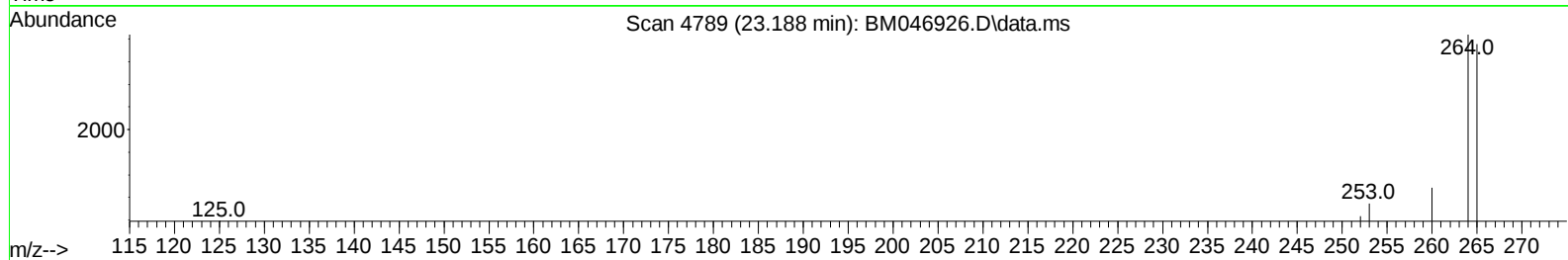
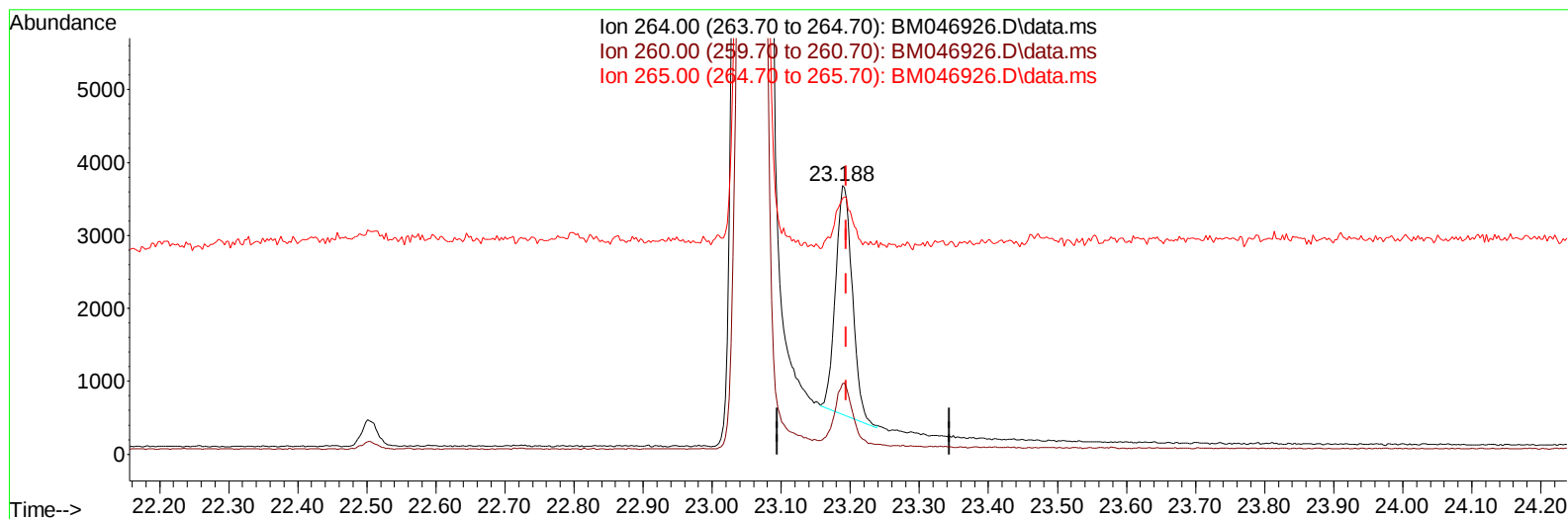
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.447	152	2201	0.400	ng/ul	0.00
4) Naphthalene-d8	10.205	136	6153	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.095	164	4059	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.854	188	8720	0.400	ng/ul	0.00
17) Chrysene-d12	21.067	240	5776	0.400	ng/ul	0.00
23) Perylene-d12	23.188	264	5332m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.084	96	6092	3.953	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.800	152	2495	0.255	ng/ul	-0.01
18) Fluoranthene-d10	18.895	212	5078	0.270	ng/ul	0.00

Target Compounds Qvalue

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

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