

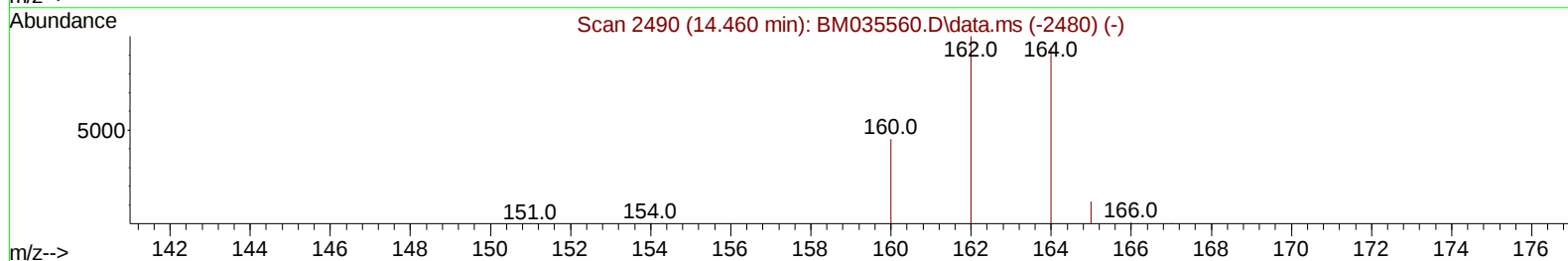
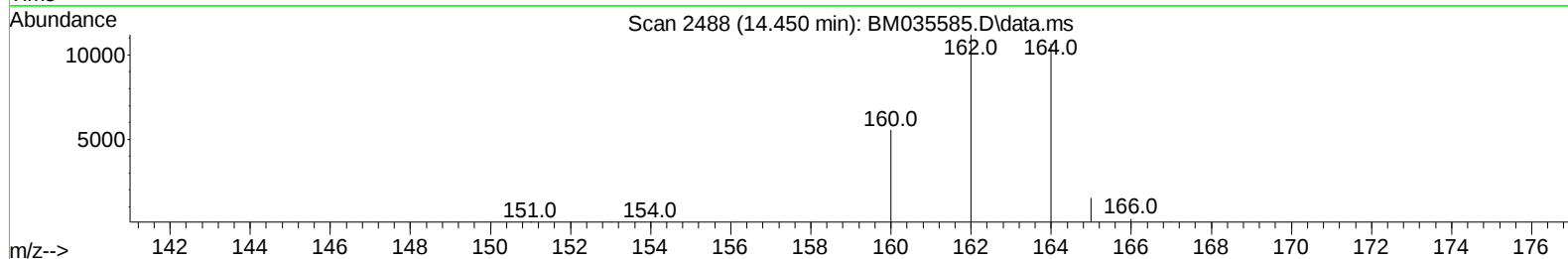
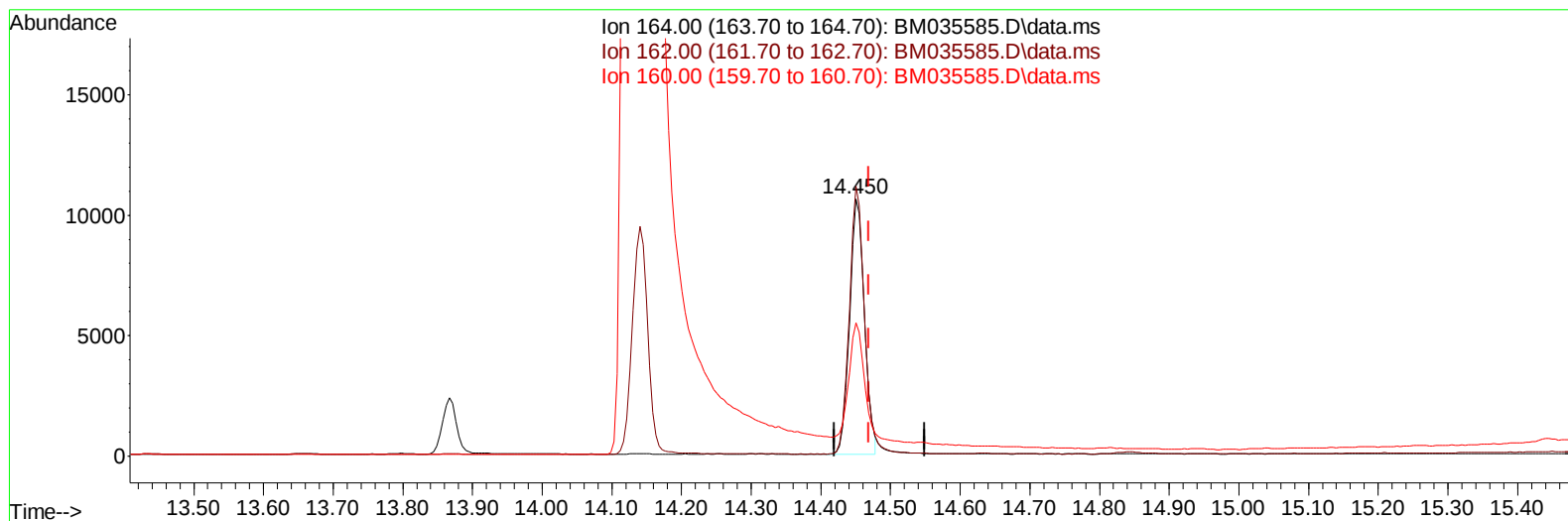
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
Data File : BM035585.D
Acq On : 21 Jun 2022 23:26
Operator : CG/JU
Sample : N3285-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4R1

Manual IntegrationsAPPROVED

Quant Time: Jun 22 01:33:36 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 20 01:56:13 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/22/2022
Supervised By :Yogesh Patel 06/25/2022



TIC: BM035585.D\data.ms

(9) Acenaphthene-d10 (I)

14.450min (-0.020) 0.40 ng/u1

response 15613

Ion	Exp%	Act%
164.00	100.00	100.00
162.00	106.40	104.76
160.00	49.90	51.86
0.00	0.00	0.00

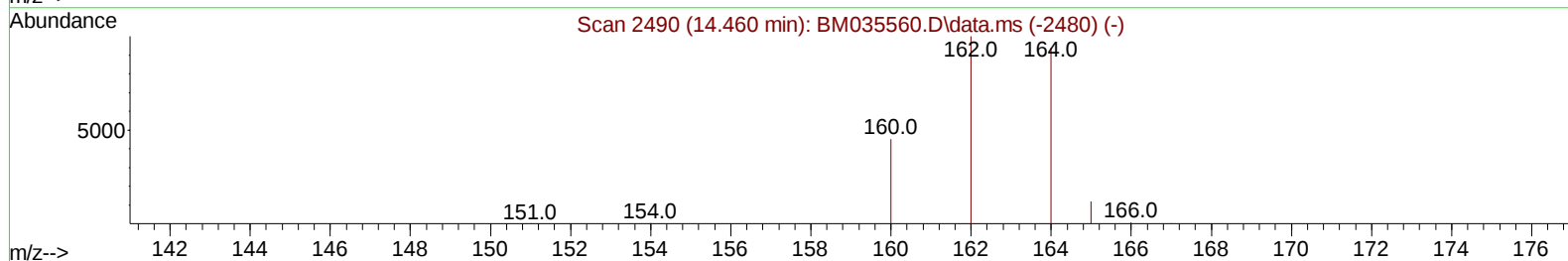
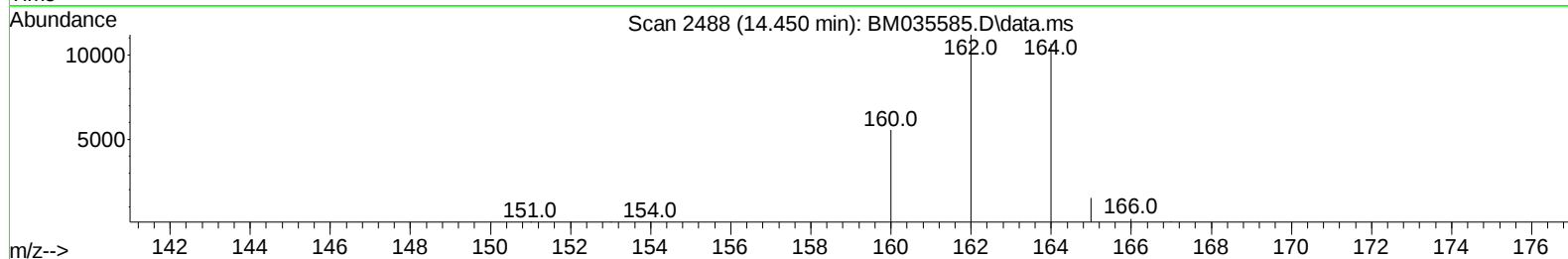
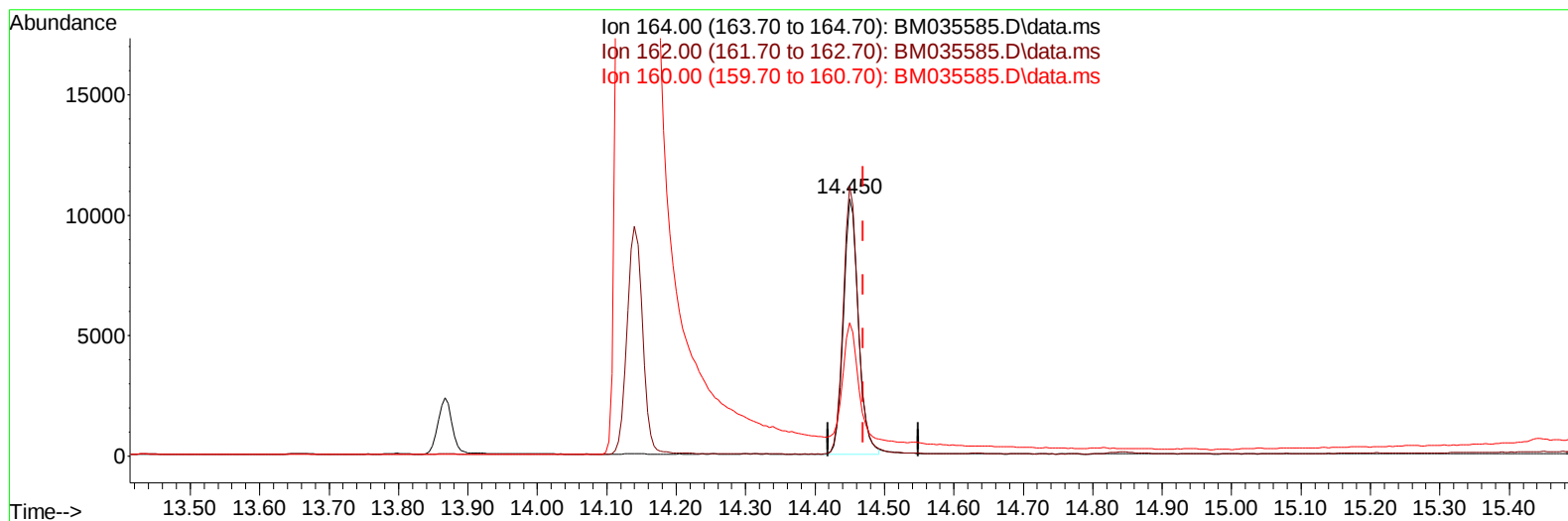
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(9) Acenaphthene-d10 (I)

14.450min (-0.020) 0.40 ng/u1 m

response 15941

Ion	Exp%	Act%
164.00	100.00	100.00
162.00	106.40	104.76
160.00	49.90	51.86
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	7.816	152		6441	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.605	136		24040	0.400	ng/ul	#-0.04
9) Acenaphthene-d10	14.450	164		15941m	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.199	188		33384	0.400	ng/ul	-0.03
17) Chrysene-d12	21.395	240		20833	0.400	ng/ul	-0.02
23) Perylene-d12	23.727	264		22891	0.400	ng/ul	#-0.04
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96		30563	3.372	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.217	152		10004	0.280	ng/ul	-0.03
18) Fluoranthene-d10	19.233	212		24728	0.353	ng/ul	-0.02

Target Compounds Qvalue

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

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