

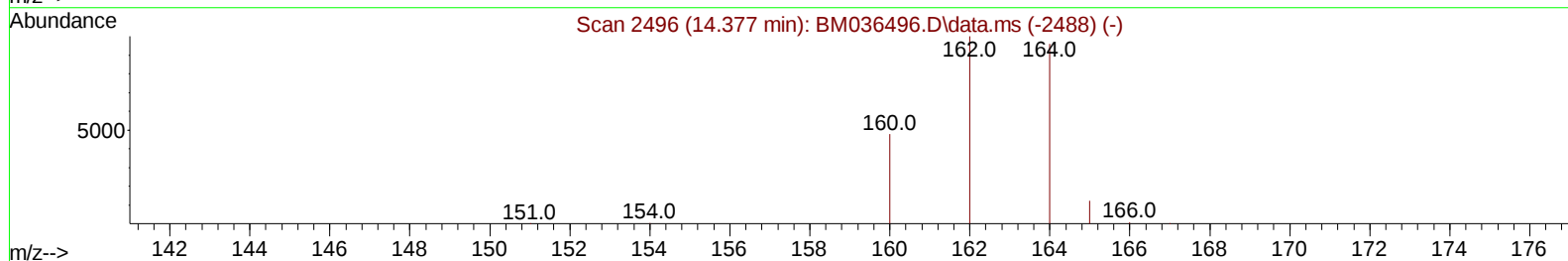
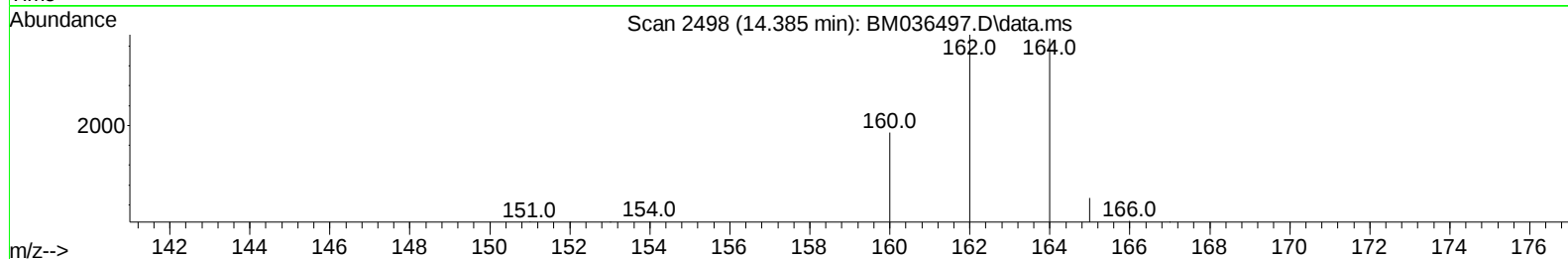
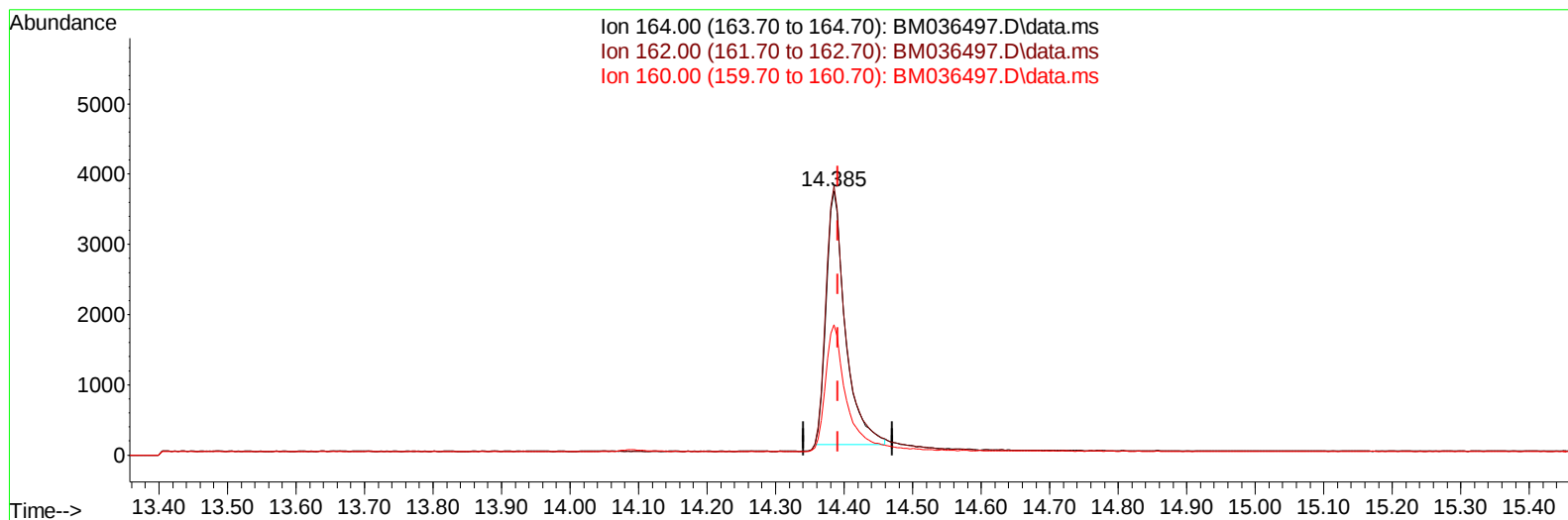
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090622\
Data File : BM036497.D
Acq On : 06 Sep 2022 11:30
Operator : CG/JU
Sample : PB146657BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK657

Manual IntegrationsAPPROVED

Quant Time: Sep 06 12:56:00 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM083022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 29 17:45:20 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/07/2022
Supervised By :mohammad ahmed 09/07/2022



TIC: BM036497.D\data.ms

(9) Acenaphthene-d10 (I)

14.385min (-0.006) 0.40 ng/u1

response 7019

Ion	Exp%	Act%
164.00	100.00	100.00
162.00	104.10	101.92
160.00	48.90	49.40
0.00	0.00	0.00

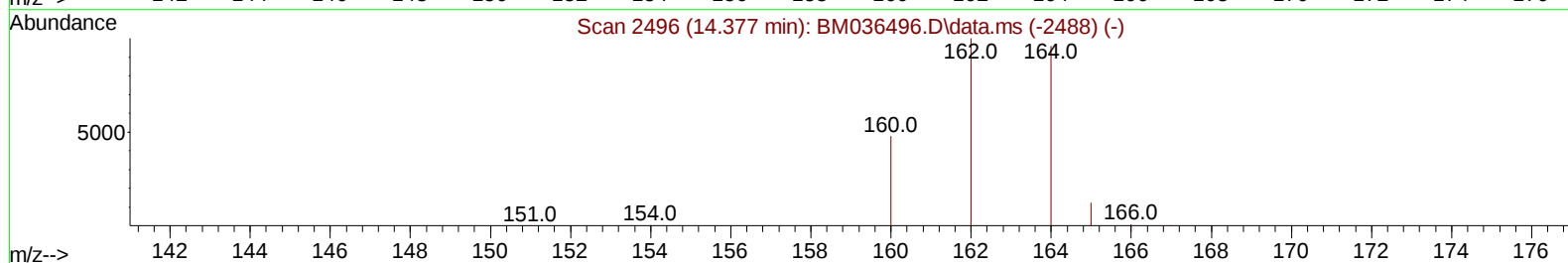
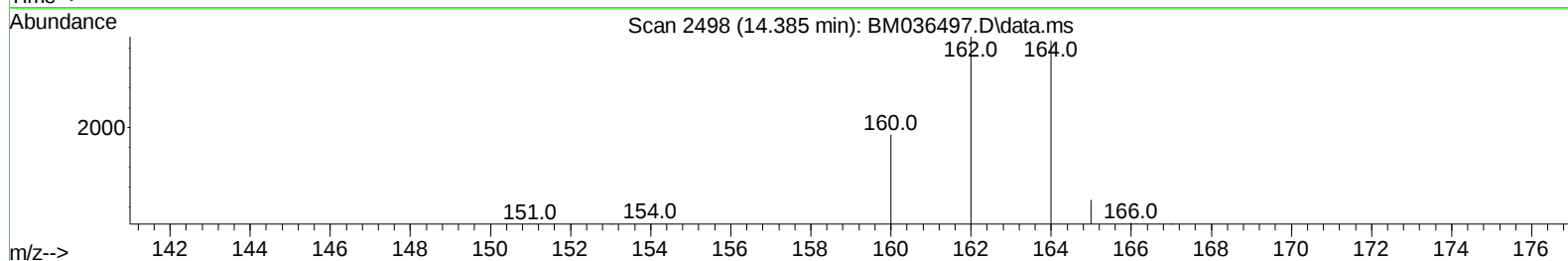
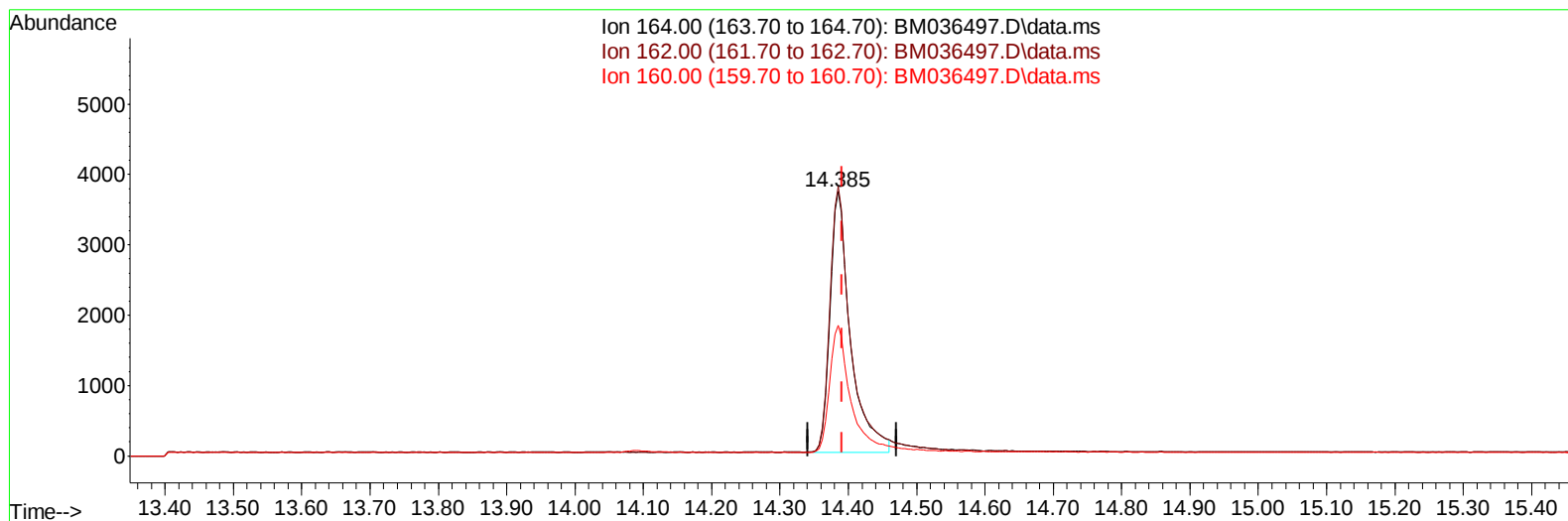
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TIC: BM036497.D\data.ms

(9) Acenaphthene-d10 (I)

14.385min (-0.006) 0.40 ng/u1 m

response 7660

Ion	Exp%	Act%
164.00	100.00	100.00
162.00	104.10	101.92
160.00	48.90	49.40
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	7.737	152		3684	0.400	ng/ul	0.00
4) Naphthalene-d8	10.545	136		13715	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.385	164		7660m	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.140	188		14561	0.400	ng/ul	0.00
17) Chrysene-d12	21.324	240		11589	0.400	ng/ul	0.00
23) Perylene-d12	23.616	264		9015	0.400	ng/ul	# 0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.134	96		4021	0.751	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.156	152		6897	0.322	ng/ul	0.02
18) Fluoranthene-d10	19.163	212		14923	0.414	ng/ul	0.00

Target Compounds Qvalue

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

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