

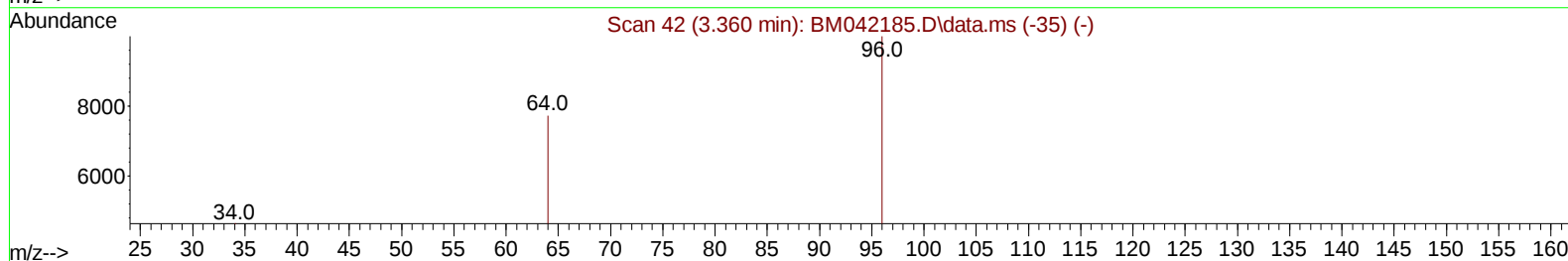
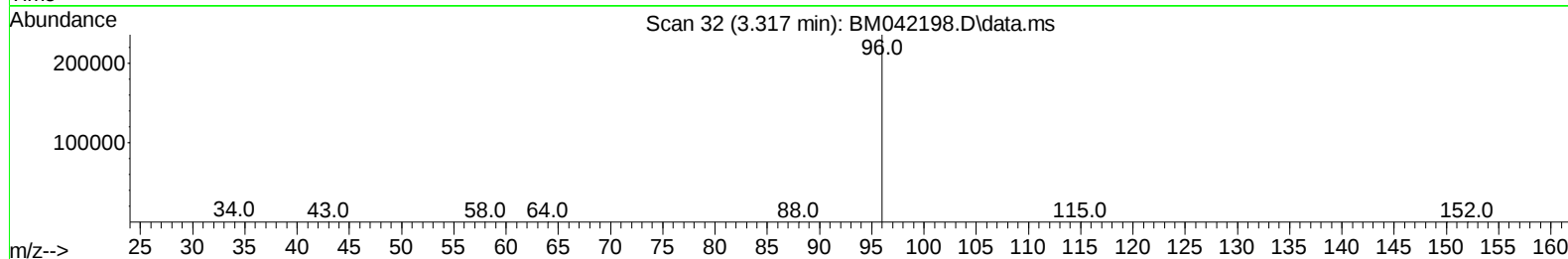
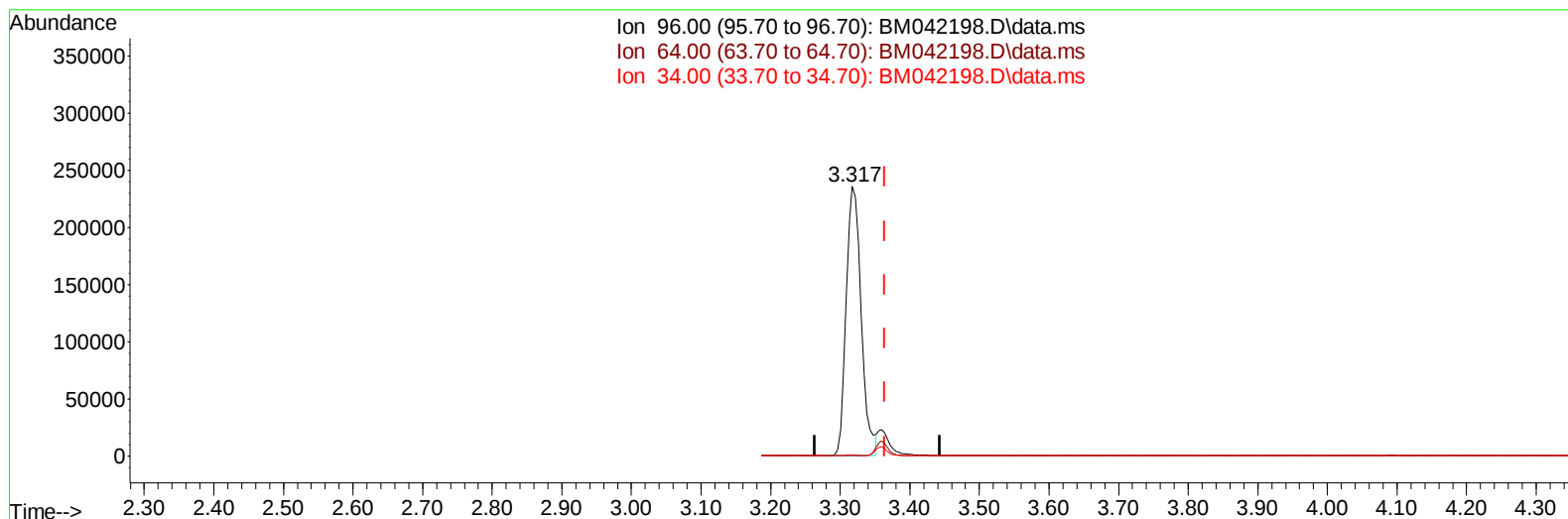
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100423\
 Data File : BM042198.D
 Acq On : 04 Oct 2023 22:28
 Operator : MA/JU
 Sample : 04588-17
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBJ33

Manual IntegrationsAPPROVED

Quant Time: Oct 05 01:55:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 02 19:28:41 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/05/2023
 Supervised By :mohammad ahmed 10/06/2023



TIC: BM042198.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.317min (-0.046) 74.42 ng/u1

response 350617

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.90	0.12#
34.00	48.60	0.45#
0.00	0.00	0.00

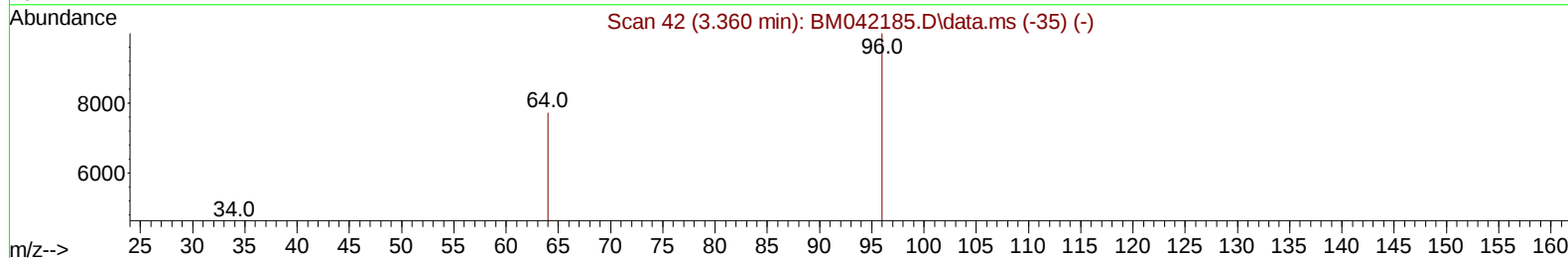
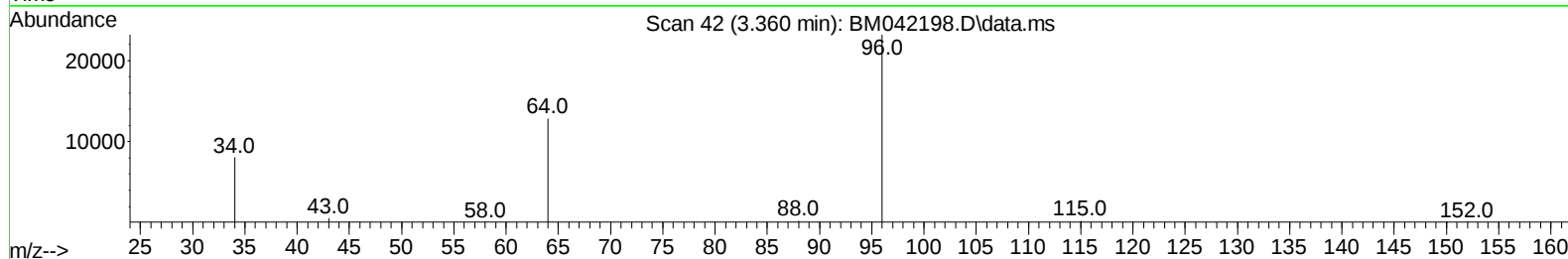
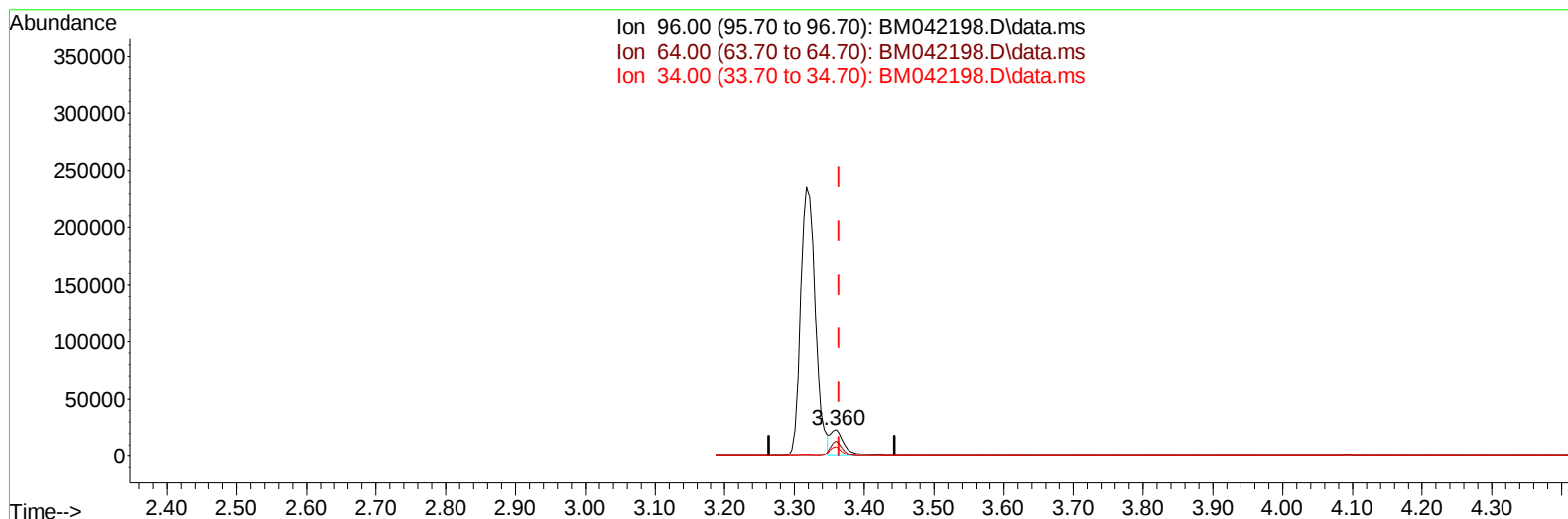
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TIC: BM042198.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.360min (-0.004) 6.65 ng/ul m

response 31313

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.90	55.56#
34.00	48.60	34.67#
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	7.933	152		3538	0.400	ng/ul	0.00
4) Naphthalene-d8	10.741	136		8908	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.578	164		4209	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.329	188		9481	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.504	240		6424	0.400	ng/ul	-0.01
23) Perylene-d12	23.901	264		6692	0.400	ng/ul	#-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.360	96		31313m	6.646	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.324	152		4228	0.346	ng/ul	-0.01
18) Fluoranthene-d10	19.352	212		8277	0.343	ng/ul	0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.397	88		1593	0.324	ng/ul	89
5) Naphthalene	10.790	128		5903	0.199	ng/ul#	94
24) Benzo(b)fluoranthene	23.164	252		947	0.024	ng/ul#	1

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

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