

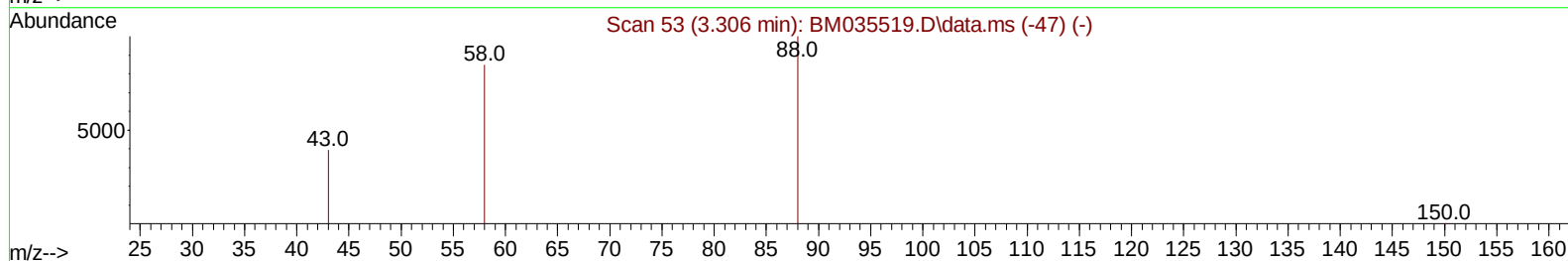
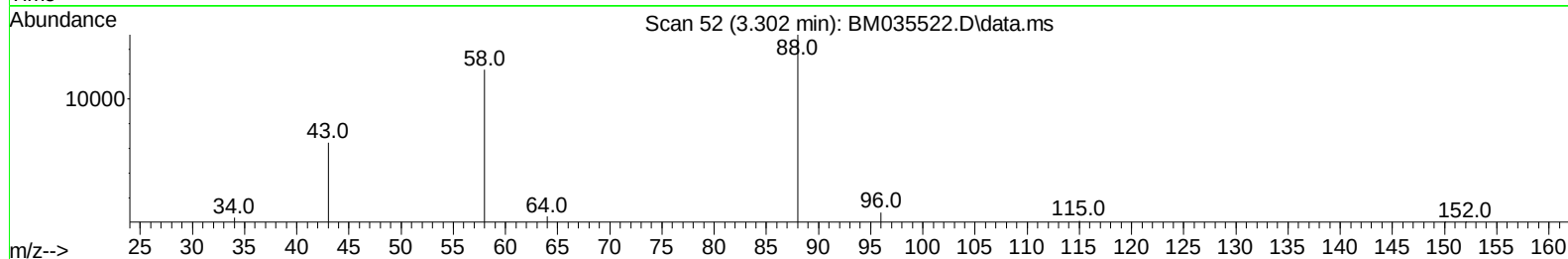
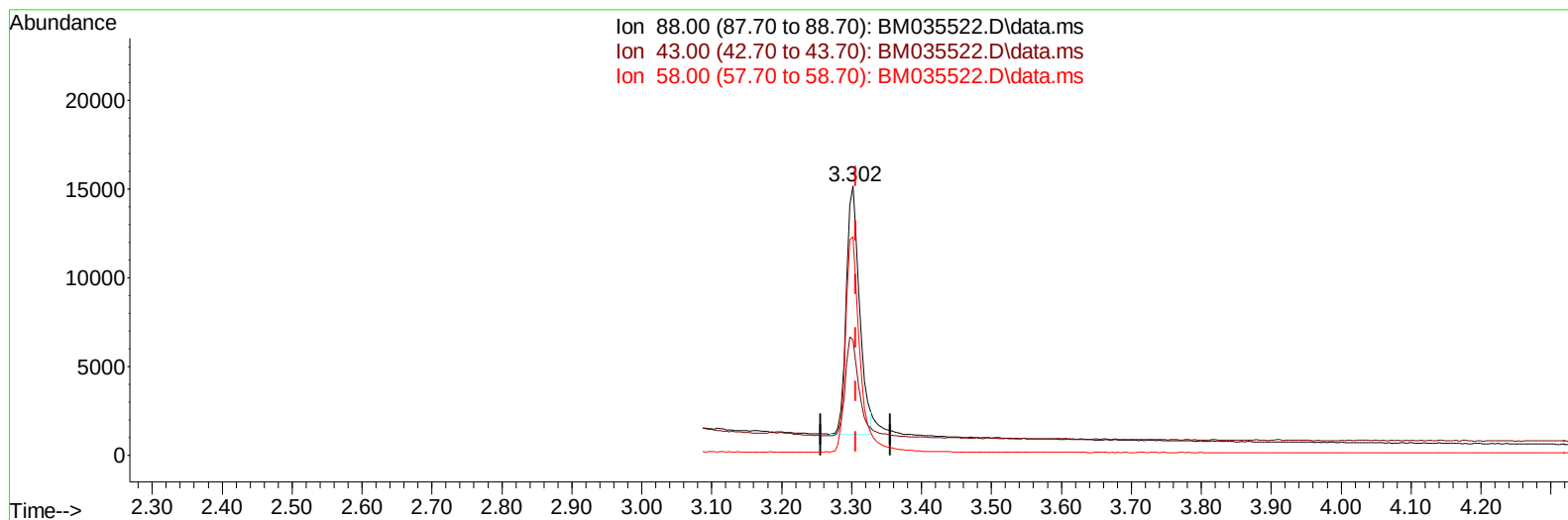
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061822\
 Data File : BM035522.D
 Acq On : 18 Jun 2022 15:12
 Operator : CG/JU
 Sample : SSTD3.248
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.2048

Manual IntegrationsAPPROVED

Quant Time: Jun 20 01:40:46 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 01:37:25 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/20/2022
 Supervised By :mohammad ahmed 06/22/2022



TIC: BM035522.D\data.ms

(2) 1,4-Dioxane

3.302min (-0.004) 3.53 ng/u1

response 18222

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	88.00	42.77#
58.00	57.00	81.33#
0.00	0.00	0.00

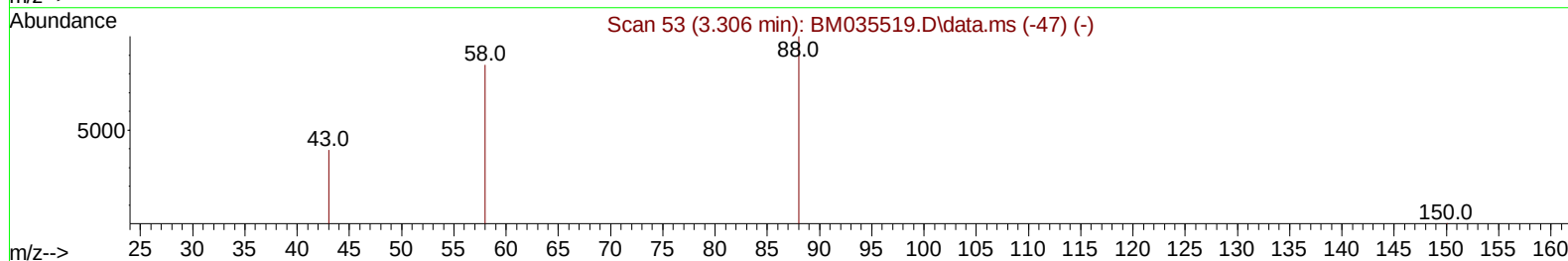
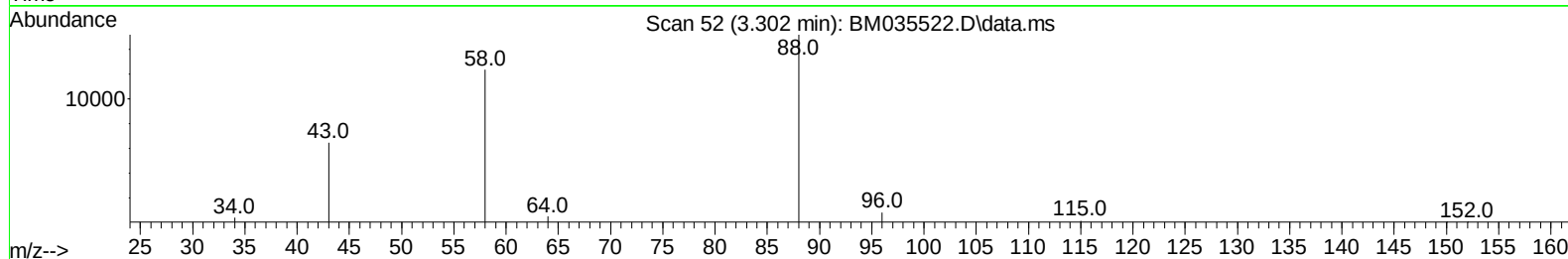
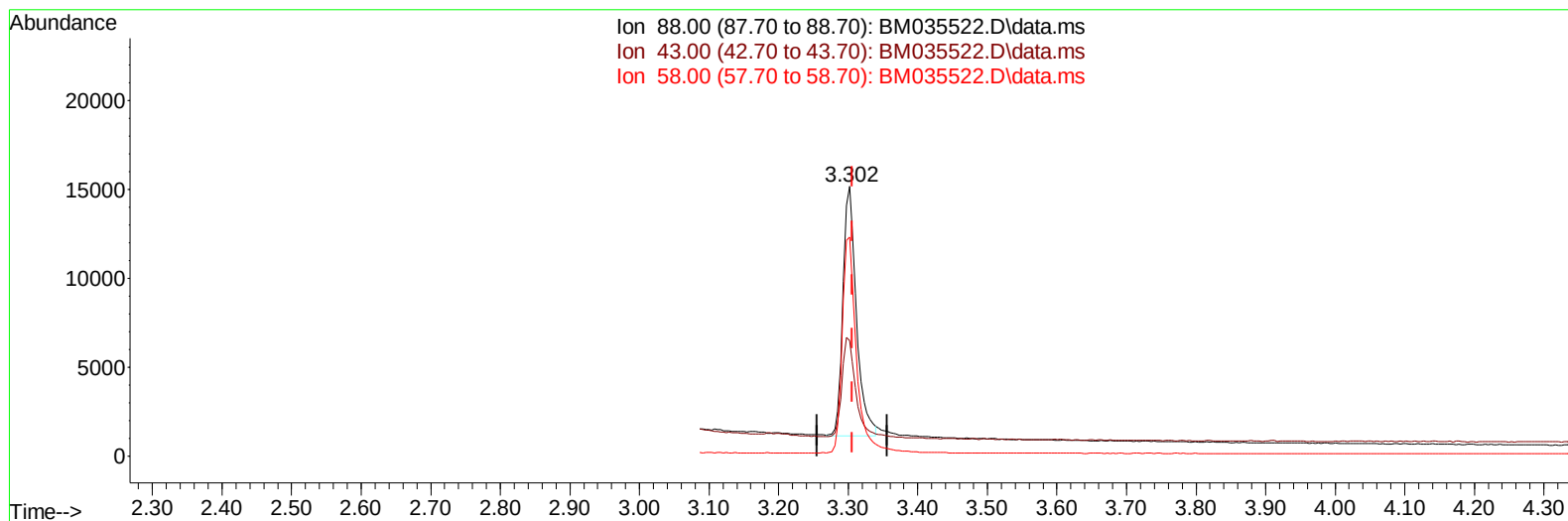
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(2) 1,4-Dioxane

3.302min (-0.004) 3.65 ng/ul m

response 18867

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	88.00	42.77#
58.00	57.00	81.33#
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	7.833	152		4291	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.616	136		14231	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.459	164		7949	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.208	188		15871	0.400	ng/ul	-0.02
17) Chrysene-d12	21.398	240		12446	0.400	ng/ul	-0.01
23) Perylene-d12	23.756	264		10689	0.400	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.268	96		17772	3.457	ng/ul	0.00
6) 2-Methylnaphthalene-d10	0.000	152		0d	0.000	ng/ul	
18) Fluoranthene-d10	0.000	212		0d	0.000	ng/ul	
Target Compounds							Qvalue
2) 1,4-Dioxane	3.302	88		18867m	3.655	ng/ul	
14) Pentachlorophenol	16.891	266		9302	1.552	ng/ul#	63

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

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