Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039458.D
Acq On : 18 Apr 2023 15:26
Operator : CG/JU
Sample : PB152151BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Apr 19 00:19:27 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 18 23:56:45 2023
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SBLK151

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/19/2023
Supervised By :Jagrut Upadhyay 04/19/2023

04/19/2023
Supervised By :Jagrut Upadhyay
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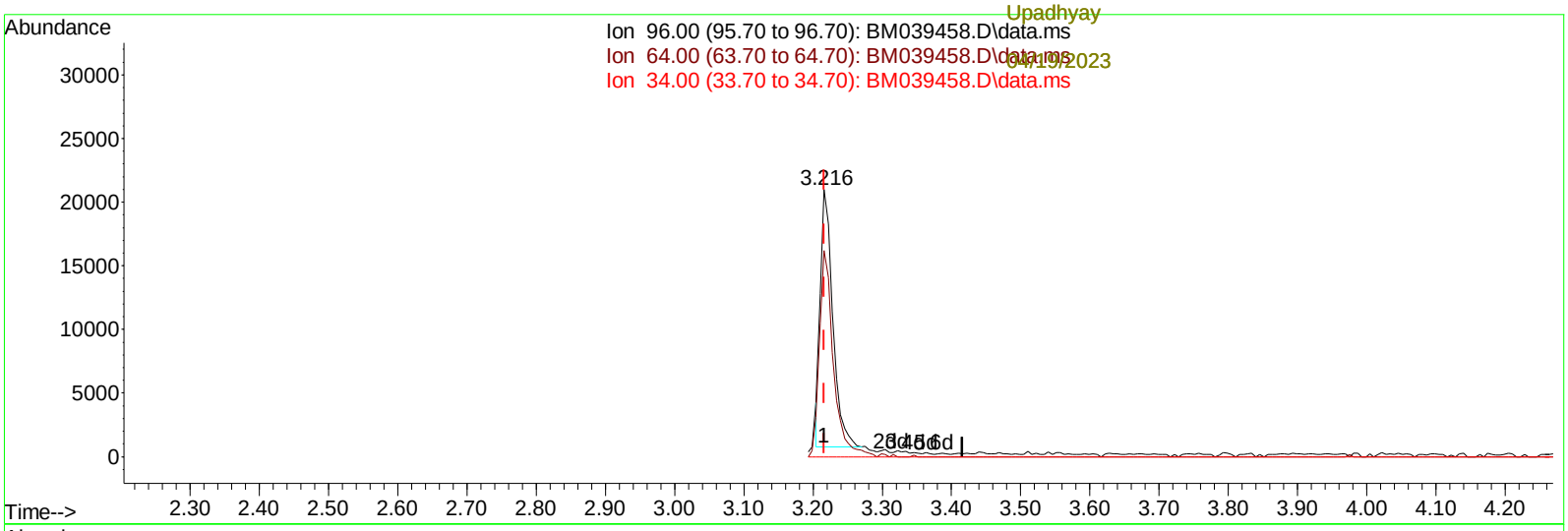

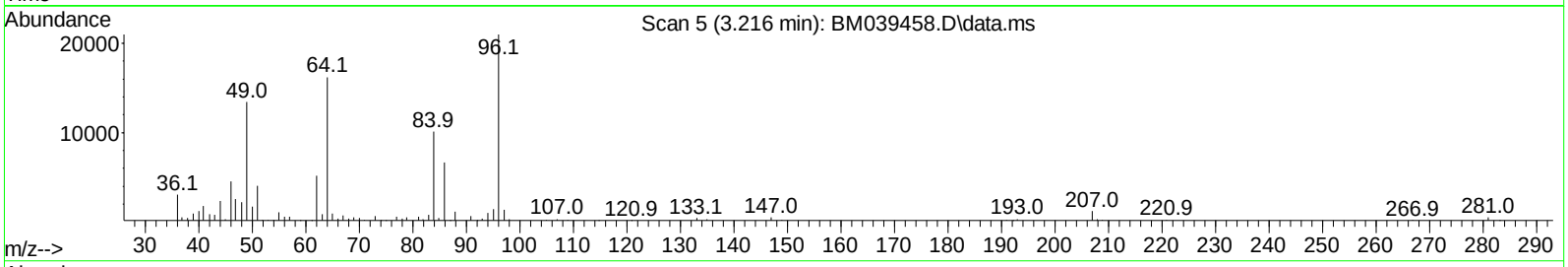

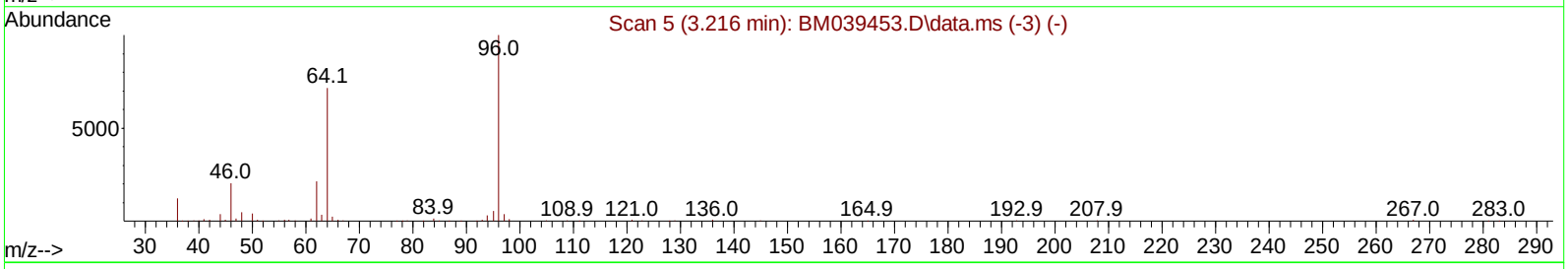

TIC: BM039458.D\data.ms

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

3.216min (+ 0.000) 4.99 ng/uL

response 25170

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.10	77.15
34.00	0.00	0.00
0.00	0.00	0.00

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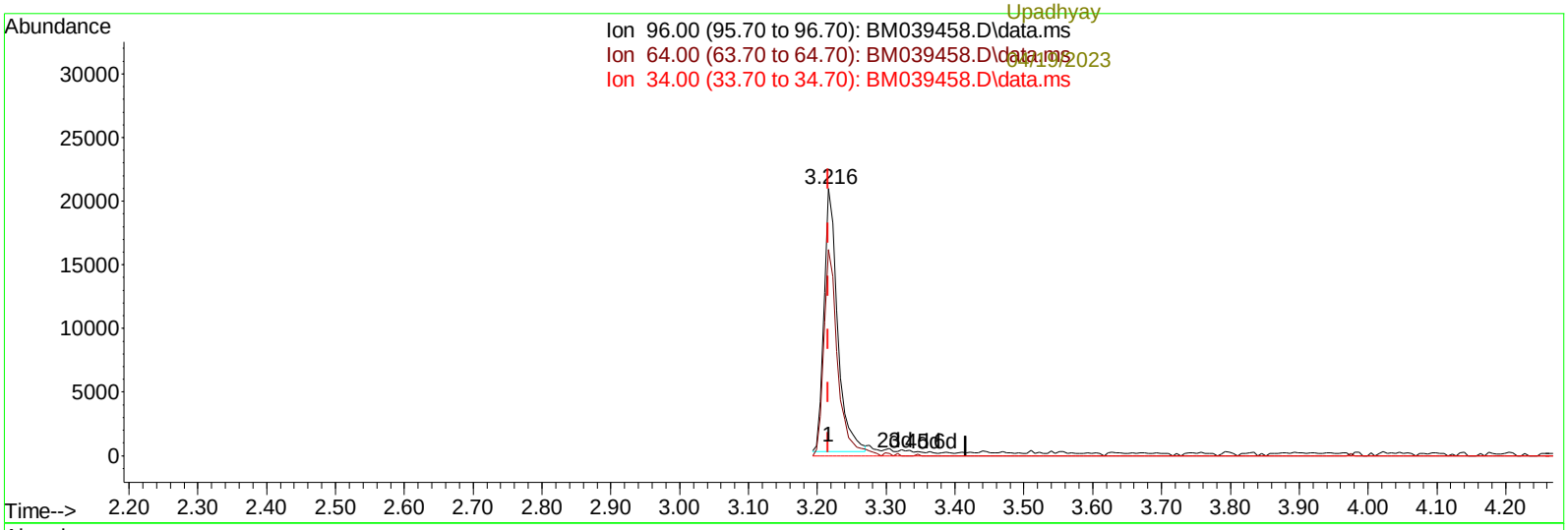

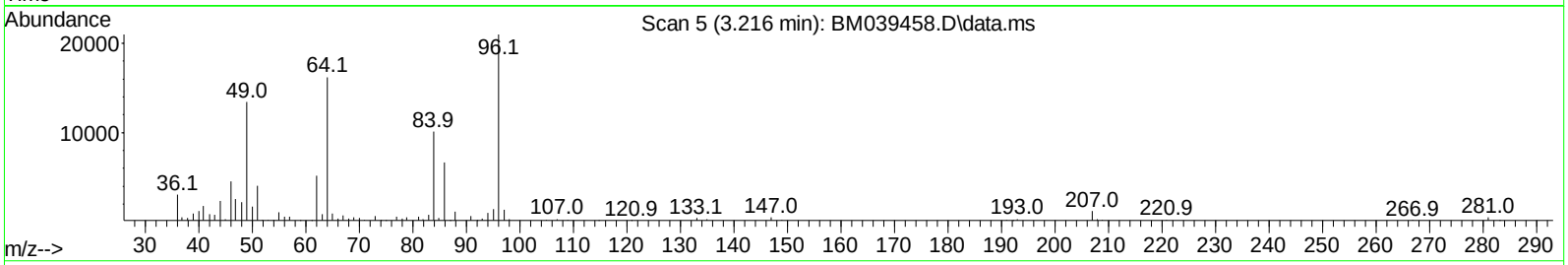

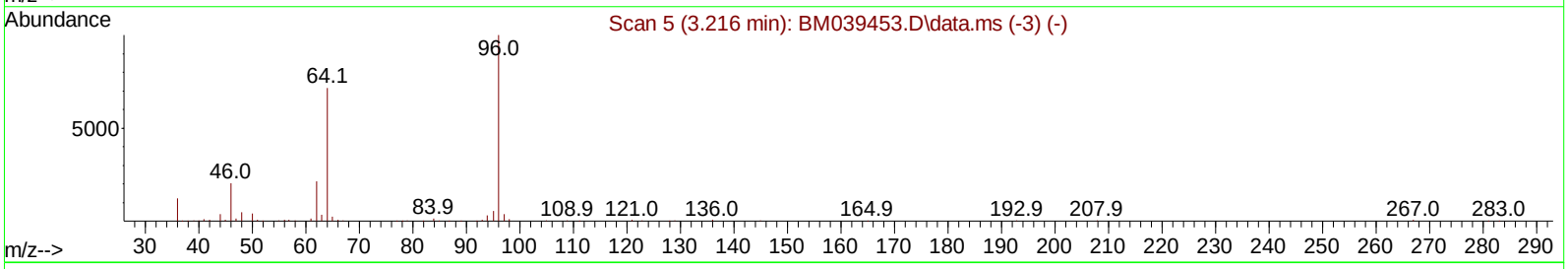

TIC: BM039458.D\data.ms

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

3.216min (+ 0.000) 5.64 ng/uL m

response 28419

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.10	77.15
34.00	0.00	0.00
0.00	0.00	0.00

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Acq On : 18 Apr 2023 15:26
Operator : CG/JU
Sample : PB152151BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Apr 19 00:19:53 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
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ClientSampleId :
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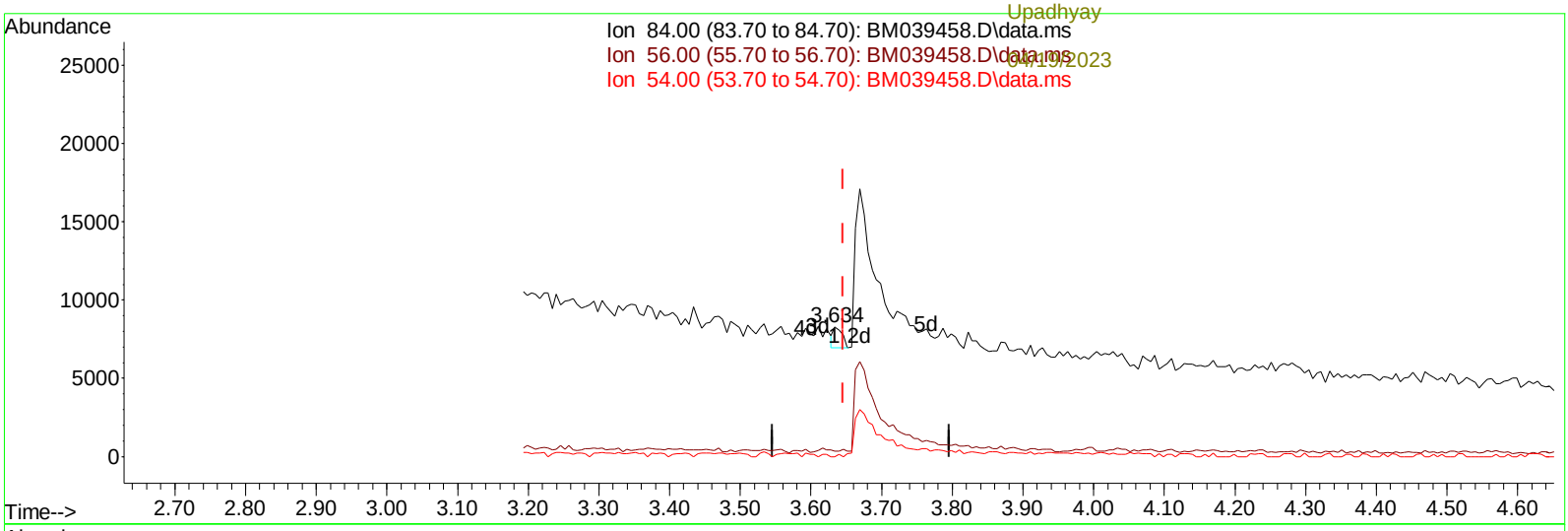

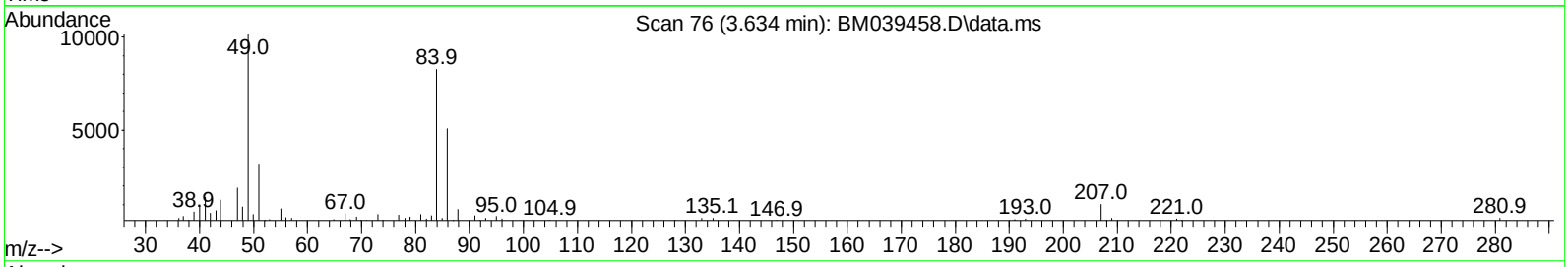

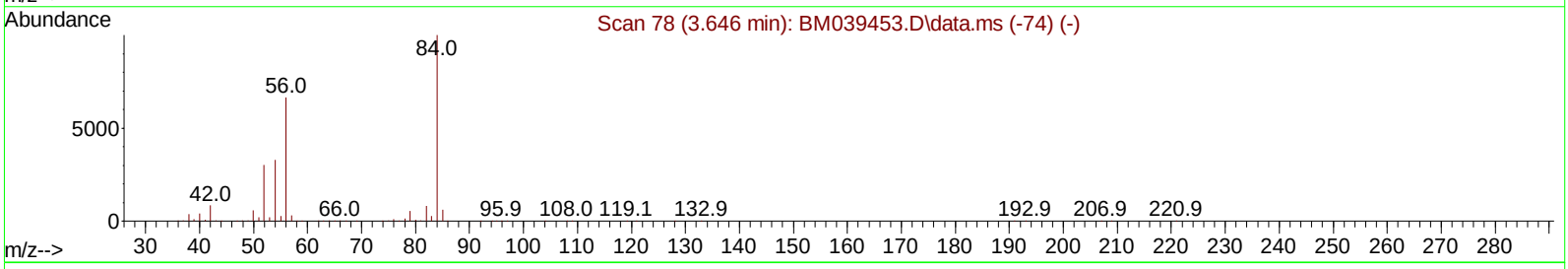

TIC: BM039458.D\data.ms

(4) Pyridine-d5 (S)

3.634min (-0.012) 0.08 ng/ul

response 1141

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	64.40	4.00#
54.00	32.00	0.00#
0.00	0.00	0.00

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Operator : CG/JU
Sample : PB152151BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Apr 19 00:19:53 2023
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Quant Title : SVOA CALIBRATION
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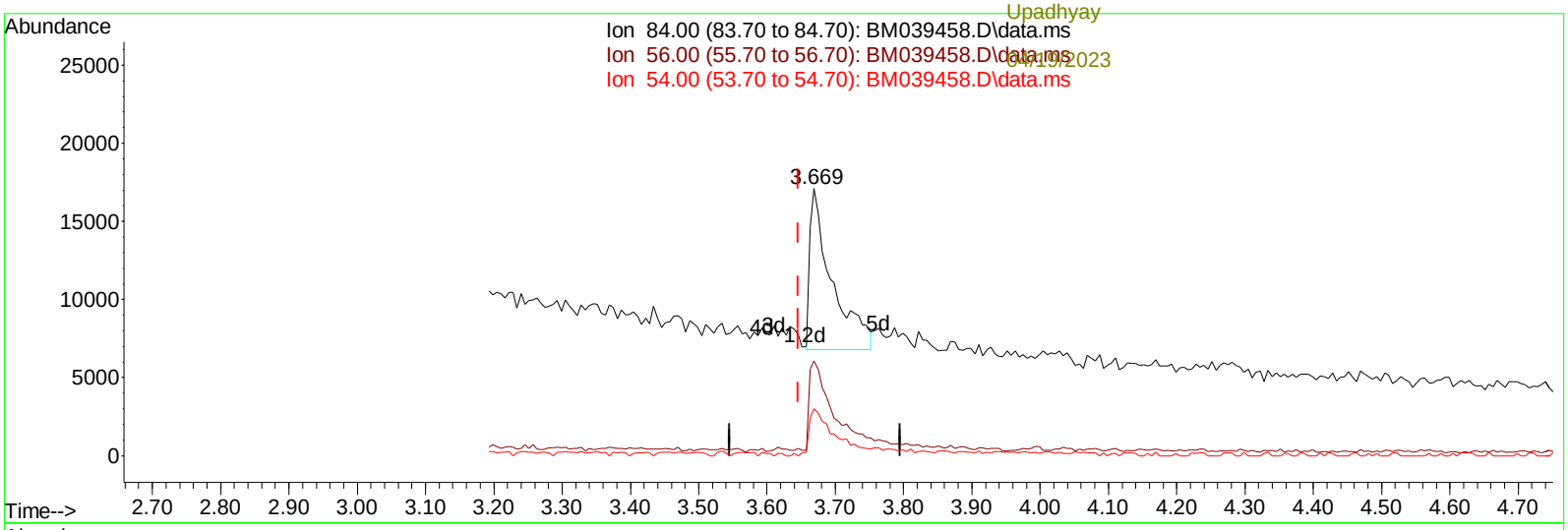

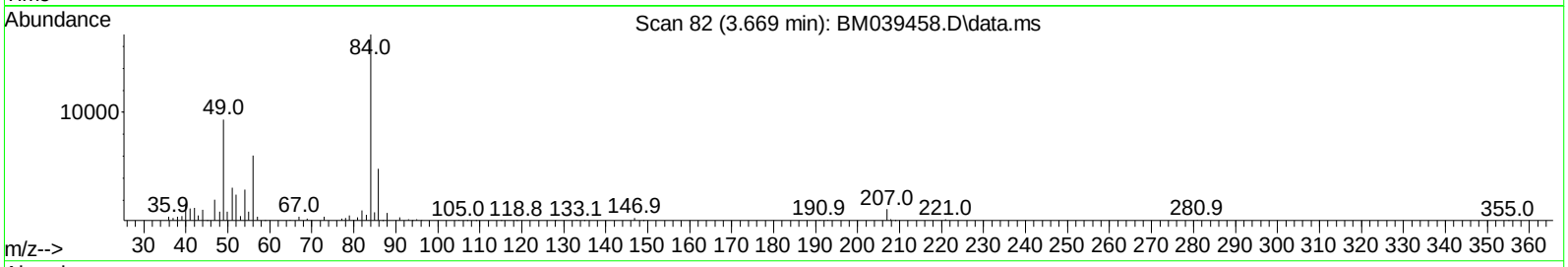

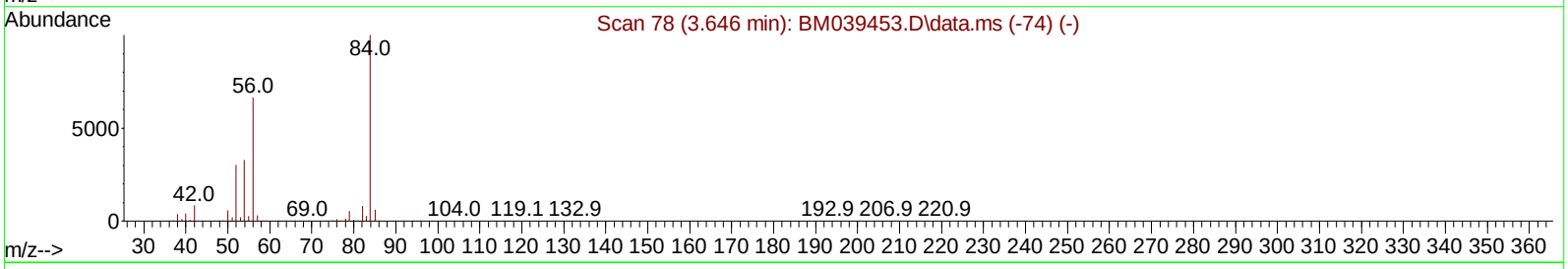

TIC: BM039458.D\data.ms

(4) Pyridine-d5 (S)

3.669min (+ 0.023) 1.63 ng/ul m

response 23153

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	64.40	35.60#
54.00	32.00	17.49#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039458.D
Acq On : 18 Apr 2023 15:26
Operator : CG/JU
Sample : PB152151BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Apr 19 00:20:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 18 23:56:45 2023
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Instrument :
BNA_M
ClientSampleId :
SBLK151

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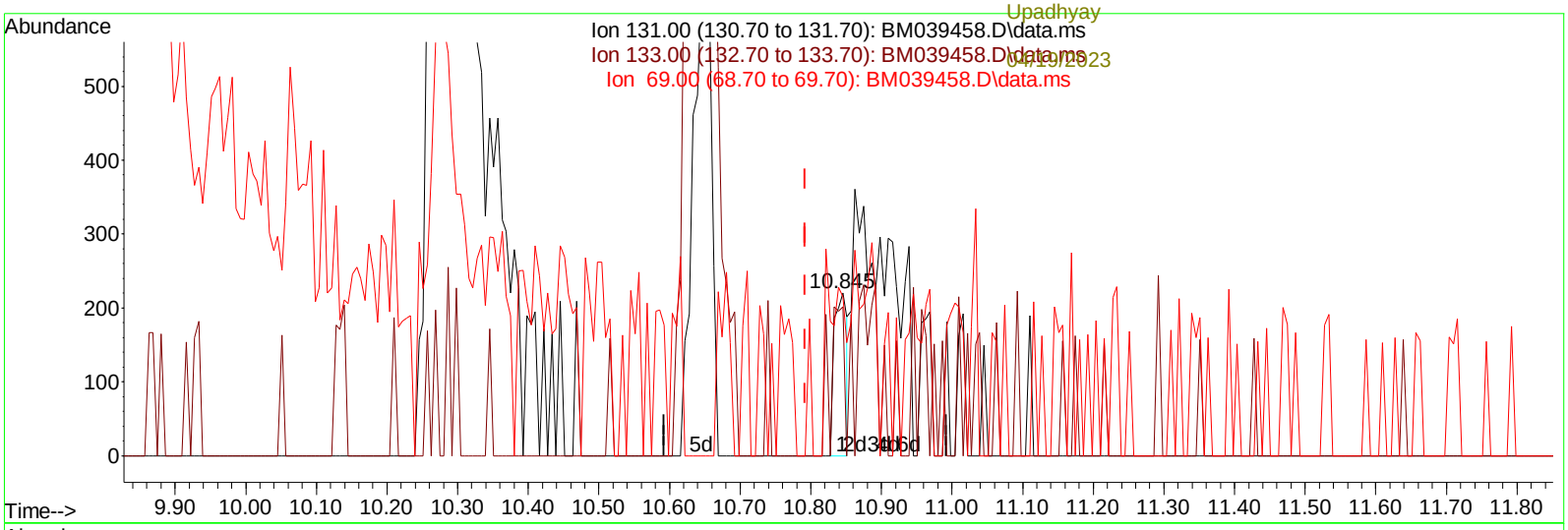

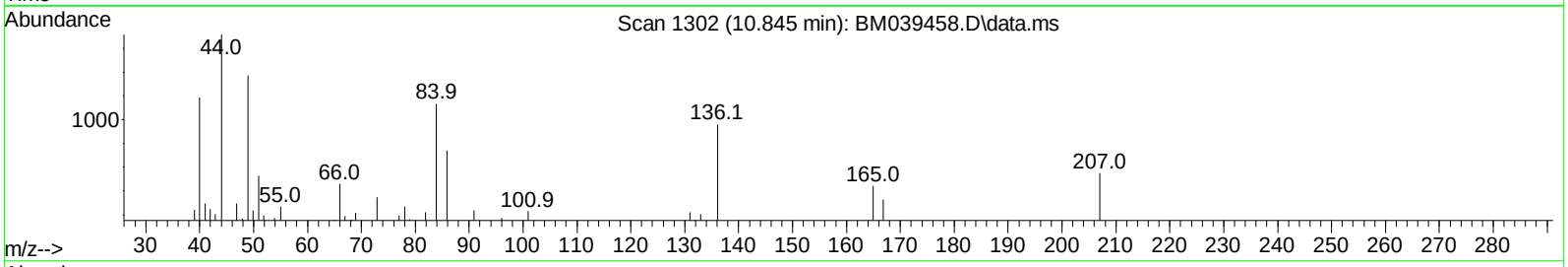

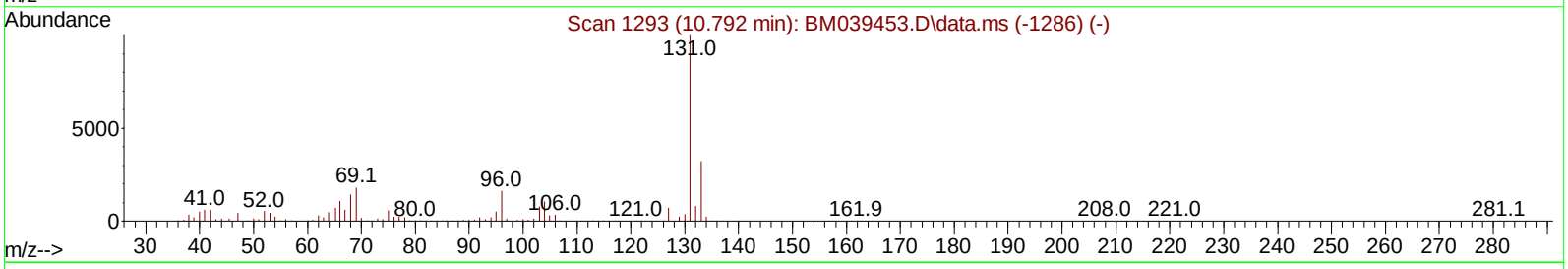

TIC: BM039458.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (+ 0.053) 0.01 ng/ul

response 280

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	91.82#
69.00	18.10	97.73#
0.00	0.00	0.00

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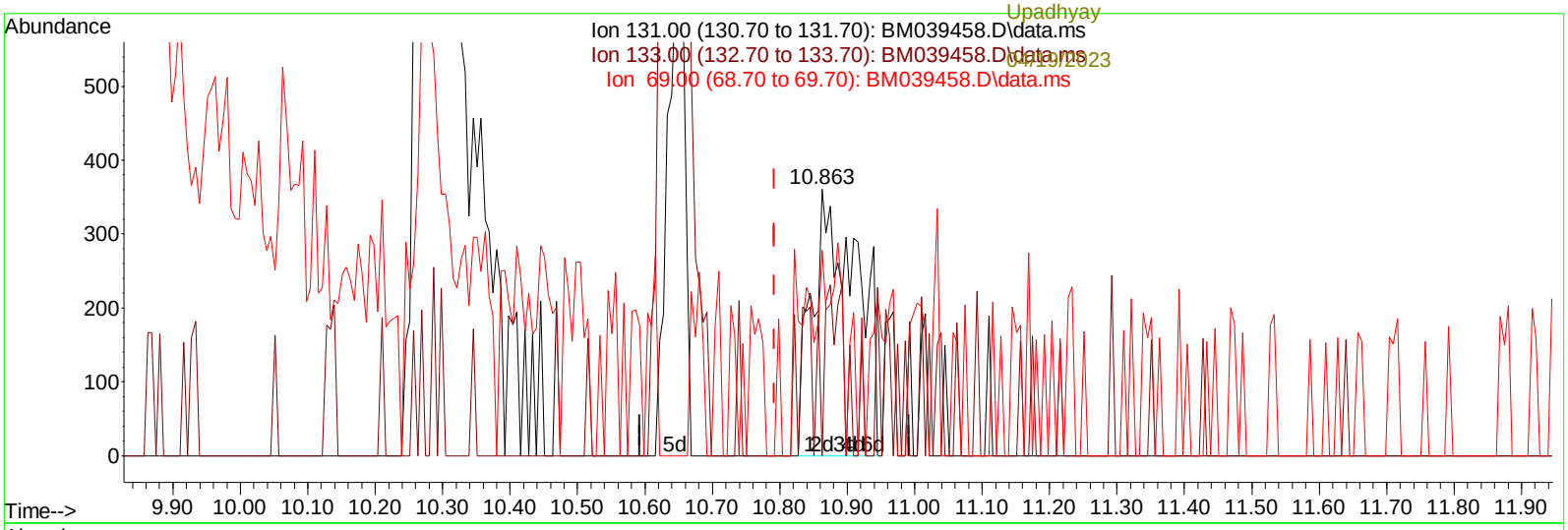

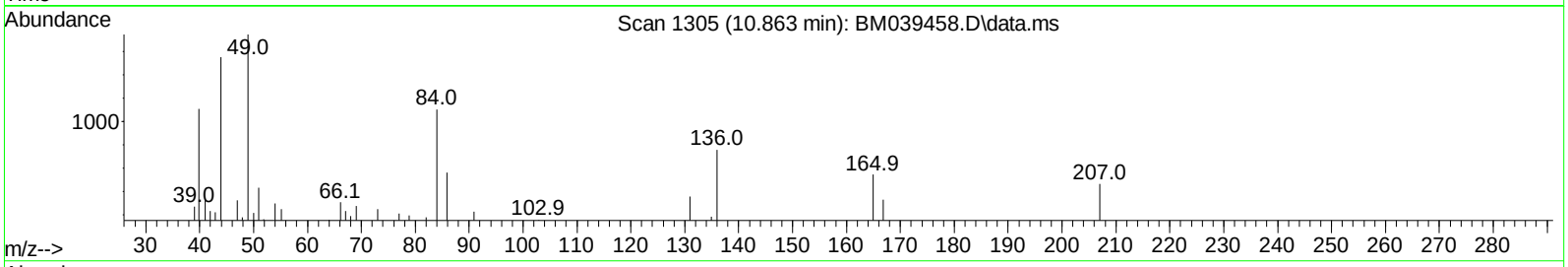

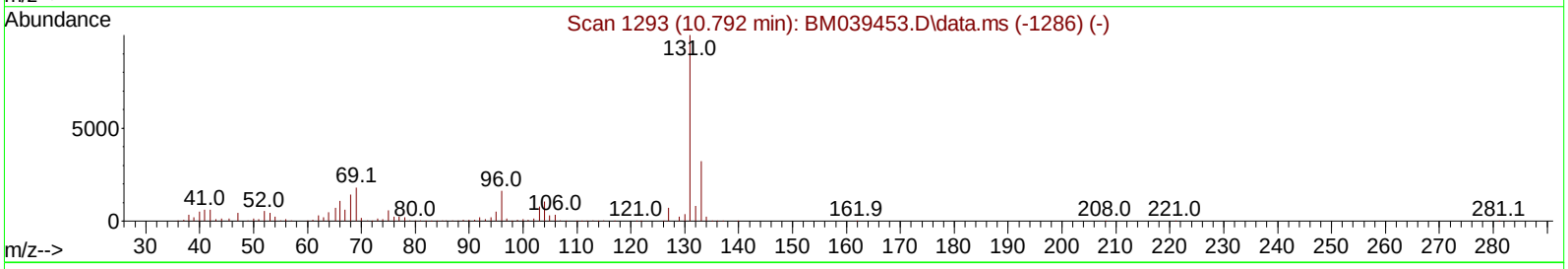

TIC: BM039458.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.863min (+ 0.070) 0.08 ng/ul m

response 1665

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	0.00#
69.00	18.10	77.01#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	193461	20.000	ng/ul	0.00
20) Naphthalene-d8	10.639	136	893497	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	601612	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	1309967	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	1189384	20.000	ng/ul	0.00
88) Perylene-d12	23.809	264	1176200	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	28419m	5.639	ng/uL	0.00
4) Pyridine-d5	3.669	84	23153m	1.635	ng/ul	0.02
7) Phenol-d5	7.022	99	316419	17.870	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	332854	28.441	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	398409	28.684	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	244968	17.175	ng/ul	0.00
21) Nitrobenzene-d5	9.004	128	202407	28.158	ng/ul	0.00
24) 2-Nitrophenol-d4	9.728	143	231097	28.630	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.280	165	345287	24.861	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	1665m	0.080	ng/ul	0.07
46) Dimethylphthalate-d6	13.904	166	1407925	29.018	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	1056327	19.379	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	141256	19.237	ng/ul	0.00
60) Fluorene-d10	15.486	176	1184317	29.793	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	205939	23.657	ng/ul	0.00
73) Anthracene-d10	17.339	188	1754335	27.847	ng/ul	0.00
81) Pyrene-d10	19.639	212	2128683	26.934	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.656	264	1267136	19.954	ng/ul	0.00

Target Compounds Qvalue

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

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