

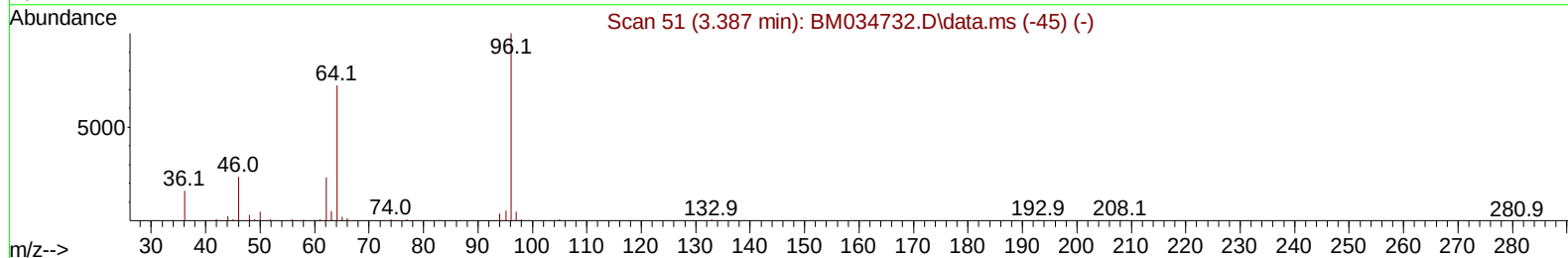
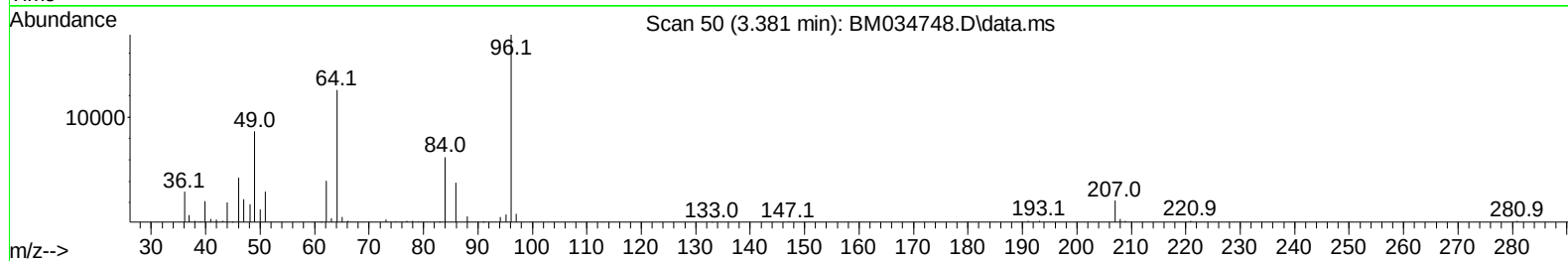
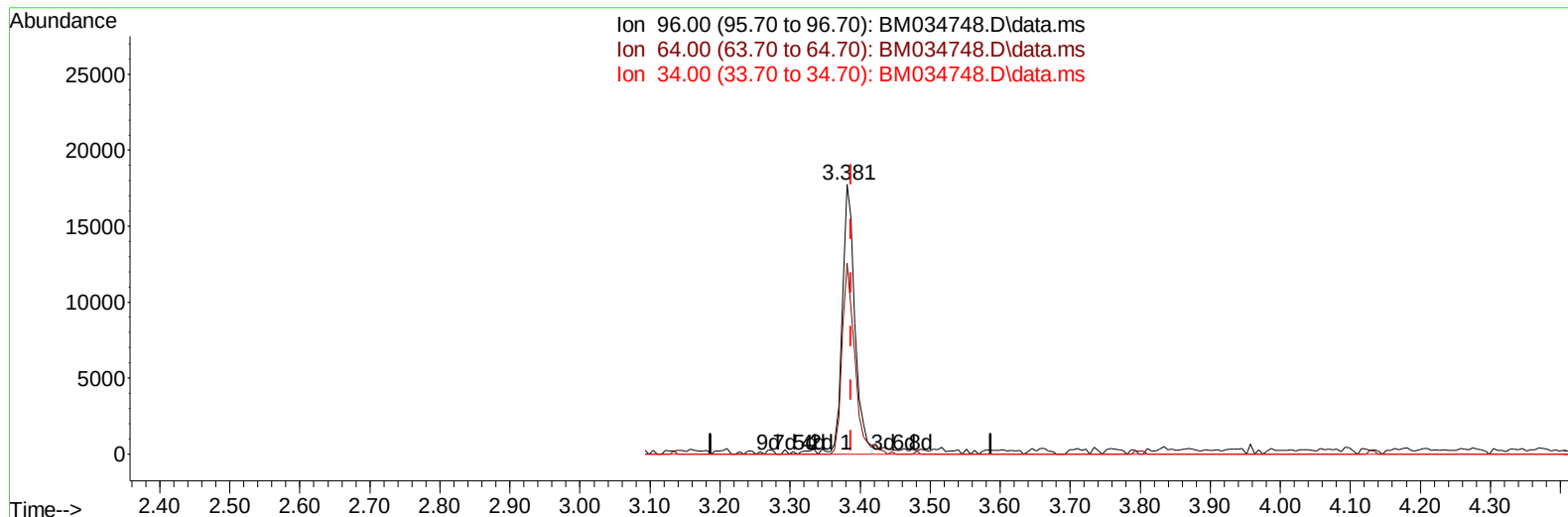
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042122\
 Data File : BM034748.D
 Acq On : 21 Apr 2022 15:42
 Operator : CG/JU
 Sample : N2473-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0J35

Manual IntegrationsAPPROVED

Quant Time: Apr 22 01:25:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 05:03:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022
 Supervised By :mohammad ahmed 04/26/2022



TIC: BM034748.D\data.ms

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

3.381min (-0.006) 5.20 ng/uL

response 22232

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.40	70.77
34.00	0.00	0.00
0.00	0.00	0.00

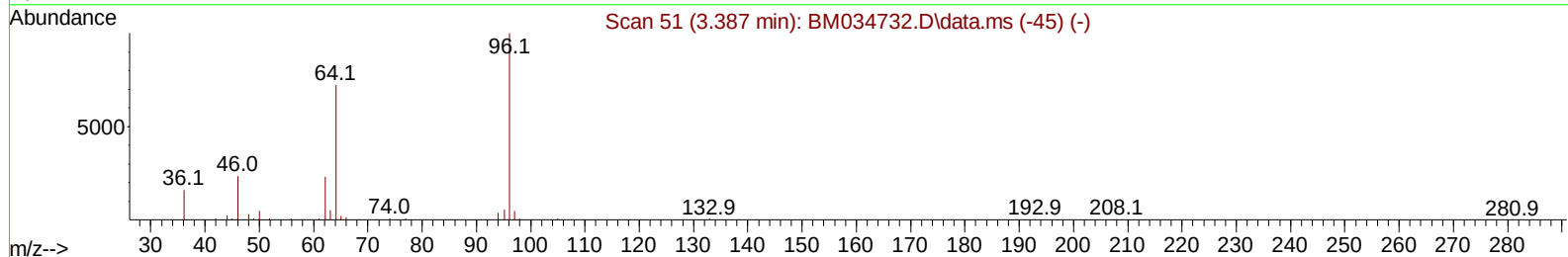
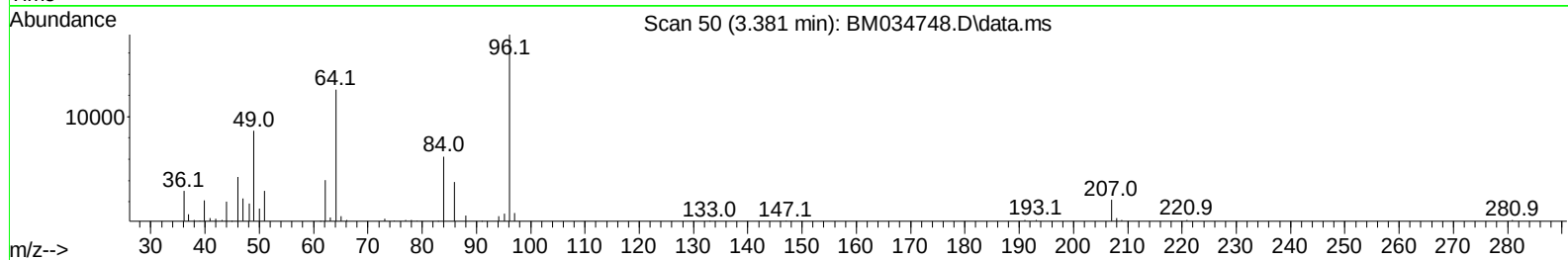
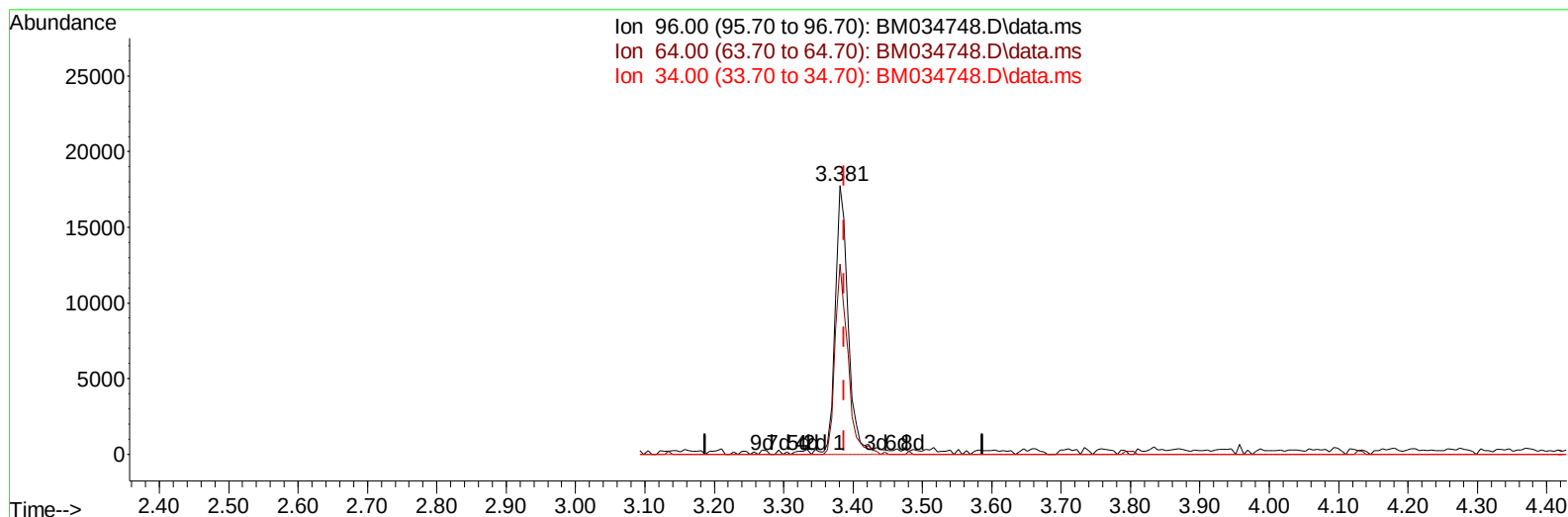
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TIC: BM034748.D\data.ms

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

3.381min (-0.006) 5.27 ng/uL m

response 22552

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.40	70.77
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	187855	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	718482	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.574	164	437239	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.315	188	872457	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	843194	20.000	ng/ul	0.00
88) Perylene-d12	23.891	264	865266	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	22552m	5.275	ng/uL	0.00
4) Pyridine-d5	3.799	84	106033	9.992	ng/ul	0.00
7) Phenol-d5	7.104	99	104151	7.341	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	288399	36.245	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	319113	26.574	ng/ul	0.00
15) 4-Methylphenol-d8	8.645	113	198267	17.465	ng/ul	-0.01
21) Nitrobenzene-d5	9.104	128	186644	35.562	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	182573	33.504	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	322954	28.135	ng/ul	0.00
31) 4-Chloroaniline-d4	10.880	131	453199	29.199	ng/ul	0.00
46) Dimethylphthalate-d6	13.986	166	1144932	35.969	ng/ul	0.00
49) Acenaphthylene-d8	14.268	160	1344007	32.832	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	34508	6.458	ng/ul	0.00
60) Fluorene-d10	15.568	176	1005487	36.506	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.674	200	151459	30.259	ng/ul	0.00
73) Anthracene-d10	17.415	188	1680447	39.555	ng/ul	0.00
81) Pyrene-d10	19.703	212	1884042	41.055	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.738	264	1852683	42.335	ng/ul	0.00

Target Compounds Qvalue

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

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