

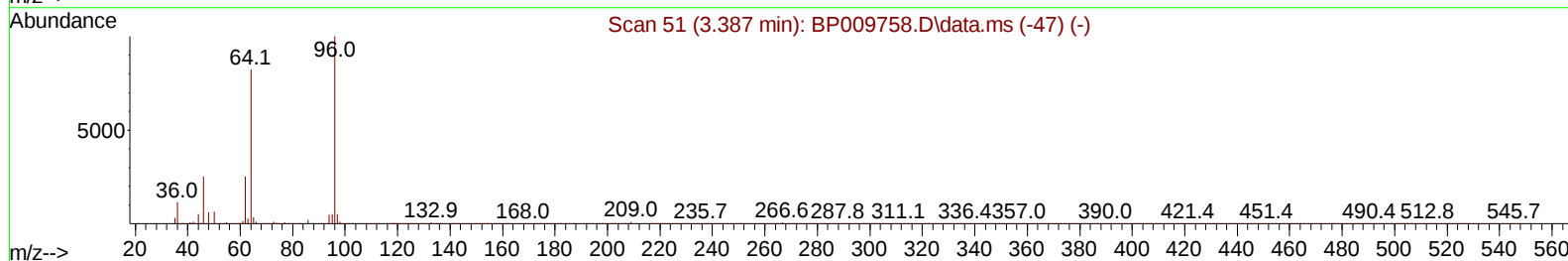
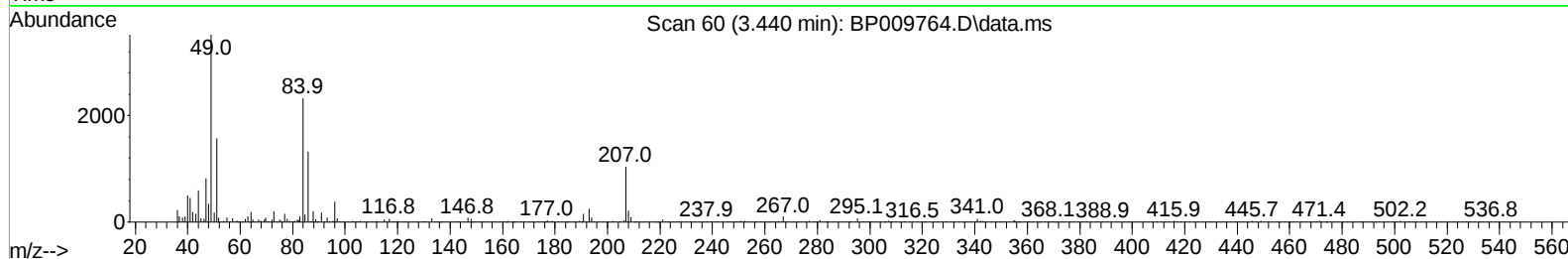
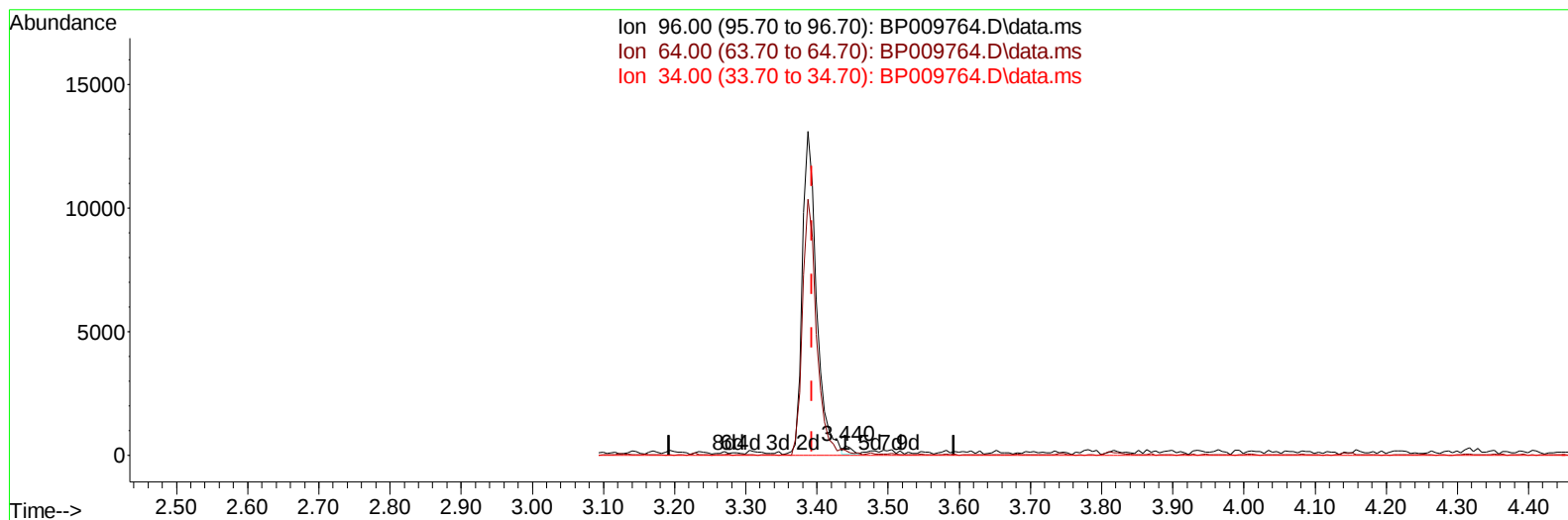
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009764.D
 Acq On : 11 Apr 2022 22:42
 Operator : CG/JU
 Sample : N2342-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY1

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:54:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009764.D\data.ms

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

3.440min (+ 0.047) 0.08 ng/uL

response 253

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	51.30
34.00	0.00	0.00
0.00	0.00	0.00

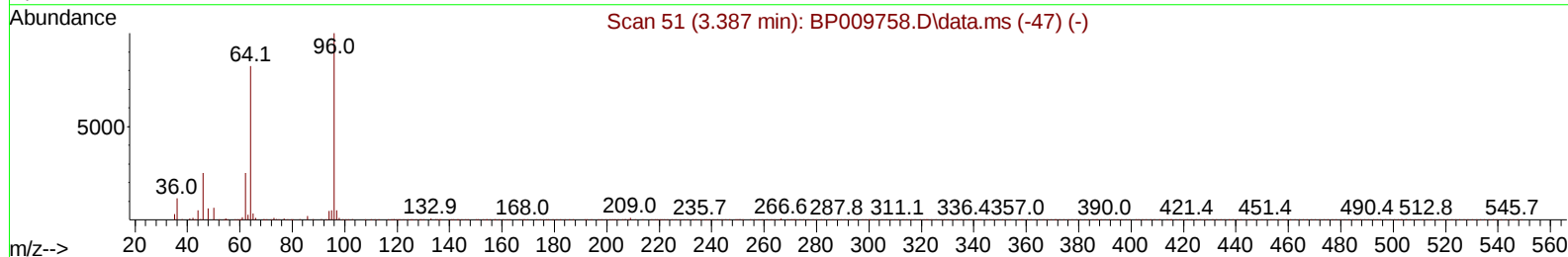
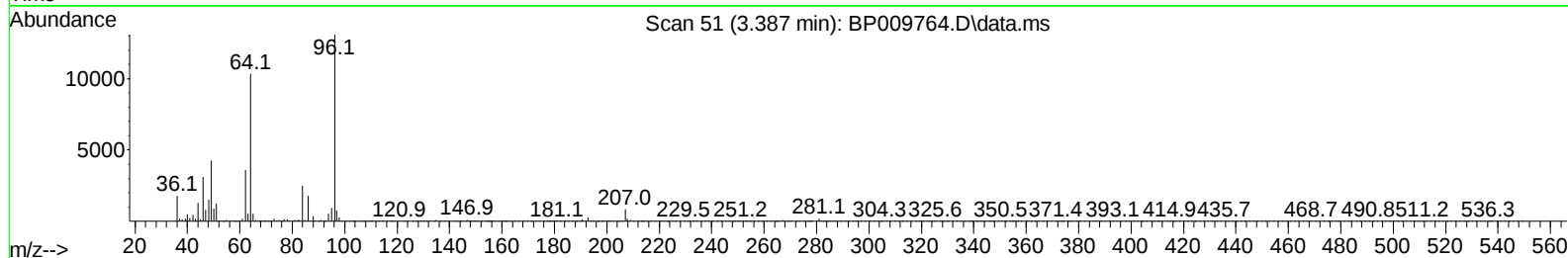
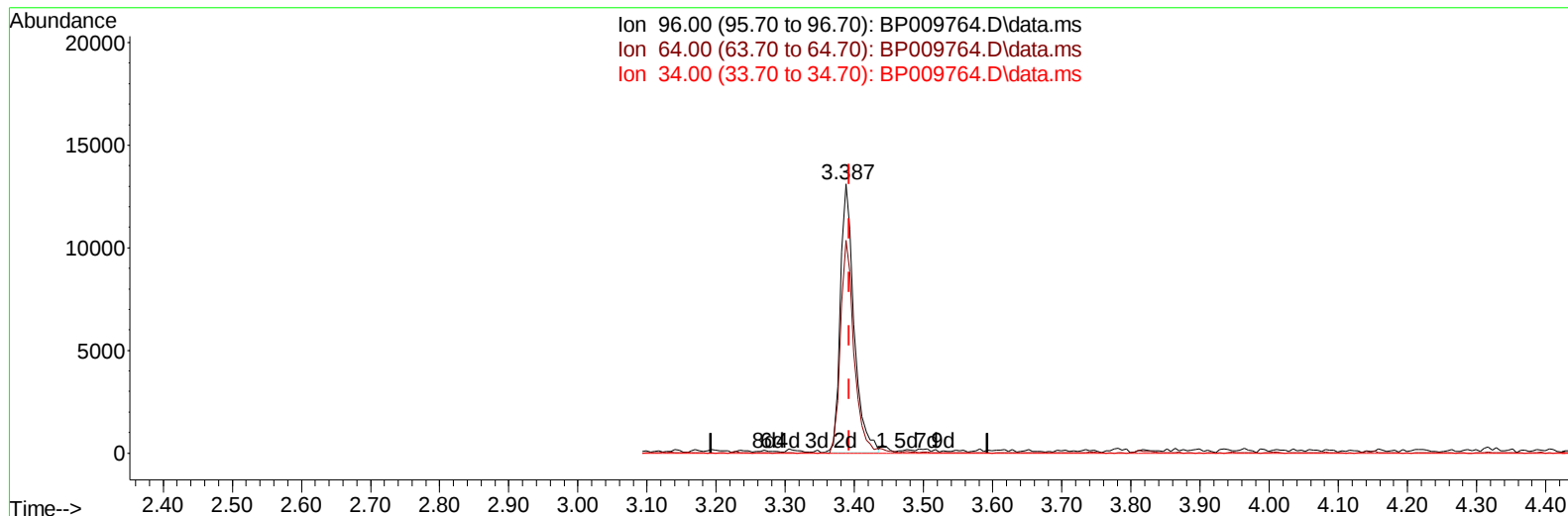
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TIC: BP009764.D\data.ms

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

3.387min (-0.006) 6.02 ng/uL m

response 18112

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.04#
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	7.969	152	134301	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	613816	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.598	164	455105	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	1006014	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.428	240	1049328	20.000	ng/ul	#-0.02
88) Perylene-d12	23.904	264	958150	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	18112m	6.019	ng/uL	0.00
4) Pyridine-d5	3.811	84	46290	5.579	ng/ul	0.00
7) Phenol-d5	7.122	99	83286	7.064	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	248329	35.377	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	230554	25.923	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	170017	17.273	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	161381	36.459	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	171761	35.394	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	291724	28.727	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.905	131	362404	25.056	ng/ul	-0.01
46) Dimethylphthalate-d6	14.010	166	1300428	36.377	ng/ul	-0.01
49) Acenaphthylene-d8	14.293	160	1378246	32.542	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	49929	7.701	ng/ul	-0.01
60) Fluorene-d10	15.593	176	1096539	36.462	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	210546	35.329	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1921553	39.893	ng/ul	-0.01
81) Pyrene-d10	19.681	212	2286641	39.900	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	2045862	41.002	ng/ul	-0.02

Target Compounds Qvalue

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

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