

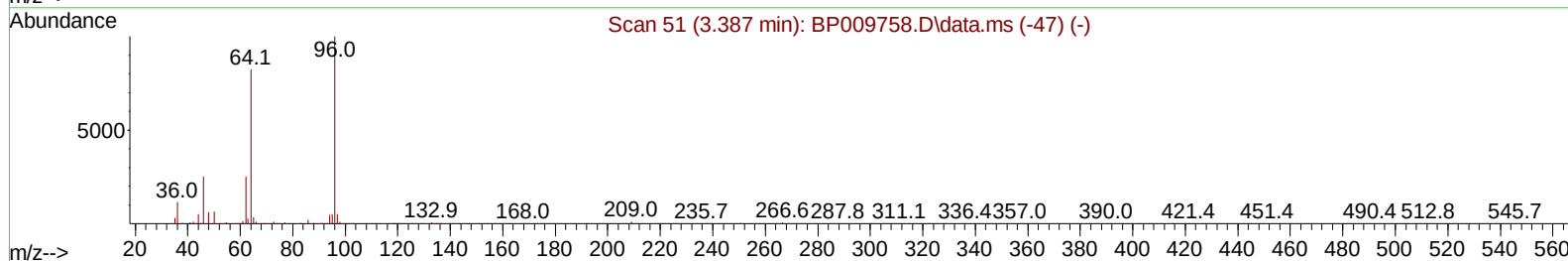
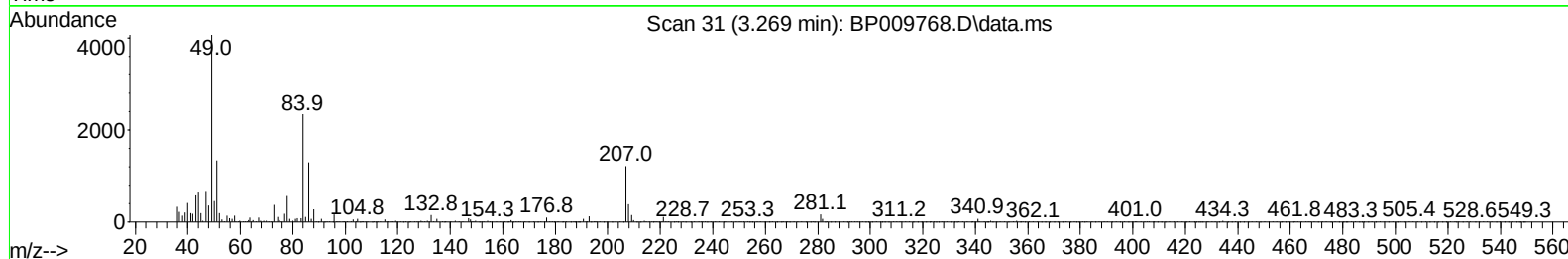
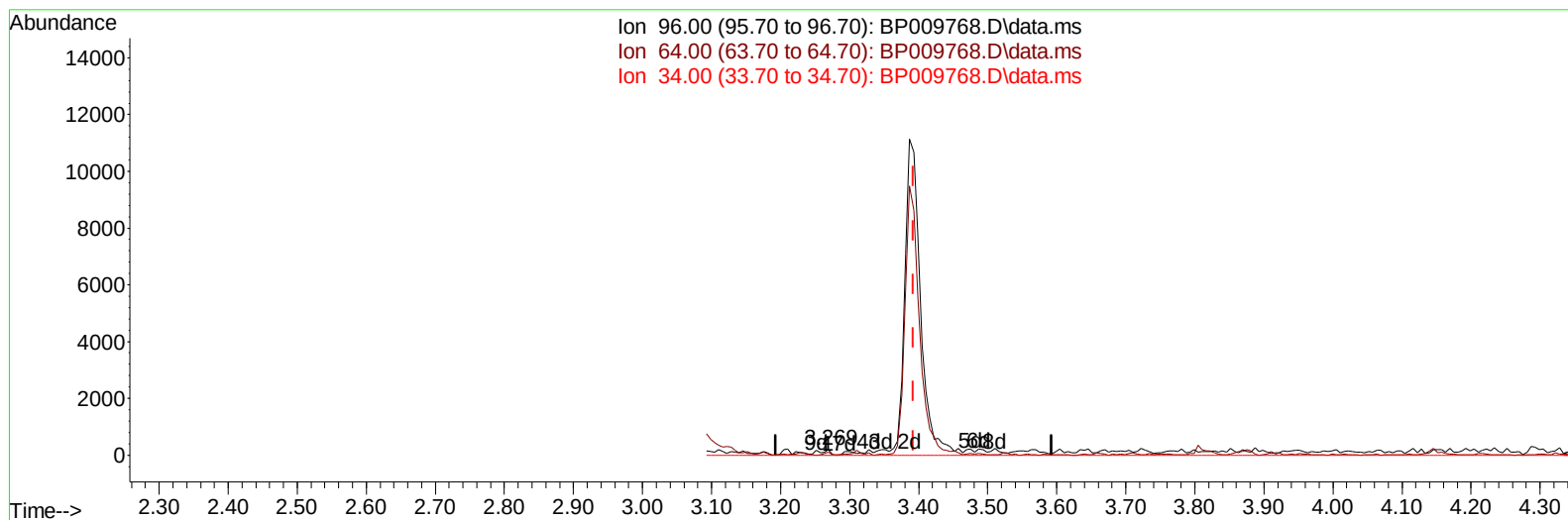
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009768.D
 Acq On : 12 Apr 2022 00:57
 Operator : CG/JU
 Sample : N2338-17
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J69

Manual IntegrationsAPPROVED

Quant Time: Apr 12 02:34:48 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: BP009768.D\data.ms

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

3.269min (-0.124) 0.04 ng/uL

response 106

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	58.29
34.00	0.00	0.00
0.00	0.00	0.00

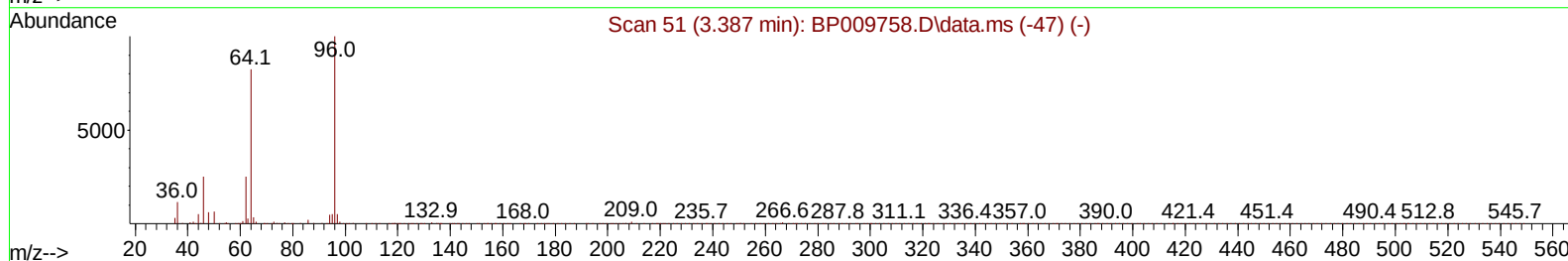
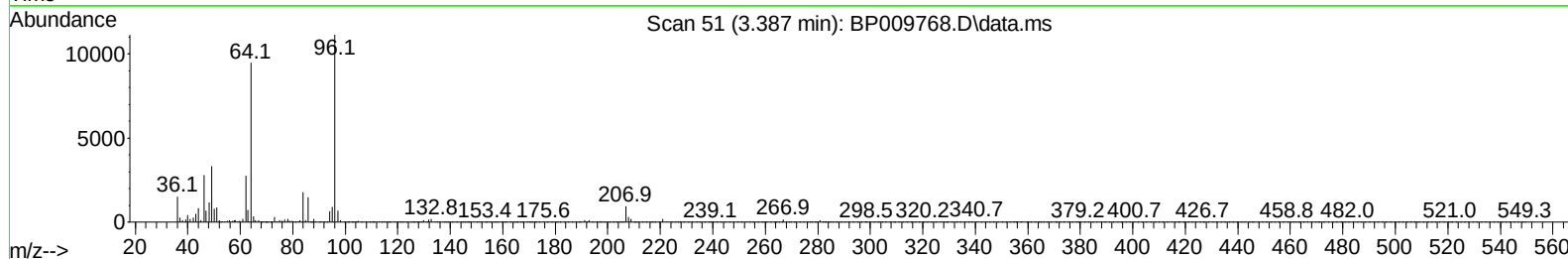
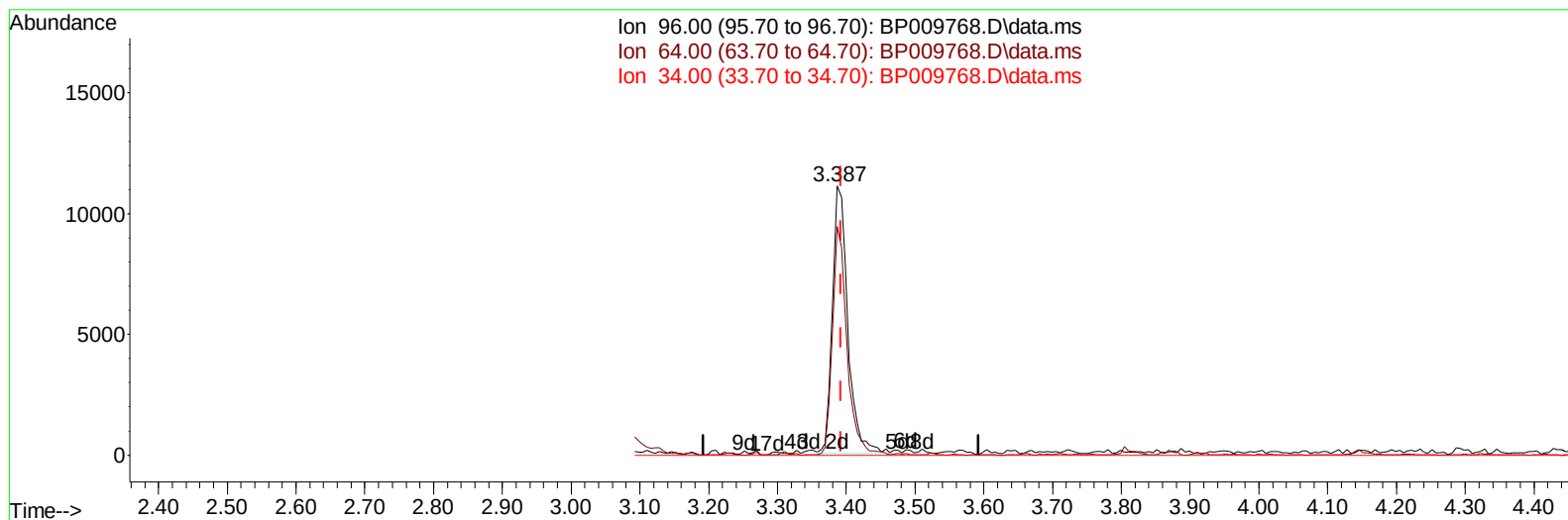
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(3) 1,4-Dioxane-d8 (S)

3.387min (-0.006) 5.79 ng/uL m

response 17212

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	85.07#
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	7.987	152	132577	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	565809	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.598	164	412993	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	920609	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.427	240	907187	20.000	ng/ul	-0.02
88) Perylene-d12	23.898	264	765442	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17212m	5.794	ng/uL	0.00
4) Pyridine-d5	3.810	84	95068	11.607	ng/ul	0.00
7) Phenol-d5	7.128	99	96452	8.287	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	350122	50.527	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	239301	27.256	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	176840	18.200	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	150670	36.927	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.851	143	165631	37.026	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.392	165	291243	31.113	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131	369578	27.720	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1153462	35.556	ng/ul	-0.01
49) Acenaphthylene-d8	14.292	160	1250269	32.531	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	56869	9.666	ng/ul	-0.01
60) Fluorene-d10	15.592	176	997455	36.549	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	191065	35.034	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1713086	38.865	ng/ul	-0.01
81) Pyrene-d10	19.674	212	2089445	42.172	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.739	264	1604585	40.254	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.157	94	49276	4.229	ng/ul#	91
12) 2-Chlorophenol	7.534	128	13885	1.569	ng/ul	94
22) Nitrobenzene	9.181	77	2274486	217.176	ng/ul#	85
32) 4-Chloroaniline	10.934	127	41026	3.110	ng/ul	97
39) 1,2,4,5-Tetrachloroben...	12.798	216	58827	4.656	ng/ul#	94
75) 1,2,3,4-Tetrachloroben...	13.404	216	71635	5.382	ng/ul#	85

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

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