

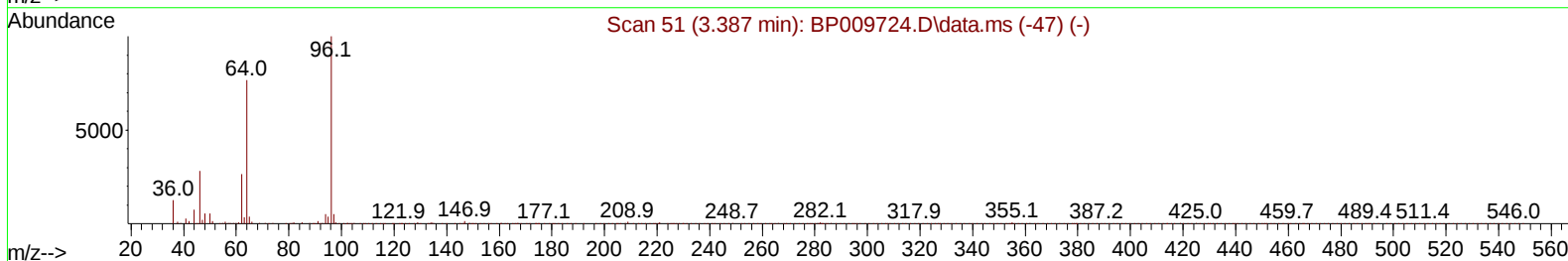
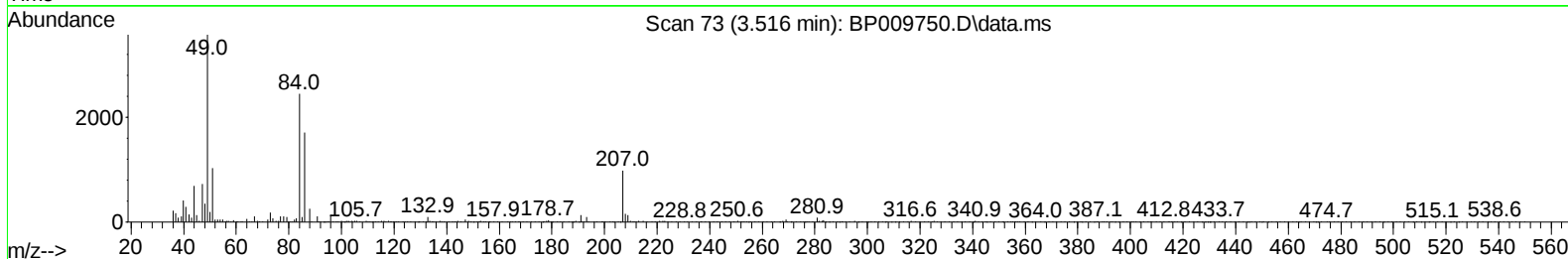
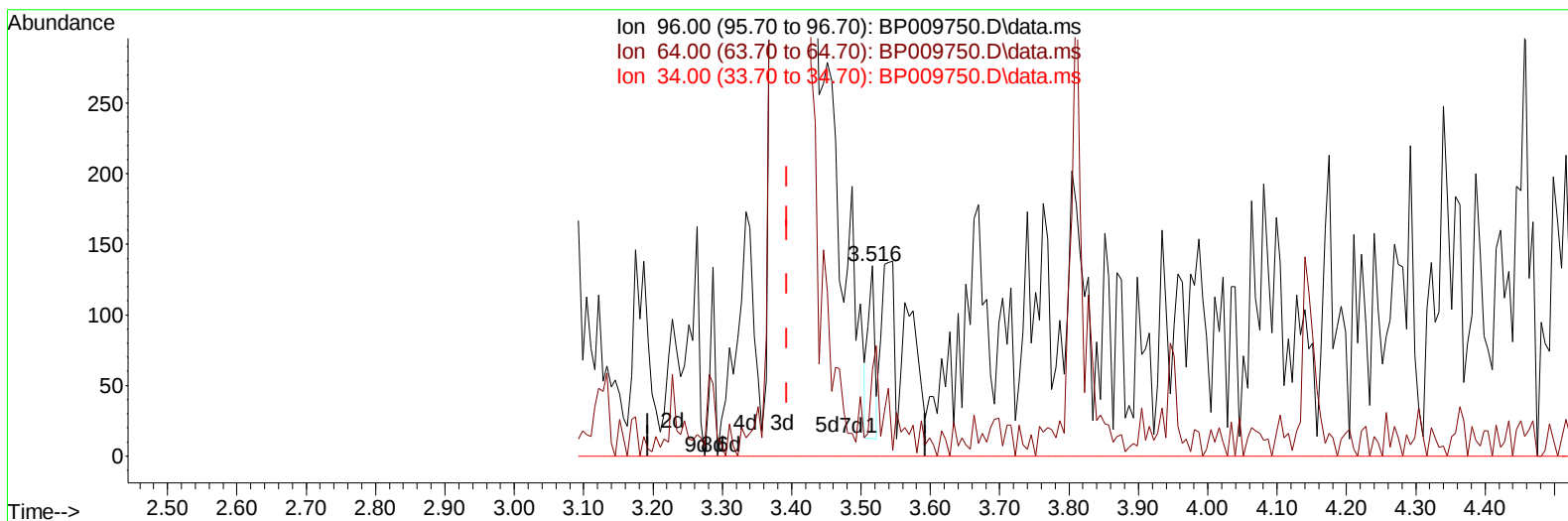
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009750.D
 Acq On : 11 Apr 2022 14:15
 Operator : CG/JU
 Sample : N2338-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J93

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:07:40 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 04/12/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009750.D\data.ms

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

3.516min (+ 0.123) 0.03 ng/uL

response 82

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.19
34.00	0.00	0.00
0.00	0.00	0.00

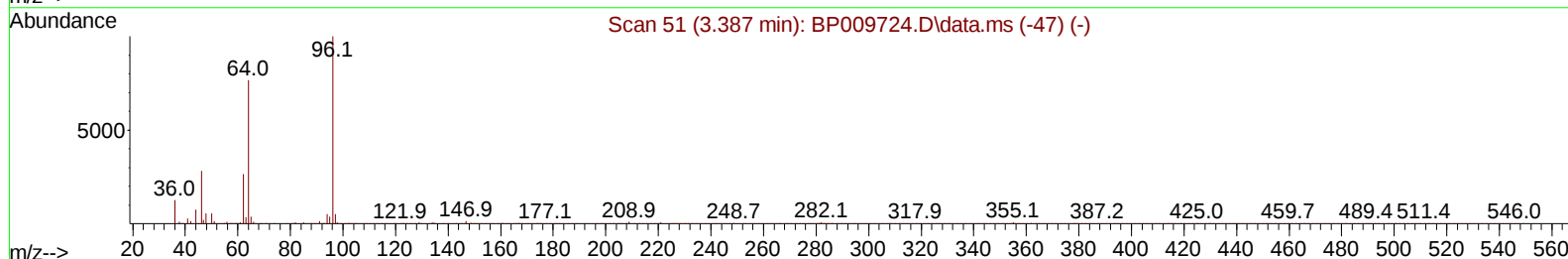
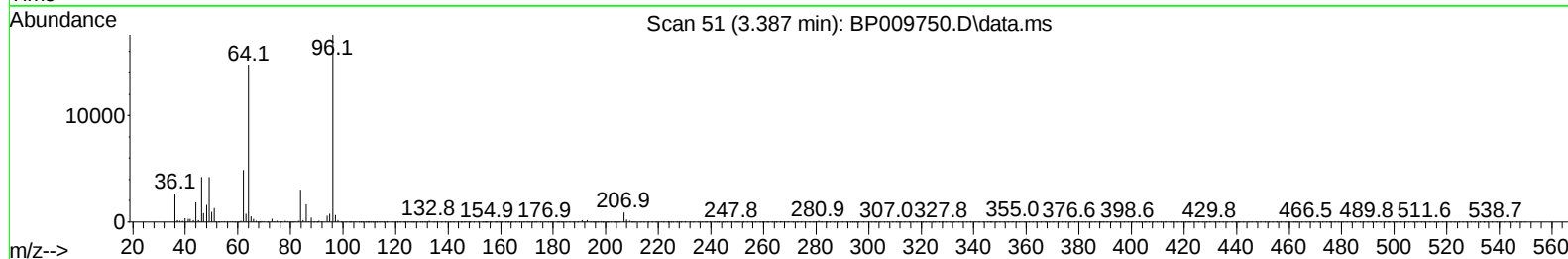
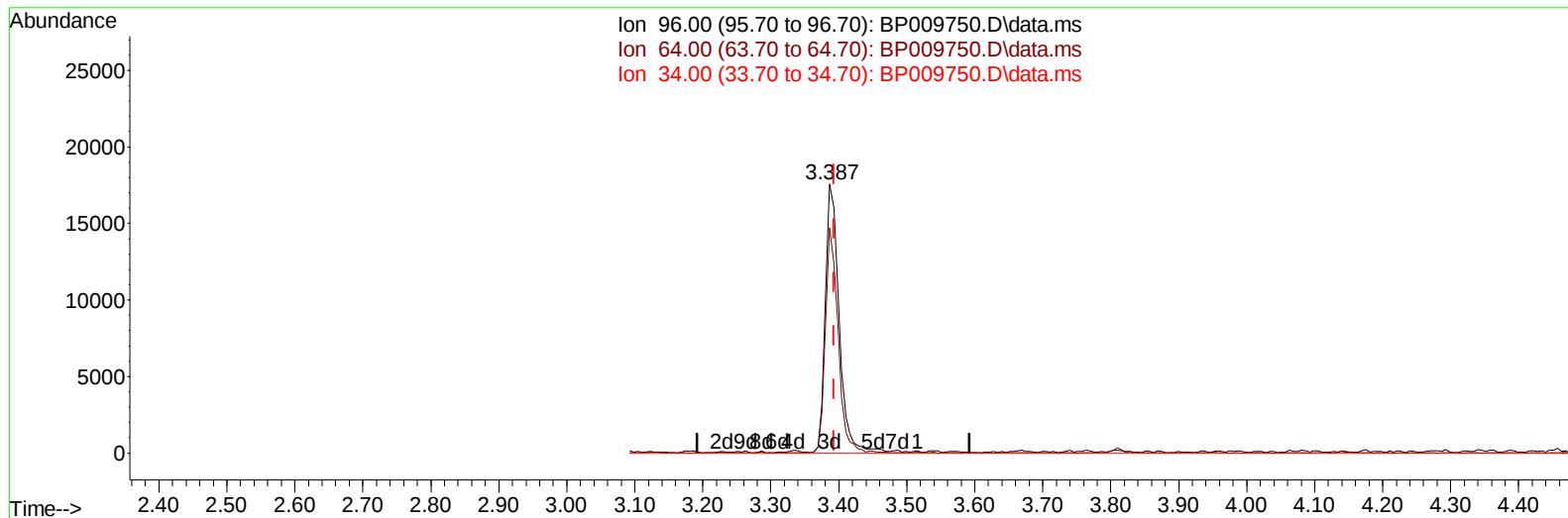
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(3) 1,4-Dioxane-d8 (S)

3.387min (-0.006) 8.07 ng/uL m

response 24859

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.70#
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	137509	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.781	136	598474	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	413941	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	918771	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240	940255	20.000	ng/ul	-0.01
88) Perylene-d12	23.909	264	838979	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	24859m	8.068	ng/uL	0.00
4) Pyridine-d5	3.804	84	93569	11.015	ng/ul	0.00
7) Phenol-d5	7.122	99	103062	8.537	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	284479	39.582	ng/ul	0.00
11) 2-Chlorophenol-d4	7.498	132	277966	30.524	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	194402	19.290	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	188628	43.707	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143	198747	42.004	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.392	165	326829	33.009	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131	483458	34.283	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1389245	42.726	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	1514842	39.324	ng/ul	0.00
54) 4-Nitrophenol-d4	14.774	143	50574	8.576	ng/ul	-0.01
60) Fluorene-d10	15.598	176	1201351	43.920	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	229841	42.228	ng/ul	0.00
73) Anthracene-d10	17.457	188	2063459	46.907	ng/ul	0.00
81) Pyrene-d10	19.680	212	2515532	48.986	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.750	264	2151804	49.250	ng/ul	-0.01

Target Compounds Qvalue

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

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