

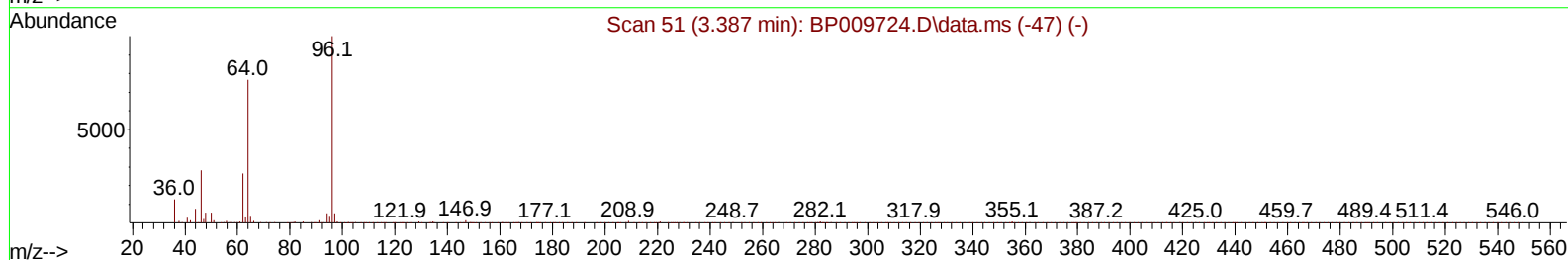
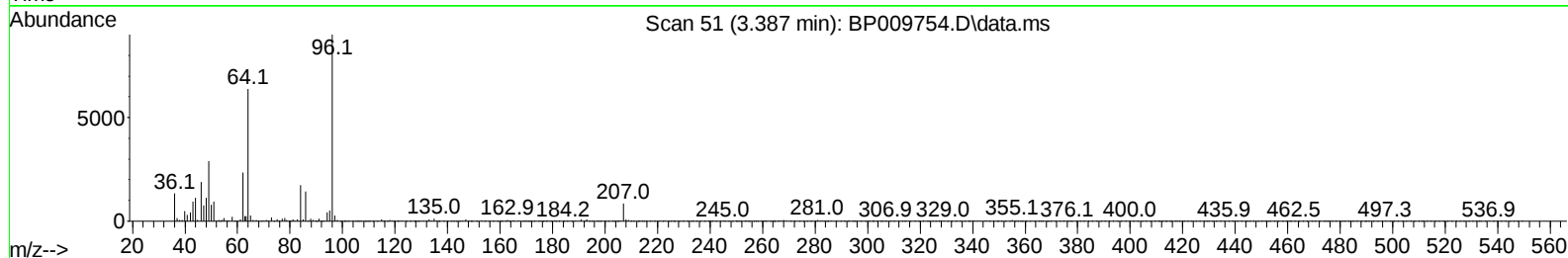
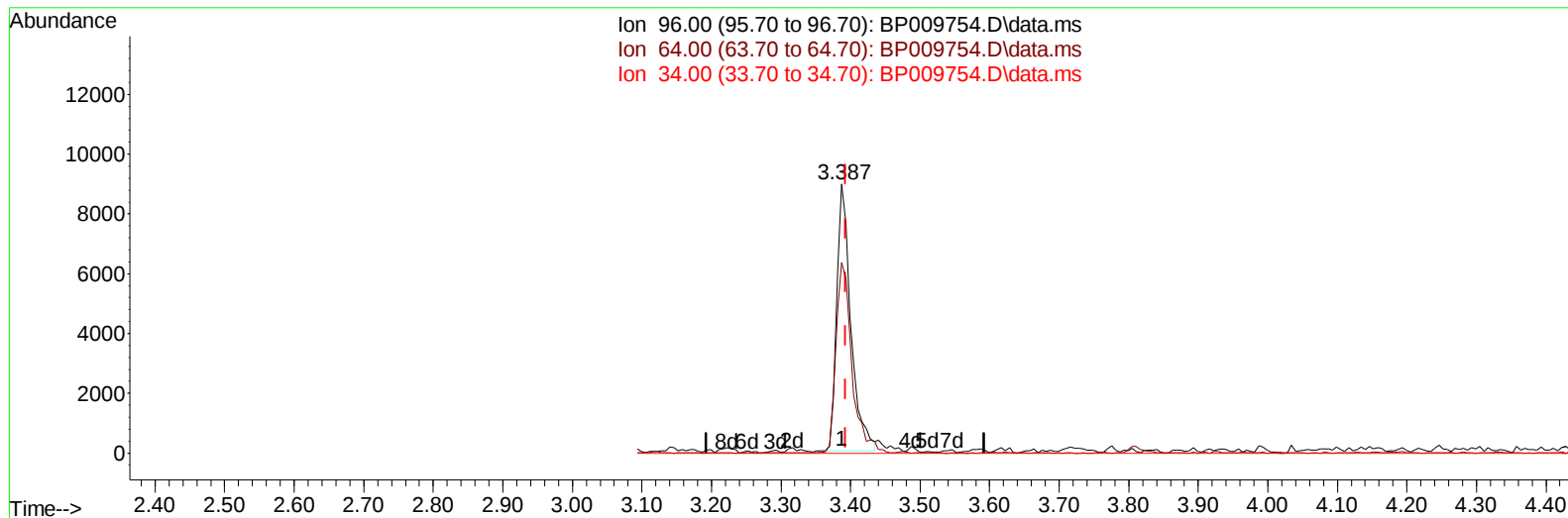
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009754.D
Acq On : 11 Apr 2022 16:30
Operator : CG/JU
Sample : N2338-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J66

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/12/2022
Supervised By :Yogesh Patel 04/13/2022

Quant Time: Apr 12 01:44:23 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



TIC: BP009754.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.387min (-0.006) 5.51 ng/uL****response 12587**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	70.92#
34.00	0.00	0.00
0.00	0.00	0.00

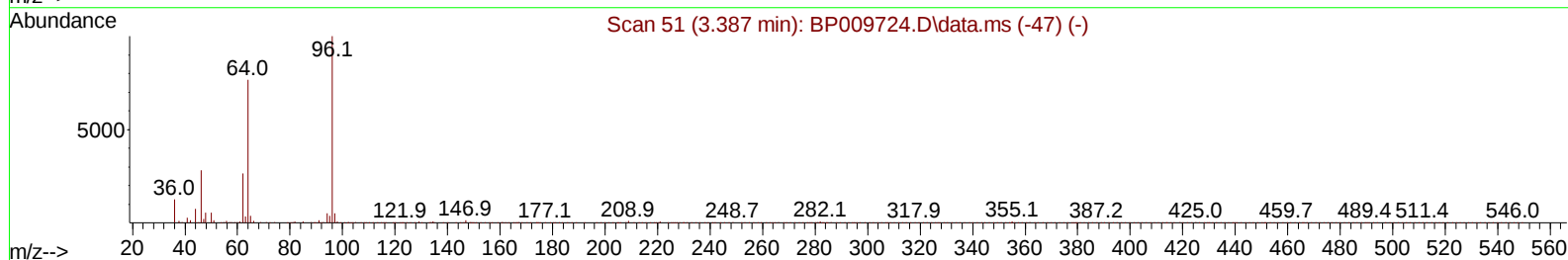
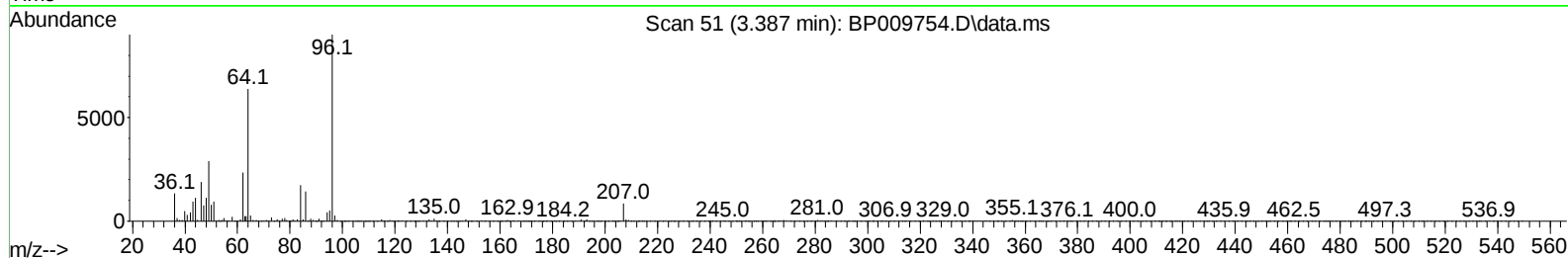
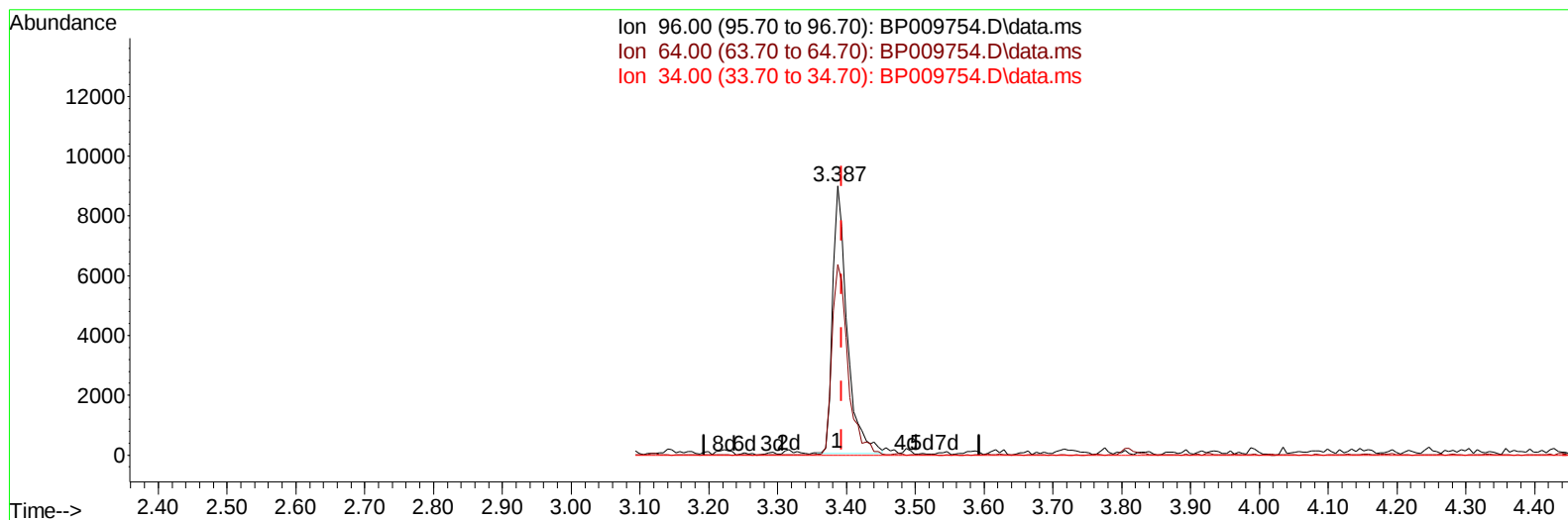
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TIC: BP009754.D\data.ms

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

3.387min (-0.006) 5.65 ng/uL m

response 12896

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	70.92#
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	101871	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.781	136	454233	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	333791	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	750134	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	801939	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264	796930	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	12896m	5.650	ng/uL	0.00
4) Pyridine-d5	3.805	84	96511	15.335	ng/ul	0.00
7) Phenol-d5	7.122	99	80808	9.035	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	230723	43.333	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	220764	32.724	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	157776	21.132	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	150979	46.092	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	156873	43.682	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	269967	35.925	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131	399917	37.364	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1176593	44.874	ng/ul	-0.01
49) Acenaphthylene-d8	14.292	160	1281132	41.243	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	40940	8.610	ng/ul	-0.01
60) Fluorene-d10	15.598	176	1015500	46.040	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	178856	40.248	ng/ul	0.00
73) Anthracene-d10	17.451	188	1723584	47.990	ng/ul	-0.01
81) Pyrene-d10	19.680	212	2132939	48.699	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	2057533	49.578	ng/ul	-0.01

Target Compounds Qvalue

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

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