

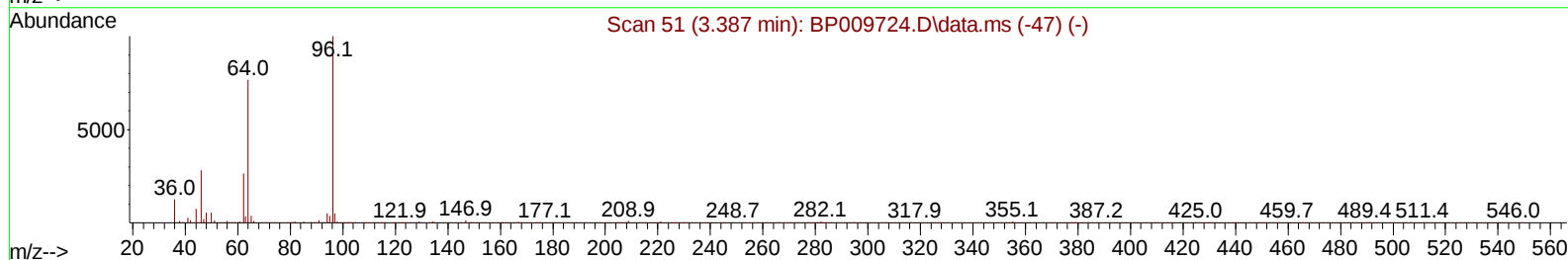
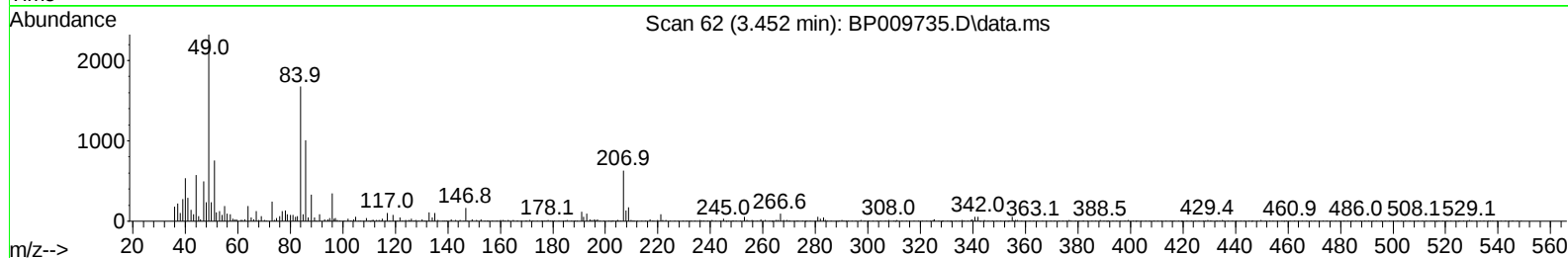
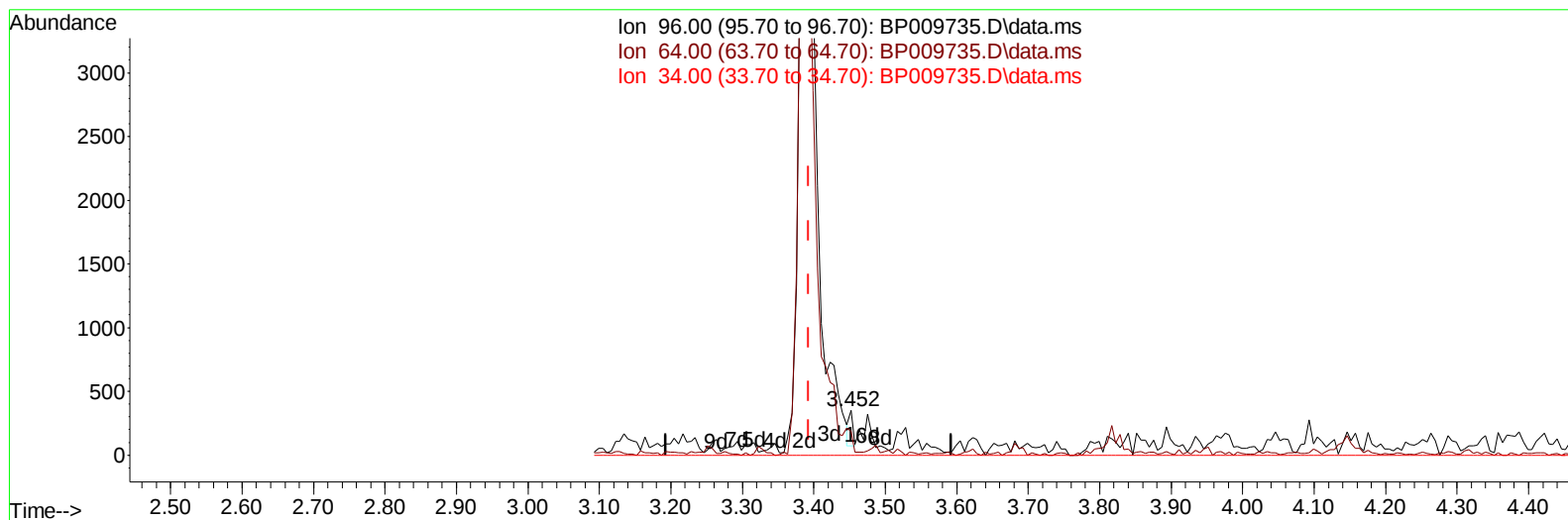
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009735.D
 Acq On : 09 Apr 2022 01:32
 Operator : CG/JU
 Sample : N2266-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YAZ18

Manual IntegrationsAPPROVED

Quant Time: Apr 09 02:18:16 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 04/11/2022
 Supervised By :Yogesh Patel 04/12/2022



TIC: BP009735.D\data.ms

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

3.452min (+ 0.059) 0.05 ng/uL

response 102

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	54.00
34.00	0.00	0.00
0.00	0.00	0.00

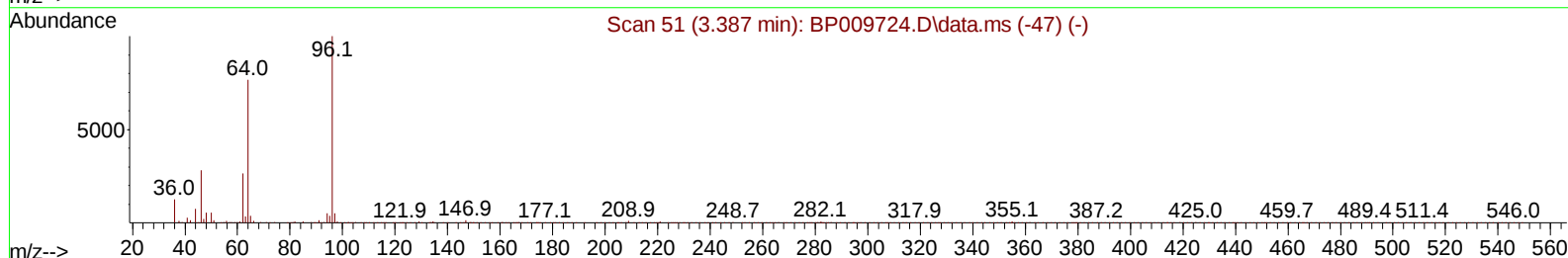
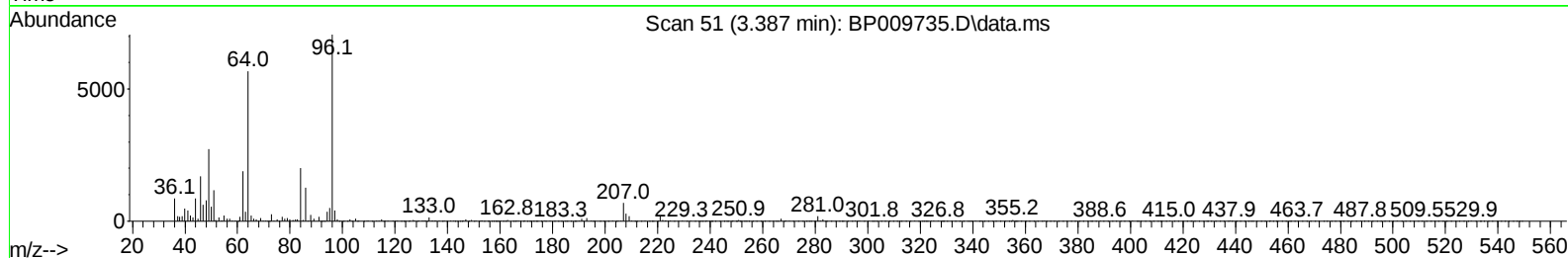
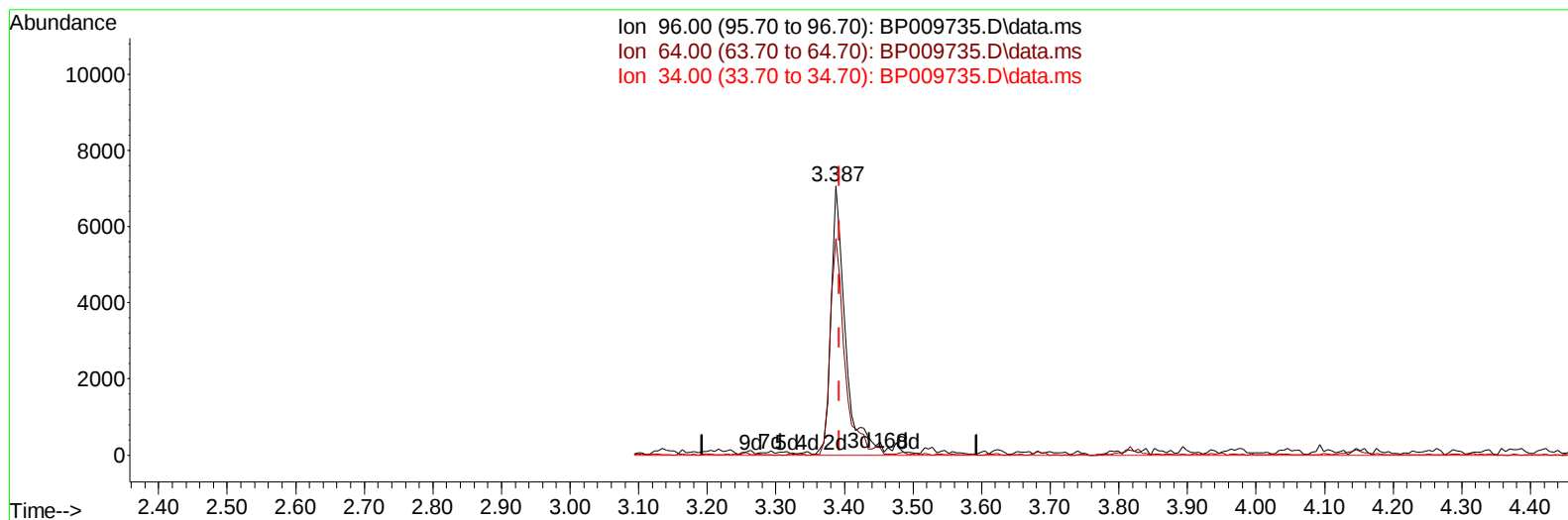
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TIC: BP009735.D\data.ms

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

3.387min (-0.006) 5.45 ng/uL m

response 10282

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.43#
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	7.975	152	84262	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.781	136	382103	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	283874	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	645966	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240	724333	20.000	ng/ul	-0.01
88) Perylene-d12	23.910	264	769674	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	10282m	5.446	ng/uL	0.00
4) Pyridine-d5	3.811	84	47809	9.184	ng/ul	0.00
7) Phenol-d5	7.122	99	65349	8.834	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	184433	41.878	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	175755	31.496	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	133092	21.552	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128	124473	45.173	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143	131839	43.642	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	227590	36.002	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.910	131	290646	32.281	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	998928	44.798	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	1078767	40.835	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	38103	9.422	ng/ul	-0.01
60) Fluorene-d10	15.598	176	845631	45.080	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	155361	40.599	ng/ul	0.00
73) Anthracene-d10	17.457	188	1428137	46.176	ng/ul	0.00
81) Pyrene-d10	19.680	212	1797288	45.432	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	1926114	48.054	ng/ul	-0.01

Target Compounds Qvalue

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

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