

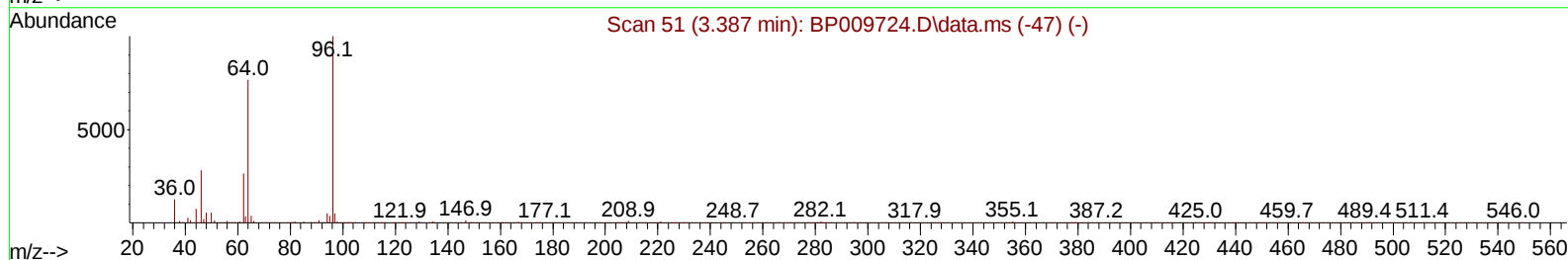
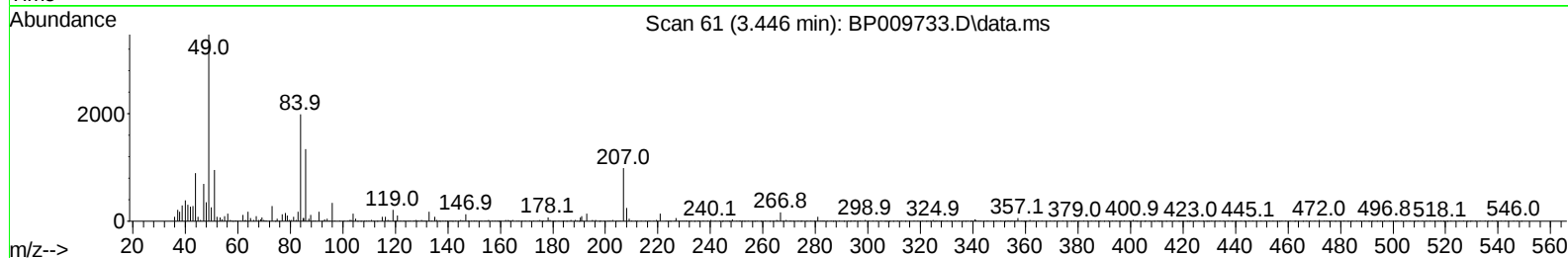
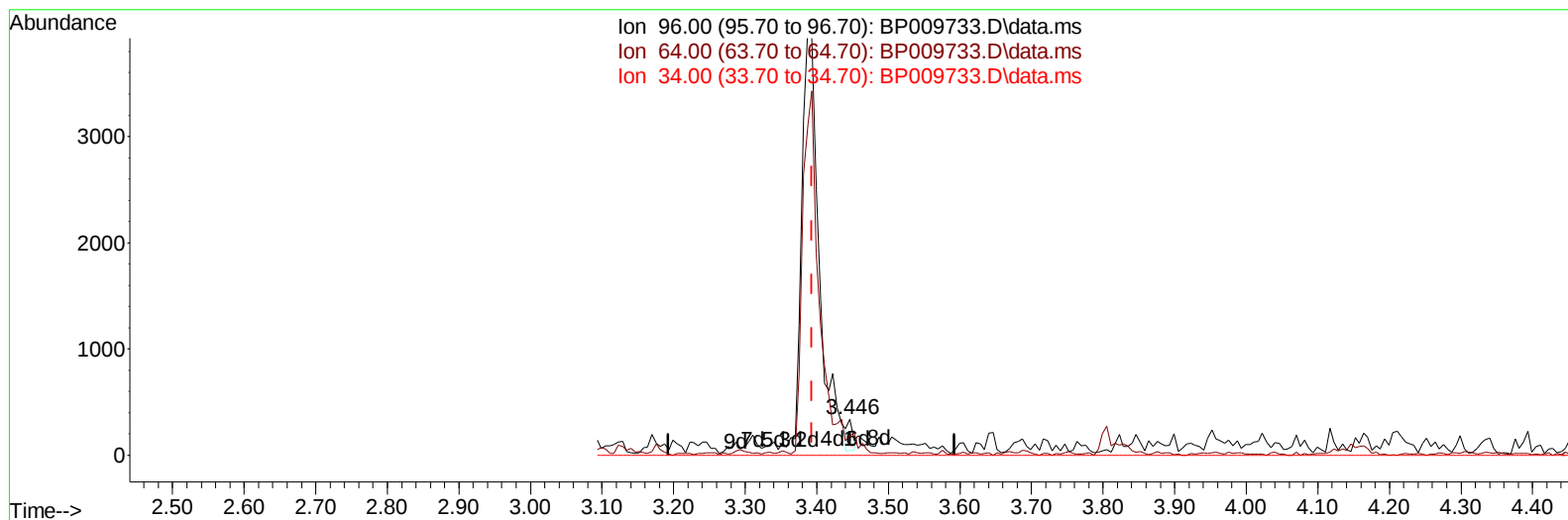
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009733.D
Acq On : 09 Apr 2022 00:24
Operator : CG/JU
Sample : N2182-13
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0R20

Manual IntegrationsAPPROVED

Quant Time: Apr 09 01:17:03 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: BP009733.D\data.ms

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

3.446min (+ 0.053) 0.10 ng/uL

response 137

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	53.08
34.00	0.00	0.00
0.00	0.00	0.00

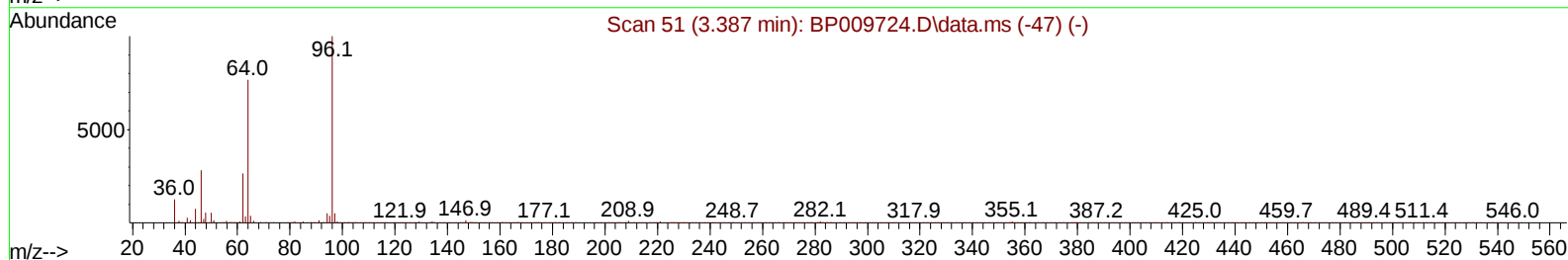
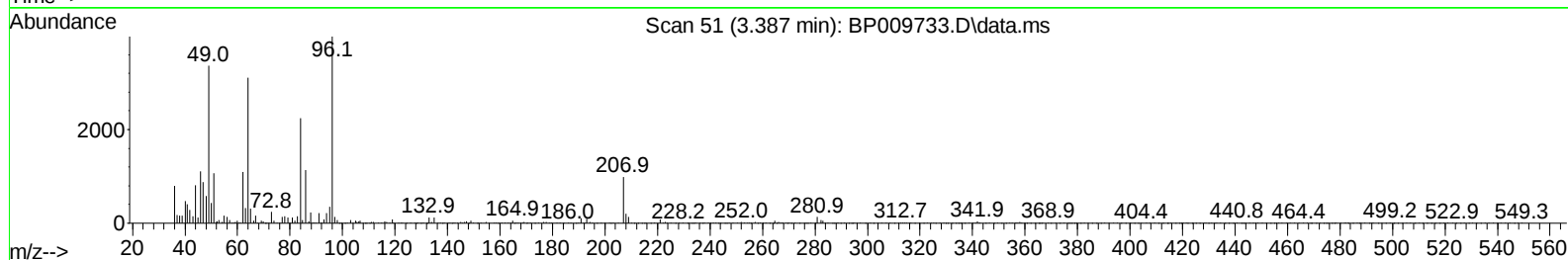
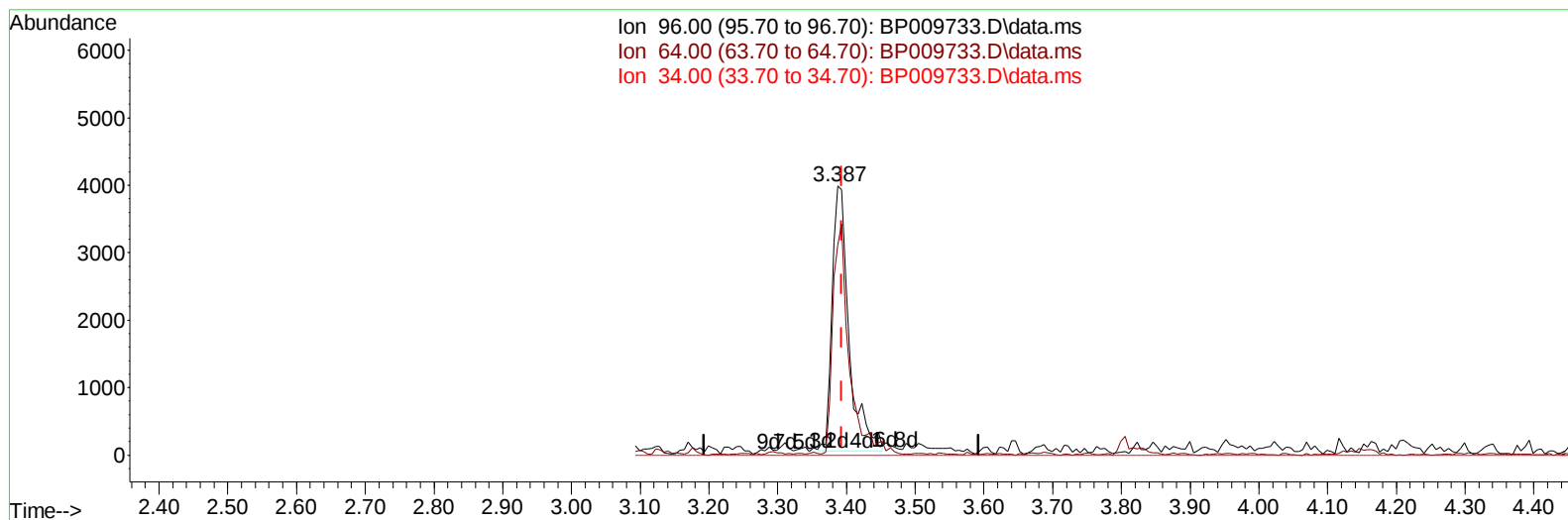
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TIC: BP009733.D\data.ms

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

3.387min (-0.006) 5.08 ng/uL m

response 6756

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.08#
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	59337	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	263569	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	194897	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	451118	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240	526839	20.000	ng/ul	#-0.01
88) Perylene-d12	23.910	264	582301	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	6756m	5.081	ng/uL	0.00
4) Pyridine-d5	3.805	84	98266	26.807	ng/ul	0.00
7) Phenol-d5	7.122	99	141979	27.254	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	90163	29.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	110814	28.200	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	128843	29.628	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	59365	31.234	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.858	143	63888	30.659	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	124930	28.650	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	188112	30.289	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	496268	32.416	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	544705	30.032	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	68071	24.517	ng/ul	-0.01
60) Fluorene-d10	15.598	176	423996	32.922	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	47639	17.826	ng/ul	0.00
73) Anthracene-d10	17.451	188	696369	32.241	ng/ul	-0.01
81) Pyrene-d10	19.681	212	912503	31.713	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	1003267	33.085	ng/ul	-0.01

Target Compounds Qvalue

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

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