

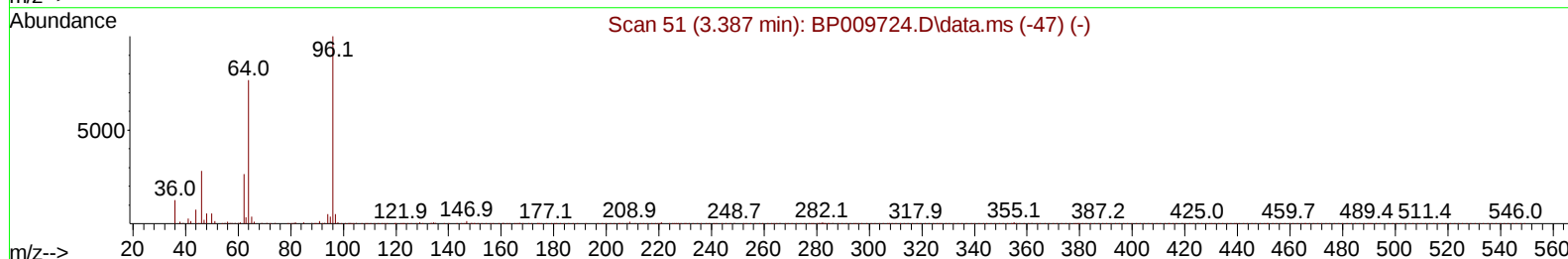
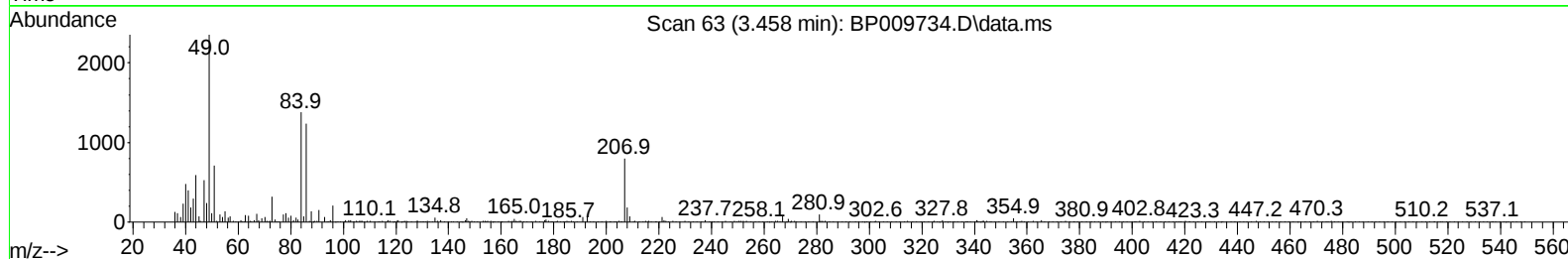
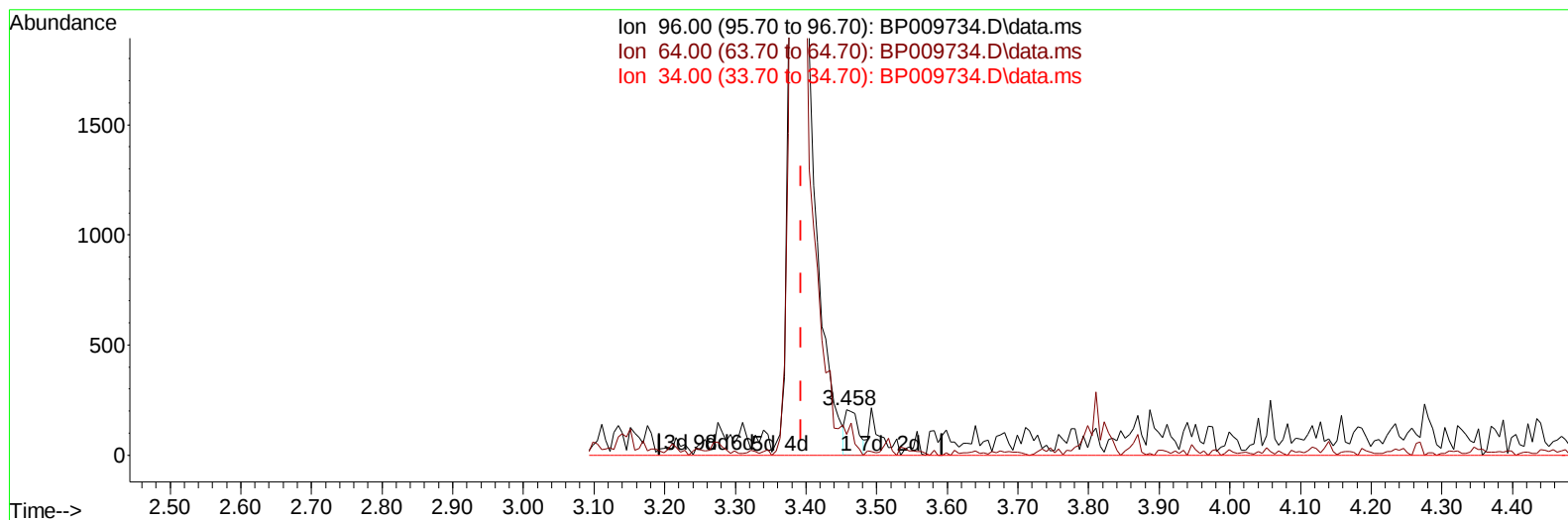
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009734.D
 Acq On : 09 Apr 2022 00:58
 Operator : CG/JU
 Sample : N2182-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R11

Manual IntegrationsAPPROVED

Quant Time: Apr 09 02:17:42 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: BP009734.D\data.ms

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

3.458min (+ 0.065) 0.14 ng/uL

response 265

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.41
34.00	0.00	0.00
0.00	0.00	0.00

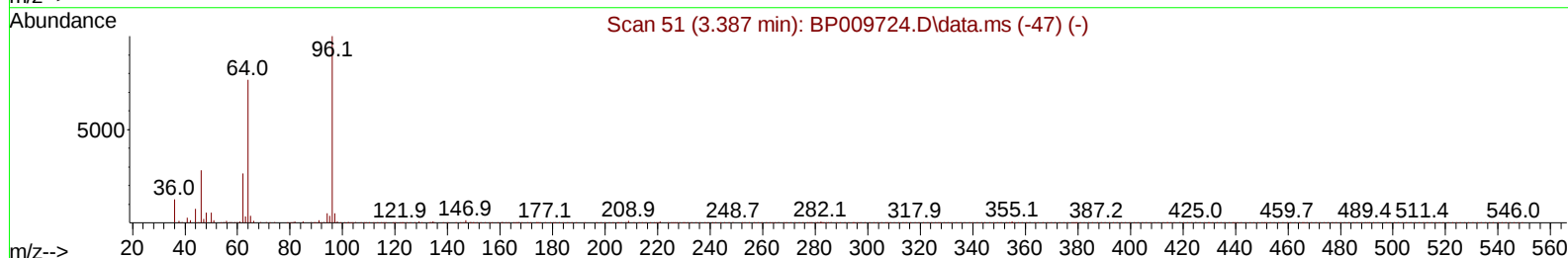
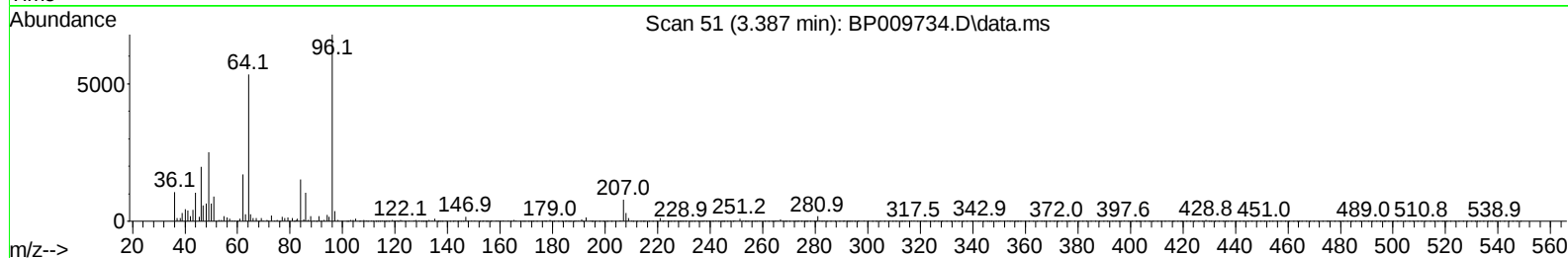
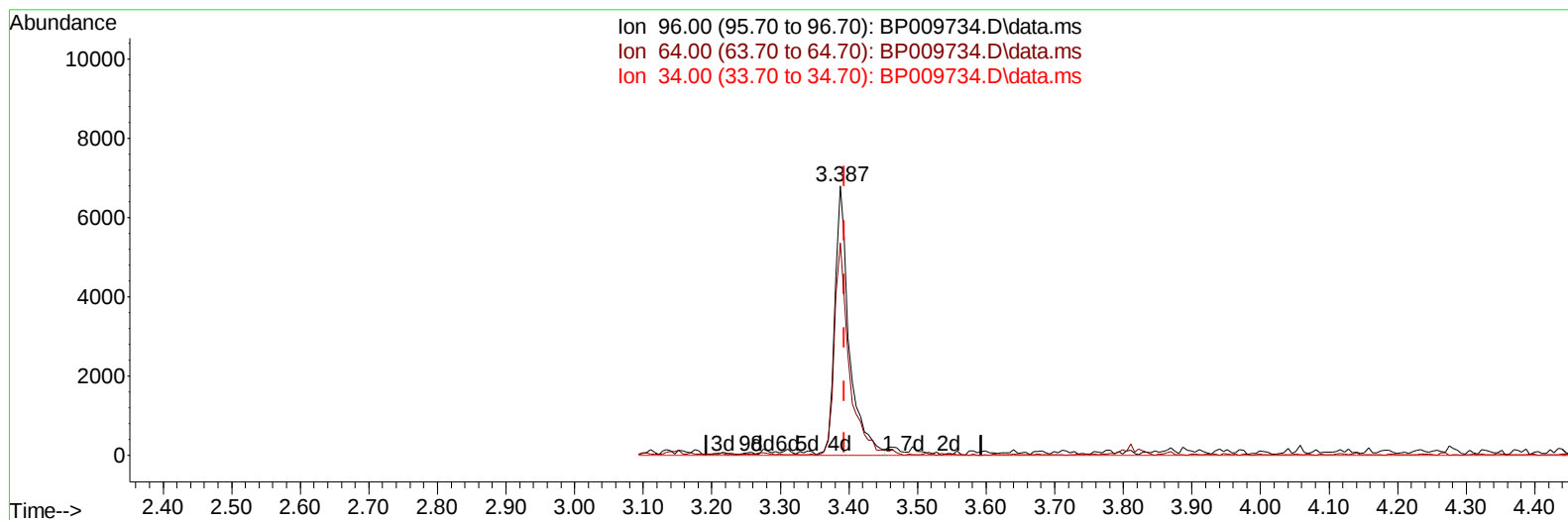
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(3) 1,4-Dioxane-d8 (S)

3.387min (-0.006) 5.09 ng/uL m

response 9930

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.80#
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	87137	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.781	136	383859	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	276617	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	638236	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	719975	20.000	ng/ul	#-0.01
88) Perylene-d12	23.910	264	790217	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	9930m	5.086	ng/uL	0.00
4) Pyridine-d5	3.805	84	101702	18.893	ng/ul	0.00
7) Phenol-d5	7.122	99	188599	24.653	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	125464	27.548	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	150553	26.090	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	155737	24.386	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	78574	28.385	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143	87726	28.906	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	164731	25.940	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	222680	24.619	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	632677	29.117	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	702580	27.293	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	100377	25.472	ng/ul	-0.01
60) Fluorene-d10	15.598	176	534450	29.239	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	79948	21.145	ng/ul	0.00
73) Anthracene-d10	17.451	188	861598	28.195	ng/ul	-0.01
81) Pyrene-d10	19.680	212	1127822	28.682	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	1193622	29.005	ng/ul	-0.01

Target Compounds Qvalue

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

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