

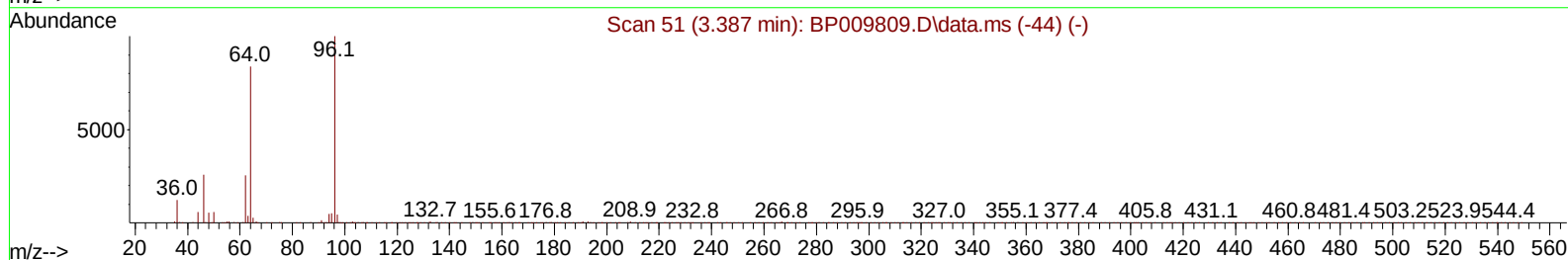
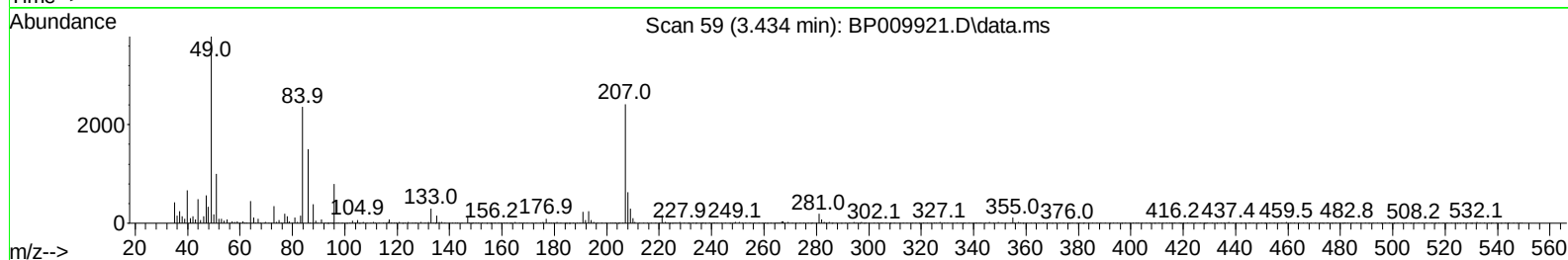
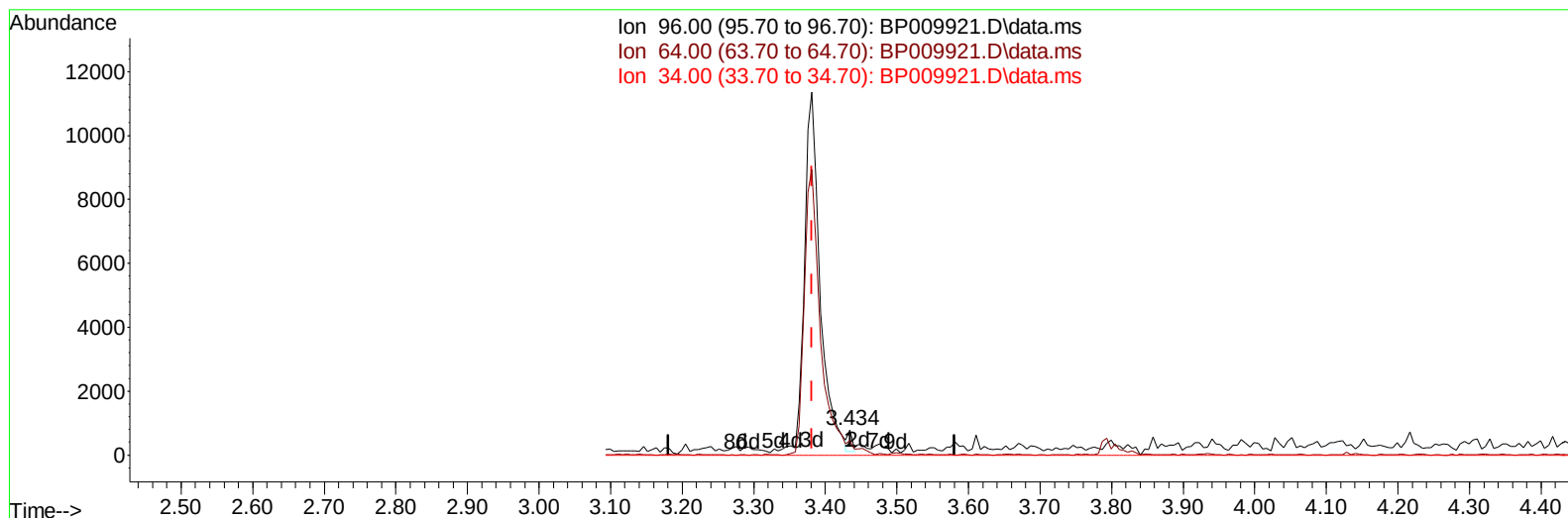
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009921.D
 Acq On : 18 Apr 2022 18:22
 Operator : CG/JU
 Sample : PB144145BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK145

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/22/2022

Quant Time: Apr 19 00:57:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration



TIC: BP009921.D\data.ms

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

3.434min (+ 0.053) 0.09 ng/uL

response	285	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.37
34.00	0.00	0.00
0.00	0.00	0.00

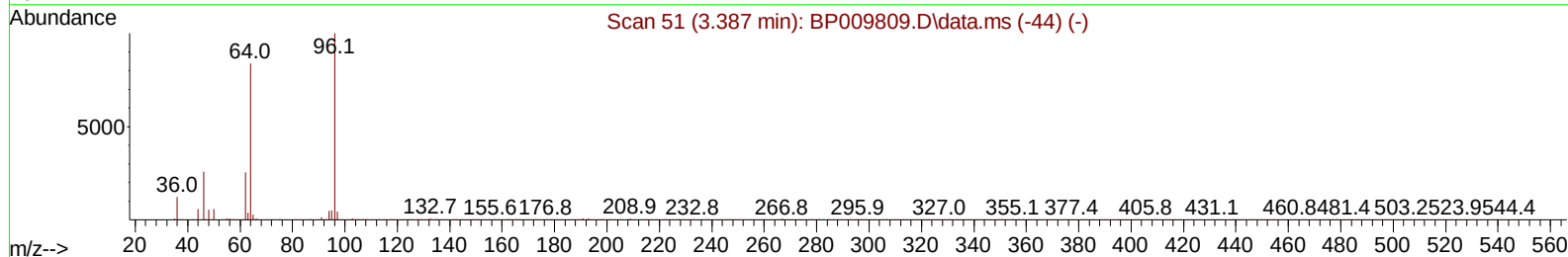
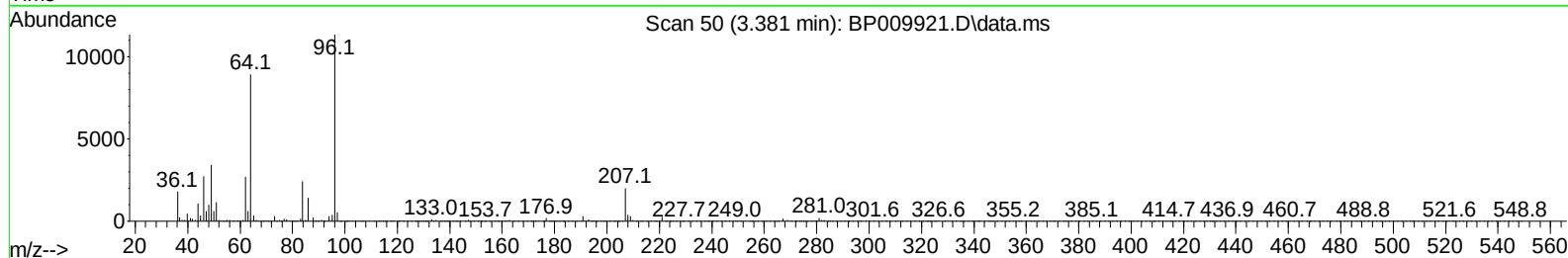
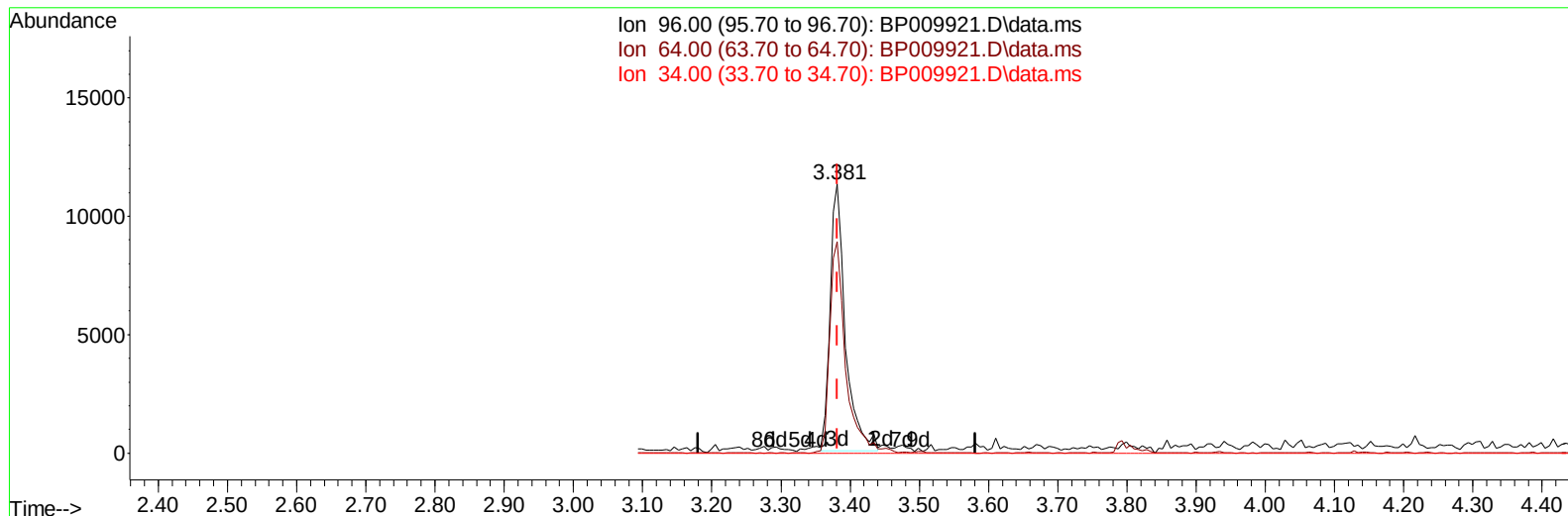
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TIC: BP009921.D\data.ms

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

3.381min (0.000) 5.48 ng/uL m

response 17171

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.70#
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	7.952	152	139834	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	652052	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.587	164	488287	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	1074387	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	1074894	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	956531	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	17171m	5.480	ng/uL	0.00
4) Pyridine-d5	3.793	84	234212	27.112	ng/ul	0.00
7) Phenol-d5	7.105	99	317858	25.891	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	219692	30.059	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	259446	28.017	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	284154	27.727	ng/ul	-0.01
21) Nitrobenzene-d5	9.110	128	142498	30.305	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	155861	30.234	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	273429	25.347	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	425827	27.715	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1077271	28.086	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1182942	26.033	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	157154	22.593	ng/ul	0.00
60) Fluorene-d10	15.575	176	915990	28.389	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	149795	23.535	ng/ul	0.00
73) Anthracene-d10	17.433	188	1486351	28.894	ng/ul	0.00
81) Pyrene-d10	19.663	212	1797547	30.620	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1467905	29.468	ng/ul	0.00

Target Compounds Qvalue

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

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