

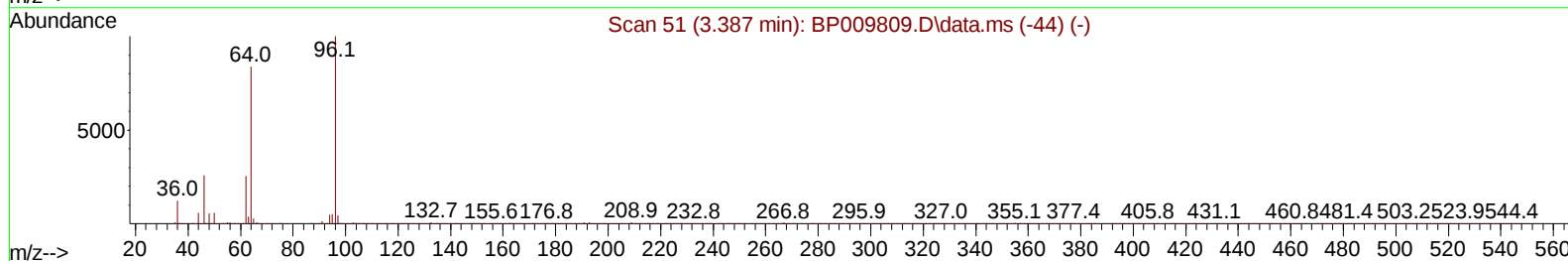
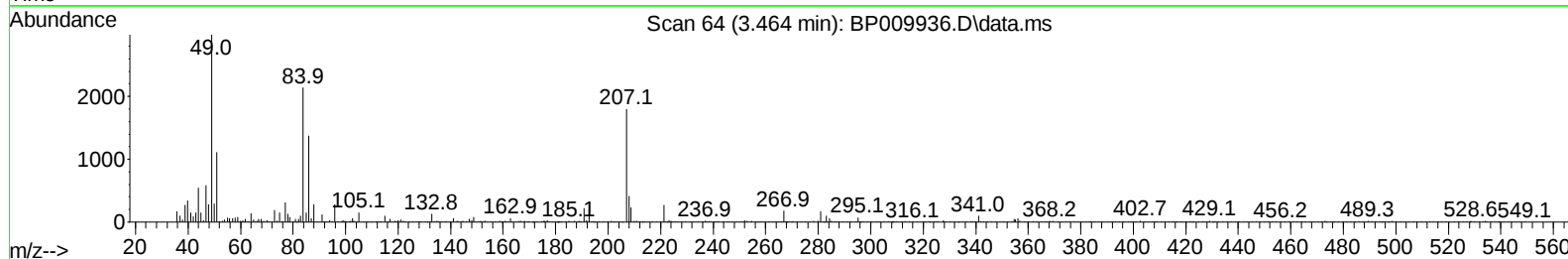
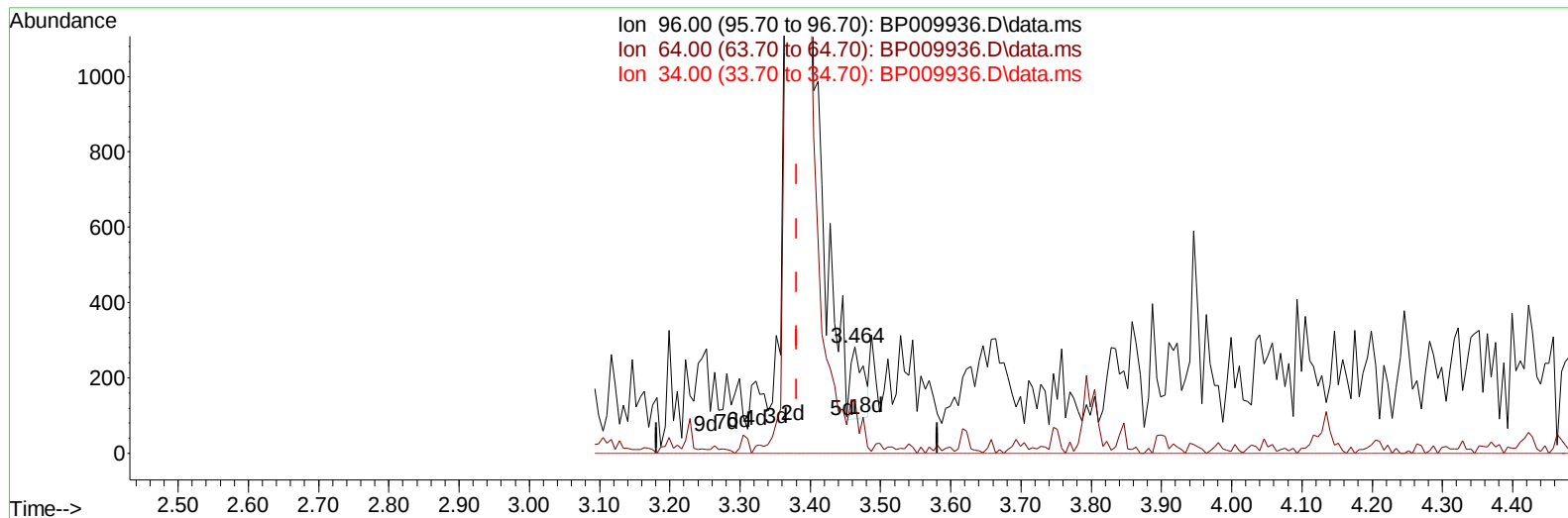
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009936.D
 Acq On : 19 Apr 2022 02:48
 Operator : CG/JU
 Sample : N2421-17
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J77

Manual IntegrationsAPPROVED

Quant Time: Apr 19 03:22:37 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

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



TIC: BP009936.D\data.ms

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

3.464min (+ 0.082) 0.08 ng/uL

response 241

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	50.71
34.00	0.00	0.00
0.00	0.00	0.00

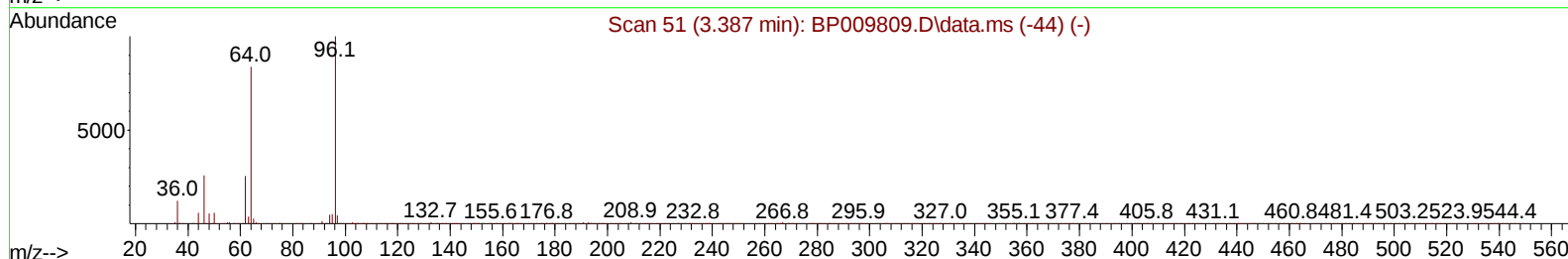
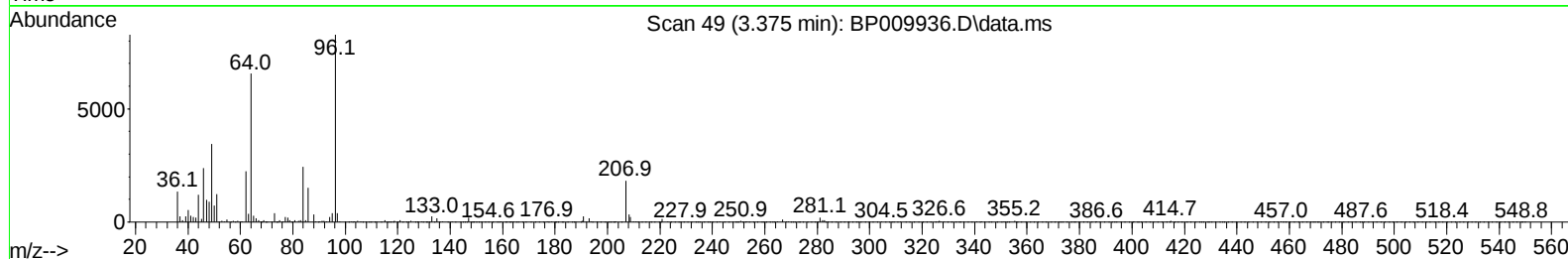
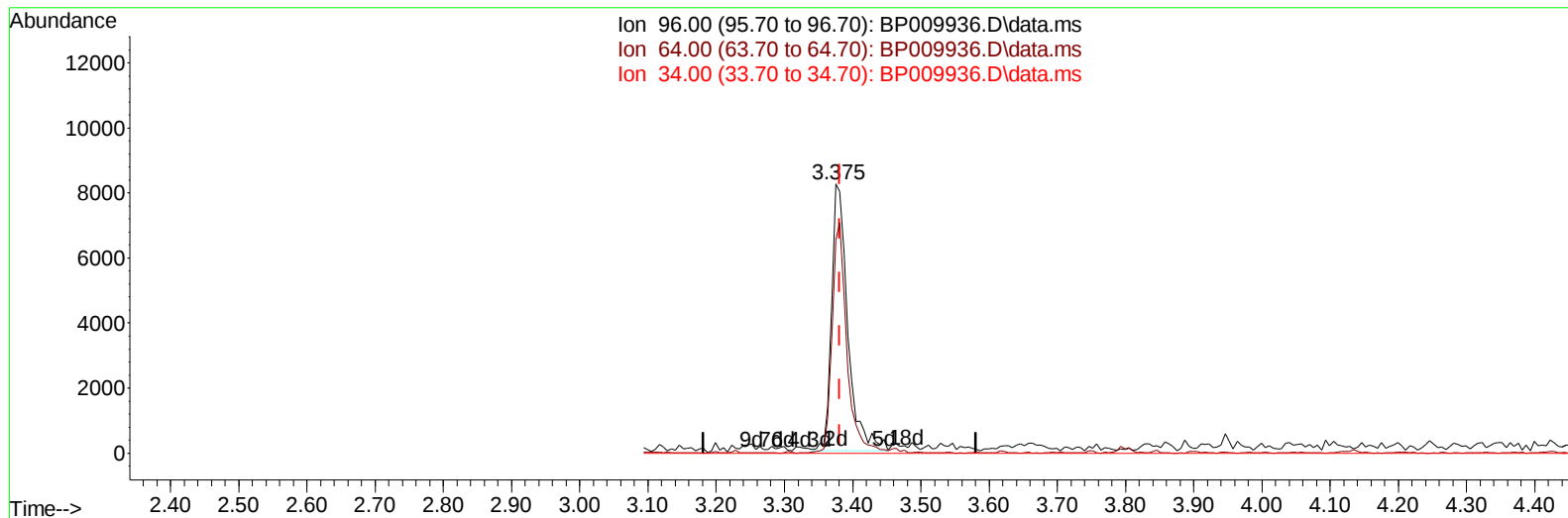
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(3) 1,4-Dioxane-d8 (S)

3.375min (-0.006) 4.76 ng/uL m

response 13604

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.31#
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	127680	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.757	136	583481	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.581	164	438201	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.334	188	971537	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	931943	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	830214	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	13604m	4.755	ng/uL	0.00
4) Pyridine-d5	3.799	84	73109	9.269	ng/ul	0.00
7) Phenol-d5	7.105	99	82926	7.398	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	217211	32.549	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	212070	25.081	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	161305	17.238	ng/ul	-0.01
21) Nitrobenzene-d5	9.105	128	146819	34.893	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.834	143	156016	33.820	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	265125	27.465	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	357913	26.032	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	1135225	32.980	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	1259516	30.886	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	38307	6.136	ng/ul	0.00
60) Fluorene-d10	15.575	176	1000093	34.538	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	182070	31.635	ng/ul	-0.01
73) Anthracene-d10	17.434	188	1762116	37.882	ng/ul	0.00
81) Pyrene-d10	19.663	212	2138057	42.006	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1734850	40.126	ng/ul	0.00

Target Compounds Qvalue

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

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