

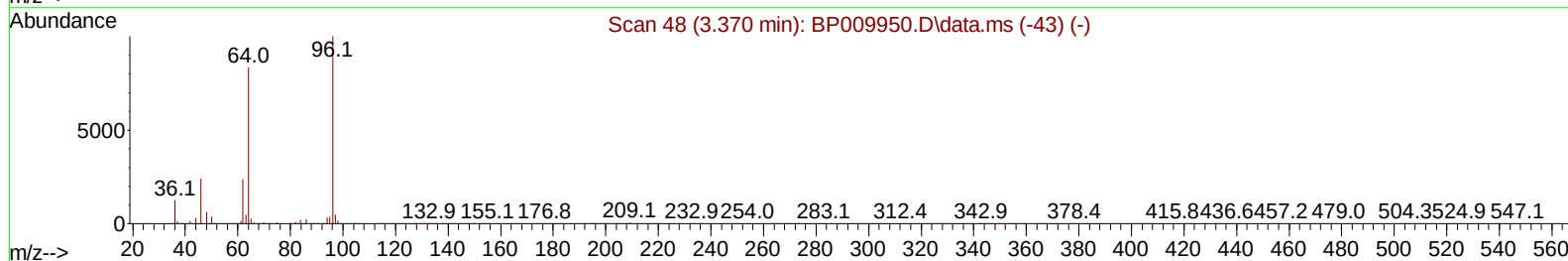
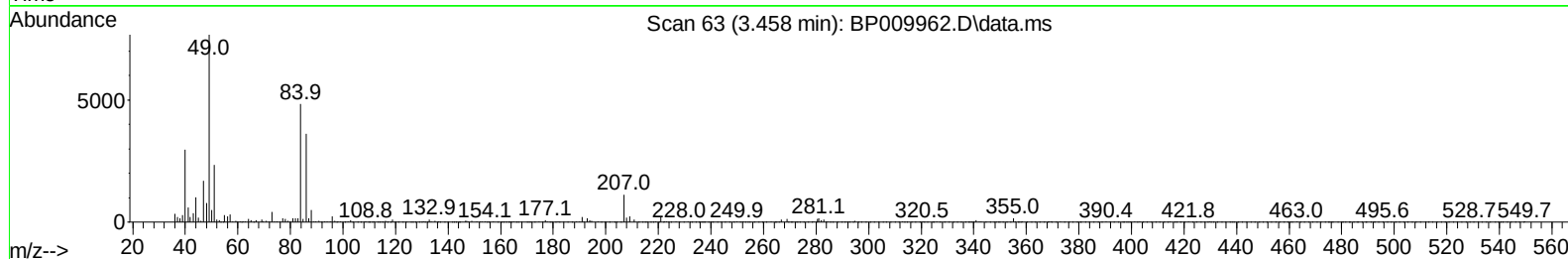
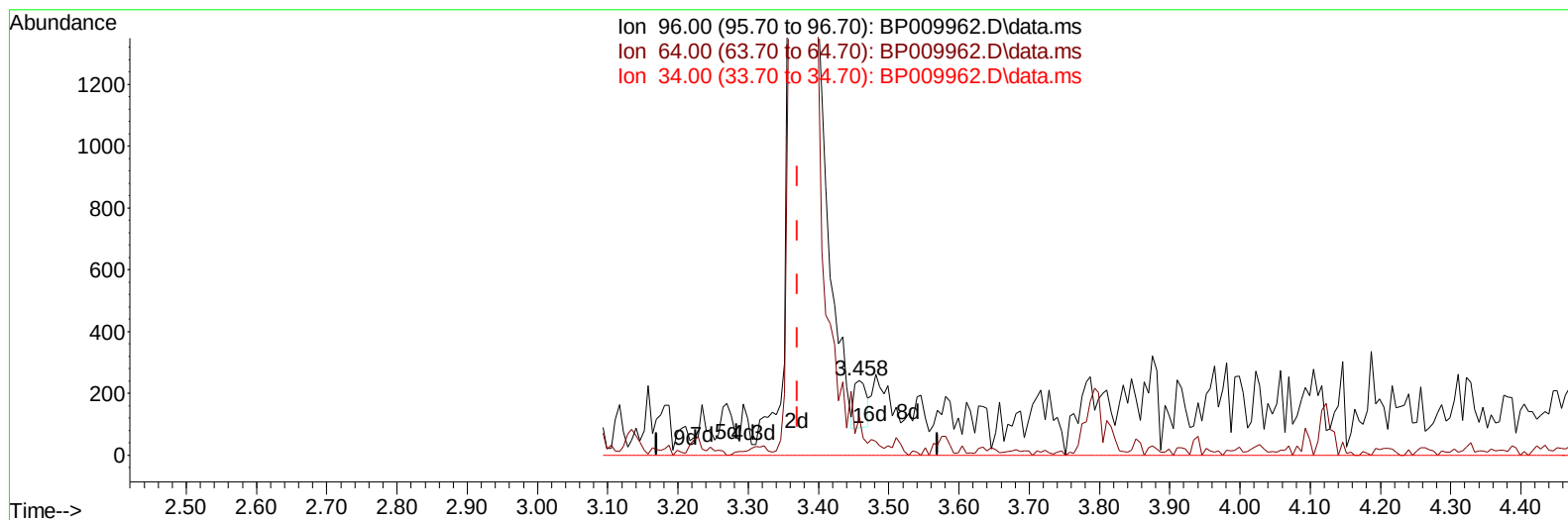
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009962.D
 Acq On : 20 Apr 2022 00:47
 Operator : CG/JU
 Sample : N2388-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFF2

Manual IntegrationsAPPROVED

Quant Time: Apr 20 02:38:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/20/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP009962.D\data.ms

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

3.458min (+ 0.088) 0.05 ng/uL

response 189

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	50.41
34.00	0.00	0.00
0.00	0.00	0.00

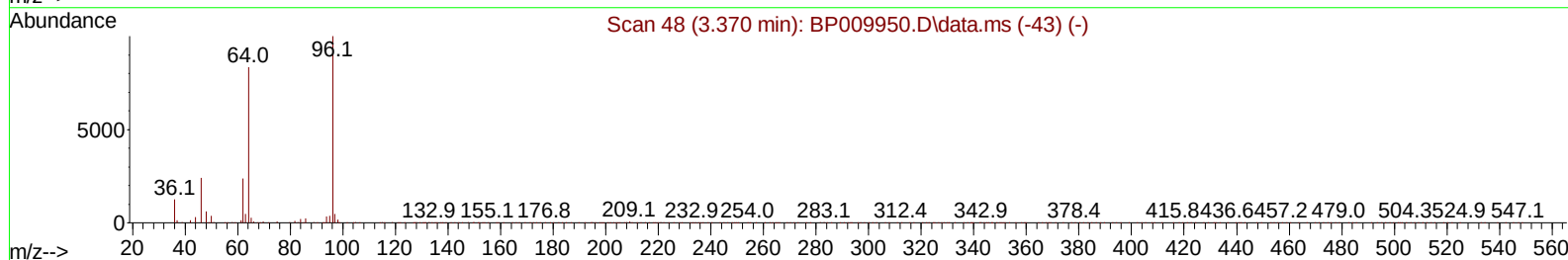
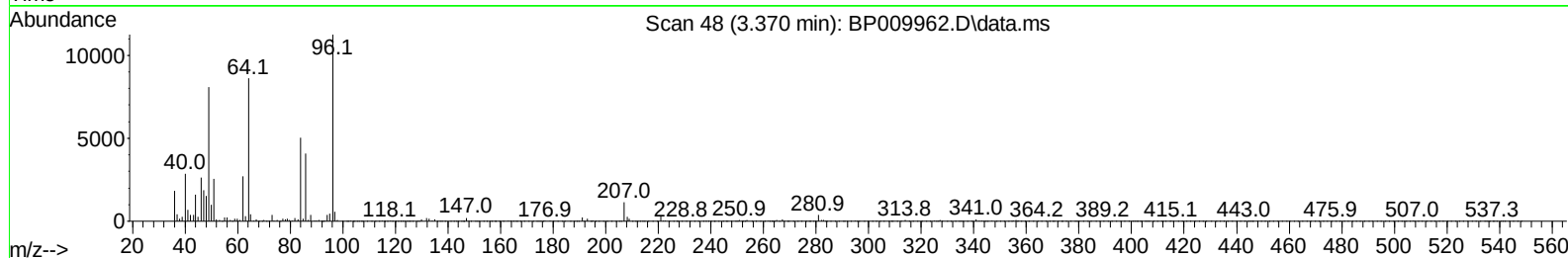
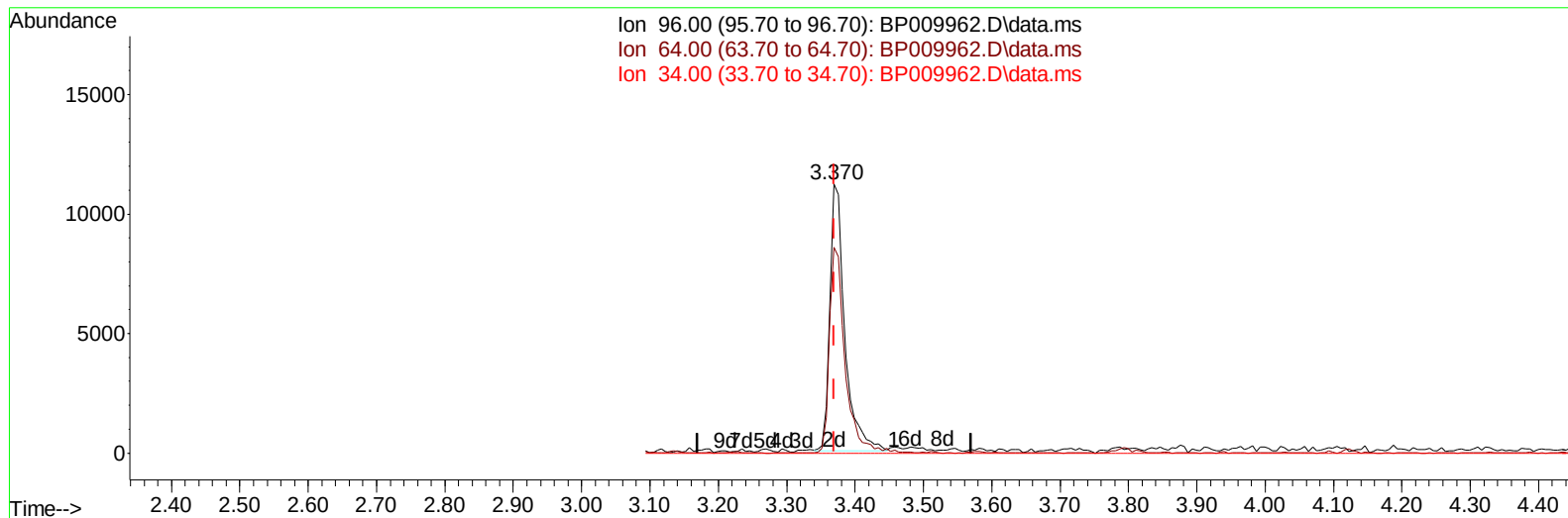
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 4.57 ng/uL m

response 16910

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	76.52#
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	7.946	152	165013	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	752462	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	536714	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1163923	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	1271880	20.000	ng/ul	# 0.00
88) Perylene-d12	23.863	264	1335354	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	16910m	4.573	ng/uL	0.00
4) Pyridine-d5	3.787	84	113691	11.152	ng/ul	0.00
7) Phenol-d5	7.099	99	103544	7.147	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	297172	34.456	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	284479	26.033	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	212767	17.593	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	199011	36.676	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	218189	36.676	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	362781	29.142	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	463148	26.121	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1450323	34.401	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1662086	33.277	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	49983	6.537	ng/ul	0.00
60) Fluorene-d10	15.569	176	1264031	35.640	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	223633	32.434	ng/ul	0.00
73) Anthracene-d10	17.428	188	1992330	35.751	ng/ul	0.00
81) Pyrene-d10	19.657	212	2501136	36.006	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	2528021	36.353	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.128	94	14794	1.020	ng/ul#	94
78) Di-n-butylphthalate	18.281	149	146598	2.094	ng/ul#	96

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

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