

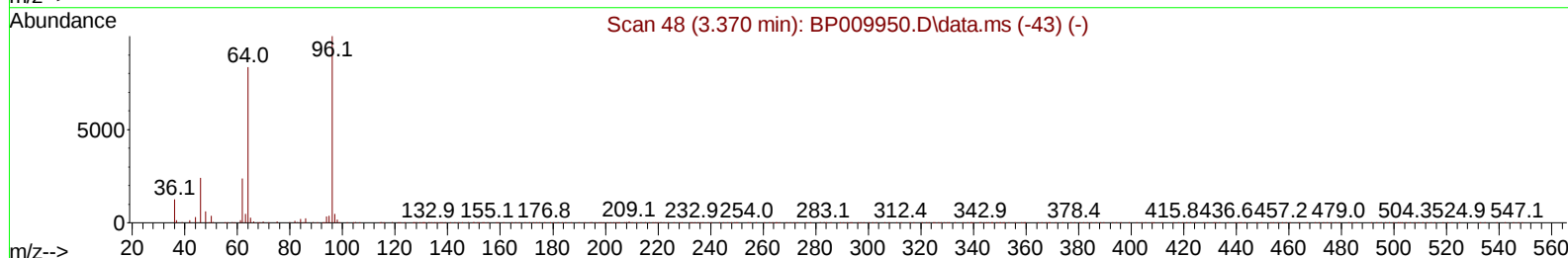
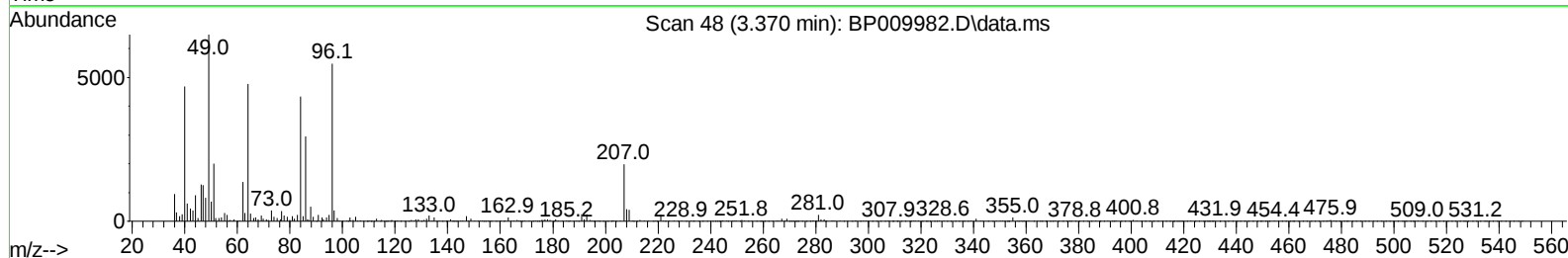
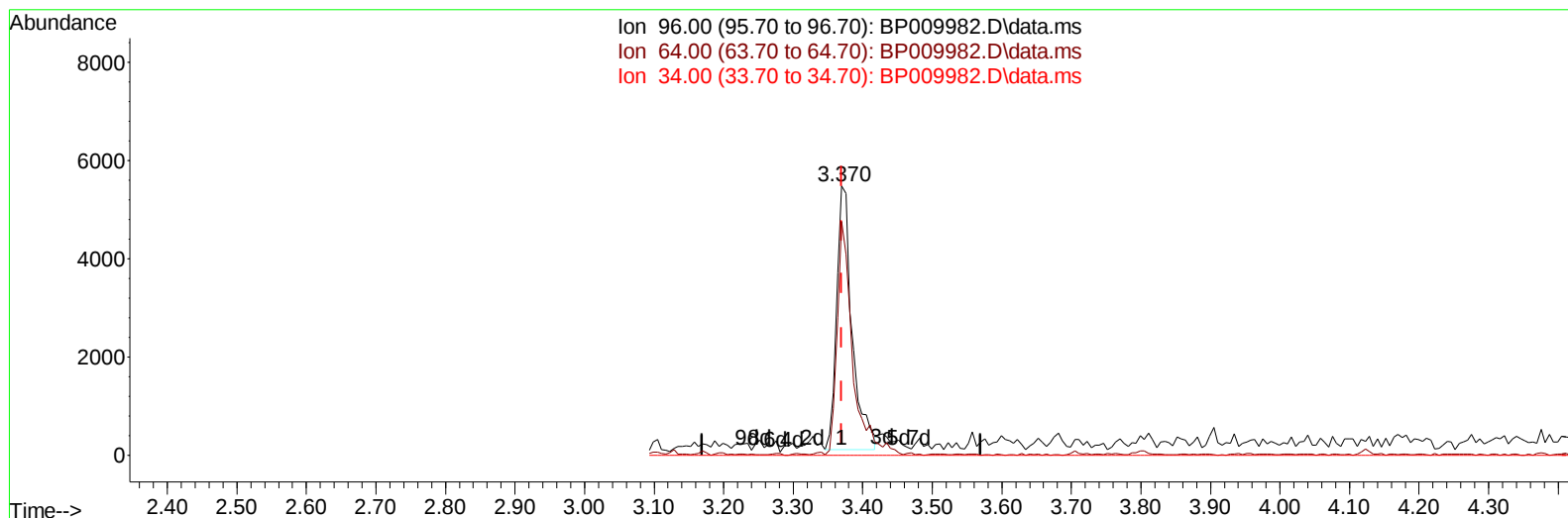
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009982.D
 Acq On : 20 Apr 2022 18:24
 Operator : CG/JU
 Sample : N2374-10
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFF4

Manual IntegrationsAPPROVED

Quant Time: Apr 21 02:13:46 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 04/21/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP009982.D\data.ms

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

3.370min (0.000) 2.60 ng/uL

response 8365

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.98#
34.00	0.00	0.00
0.00	0.00	0.00

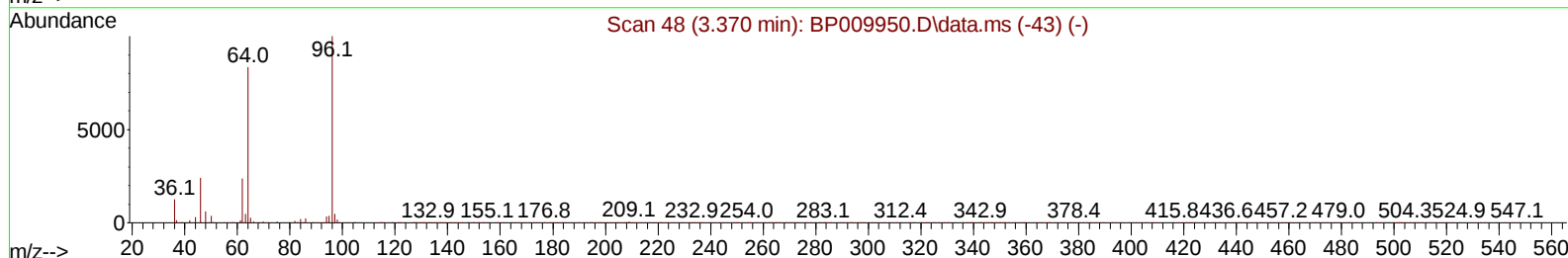
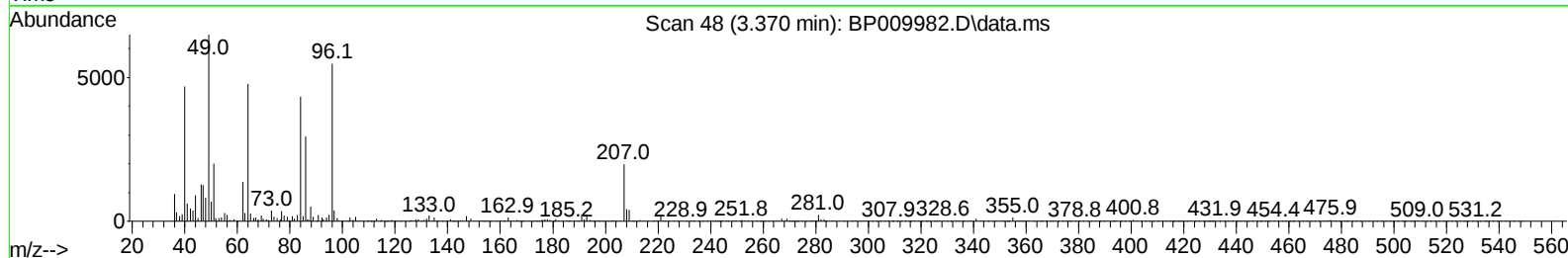
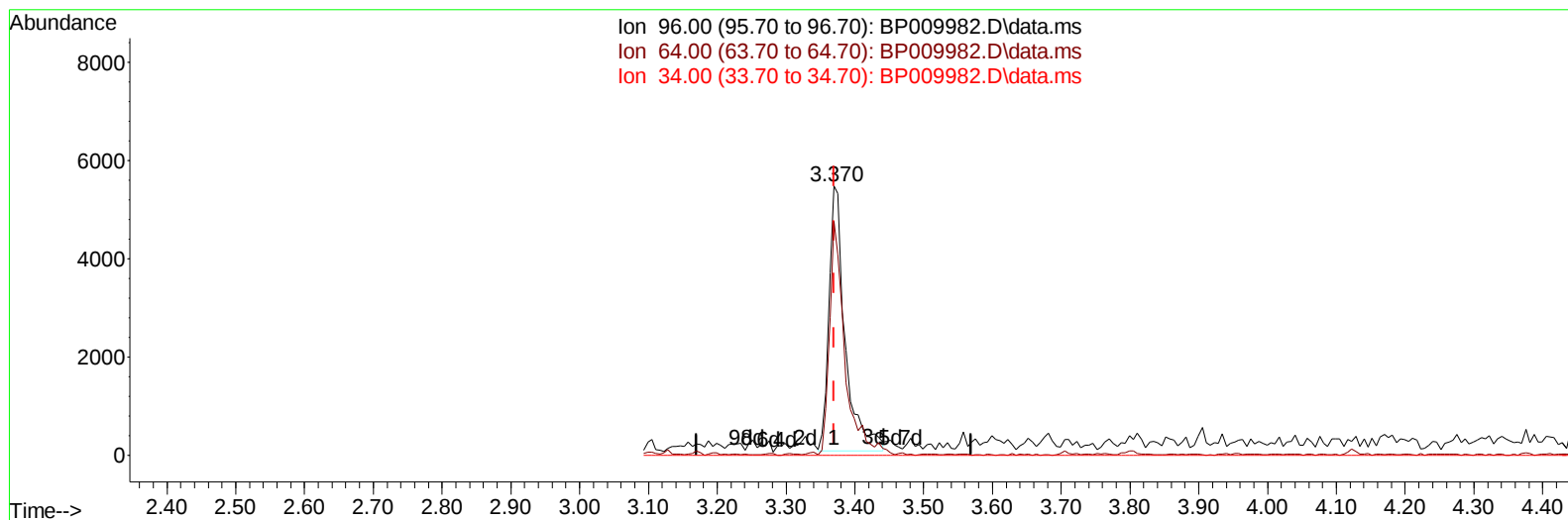
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TIC: BP009982.D\data.ms

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

3.370min (0.000) 2.75 ng/uL m

response 8834

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.98#
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	143389	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	649332	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	458842	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	992892	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	952402	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	837334	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	8834m	2.750	ng/uL	0.00
4) Pyridine-d5	3.799	84	24632	2.781	ng/ul	0.01
7) Phenol-d5	7.105	99	170757	13.564	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	114677	15.302	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	142778	15.036	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	136900	13.027	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	73097	15.611	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	79164	15.420	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	143942	13.399	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	186648	12.199	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	518874	14.396	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	592675	13.880	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	68225	10.437	ng/ul	0.00
60) Fluorene-d10	15.563	176	452536	14.925	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	48750	8.288	ng/ul	0.00
73) Anthracene-d10	17.422	188	738294	15.530	ng/ul	0.00
81) Pyrene-d10	19.651	212	989149	19.016	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	745403	17.094	ng/ul	-0.01

Target Compounds Qvalue

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

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