

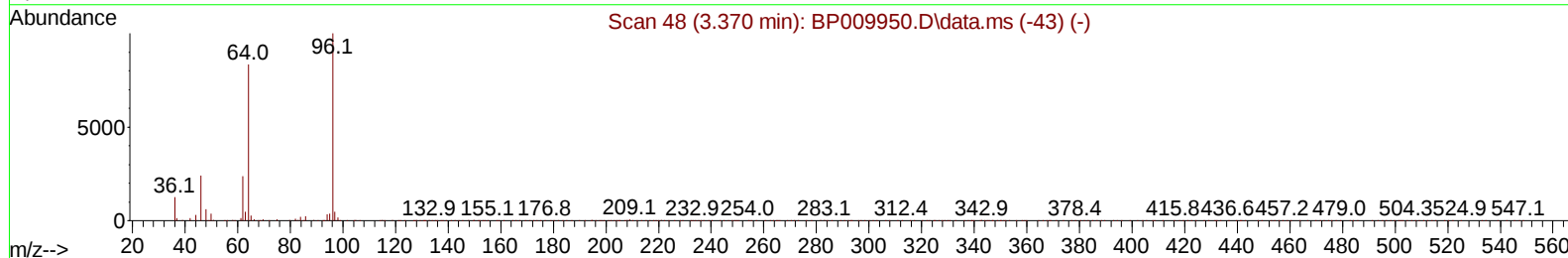
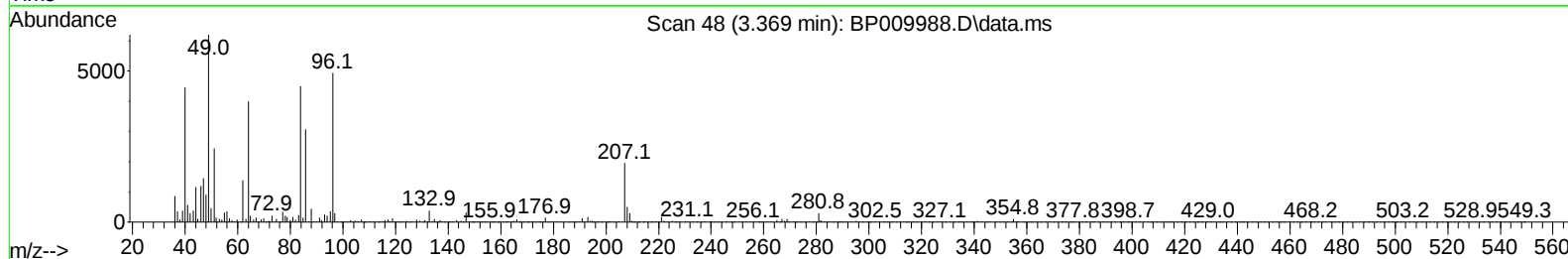
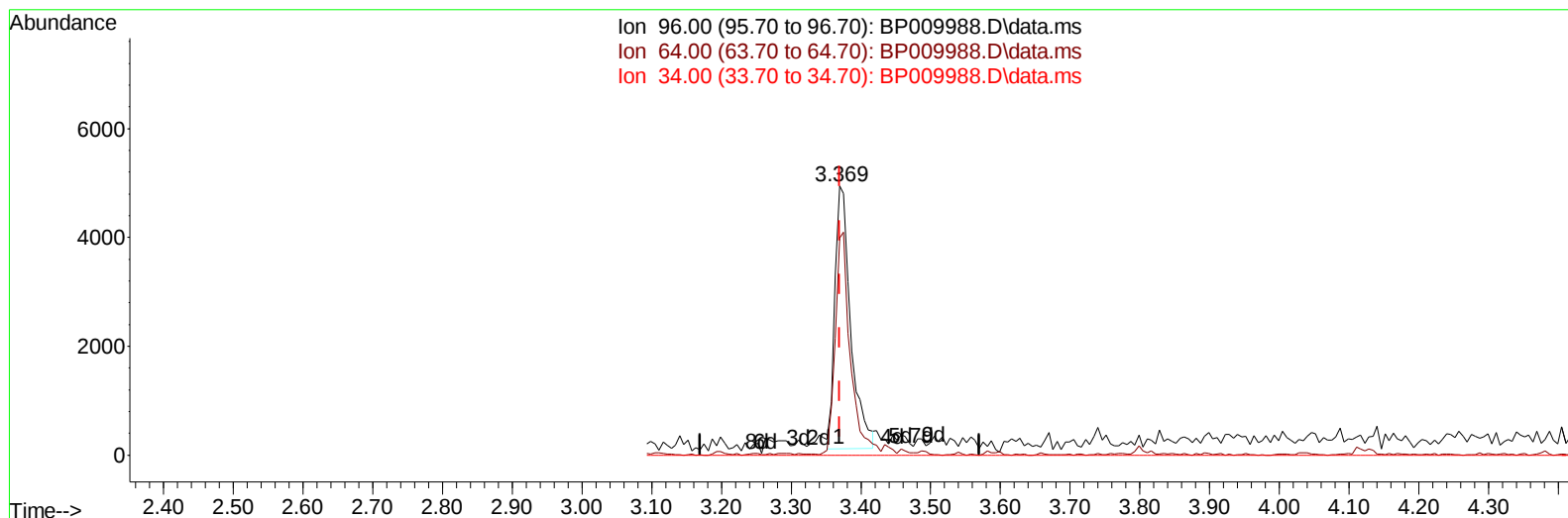
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009988.D
 Acq On : 20 Apr 2022 21:45
 Operator : CG/JU
 Sample : N2472-11
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J31

Manual IntegrationsAPPROVED

Quant Time: Apr 21 01:09:21 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: BP009988.D\data.ms

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

3.369min (-0.000) 2.78 ng/uL

response 7607

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.78#
34.00	0.00	0.00
0.00	0.00	0.00

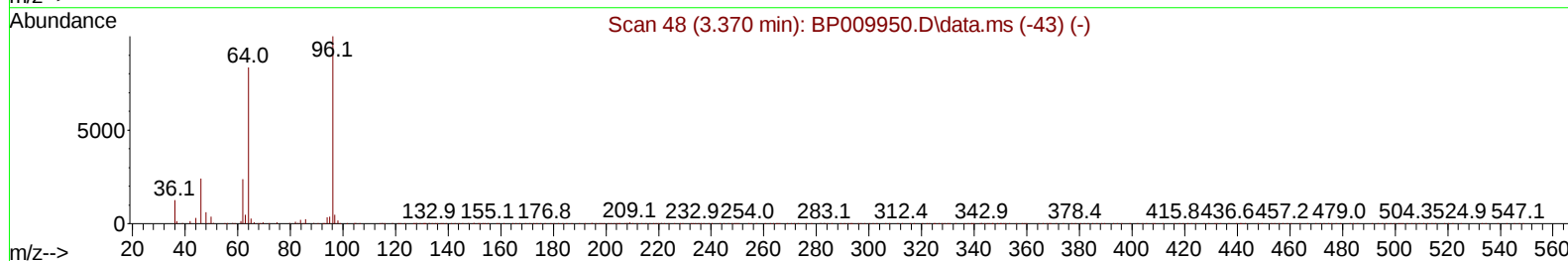
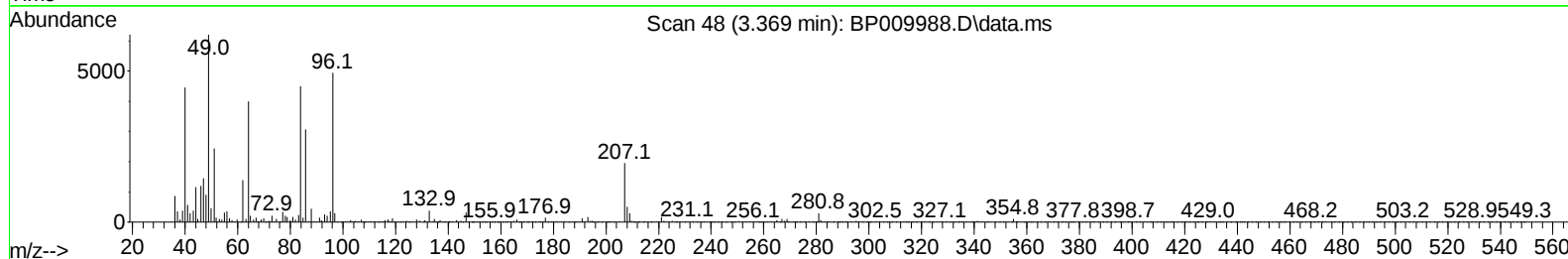
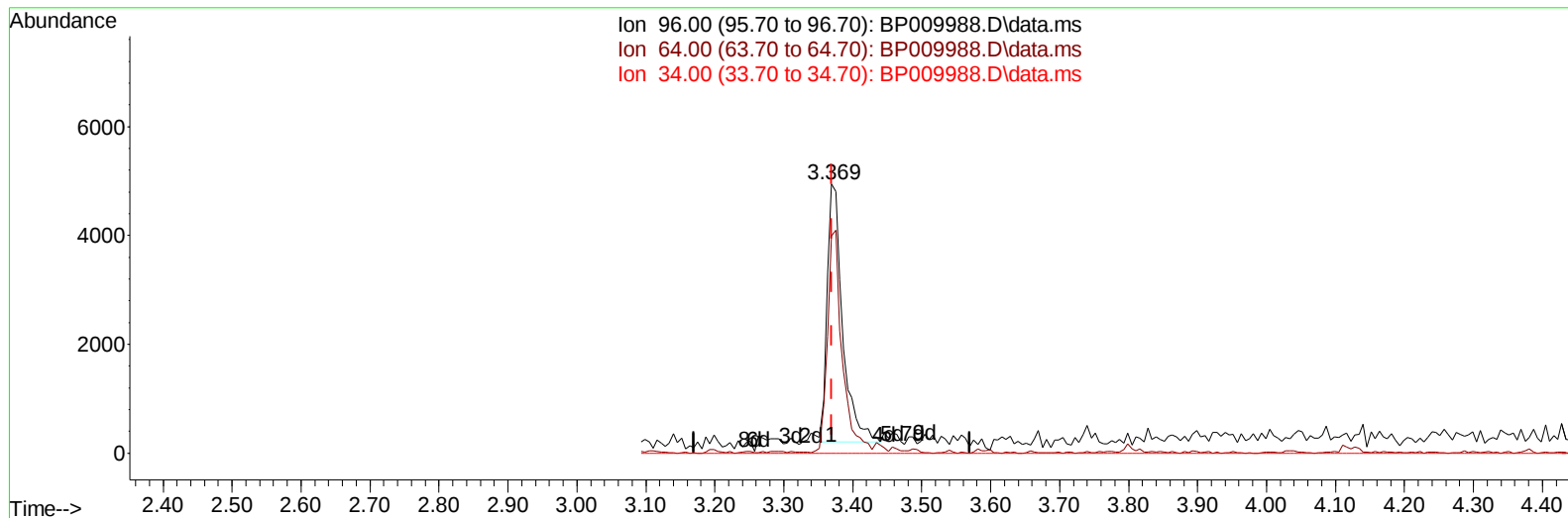
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TIC: BP009988.D\data.ms

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

3.369min (-0.000) 2.73 ng/uL m

response 7450

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.78#
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	121967	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	558634	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	395799	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.327	188	884723	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	875609	20.000	ng/ul	# 0.00
88) Perylene-d12	23.856	264	766290	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	7450m	2.726	ng/uL	0.00
4) Pyridine-d5	3.787	84	80353	10.664	ng/ul	0.00
7) Phenol-d5	7.104	99	83863	7.832	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	247403	38.810	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	241646	29.917	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	167099	18.694	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	166001	41.207	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	176607	39.987	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	290797	31.465	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	399913	30.381	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1226497	39.449	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1352742	36.726	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	41458	7.353	ng/ul	0.00
60) Fluorene-d10	15.569	176	1122894	42.933	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.680	200	191942	36.622	ng/ul	0.00
73) Anthracene-d10	17.422	188	1941877	45.842	ng/ul	0.00
81) Pyrene-d10	19.657	212	2365460	49.464	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1848291	46.317	ng/ul	0.00

Target Compounds Qvalue

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

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