

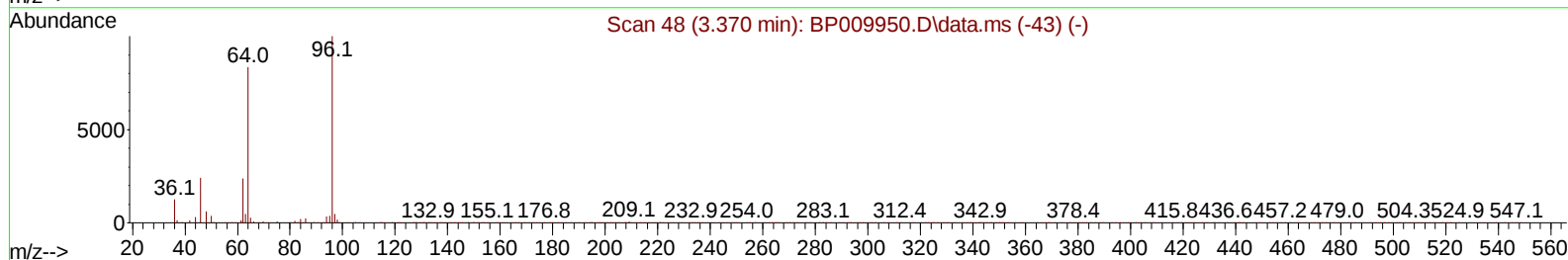
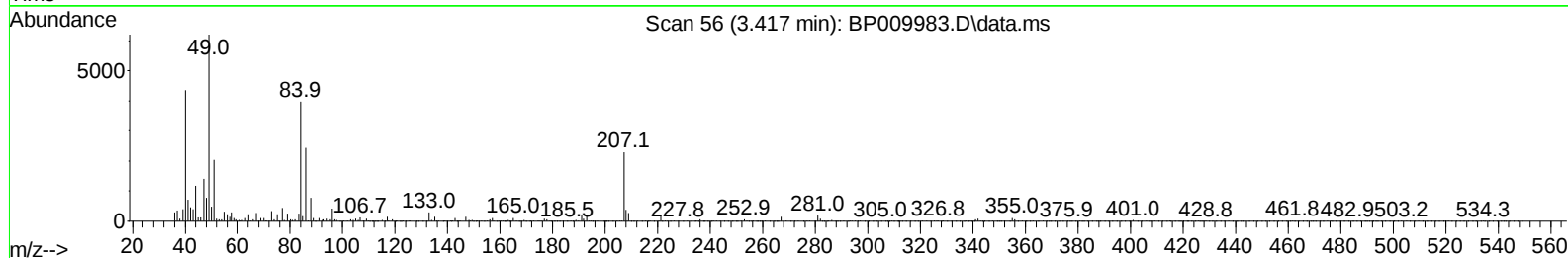
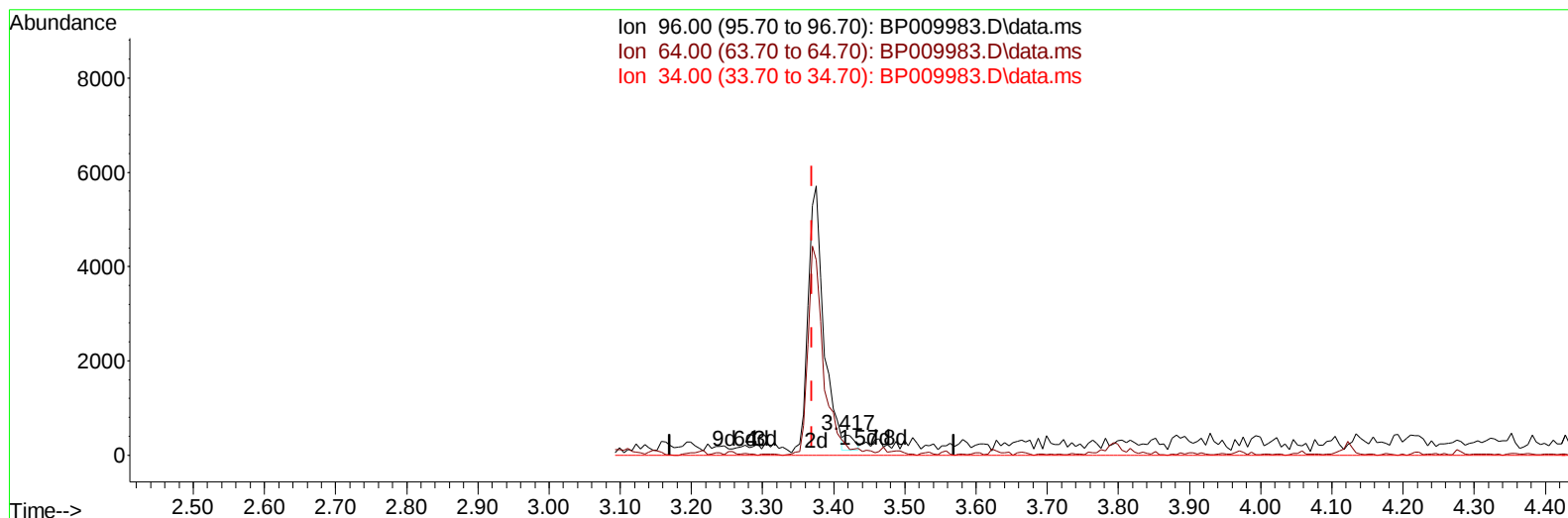
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009983.D
Acq On : 20 Apr 2022 18:57
Operator : CG/JU
Sample : N2472-10
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J10

Manual IntegrationsAPPROVED

Quant Time: Apr 21 01:08:11 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: BP009983.D\data.ms

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

3.417min (+ 0.047) 0.08 ng/uL

response 320

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	55.06
34.00	0.00	0.00
0.00	0.00	0.00

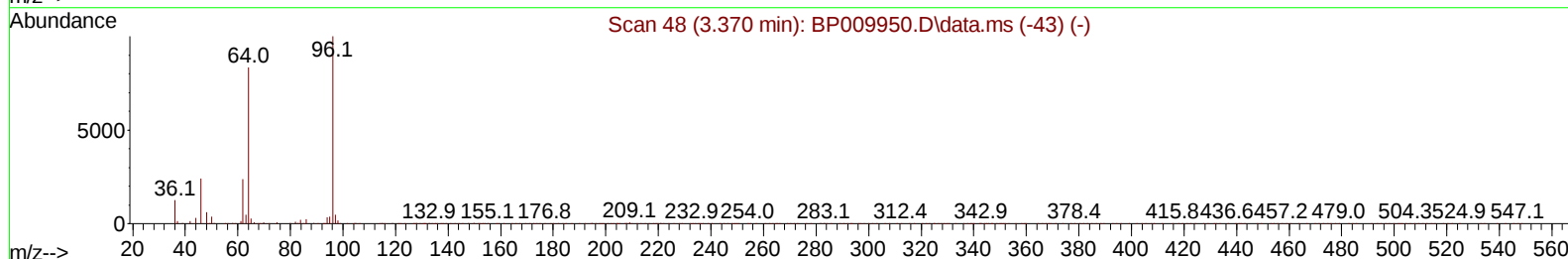
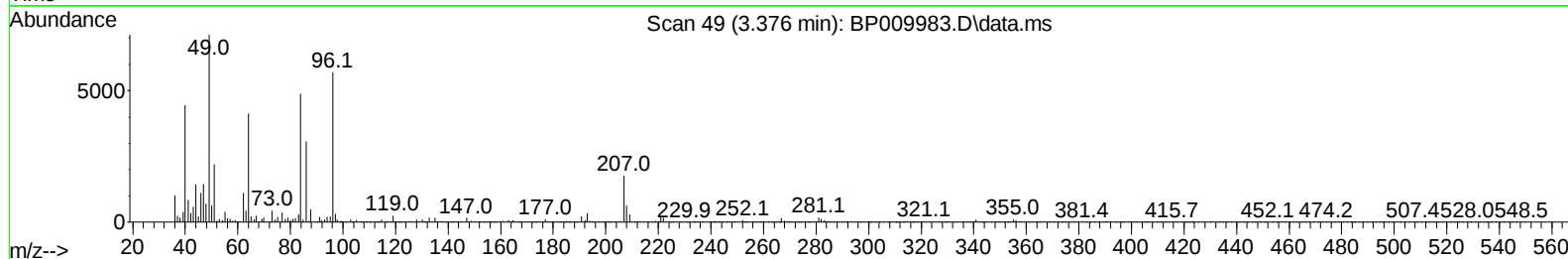
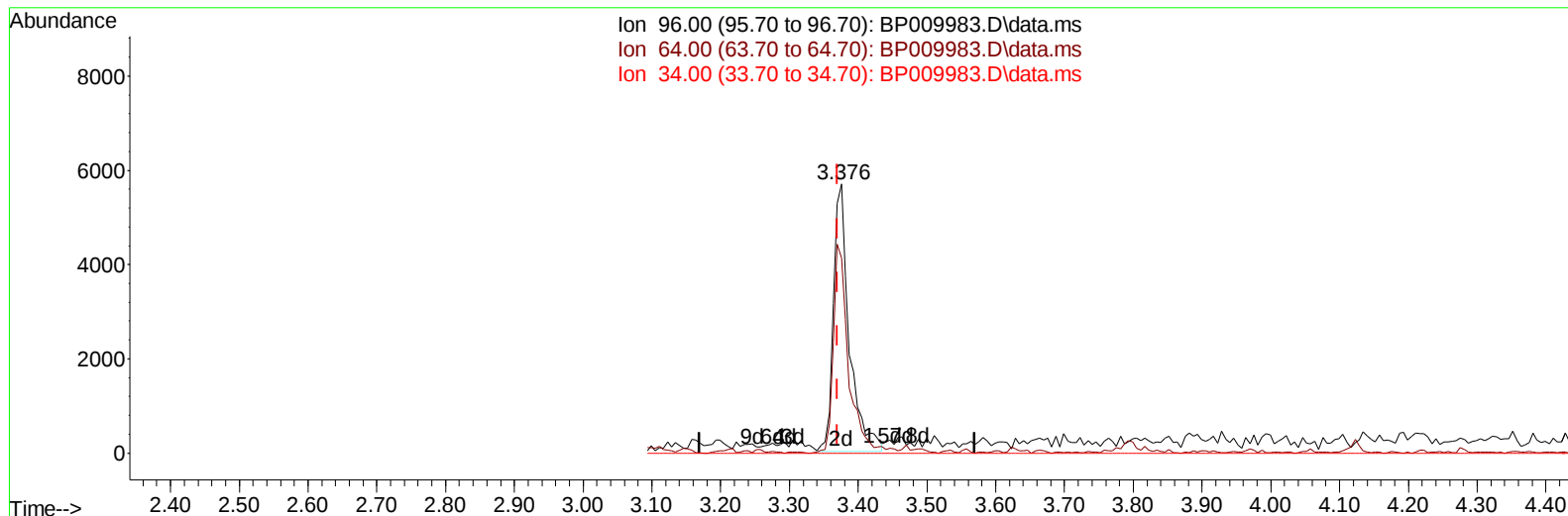
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(3) 1,4-Dioxane-d8 (S)

3.376min (+ 0.006) 2.39 ng/uL m

response 9121

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.56#
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	170070	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	761357	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	538362	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1138563	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	1033426	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	901068	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	9121m	2.394	ng/uL	0.00
4) Pyridine-d5	3.787	84	111521	10.614	ng/ul	0.00
7) Phenol-d5	7.105	99	110280	7.386	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	280720	31.581	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	297893	26.449	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	224877	18.042	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	195744	35.652	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	208961	34.715	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	402233	31.934	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	472594	26.343	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1492041	35.282	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1676326	33.459	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	44676	5.825	ng/ul	0.00
60) Fluorene-d10	15.569	176	1263661	35.521	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	207303	30.735	ng/ul	0.00
73) Anthracene-d10	17.428	188	1932600	35.452	ng/ul	0.00
81) Pyrene-d10	19.651	212	2302067	40.787	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1703964	36.313	ng/ul	-0.01

Target Compounds Qvalue

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

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