

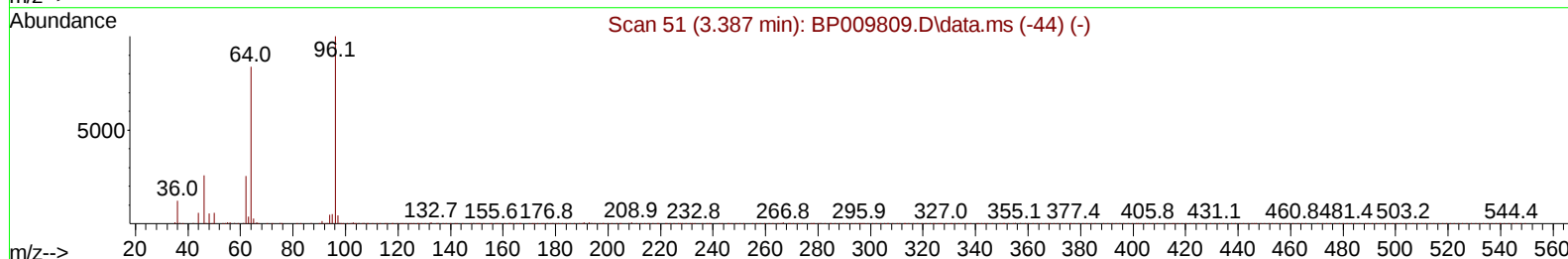
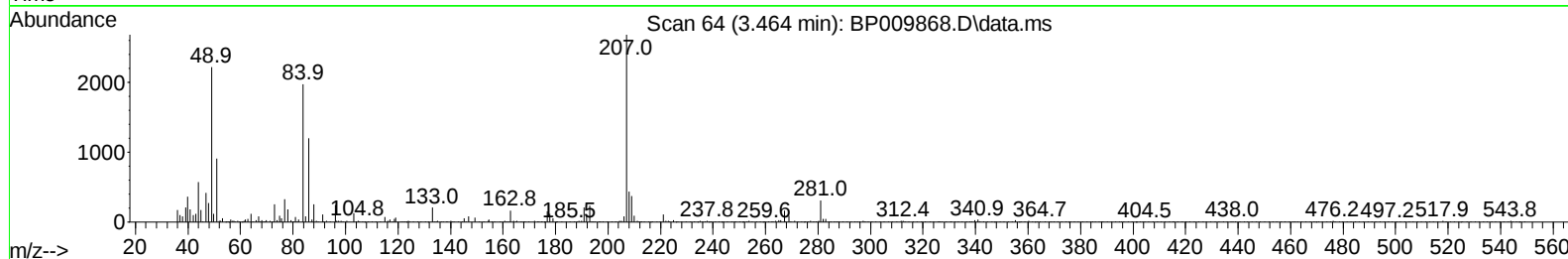
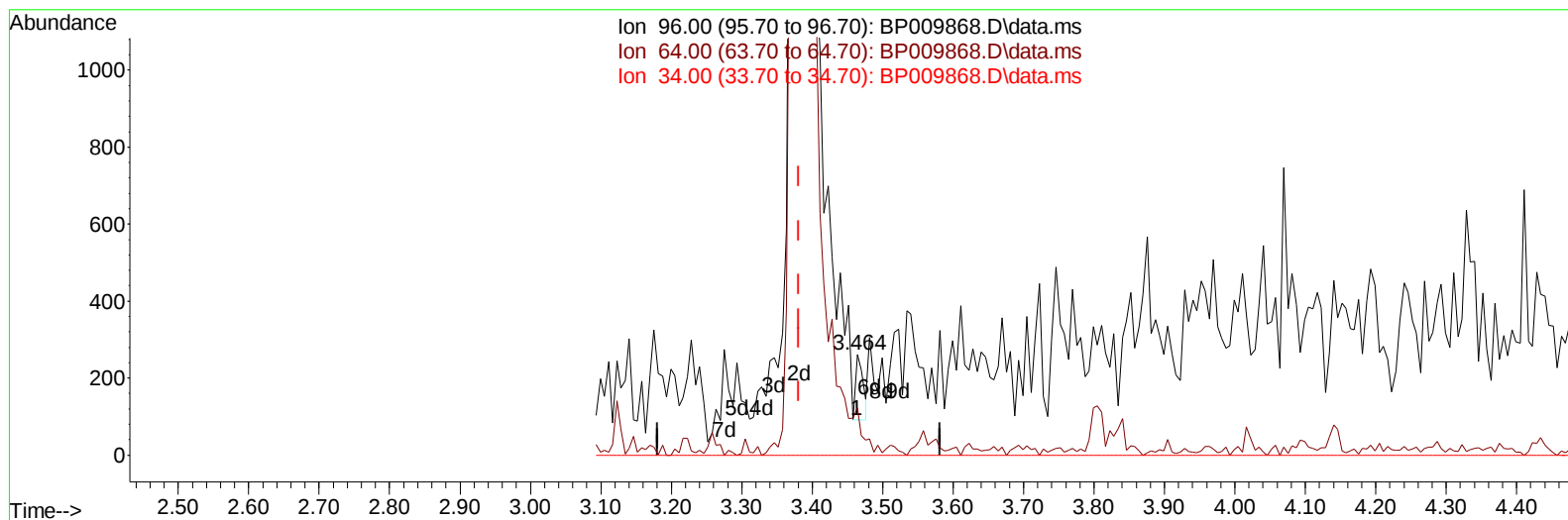
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009868.D
 Acq On : 15 Apr 2022 18:59
 Operator : CG/JU
 Sample : N2422-16
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J78

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:03:36 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 04/18/2022
 Supervised By :Yogesh Patel 04/18/2022



TIC: BP009868.D\data.ms

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

3.464min (+ 0.082) 0.04 ng/uL

response 125

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	47.13
34.00	0.00	0.00
0.00	0.00	0.00

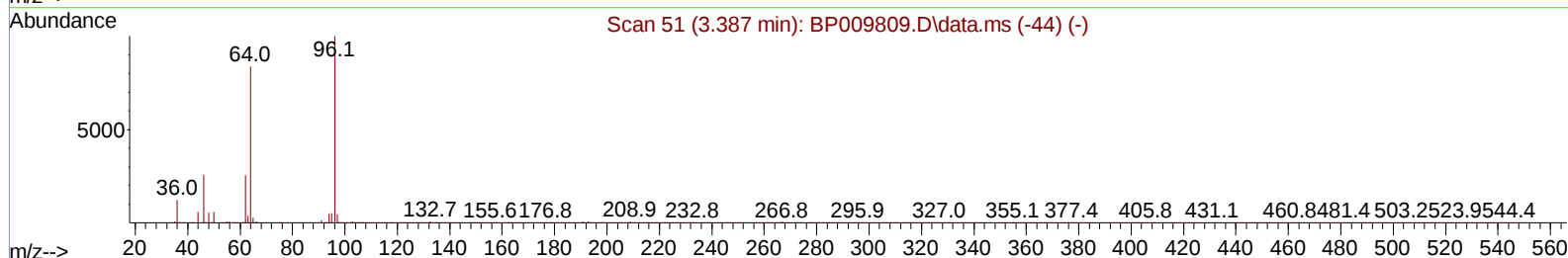
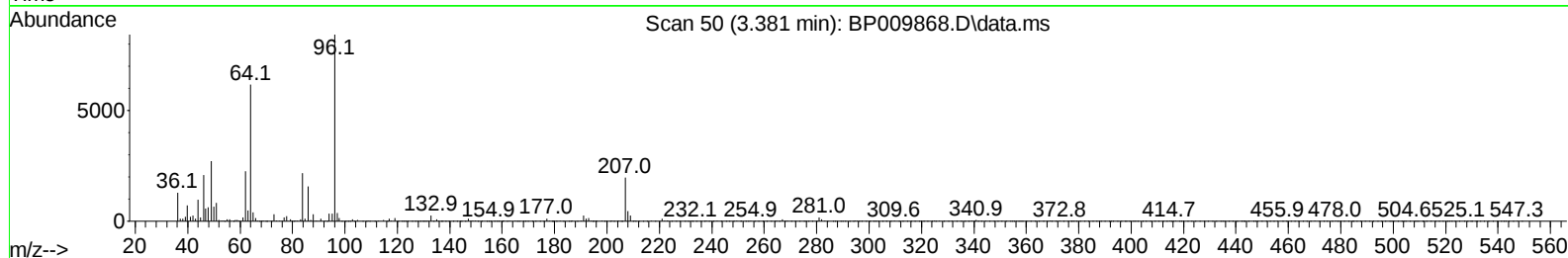
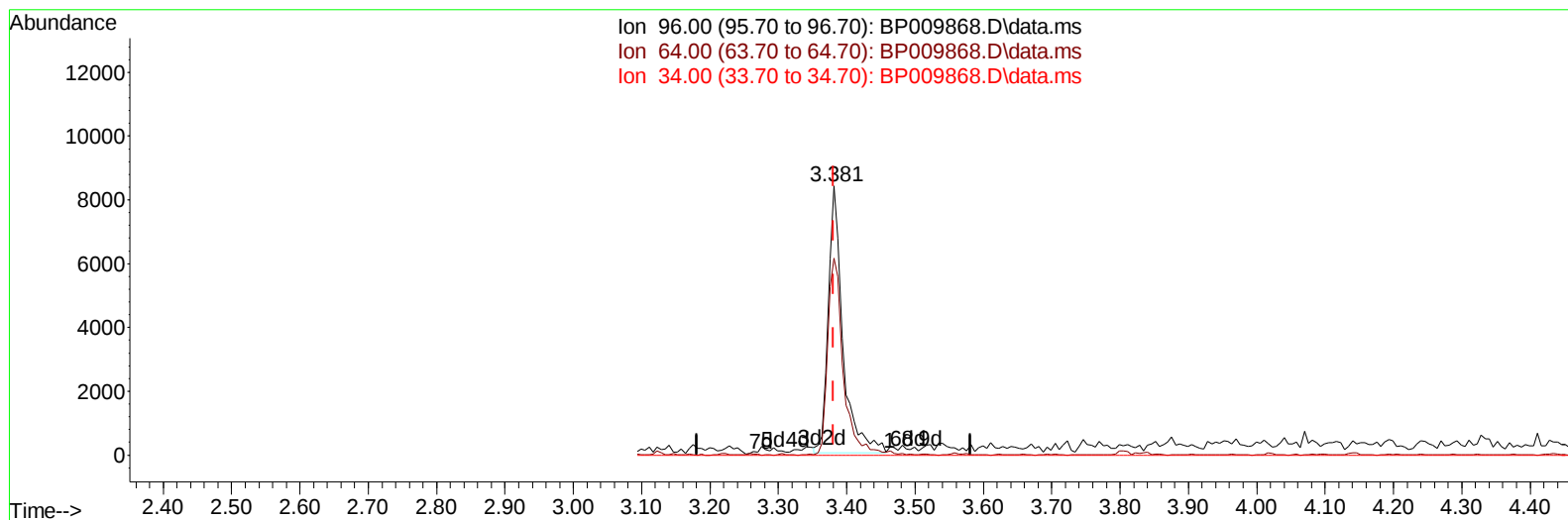
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TIC: BP009868.D\data.ms

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

3.381min (+ 0.000) 4.28 ng/uL m

response 12401

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.17#
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	7.958	152	129404	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	594881	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	441971	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1002298	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	1024864	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	955158	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12401m	4.277	ng/uL	0.00
4) Pyridine-d5	3.805	84	57531	7.196	ng/ul	0.00
7) Phenol-d5	7.111	99	68184	6.002	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	211067	31.207	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	195148	22.772	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	139276	14.685	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	140239	32.691	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	145510	30.939	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	255342	25.945	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	356765	25.452	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1123897	32.373	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1216744	29.583	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	36741	5.835	ng/ul	0.00
60) Fluorene-d10	15.581	176	980207	33.562	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	180931	30.472	ng/ul	0.00
73) Anthracene-d10	17.439	188	1745713	36.377	ng/ul	0.00
81) Pyrene-d10	19.669	212	2140580	38.243	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1948510	39.173	ng/ul	0.00

Target Compounds Qvalue

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

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