

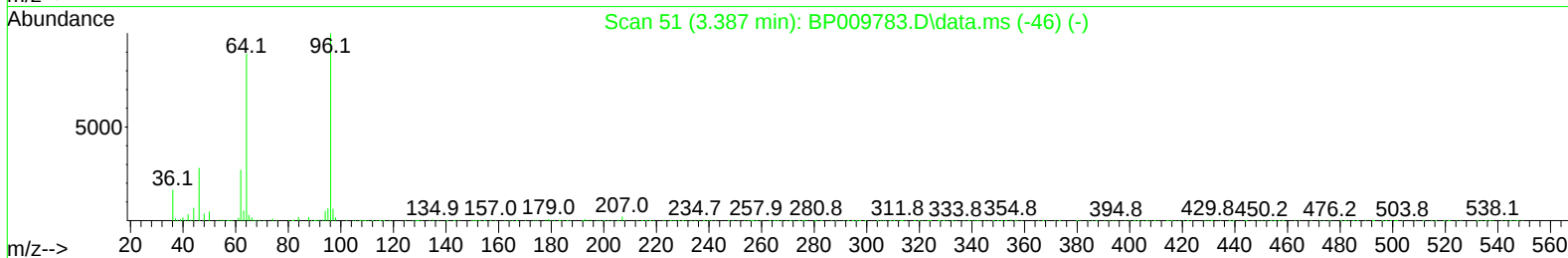
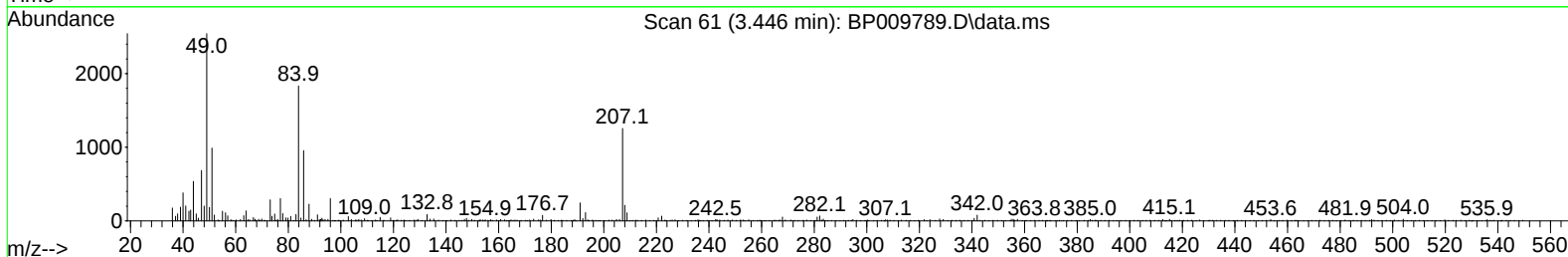
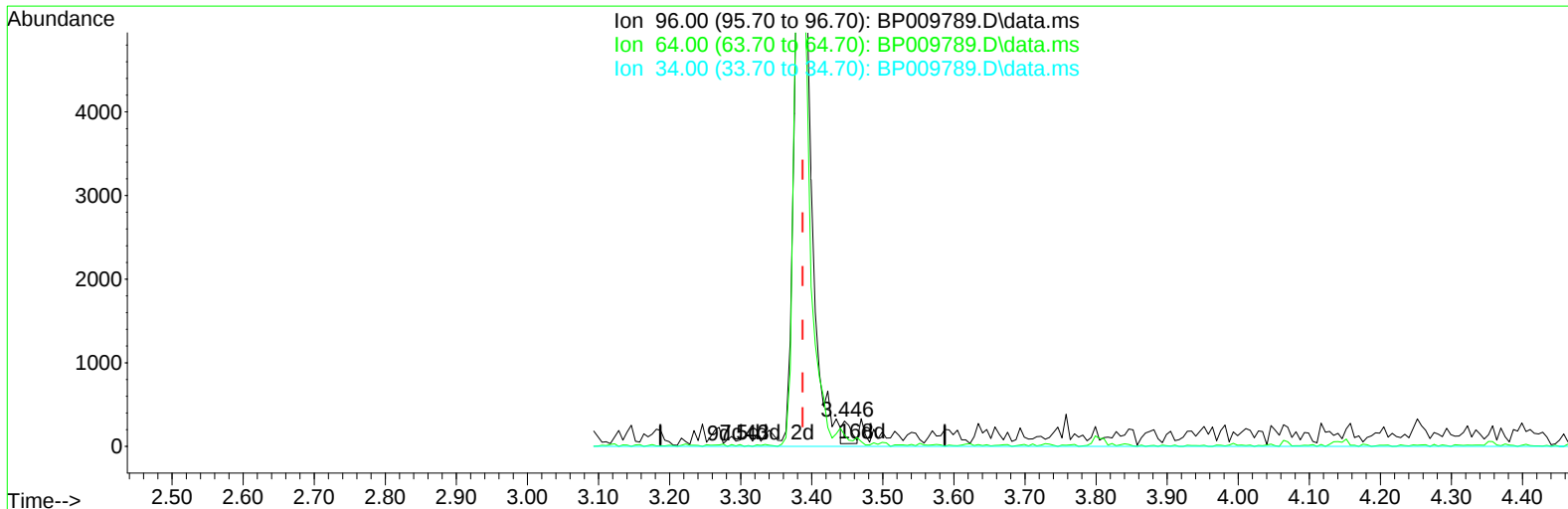
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009789.D
Acq On : 12 Apr 2022 18:42
Operator : CG/JU
Sample : N2316-17
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNR5

Manual IntegrationsAPPROVED

Quant Time: Apr 13 10:44:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
Supervised By :Yogesh Patel 04/13/2022



TIC: BP009789.D\data.ms

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

3.446min (+ 0.059) 0.10 ng/uL

response 225

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	44.88
34.00	0.00	0.00
0.00	0.00	0.00

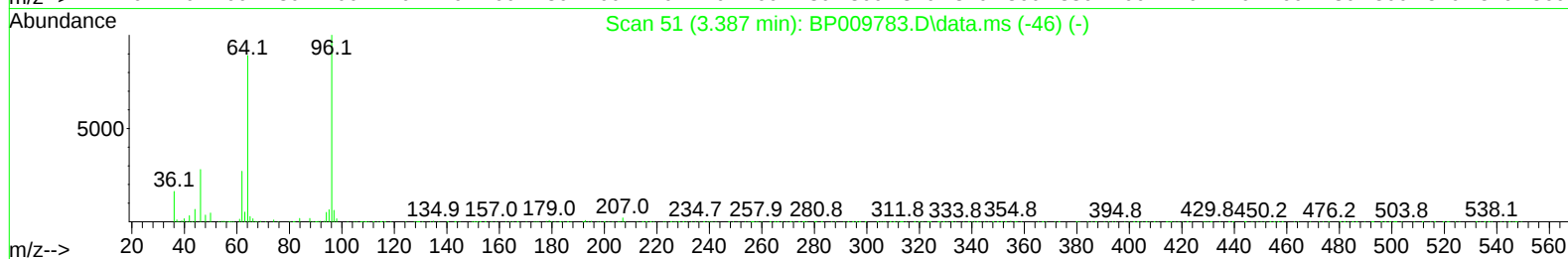
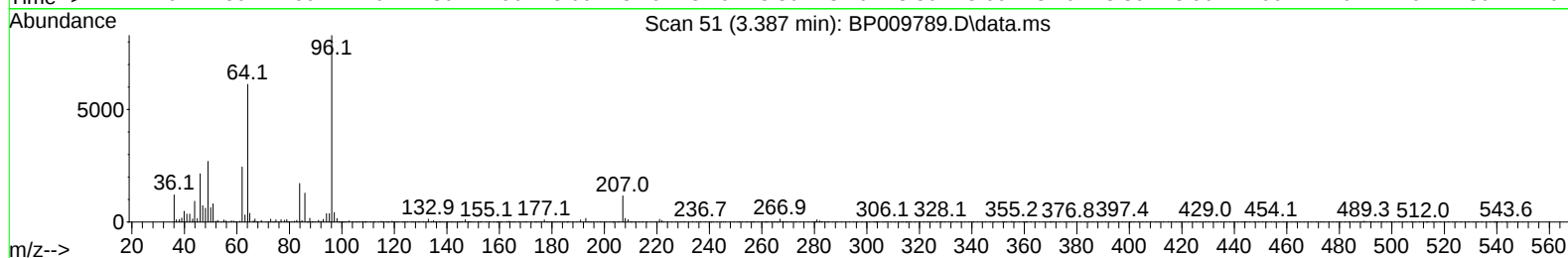
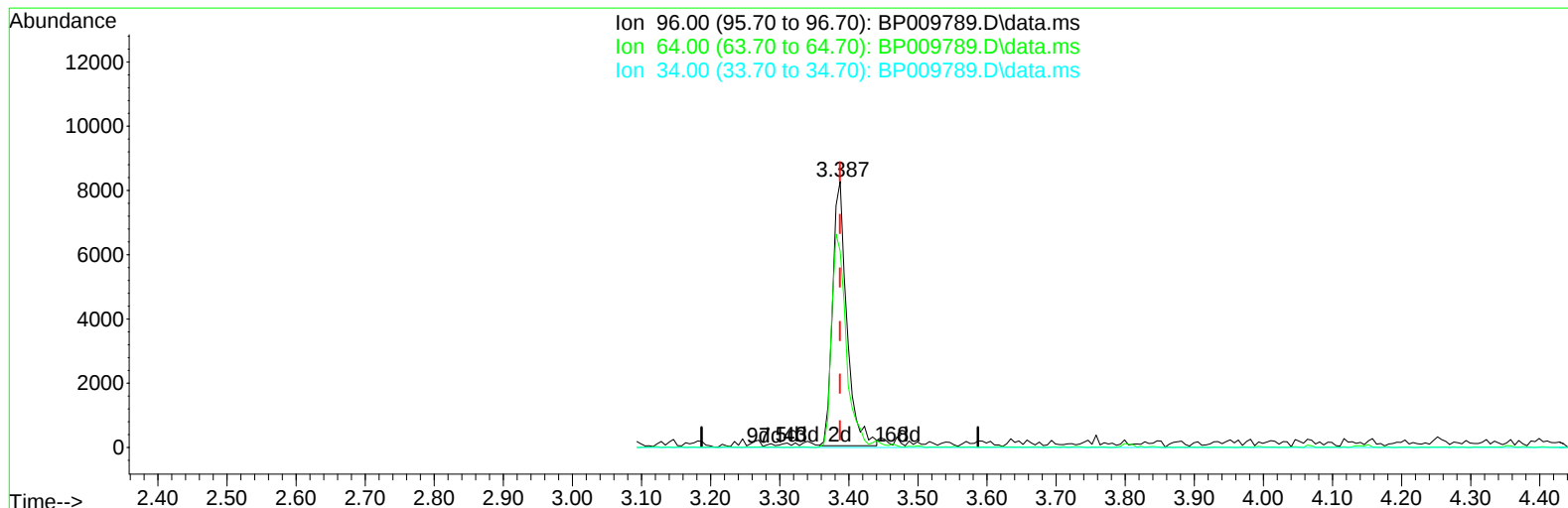
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TIC: BP009789.D\data.ms

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

3.387min (+ 0.000) 5.07 ng/uL m

response 11804

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.82#
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	7.969	152	103979	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	467073	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	337516	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	755039	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.428	240	798650	20.000	ng/u1	# 0.00
88) Perylene-d12	23.892	264	754078	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	11804m	5.066	ng/uL	0.00
4) Pyridine-d5	3.805	84	56297	8.764	ng/u1	0.00
7) Phenol-d5	7.117	99	55563	6.087	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	164565	30.281	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	155391	22.567	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	108491	14.237	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	104594	31.053	ng/u1	0.00
24) 2-Nitrophenol-d4	9.852	143	105837	28.661	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	185665	24.027	ng/u1	0.00
31) 4-Chloroaniline-d4	10.905	131	280056	25.446	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	825955	31.154	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	867834	27.630	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	31747	6.603	ng/u1	-0.01
60) Fluorene-d10	15.587	176	740138	33.185	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	136522	30.522	ng/u1	0.00
73) Anthracene-d10	17.445	188	1341665	37.113	ng/u1	0.00
81) Pyrene-d10	19.675	212	1663309	38.133	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	1494582	38.059	ng/u1	0.00

Target Compounds Qvalue

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

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