

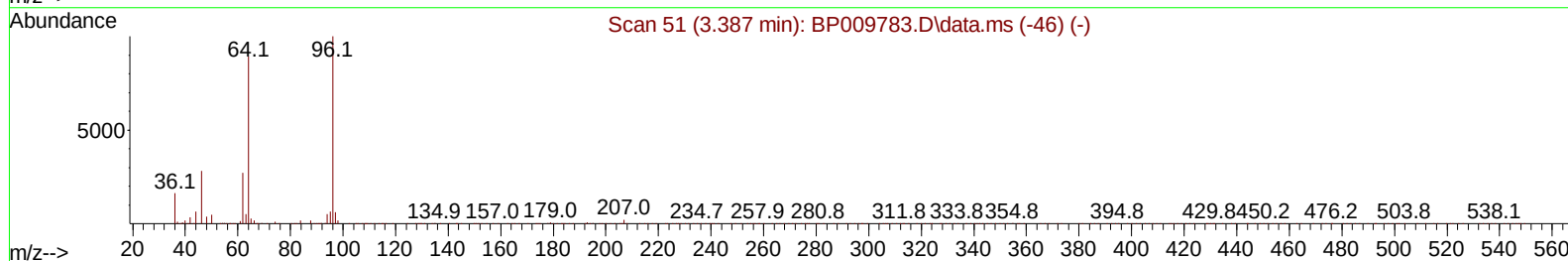
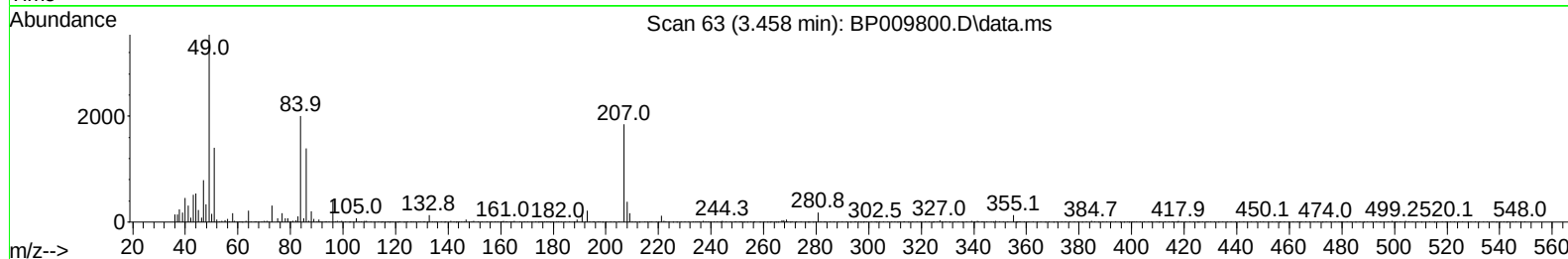
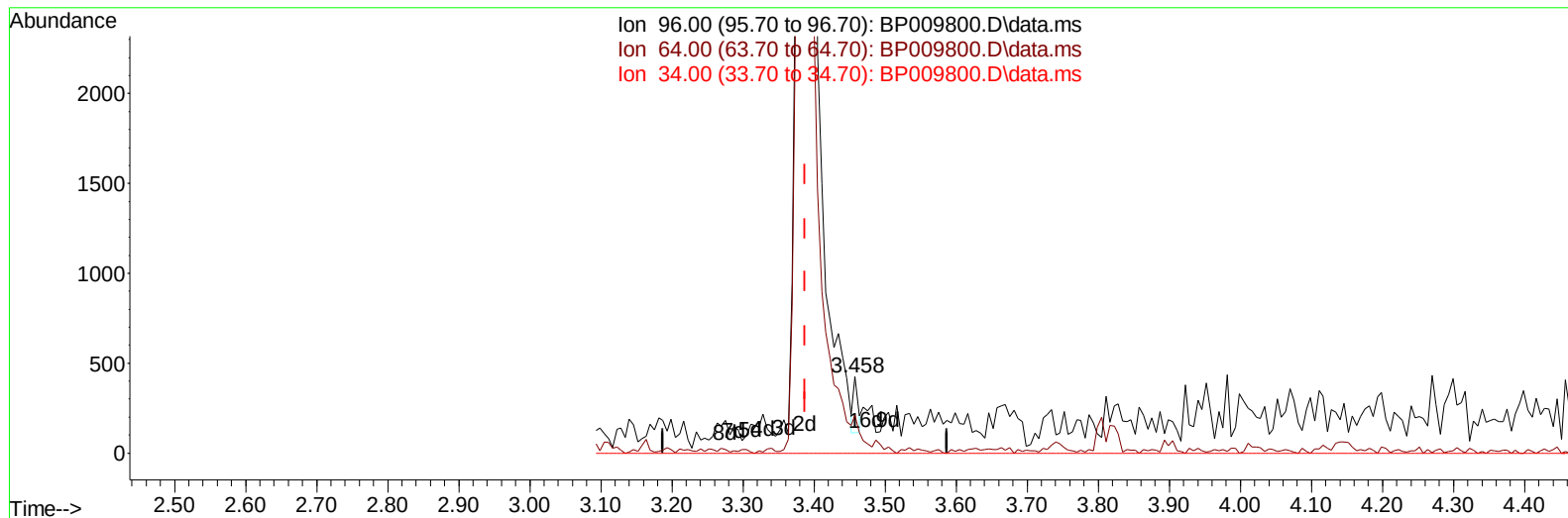
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009800.D
 Acq On : 13 Apr 2022 01:29
 Operator : CG/JU
 Sample : N2342-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY4

Manual IntegrationsAPPROVED

Quant Time: Apr 13 04:27:46 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: BP009800.D\data.ms

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

3.458min (+ 0.071) 0.07 ng/uL

response 143

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	49.76
34.00	0.00	0.00
0.00	0.00	0.00

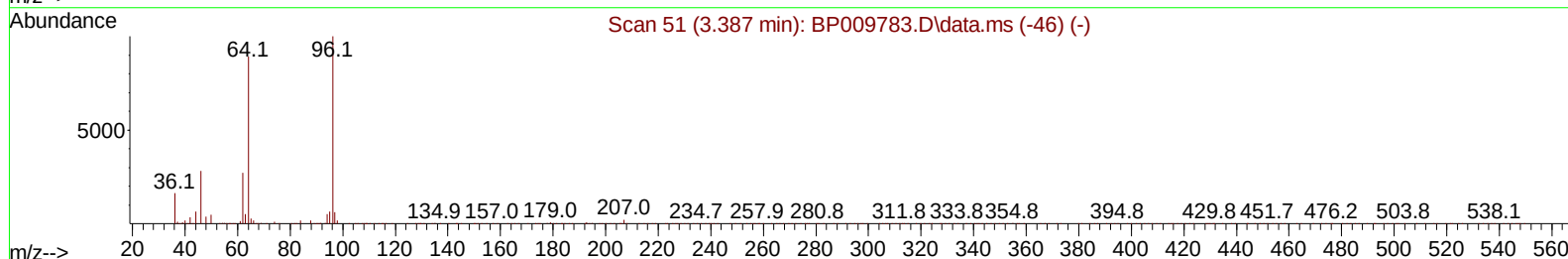
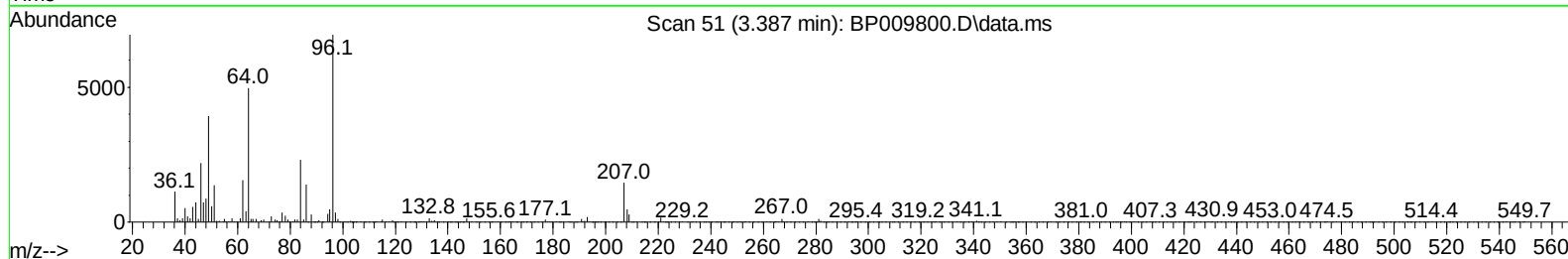
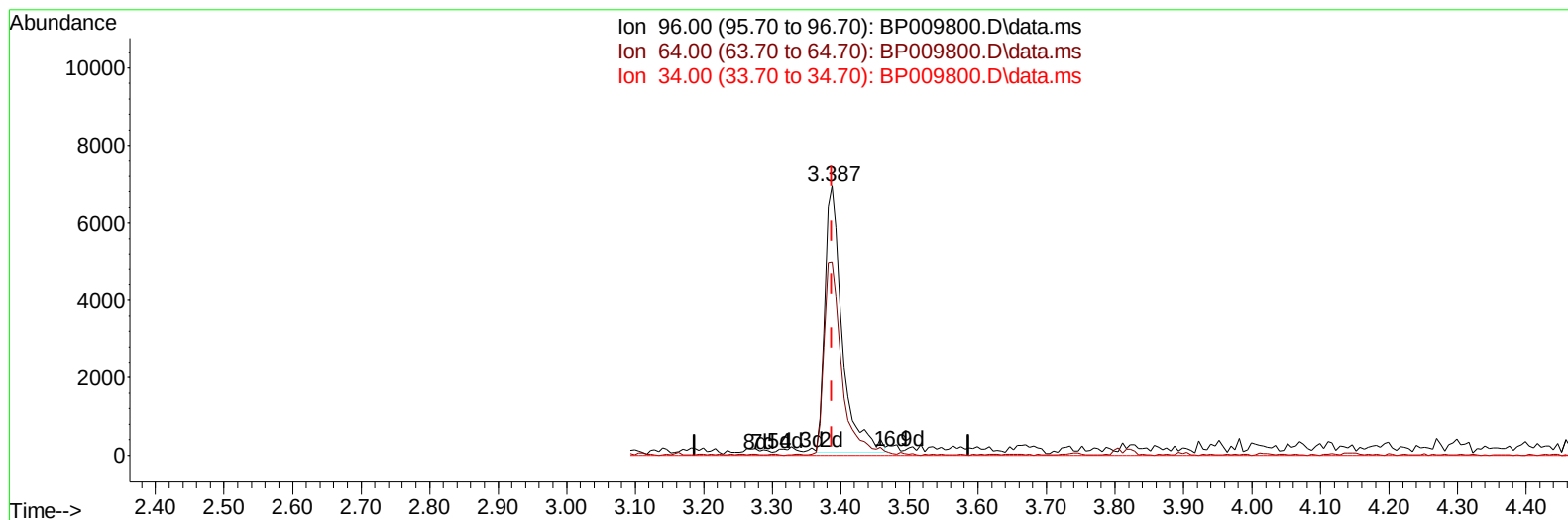
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TIC: BP009800.D\data.ms

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

3.387min (+ 0.000) 5.49 ng/uL m

response 12078

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	71.49#
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	7.963	152	98120	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	463332	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	354372	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	807046	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.421	240	946814	20.000	ng/ul	#-0.01
88) Perylene-d12	23.892	264	1011928	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	12078m	5.494	ng/uL	0.00
4) Pyridine-d5	3.805	84	73475	12.121	ng/ul	0.00
7) Phenol-d5	7.116	99	61061	7.088	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	162070	31.602	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	158126	24.335	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	120190	16.714	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	106194	31.783	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	111152	30.343	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	194364	25.356	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	288213	26.399	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	863547	31.022	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	936666	28.402	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	35199	6.972	ng/ul	-0.01
60) Fluorene-d10	15.586	176	770810	32.917	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	137775	28.817	ng/ul	0.00
73) Anthracene-d10	17.445	188	1378457	35.674	ng/ul	0.00
81) Pyrene-d10	19.669	212	1763112	34.096	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1935548	36.729	ng/ul	0.00

Target Compounds Qvalue

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

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