

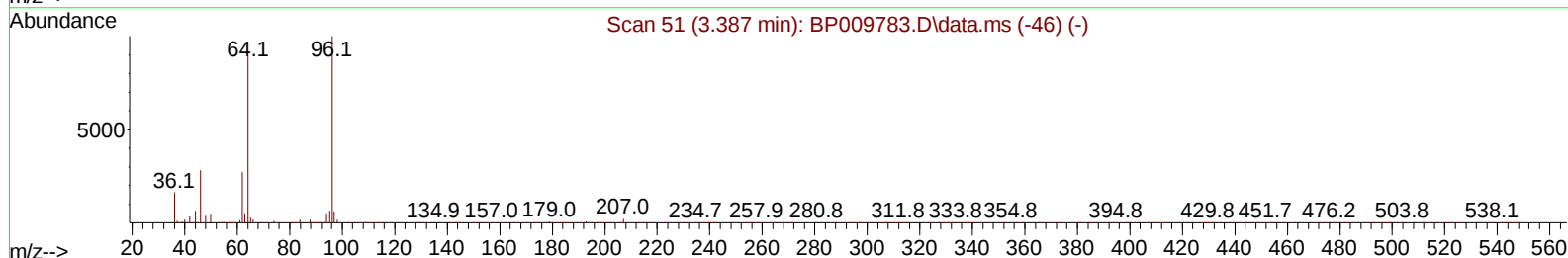
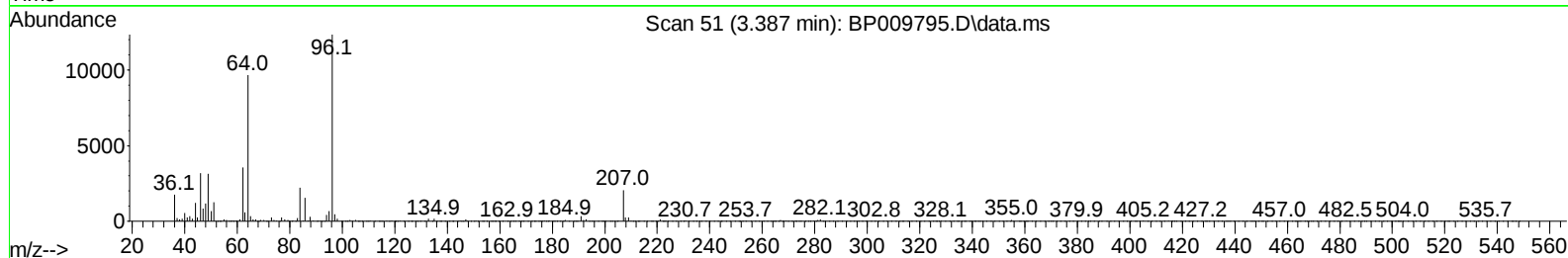
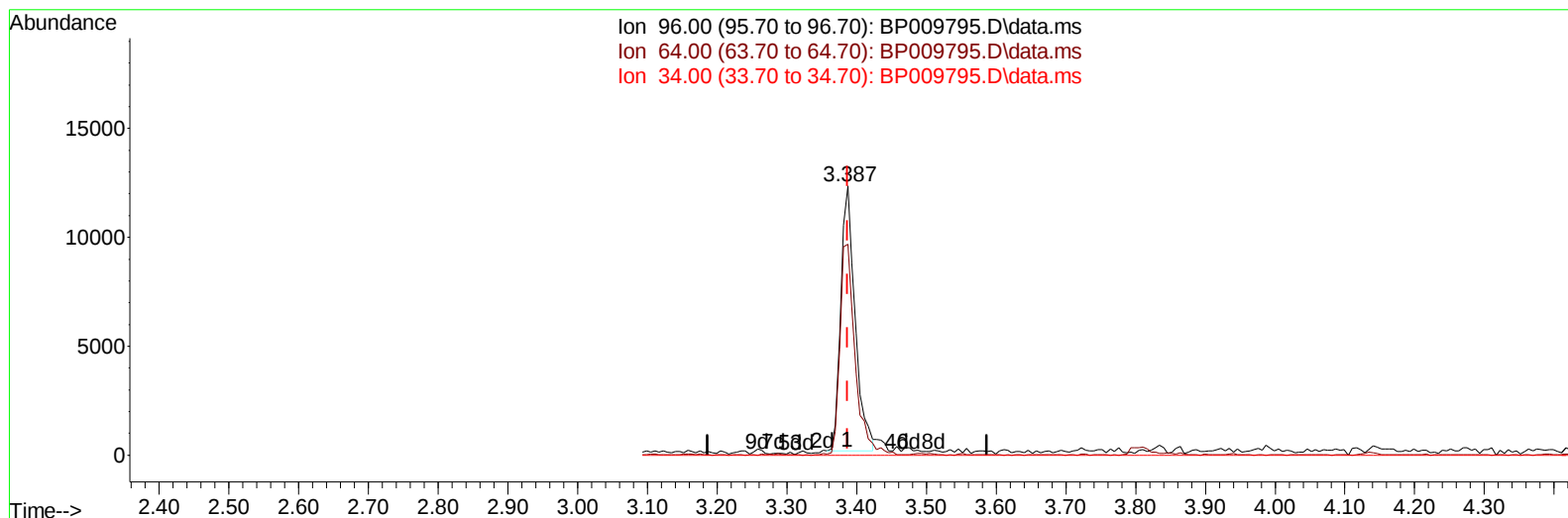
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009795.D
 Acq On : 12 Apr 2022 22:40
 Operator : CG/JU
 Sample : PB144055BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK055

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:26:12 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: BP009795.D\data.ms

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

3.387min (+ 0.000) 5.71 ng/uL

response 17314

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.47#
34.00	0.00	0.00
0.00	0.00	0.00

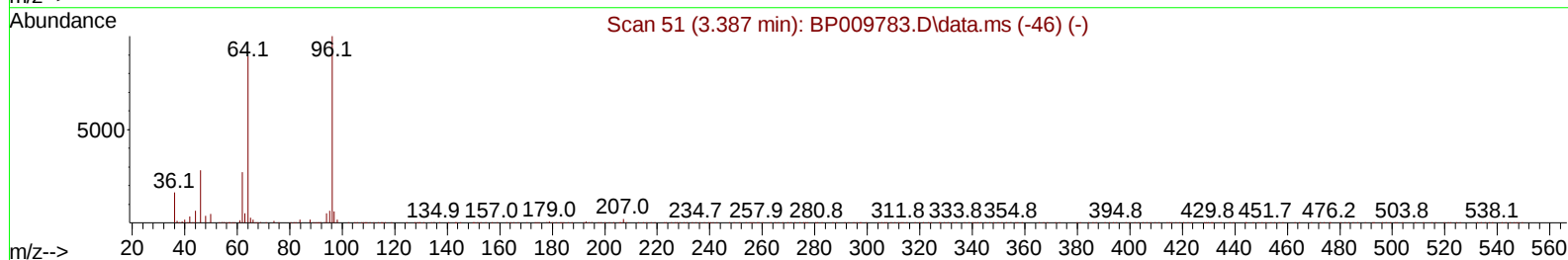
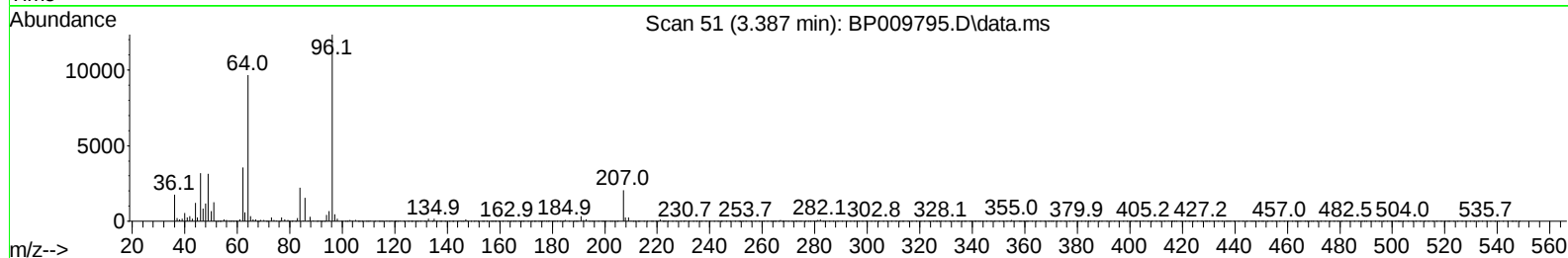
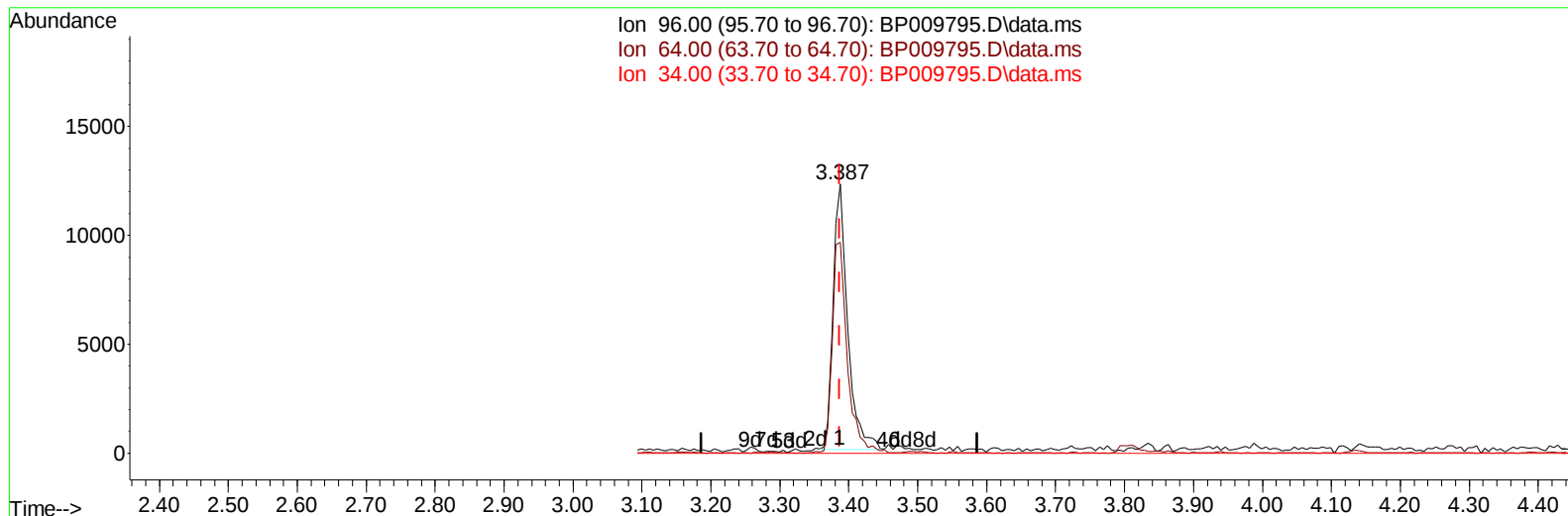
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TIC: BP009795.D\data.ms

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

3.387min (+ 0.000) 5.89 ng/uL m

response 17866

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.47#
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	135353	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	602861	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	440495	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	966971	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	1028870	20.000	ng/ul	#-0.01
88) Perylene-d12	23.892	264	971220	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17866m	5.891	ng/uL	0.00
4) Pyridine-d5	3.799	84	254873	30.480	ng/ul	0.00
7) Phenol-d5	7.116	99	320490	26.970	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	213436	30.170	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	263881	29.439	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	279106	28.136	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	140035	32.211	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	150209	31.515	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	265947	26.665	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	414503	29.179	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	1063866	30.746	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1147890	28.002	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	178360	28.423	ng/ul	0.00
60) Fluorene-d10	15.587	176	877157	30.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	149738	26.140	ng/ul	0.00
73) Anthracene-d10	17.445	188	1439740	31.097	ng/ul	0.00
81) Pyrene-d10	19.675	212	1751304	31.166	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1616229	31.955	ng/ul	0.00

Target Compounds Qvalue

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

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