

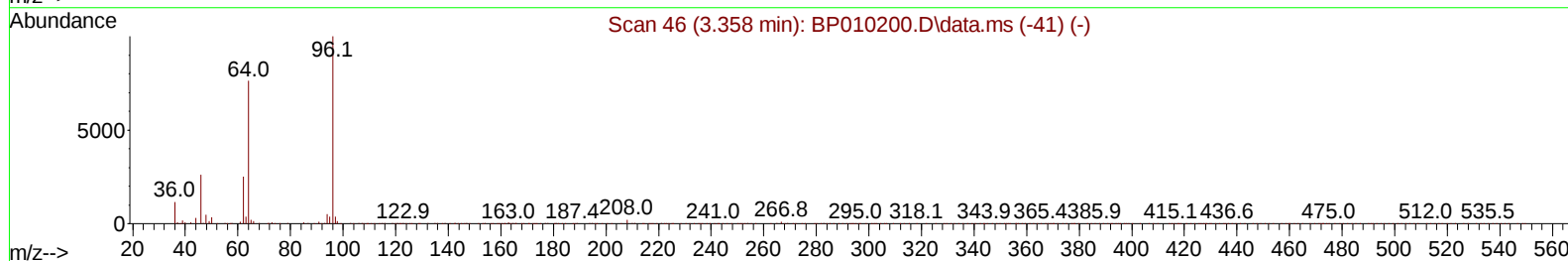
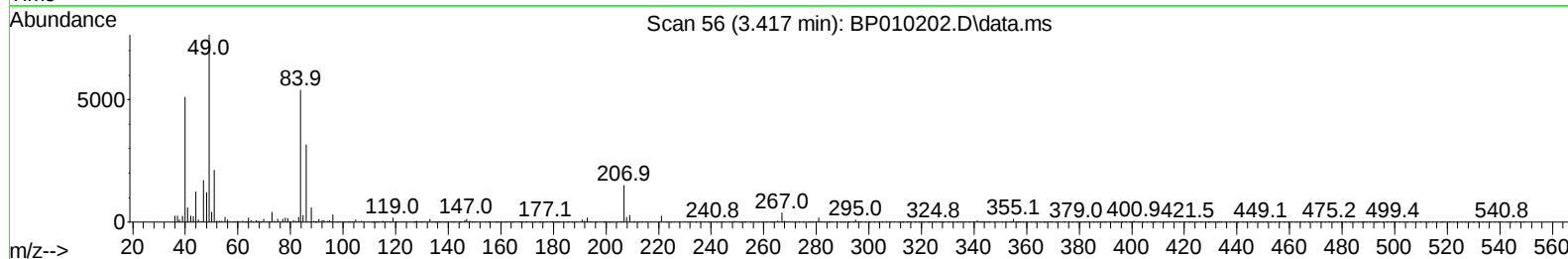
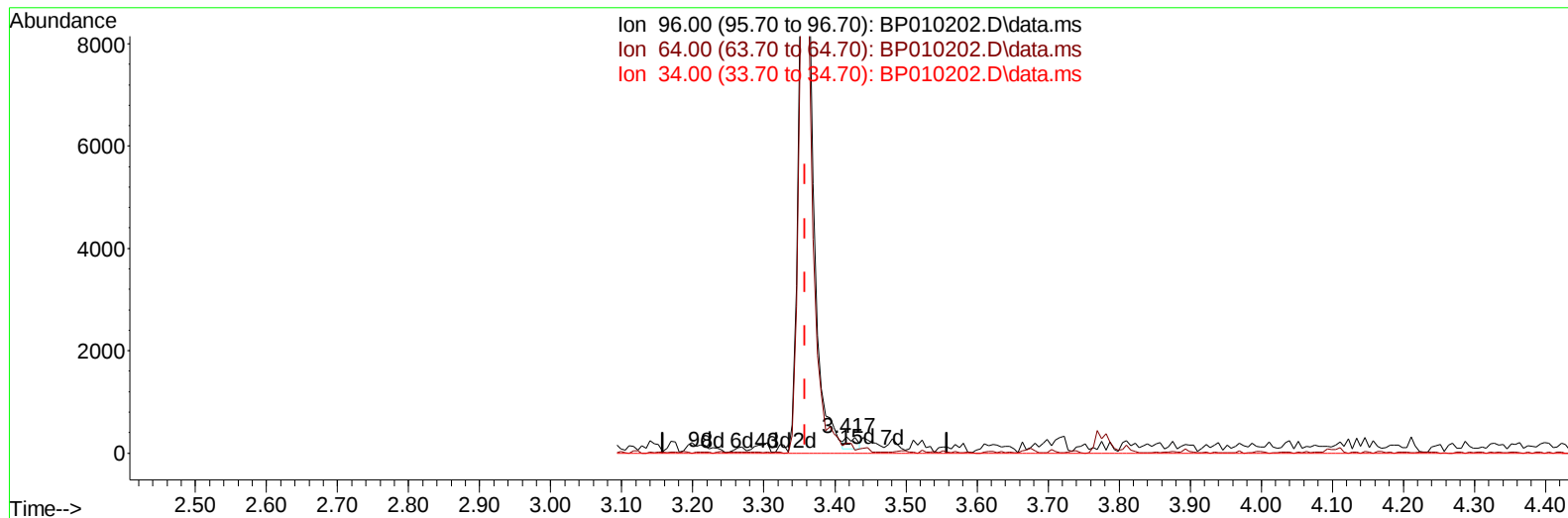
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010202.D
 Acq On : 03 May 2022 10:20
 Operator : CG/JU
 Sample : PB144510BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK510

Manual IntegrationsAPPROVED

Quant Time: May 04 01:08:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 00:58:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/04/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010202.D\data.ms

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

3.417min (+ 0.059) 0.09 ng/uL

response 256

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	61.54
34.00	0.00	0.00
0.00	0.00	0.00

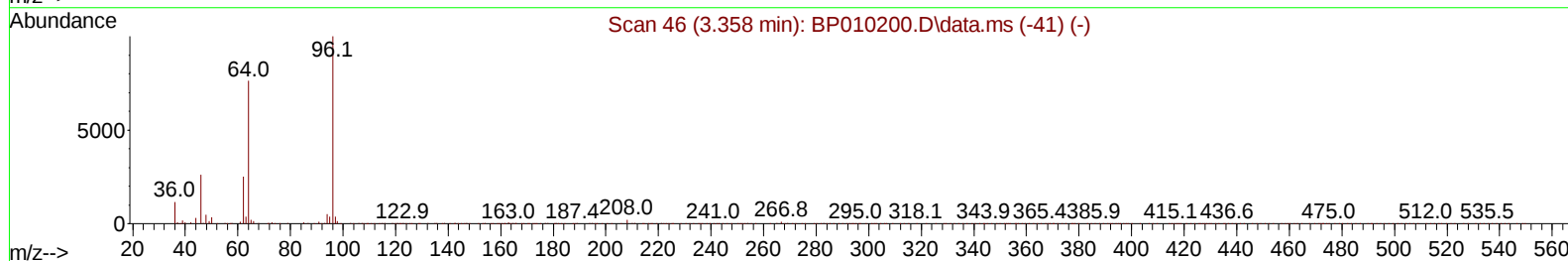
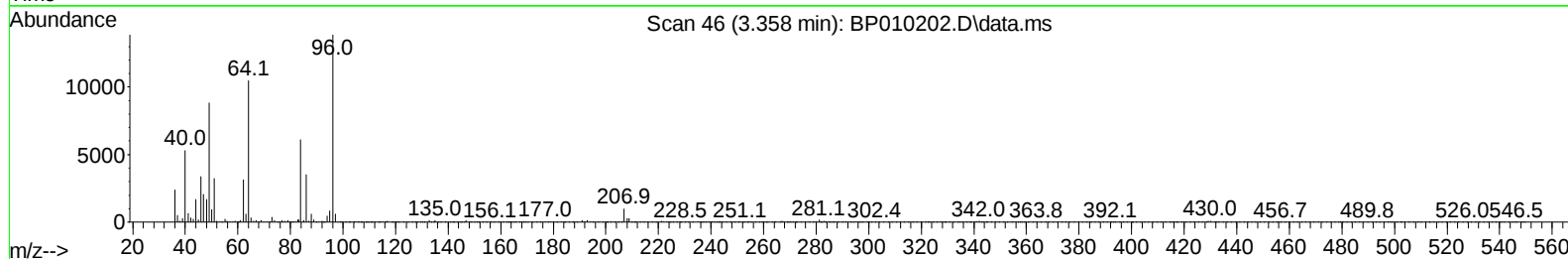
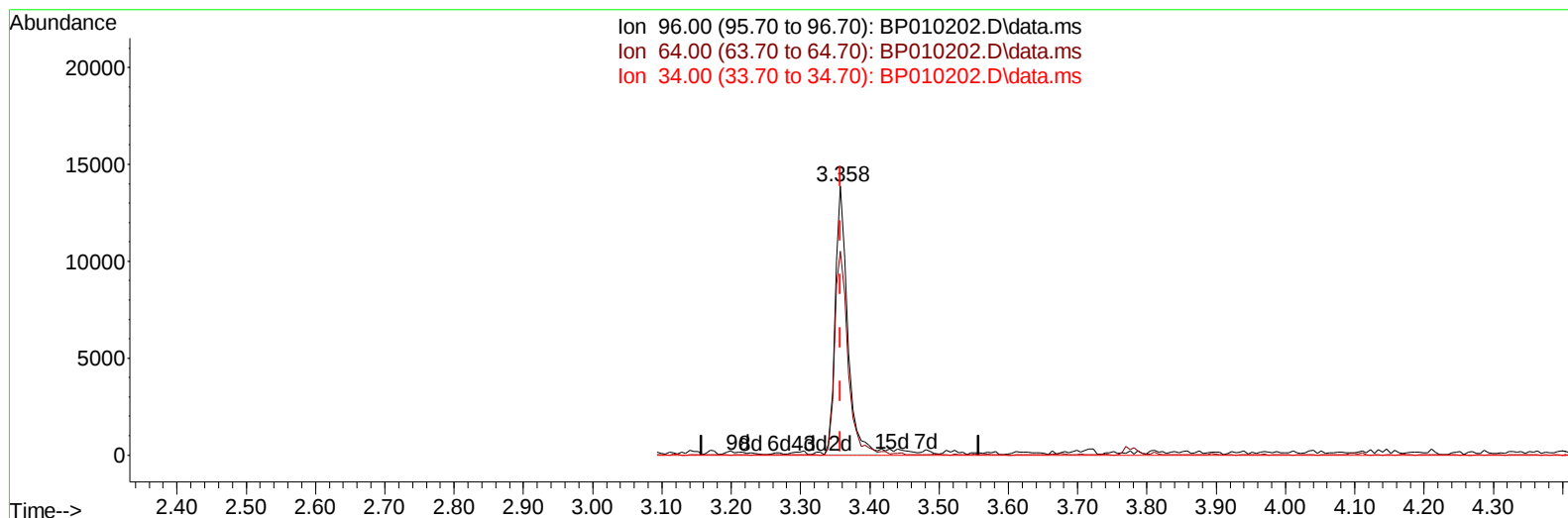
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TIC: BP010202.D\data.ms

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

3.358min (-0.000) 5.97 ng/uL m

response 17275

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	75.72#
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	7.922	152	126237	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	517093	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	328580	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	671999	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	627446	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	595094	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	17275m	5.967	ng/uL	0.00
4) Pyridine-d5	3.775	84	232305	28.525	ng/ul	0.00
7) Phenol-d5	7.081	99	278652	24.268	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	201747	29.281	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	236352	27.794	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	233080	24.500	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	116619	28.356	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	131400	29.255	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	211225	24.156	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	304282	24.270	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	721941	28.245	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	831694	26.886	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	91977	19.547	ng/ul	0.00
60) Fluorene-d10	15.545	176	630472	28.675	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	84152	19.804	ng/ul	0.00
73) Anthracene-d10	17.404	188	974198	29.966	ng/ul	0.00
81) Pyrene-d10	19.633	212	1140297	31.956	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	976628	31.304	ng/ul	0.00

Target Compounds Qvalue

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

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