

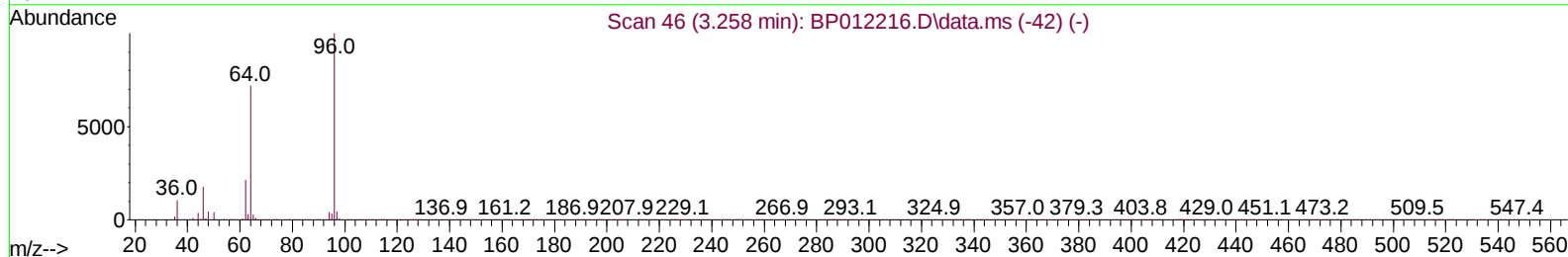
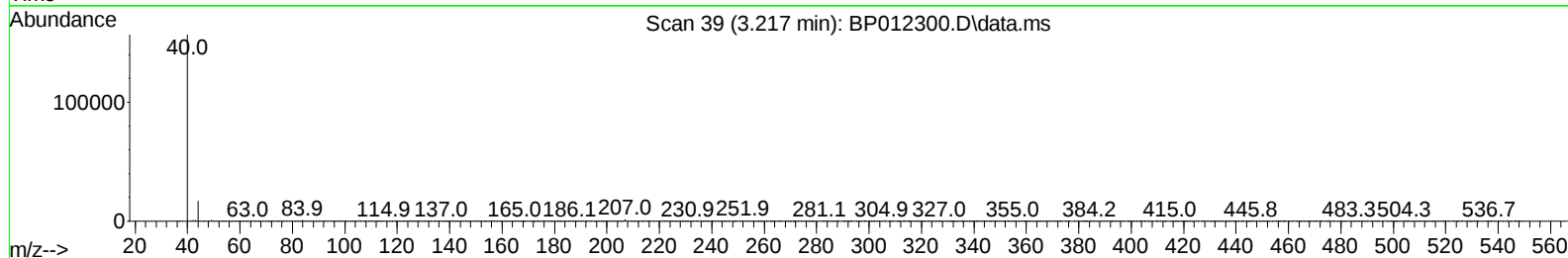
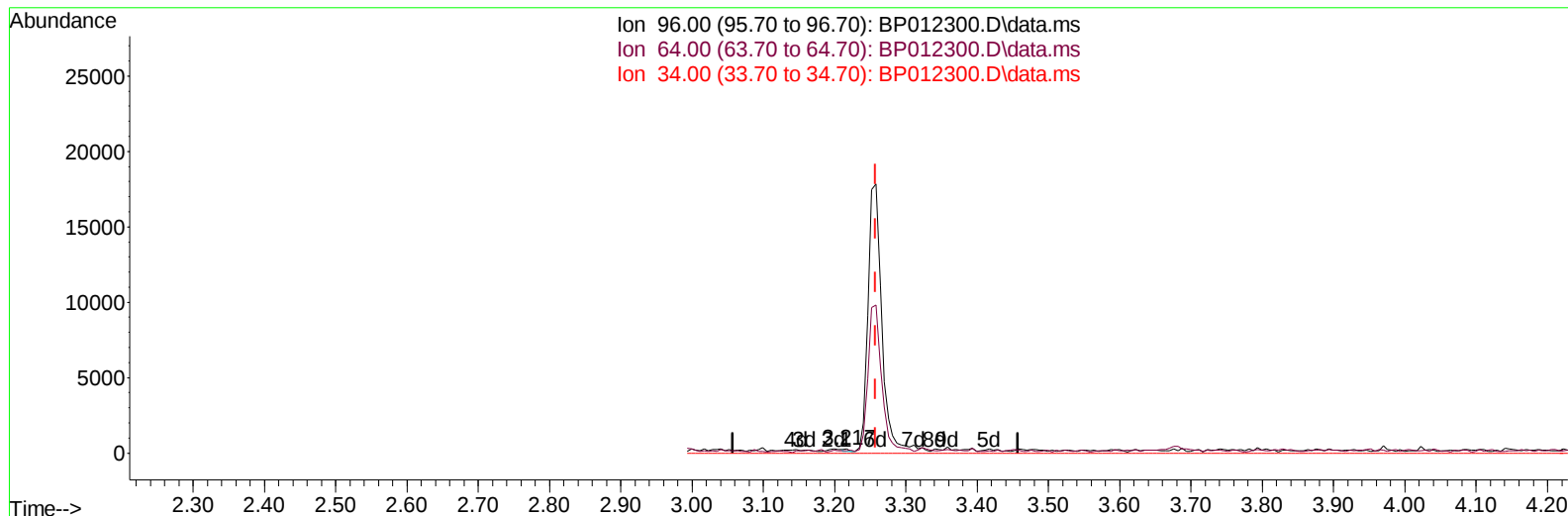
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012300.D
 Acq On : 25 Oct 2022 00:13
 Operator : CG/JU
 Sample : N5015-14
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BF0

Manual IntegrationsAPPROVED

Quant Time: Oct 25 07:06:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 22 01:24:49 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/25/2022
 Supervised By :mohammad ahmed 10/27/2022



TIC: BP012300.D\data.ms

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

3.217min (-0.041) 0.01 ng/uL

response 71

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	63.64
34.00	0.00	0.00
0.00	0.00	0.00

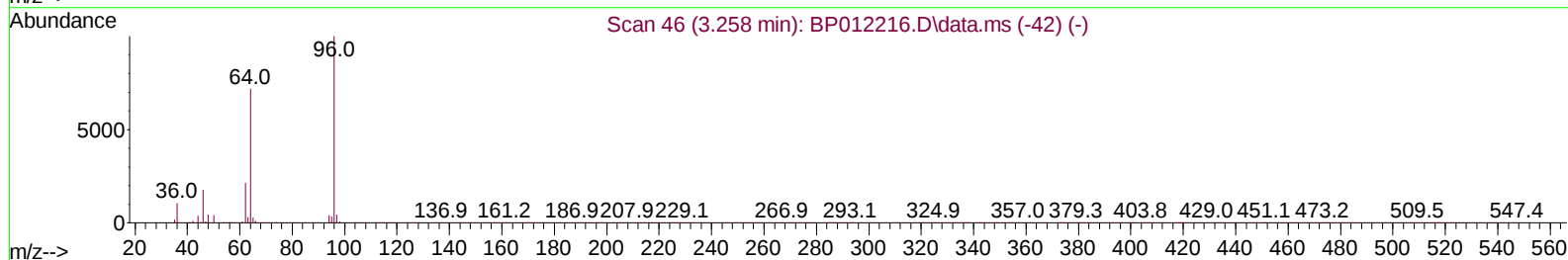
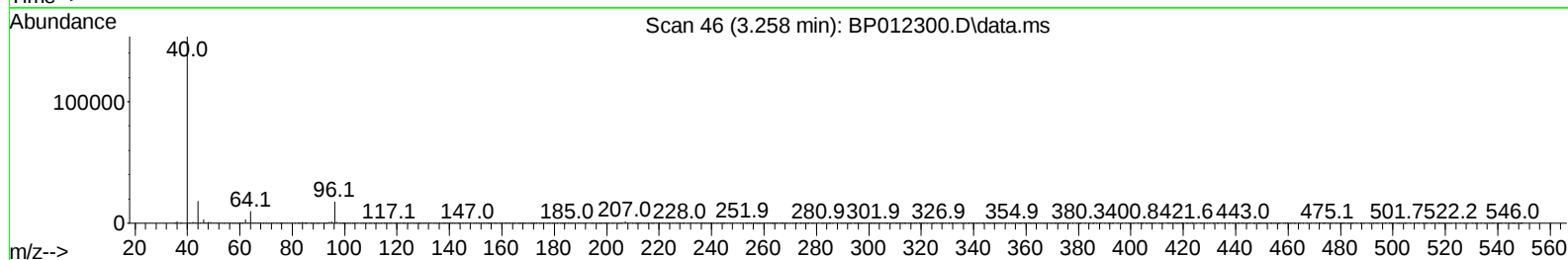
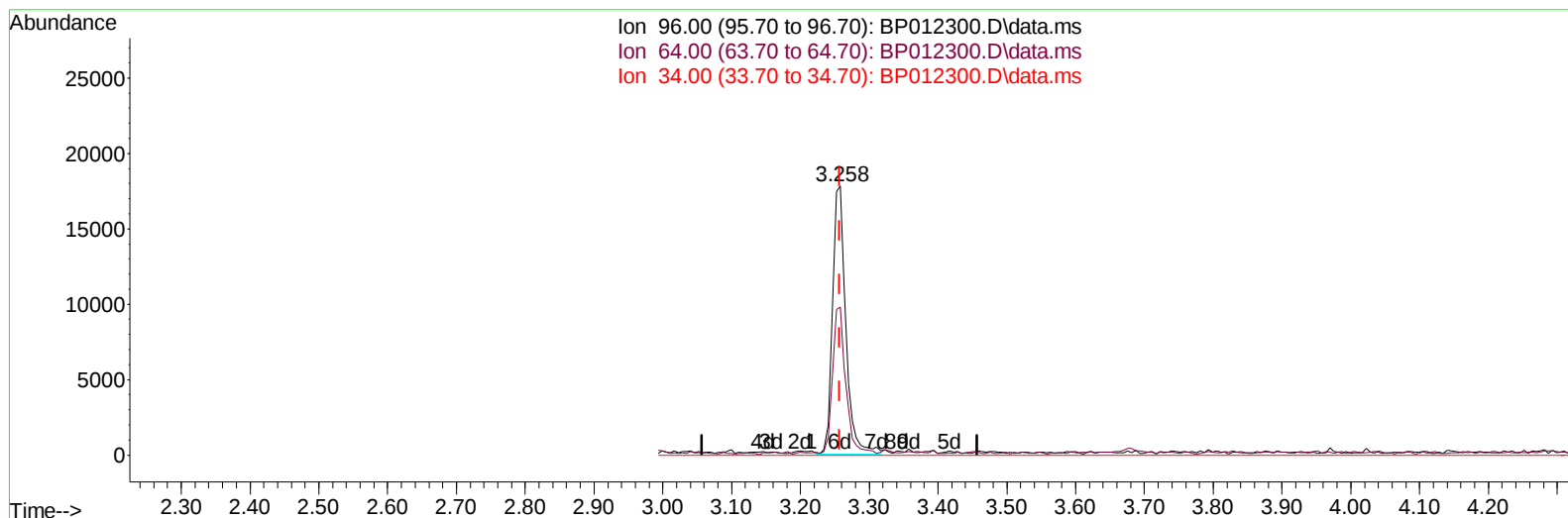
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TIC: BP012300.D\data.ms

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

3.258min (+ 0.000) 2.75 ng/uL m

response 24019

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	55.15#
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	7.816	152	353835	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	1528881	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	798533	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1418177	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1438685	20.000	ng/ul	0.00
88) Perylene-d12	23.716	264	1539861	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	24019m	2.747	ng/uL	0.00
4) Pyridine-d5	3.676	84	171390	6.803	ng/ul	0.00
7) Phenol-d5	6.987	99	1521337	49.803	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	953908	51.491	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	1128962	48.554	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	1151531	48.503	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	587552	51.918	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	597901	48.968	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	1067167	46.595	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	1165037	35.803	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	2947841	53.069	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	3427953	54.065	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	375903	36.122	ng/ul	0.00
60) Fluorene-d10	15.463	176	2434650	50.947	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	192326	23.839	ng/ul	0.00
73) Anthracene-d10	17.328	188	3535329	56.872	ng/ul	0.00
81) Pyrene-d10	19.563	212	4101393	50.071	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.563	264	4572389	59.211	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	41256	1.291	ng/ul#	81
16) Acetophenone	8.652	105	45288	1.235	ng/ul	96

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

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