

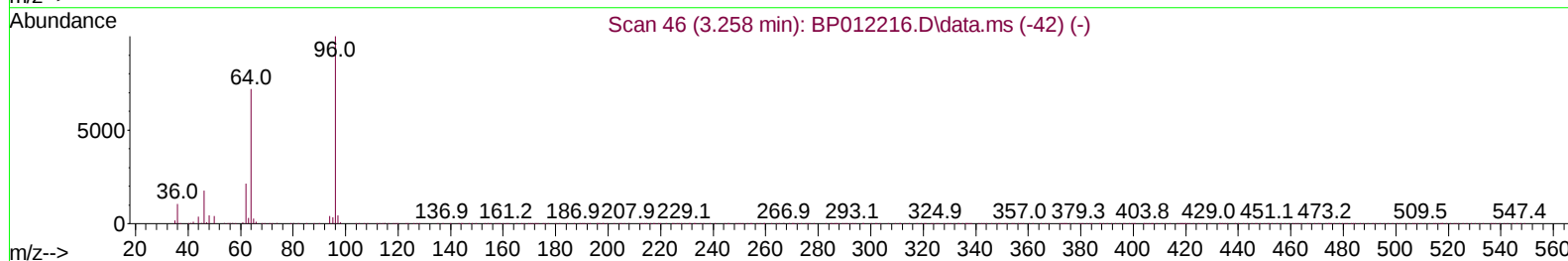
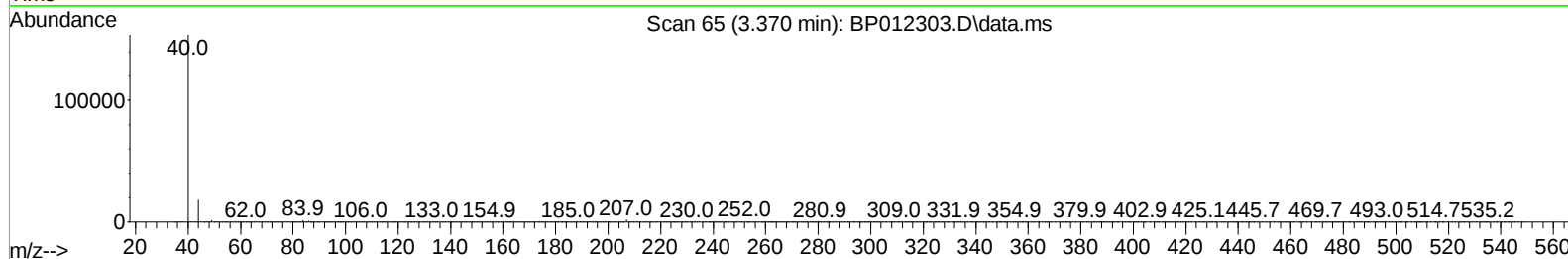
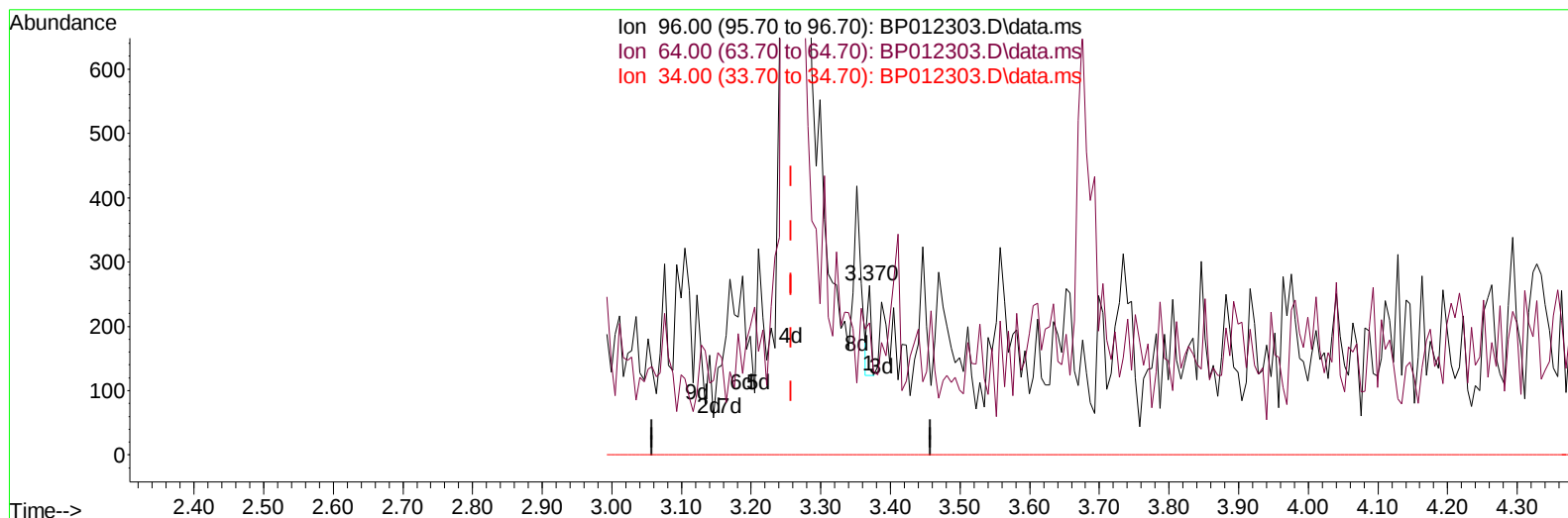
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012303.D
 Acq On : 25 Oct 2022 01:56
 Operator : CG/JU
 Sample : N5015-08
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BG0

Manual IntegrationsAPPROVED

Quant Time: Oct 25 07:16:26 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: BP012303.D\data.ms

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

3.370min (+ 0.112) 0.01 ng/uL

response 49

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	77.65
34.00	0.00	0.00
0.00	0.00	0.00

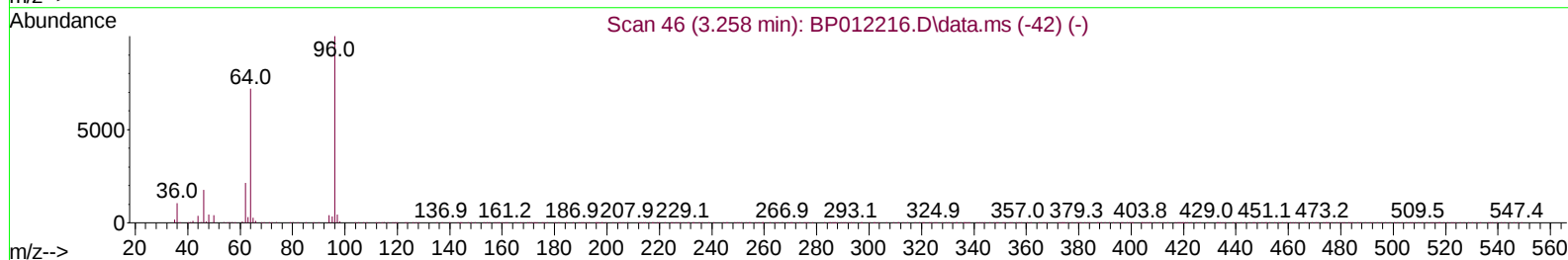
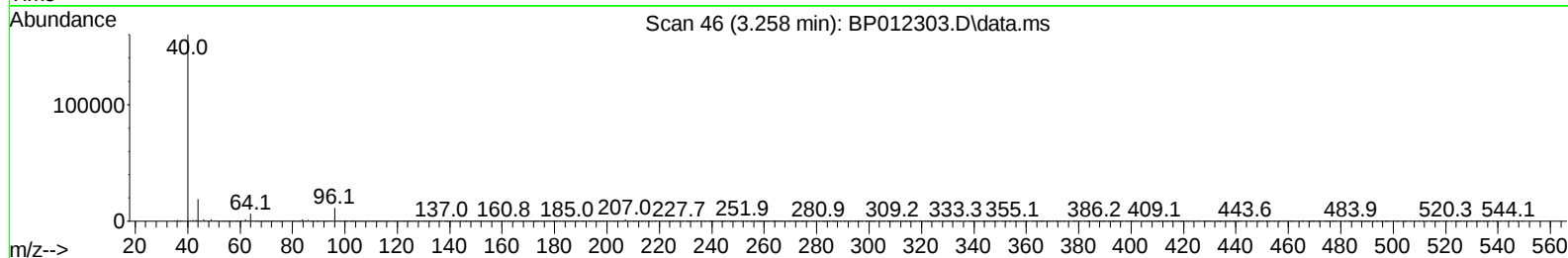
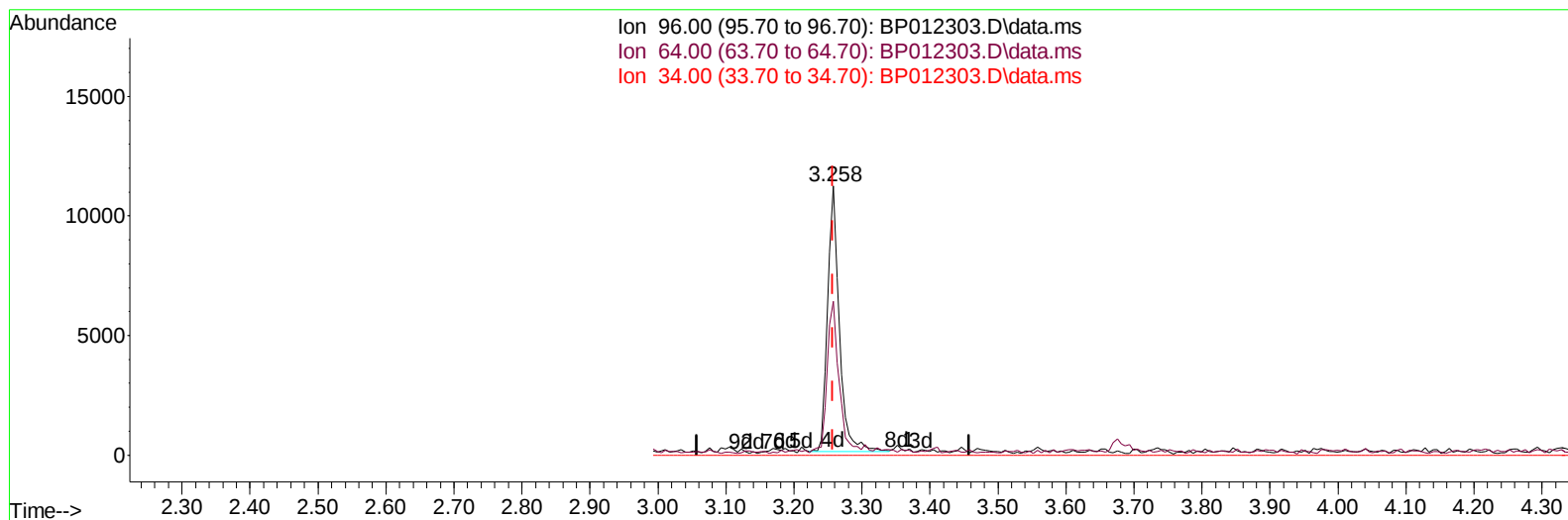
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TIC: BP012303.D\data.ms

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

3.258min (+ 0.000) 1.85 ng/uL m

response 13385

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	57.10#
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	293239	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	1313845	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	762464	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1571353	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1590630	20.000	ng/ul	0.00
88) Perylene-d12	23.710	264	1674701	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	13385m	1.847	ng/uL	0.00
4) Pyridine-d5	3.675	84	208632	9.992	ng/ul	0.00
7) Phenol-d5	6.987	99	894643	35.339	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	578783	37.698	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	696488	36.144	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	726675	36.933	ng/ul	-0.01
21) Nitrobenzene-d5	8.987	128	355706	36.576	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	350709	33.424	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	676582	34.376	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	964536	34.493	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1949241	36.752	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	2391466	39.502	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	237622	23.914	ng/ul	0.00
60) Fluorene-d10	15.463	176	1731049	37.938	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	104414	11.680	ng/ul	0.00
73) Anthracene-d10	17.322	188	2644073	38.388	ng/ul	0.00
81) Pyrene-d10	19.563	212	3066213	33.857	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	3201442	38.120	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	37522	1.417	ng/ul#	83

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

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