

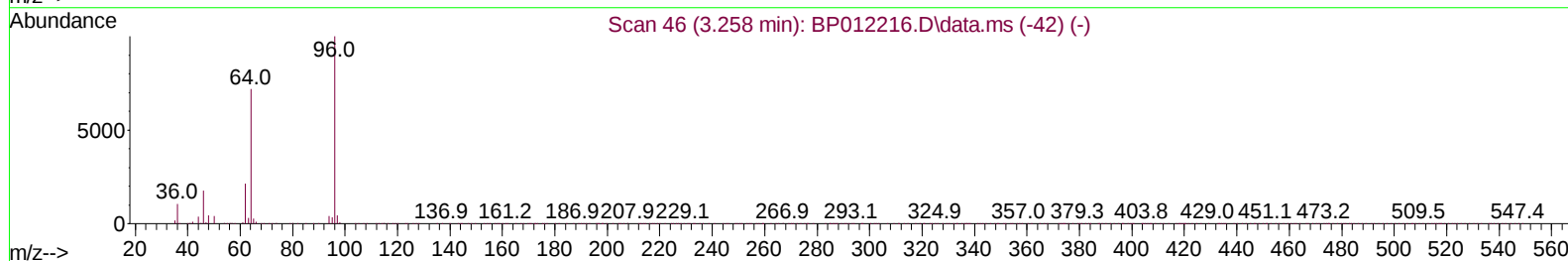
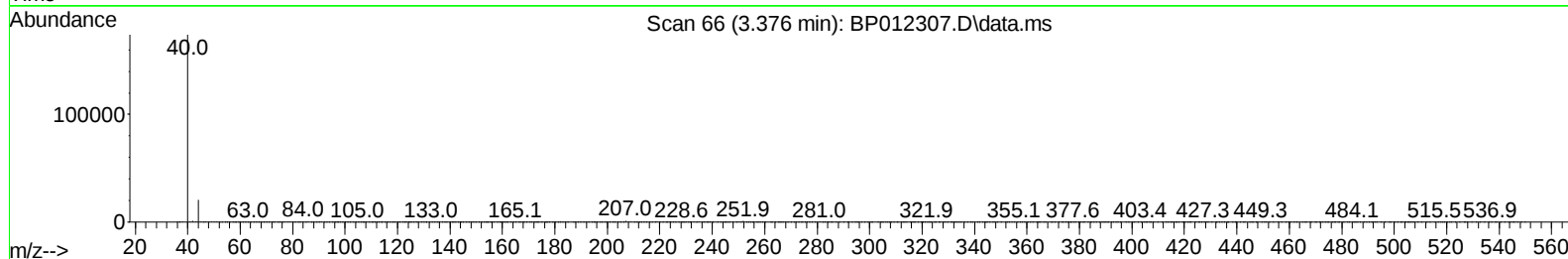
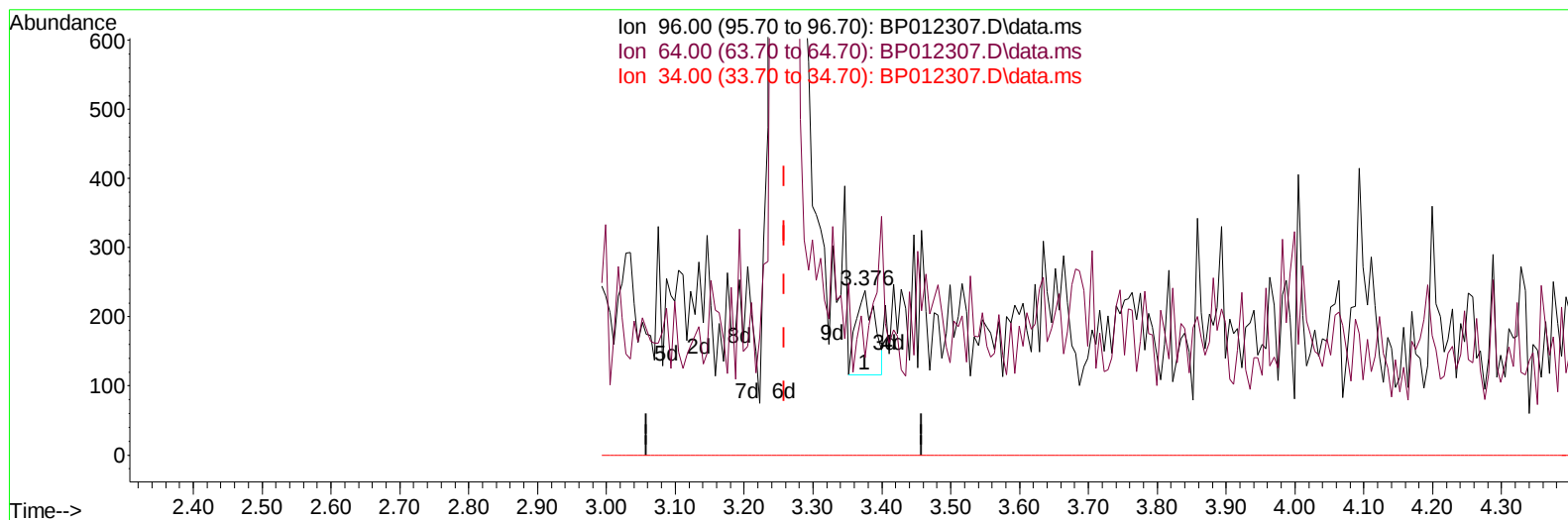
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012307.D
 Acq On : 25 Oct 2022 04:13
 Operator : CG/JU
 Sample : N5015-16
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BG1

Manual IntegrationsAPPROVED

Quant Time: Oct 25 07:26:17 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: BP012307.D\data.ms

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

3.376min (+ 0.118) 0.03 ng/uL

response 228

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	59.24
34.00	0.00	0.00
0.00	0.00	0.00

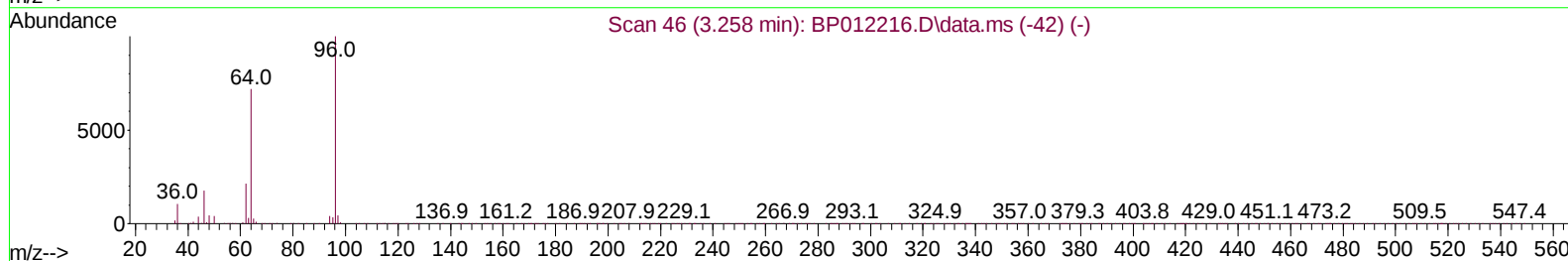
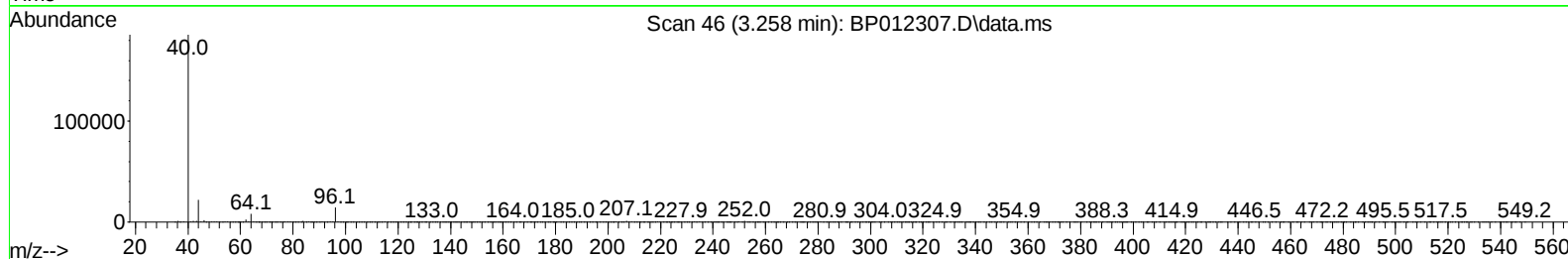
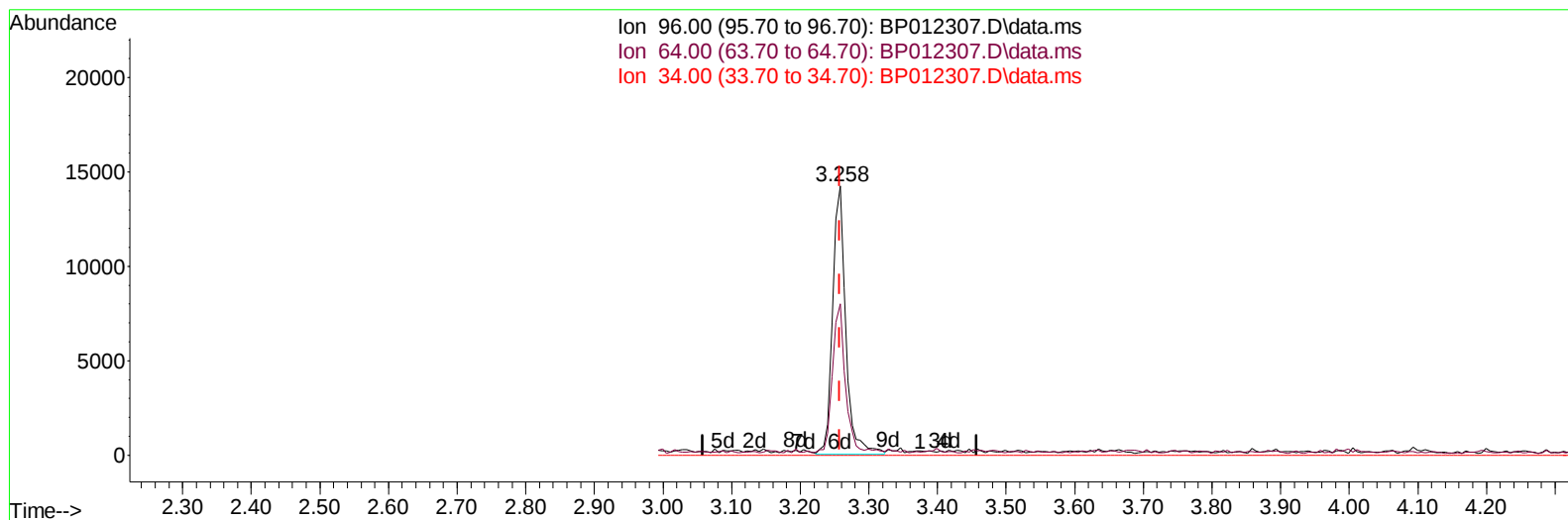
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TIC: BP012307.D\data.ms

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

3.258min (+ 0.000) 2.15 ng/uL m

response 18710

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	56.26#
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	351482	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	1511095	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	772467	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	1369673	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1404786	20.000	ng/ul	0.00
88) Perylene-d12	23.716	264	1512404	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	18710m	2.154	ng/uL	0.00
4) Pyridine-d5	3.676	84	92279	3.687	ng/ul	0.00
7) Phenol-d5	6.987	99	1151819	37.959	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	769601	41.821	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	882160	38.194	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	857157	36.346	ng/ul	-0.01
21) Nitrobenzene-d5	8.987	128	445629	39.841	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	407897	33.800	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	781652	34.531	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	1057480	32.880	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	2158517	40.171	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	2625755	42.810	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	181137	17.993	ng/ul	0.00
60) Fluorene-d10	15.463	176	1840971	39.824	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	76930	9.873	ng/ul	0.00
73) Anthracene-d10	17.322	188	2642882	44.021	ng/ul	0.00
81) Pyrene-d10	19.563	212	3080555	38.516	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	3319076	43.761	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.011	94	43246	1.362	ng/ul#	78

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

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