

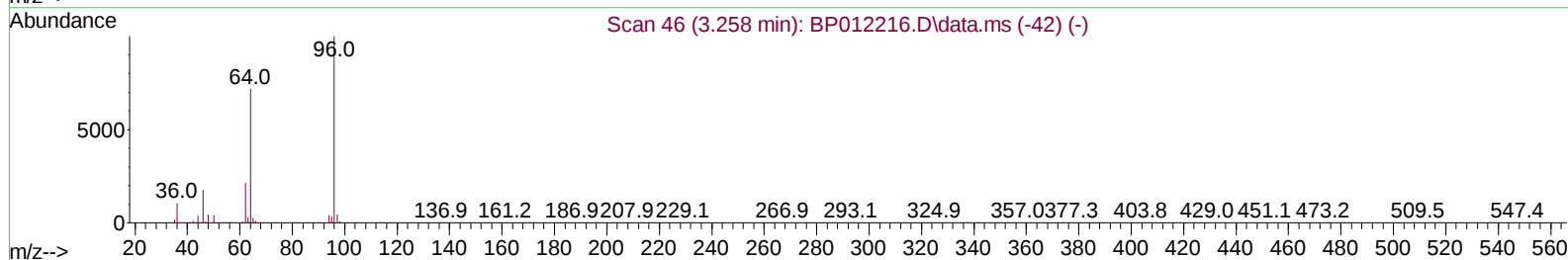
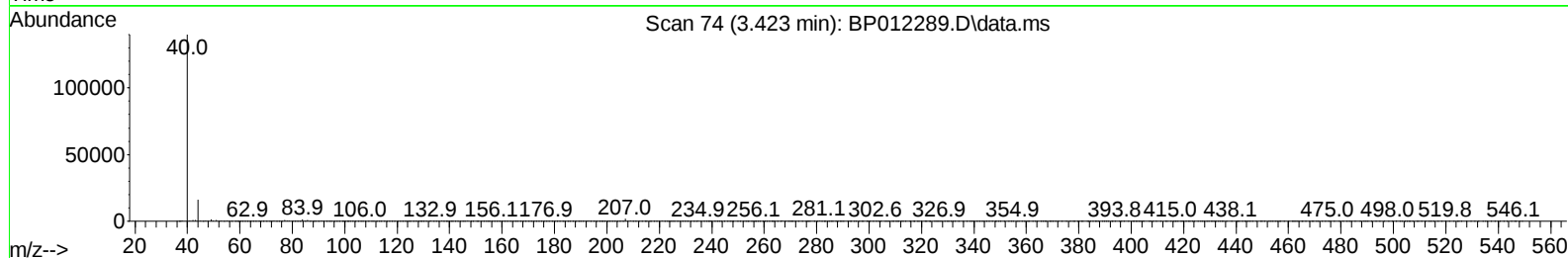
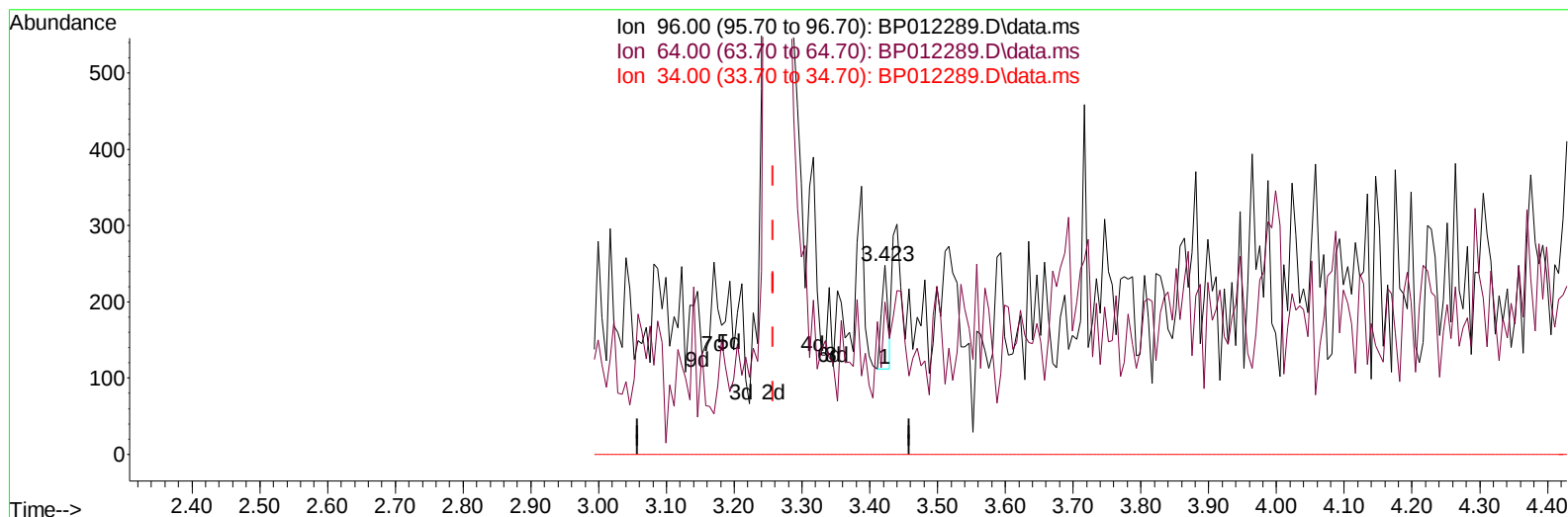
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012289.D
 Acq On : 24 Oct 2022 17:22
 Operator : CG/JU
 Sample : N5070-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BS5

Manual IntegrationsAPPROVED

Quant Time: Oct 25 06:25:29 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: BP012289.D\data.ms

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

3.423min (+ 0.165) 0.01 ng/uL

response 89

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	80.65
34.00	0.00	0.00
0.00	0.00	0.00

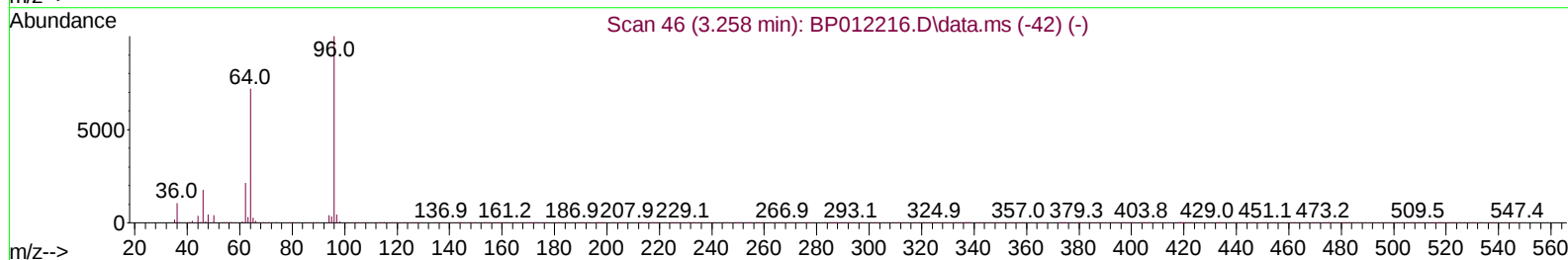
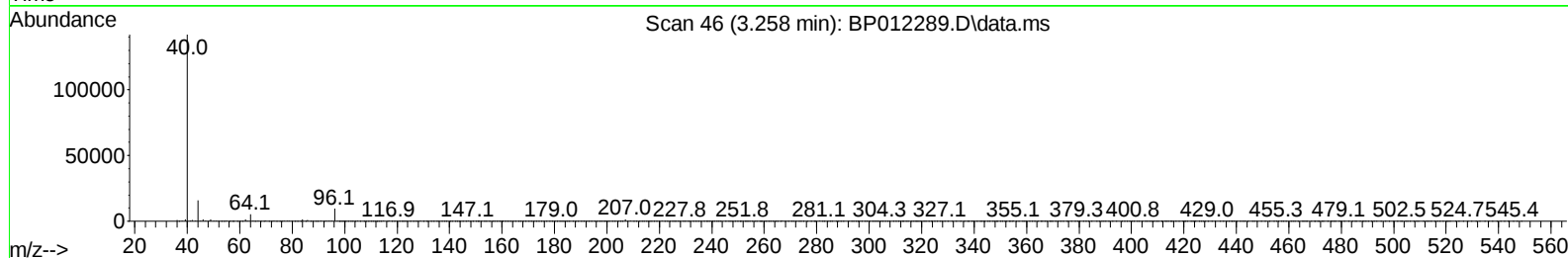
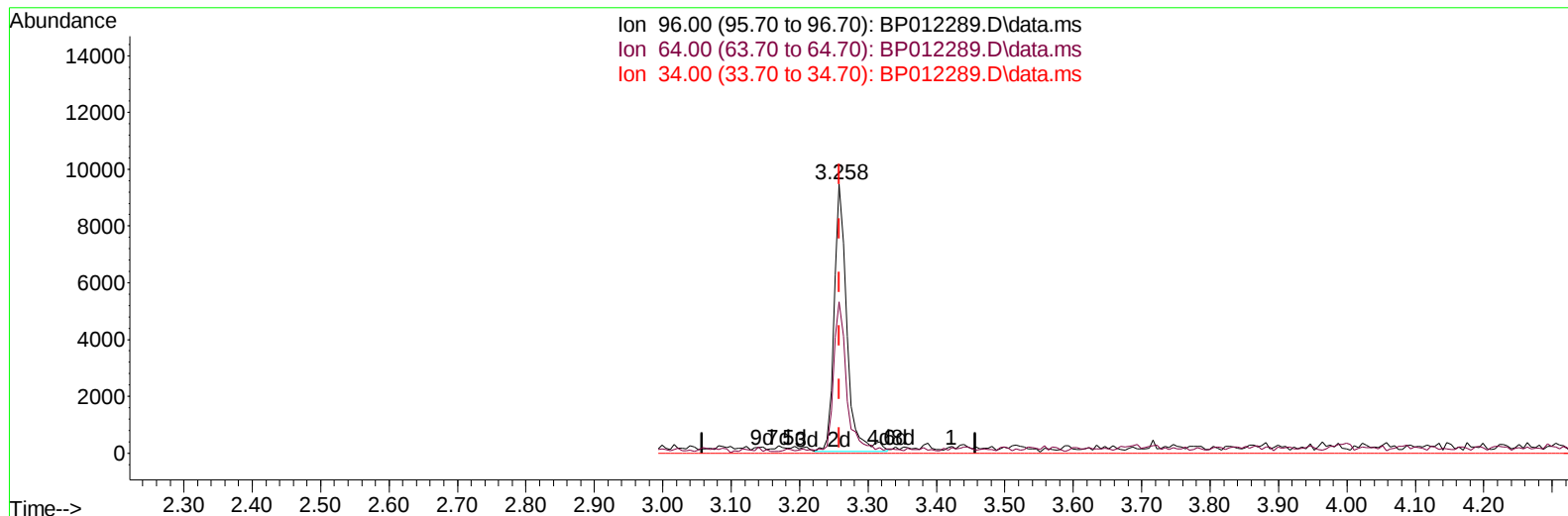
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TIC: BP012289.D\data.ms

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

3.258min (+ 0.000) 1.44 ng/uL m

response 12109

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	56.35#
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	339281	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1488277	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	783749	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1399596	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1399864	20.000	ng/ul	0.00
88) Perylene-d12	23.716	264	1515415	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	12109m	1.444	ng/uL	0.00
4) Pyridine-d5	3.687	84	50390	2.086	ng/ul	0.00
7) Phenol-d5	6.987	99	508595	17.364	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	386759	21.772	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	438257	19.657	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	362658	15.931	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	247536	22.470	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	226516	19.058	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	405162	18.173	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	419316	13.238	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1036443	19.011	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1429041	22.964	ng/ul	0.00
54) 4-Nitrophenol-d4	14.693	143	76282	7.468	ng/ul	0.00
60) Fluorene-d10	15.463	176	976938	20.829	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	41573	5.221	ng/ul	0.00
73) Anthracene-d10	17.328	188	1314528	21.427	ng/ul	0.00
81) Pyrene-d10	19.563	212	1508239	18.924	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.563	264	1562543	20.561	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.017	94	33142	1.082	ng/ul#	85

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

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