

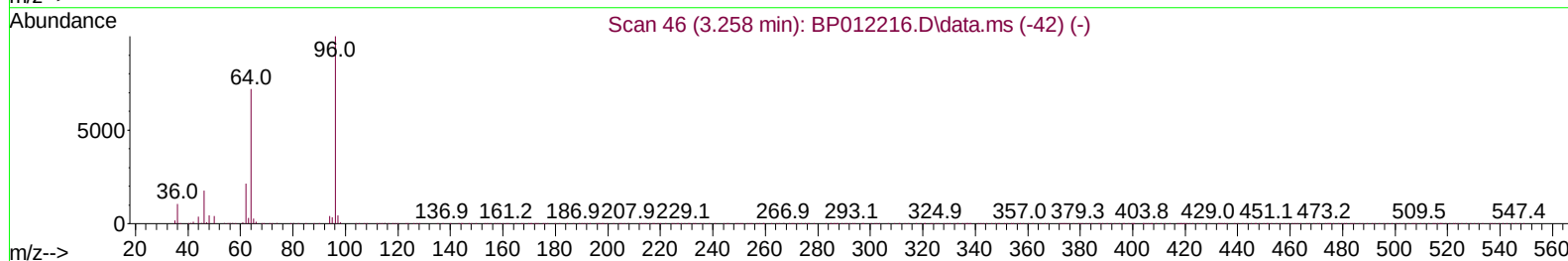
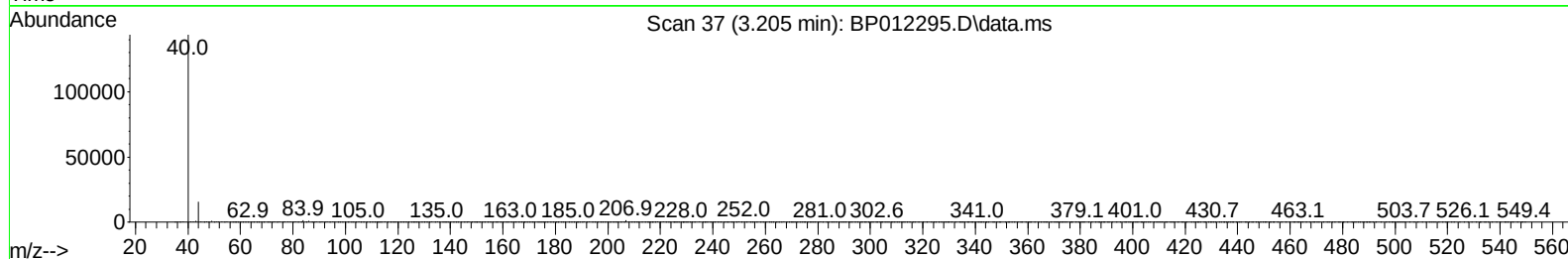
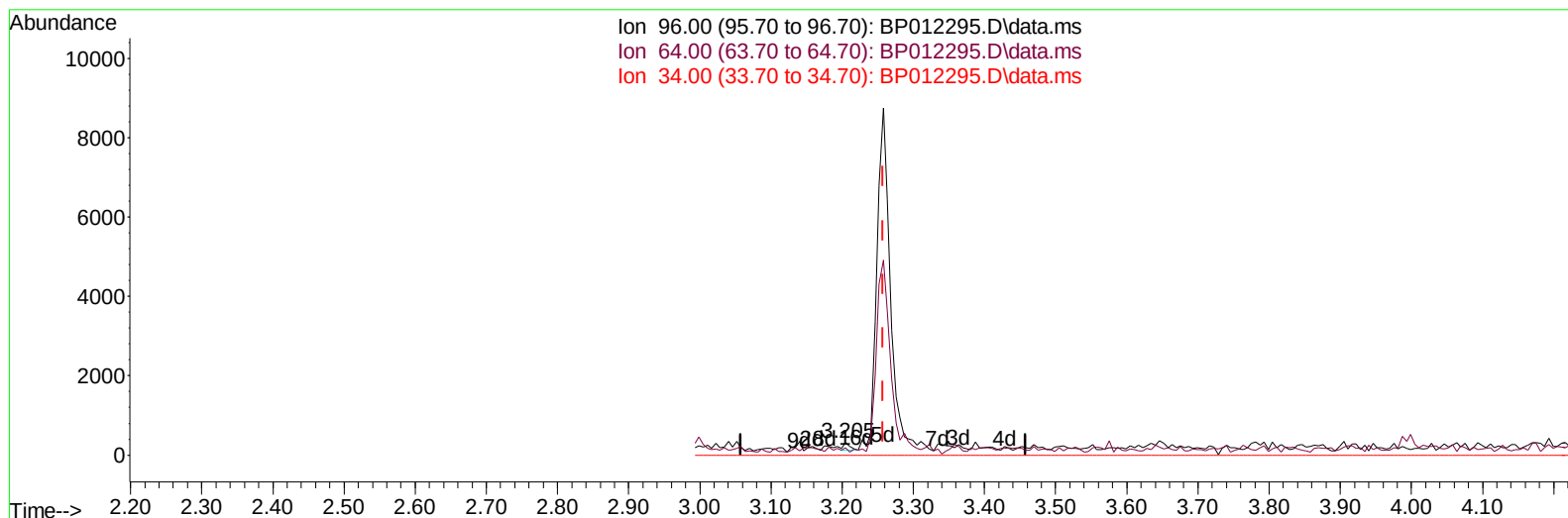
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012295.D
 Acq On : 24 Oct 2022 21:22
 Operator : CG/JU
 Sample : N5070-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BL1

Manual IntegrationsAPPROVED

Quant Time: Oct 25 06:50:12 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: BP012295.D\data.ms

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

3.205min (-0.053) 0.01 ng/uL

response	103	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	59.93
34.00	0.00	0.00
0.00	0.00	0.00

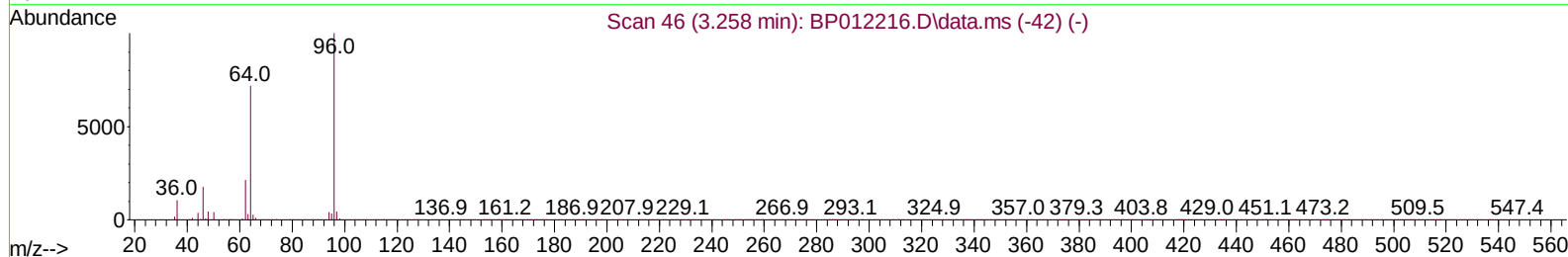
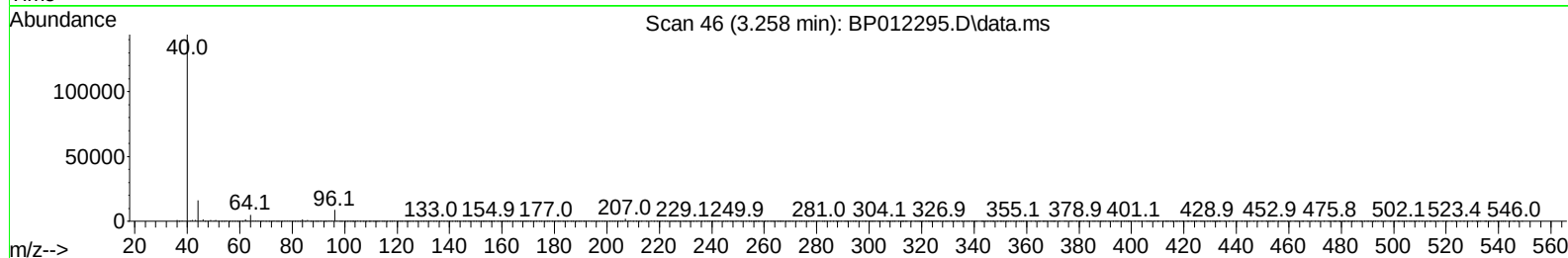
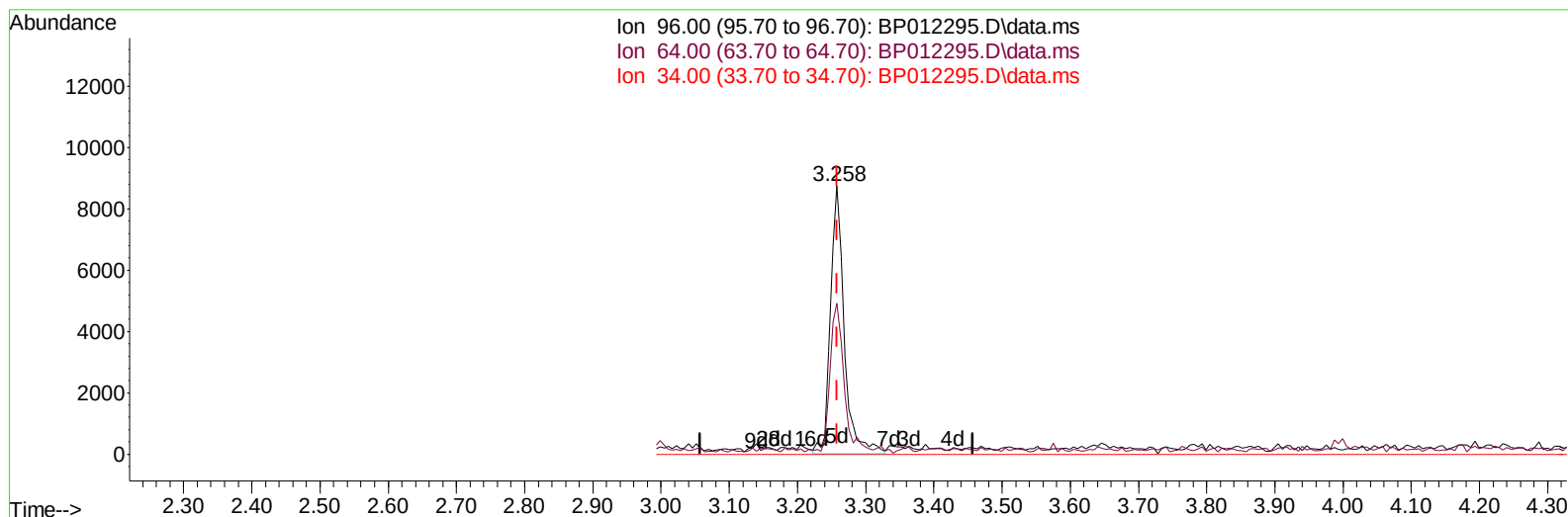
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TIC: BP012295.D\data.ms

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

3.258min (-0.000) 1.25 ng/uL m

response 12053

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	56.15#
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	390791	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1696945	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	930381	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1645558	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1414865	20.000	ng/ul	0.00
88) Perylene-d12	23.716	264	1557542	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	12053m	1.248	ng/uL	0.00
4) Pyridine-d5	3.705	84	5914	0.213	ng/ul	0.02
7) Phenol-d5	6.987	99	539767	15.999	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	404819	19.785	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	455417	17.734	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	224789	8.573	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	244999	19.505	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	241300	17.805	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	423999	16.679	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	322102	8.918	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1306177	20.182	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1541751	20.870	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	80999	6.680	ng/ul	0.00
60) Fluorene-d10	15.463	176	1100450	19.765	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	43267	4.622	ng/ul	0.00
73) Anthracene-d10	17.328	188	1413513	19.597	ng/ul	0.00
81) Pyrene-d10	19.563	212	1621115	20.124	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	1633205	20.909	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	41544	1.177	ng/ul#	88

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

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