

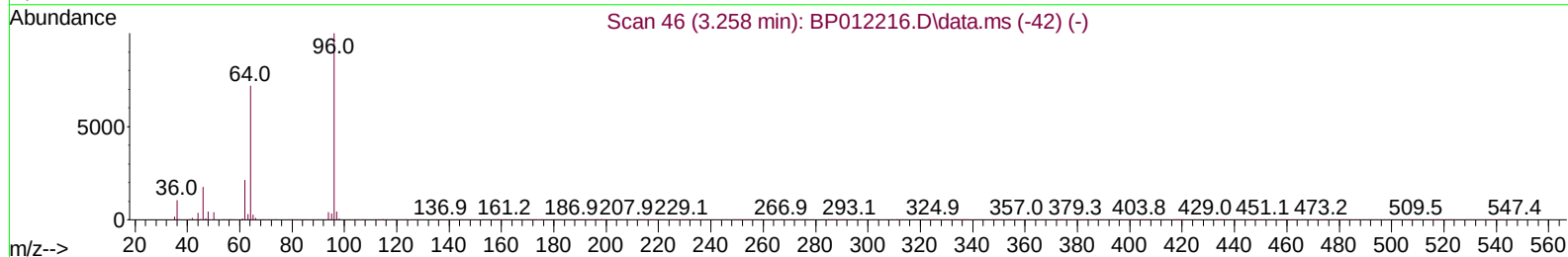
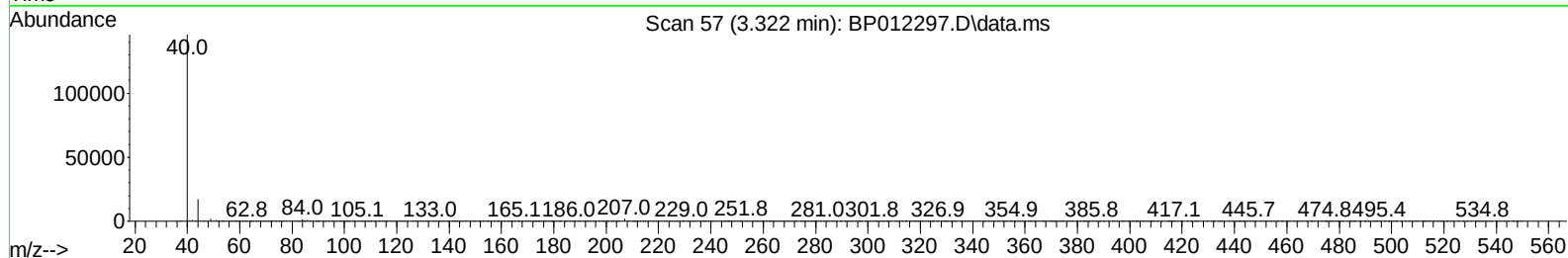
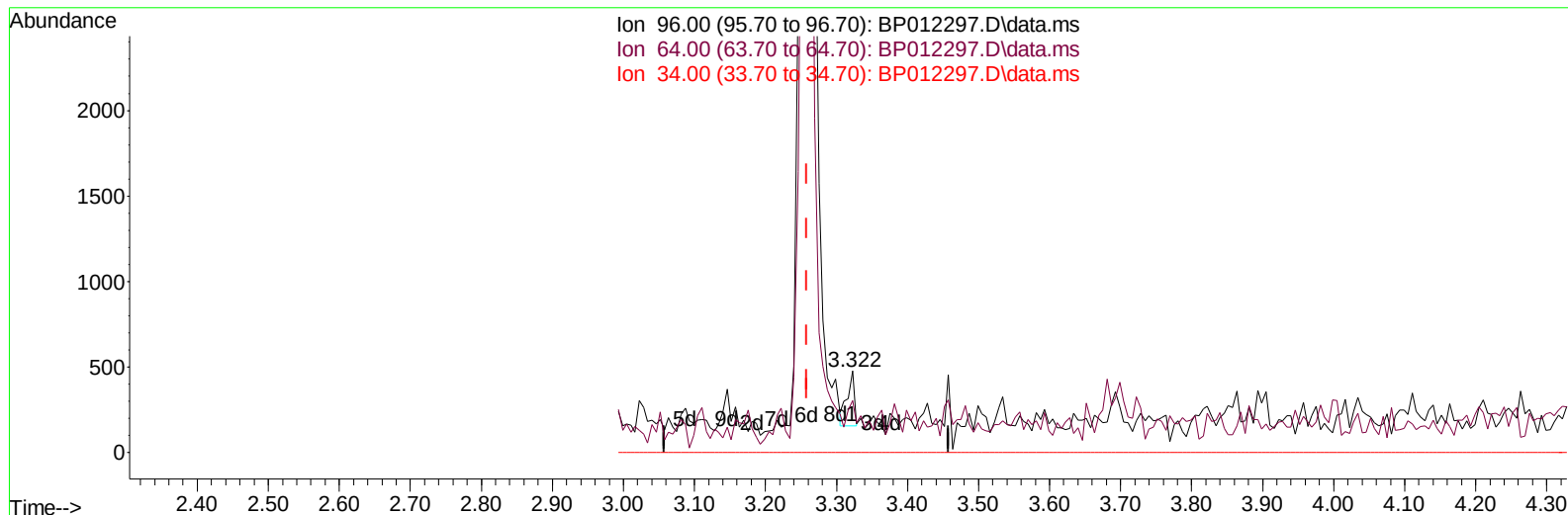
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012297.D
 Acq On : 24 Oct 2022 22:31
 Operator : CG/JU
 Sample : N5103-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BX1

Manual IntegrationsAPPROVED

Quant Time: Oct 25 06:58:42 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: BP012297.D\data.ms

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

3.322min (+ 0.065) 0.02 ng/uL

response 224

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	63.87
34.00	0.00	0.00
0.00	0.00	0.00

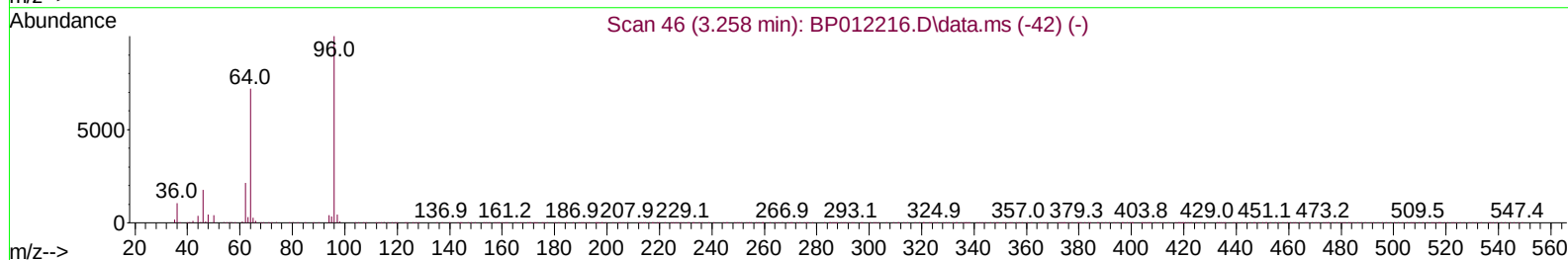
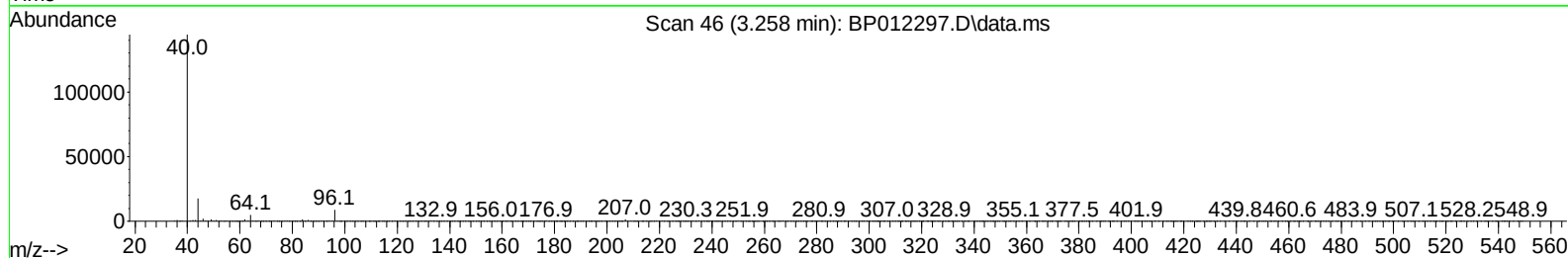
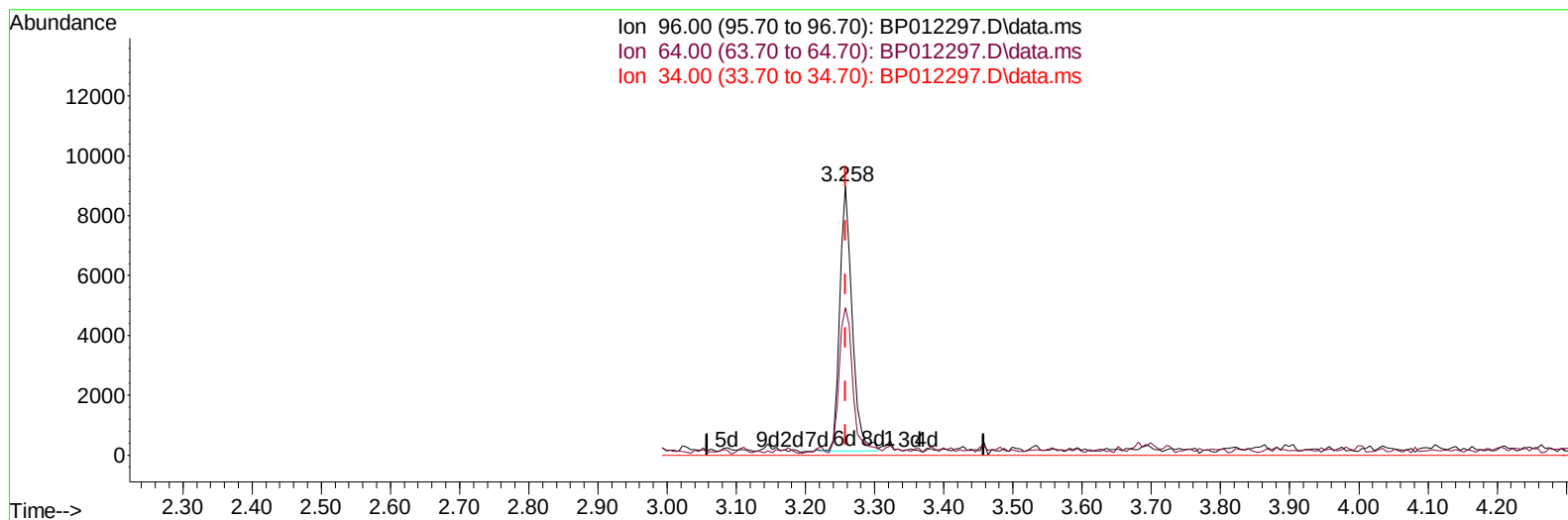
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(3) 1,4-Dioxane-d8 (S)

3.258min (-0.000) 1.02 ng/uL m

response 11131

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	54.81#
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	442640	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1936062	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	1040497	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1890502	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1816374	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1935943	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	11131m	1.018	ng/uL	0.00
4) Pyridine-d5	3.681	84	66147	2.099	ng/ul	0.00
7) Phenol-d5	6.987	99	190020	4.973	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	632075	27.274	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	564659	19.412	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	349465	11.767	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	392097	27.360	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	367329	23.757	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	613477	21.153	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	563847	13.683	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1993469	27.542	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	2315152	28.023	ng/ul	0.00
54) 4-Nitrophenol-d4	14.692	143	23778	1.754	ng/ul	0.00
60) Fluorene-d10	15.463	176	1664683	26.734	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	114838	10.678	ng/ul	0.00
73) Anthracene-d10	17.328	188	2567361	30.982	ng/ul	0.00
81) Pyrene-d10	19.563	212	2966776	28.688	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	3091213	31.840	ng/ul	0.00

Target Compounds Qvalue

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

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