

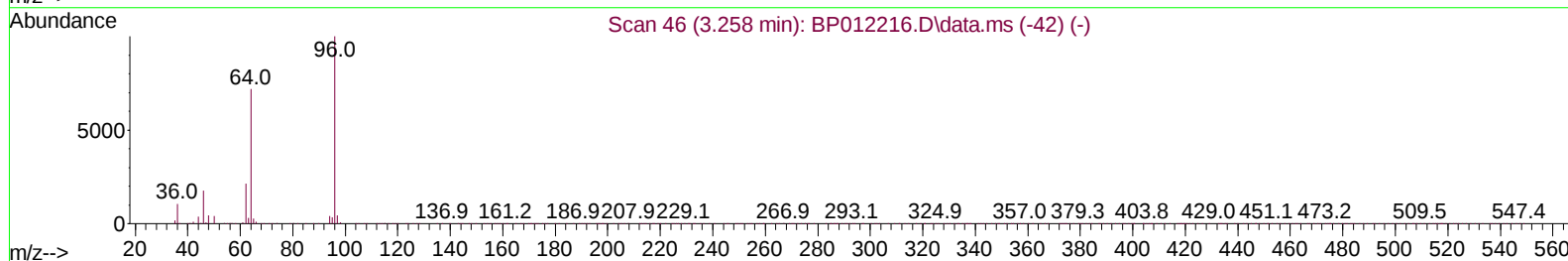
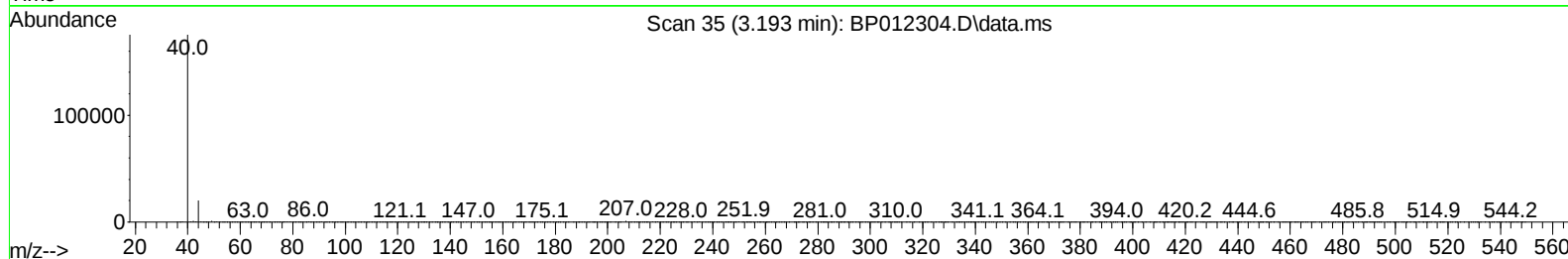
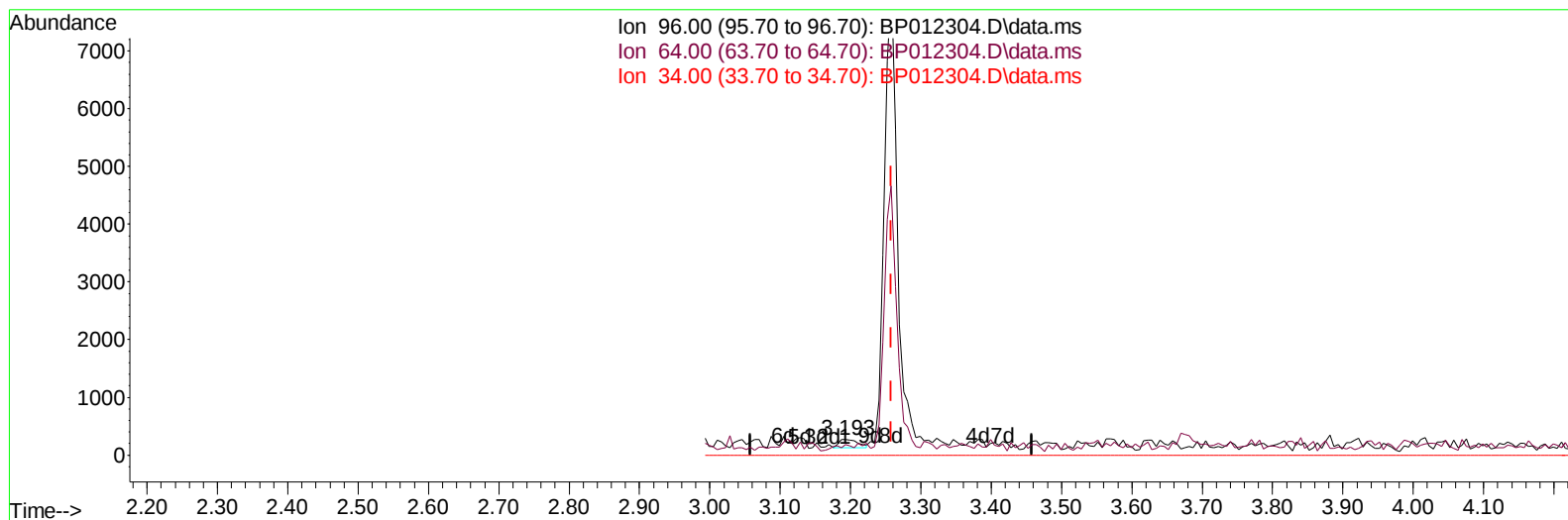
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012304.D
 Acq On : 25 Oct 2022 02:30
 Operator : CG/JU
 Sample : N5015-09
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BF8

Manual IntegrationsAPPROVED

Quant Time: Oct 25 07:20:09 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: BP012304.D\data.ms

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

3.193min (-0.065) 0.04 ng/uL

response 277

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	66.42
34.00	0.00	0.00
0.00	0.00	0.00

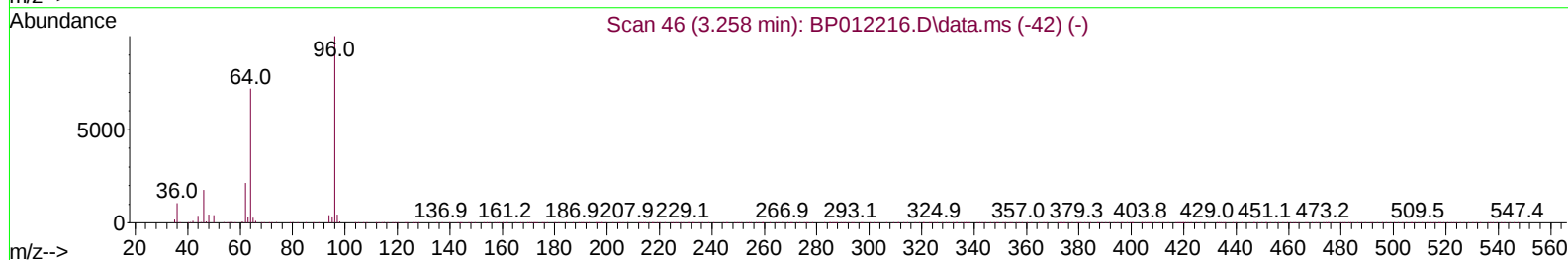
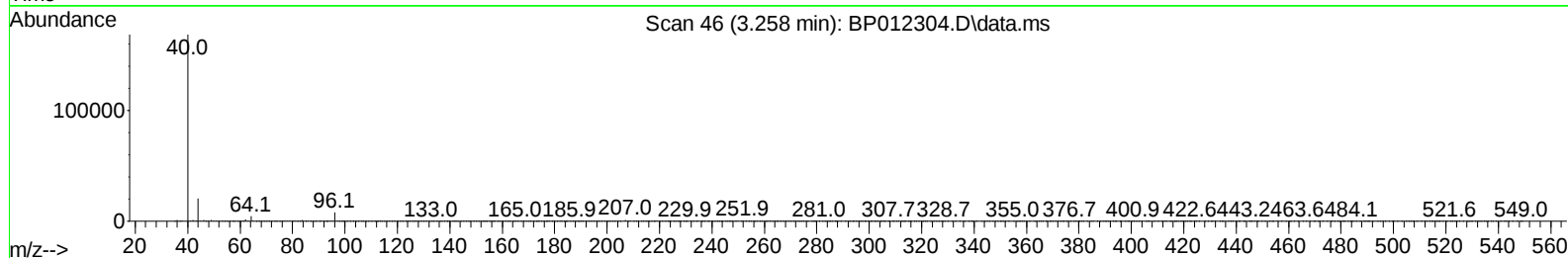
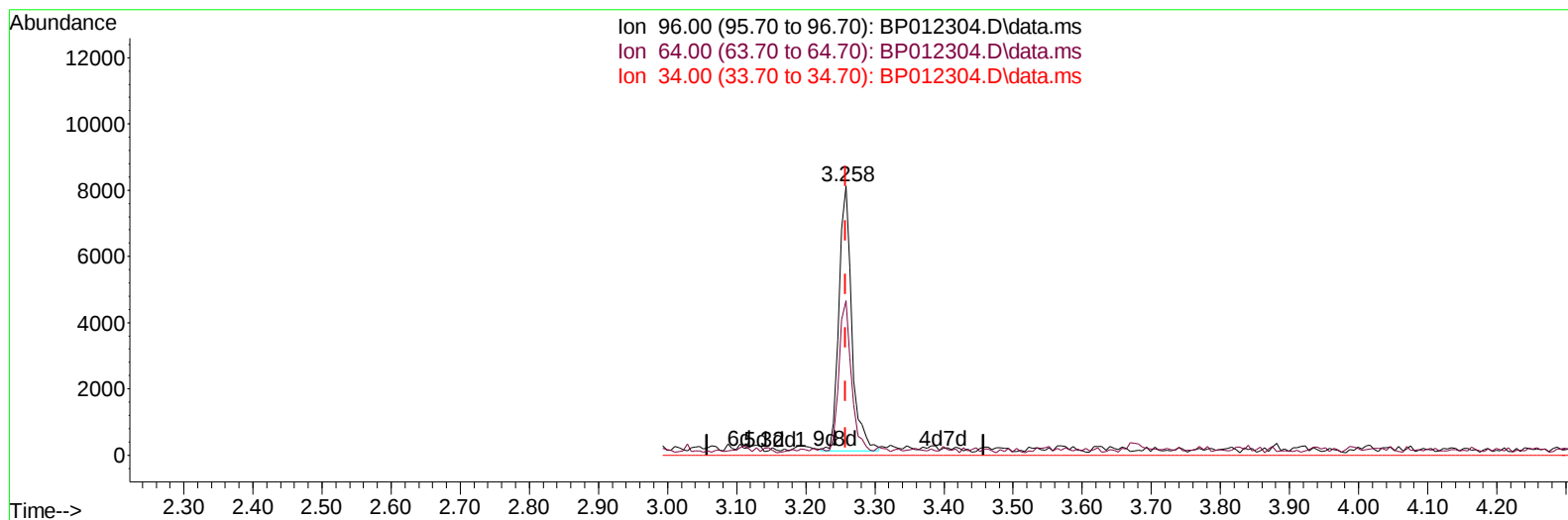
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(3) 1,4-Dioxane-d8 (S)

3.258min (-0.000) 1.35 ng/uL m

response 10339

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	57.36#
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	310421	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	1270307	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	659776	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	1267768	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1289221	20.000	ng/ul	0.00
88) Perylene-d12	23.710	264	1382873	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	10339m	1.348	ng/uL	0.00
4) Pyridine-d5	3.675	84	143996	6.515	ng/ul	0.00
7) Phenol-d5	6.987	99	581997	21.717	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	402351	24.756	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	459176	22.510	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	463461	22.251	ng/ul	-0.01
21) Nitrobenzene-d5	8.987	128	221990	23.609	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	202532	19.964	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	399848	21.012	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	571508	21.138	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1148632	25.027	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1413990	26.991	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	87115	10.132	ng/ul	0.00
60) Fluorene-d10	15.463	176	1023941	25.933	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	37063	5.139	ng/ul	0.00
73) Anthracene-d10	17.322	188	1531692	27.563	ng/ul	0.00
81) Pyrene-d10	19.563	212	1752060	23.869	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	1821147	26.261	ng/ul	-0.01

Target Compounds Qvalue

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

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