

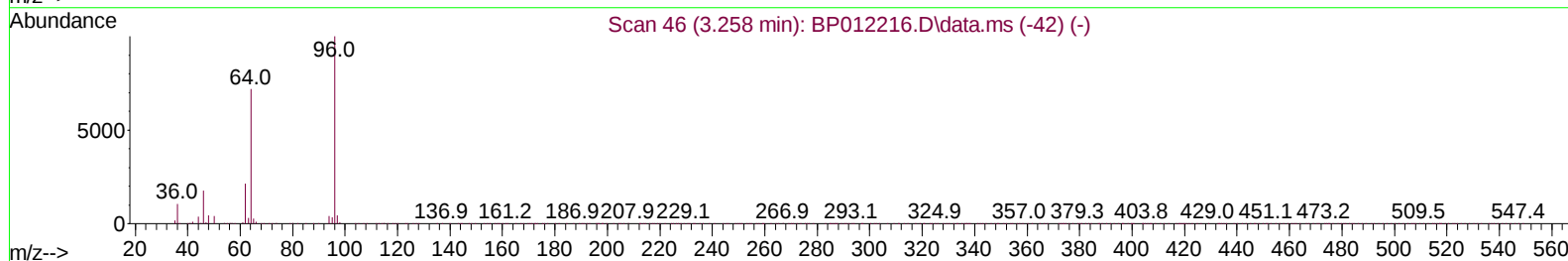
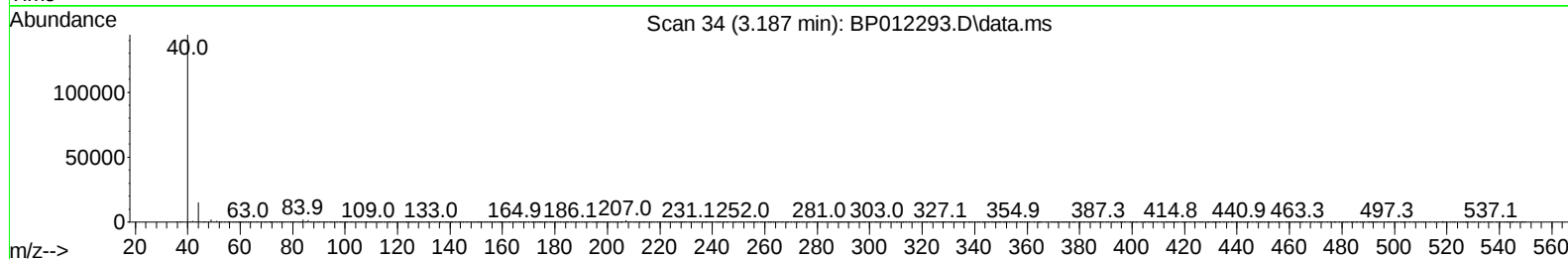
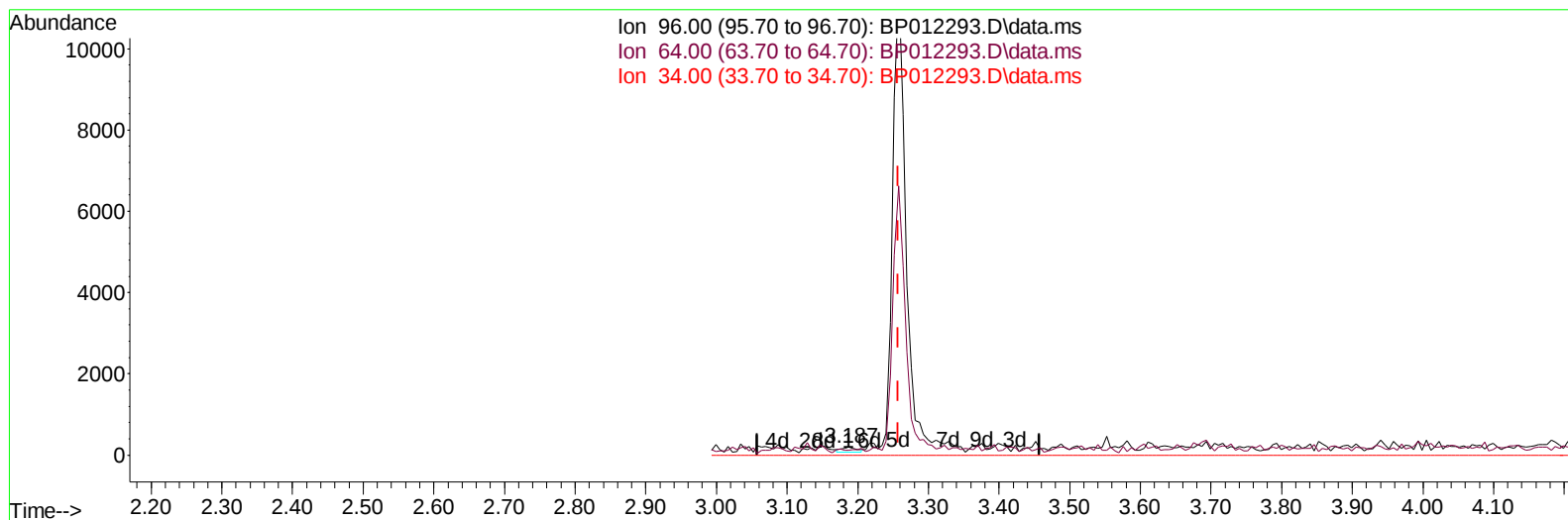
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012293.D
 Acq On : 24 Oct 2022 20:13
 Operator : CG/JU
 Sample : PB148414BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK414

Manual IntegrationsAPPROVED

Quant Time: Oct 25 06:43:49 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: BP012293.D\data.ms

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

3.187min (-0.071) 0.02 ng/uL

response	191	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	63.21
34.00	0.00	0.00
0.00	0.00	0.00

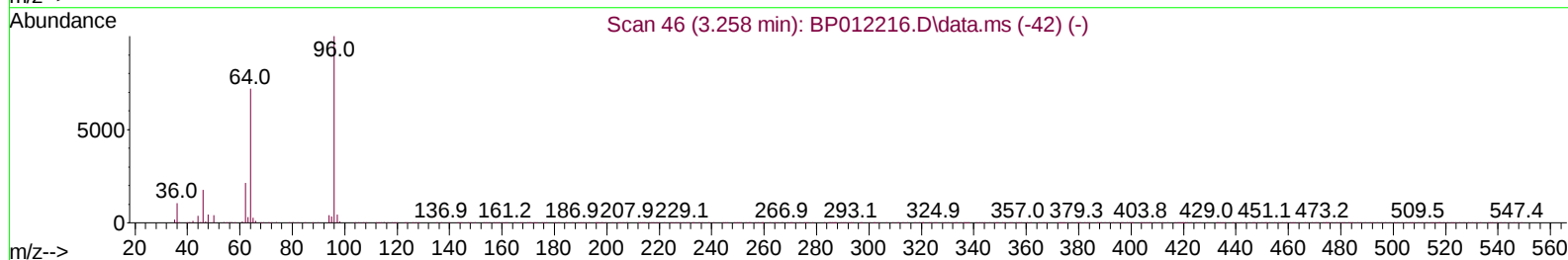
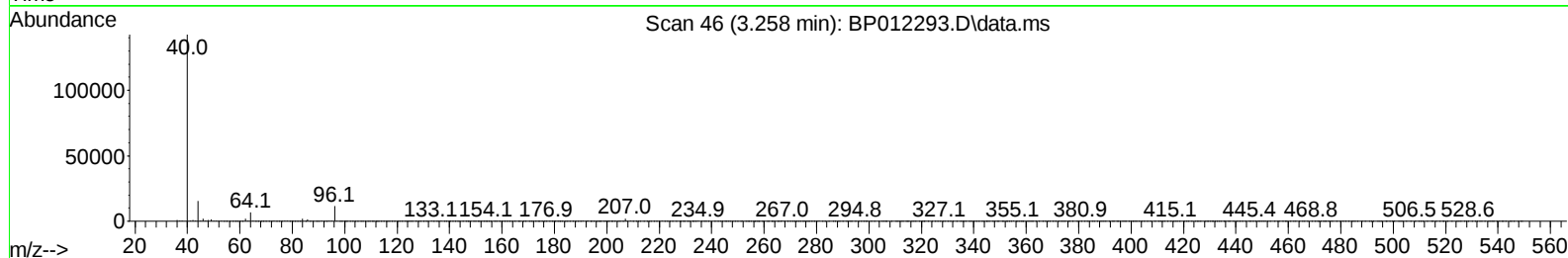
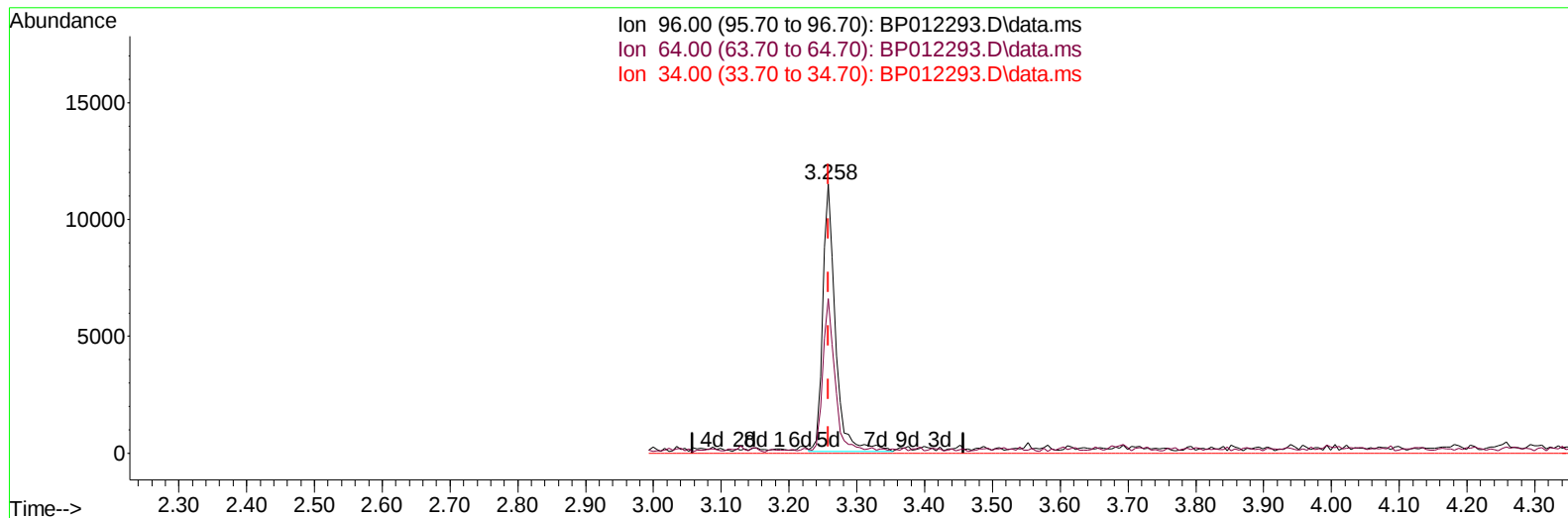
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TIC: BP012293.D\data.ms

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

3.258min (+ 0.000) 1.74 ng/uL m

response 15083

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.60	57.49#
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	351055	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1539768	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	878171	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1812964	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1806215	20.000	ng/ul	0.00
88) Perylene-d12	23.716	264	1916763	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	15083m	1.739	ng/uL	0.00
4) Pyridine-d5	3.681	84	91761	3.671	ng/ul	0.00
7) Phenol-d5	6.987	99	697515	23.015	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	445035	24.213	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	555677	24.088	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	557195	23.655	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	276245	24.237	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	283168	23.028	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	519544	22.524	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	641506	19.575	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1487443	24.350	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1788059	25.643	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	204149	17.838	ng/ul	0.00
60) Fluorene-d10	15.463	176	1305863	24.848	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	99524	9.650	ng/ul	0.00
73) Anthracene-d10	17.328	188	2124516	26.734	ng/ul	0.00
81) Pyrene-d10	19.563	212	2415557	23.489	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.563	264	2526127	26.280	ng/ul	0.00

Target Compounds Qvalue

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

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