

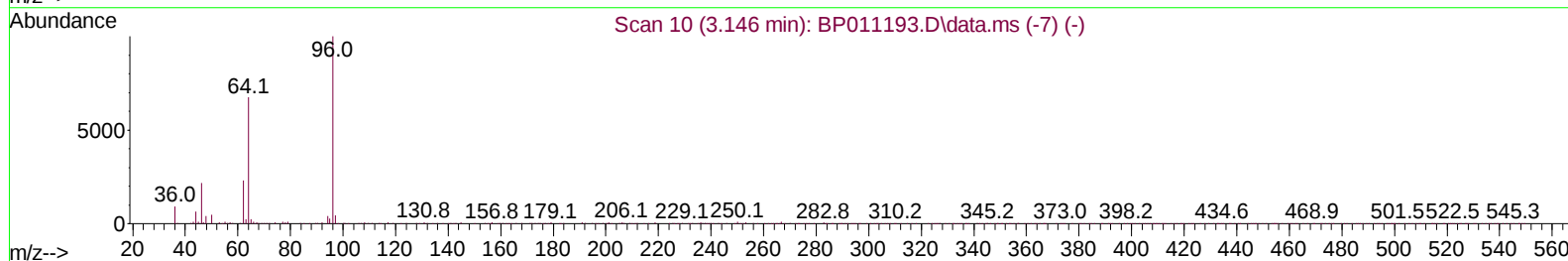
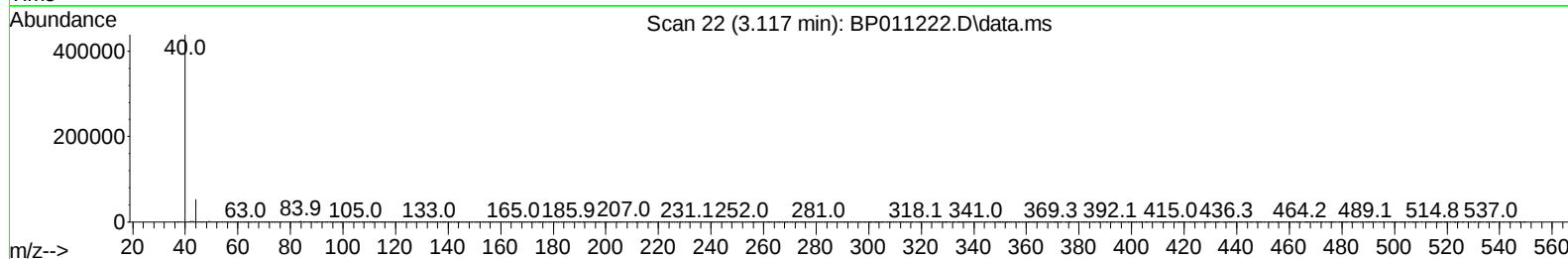
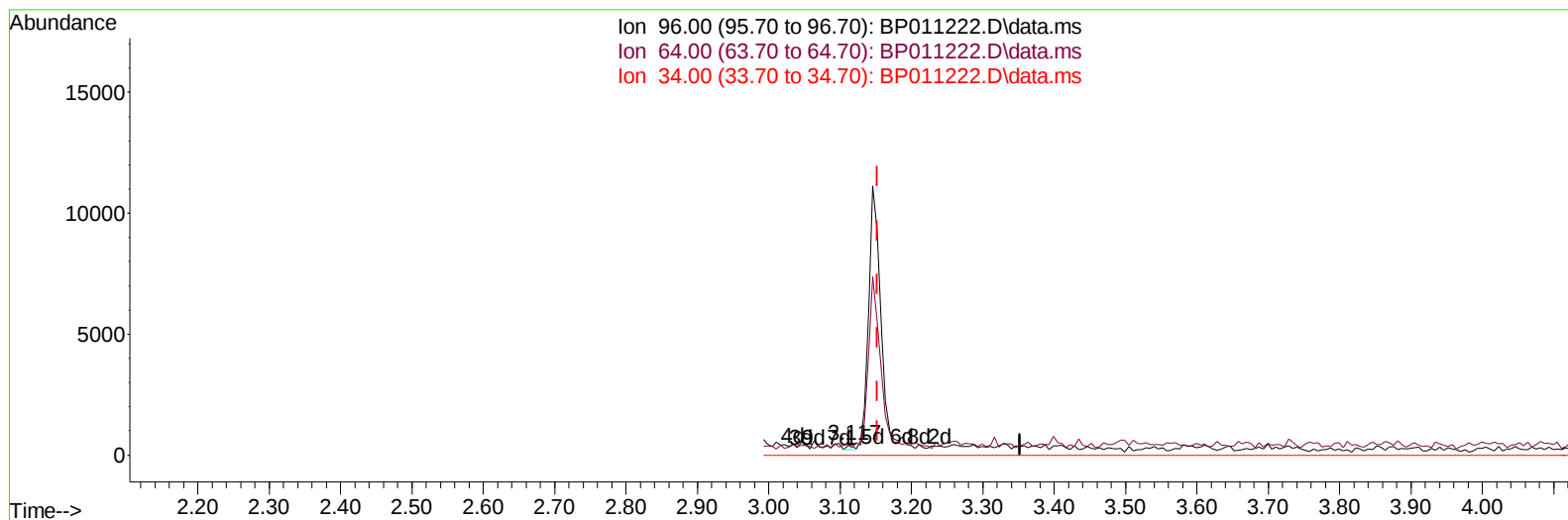
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011222.D
 Acq On : 02 Aug 2022 09:31
 Operator : CG/JU
 Sample : N3900-06
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0C98

Manual IntegrationsAPPROVED

Quant Time: Aug 02 16:09:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 16:05:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/02/2022
 Supervised By :mohammad ahmed 08/03/2022



TIC: BP011222.D\data.ms

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

3.117min (-0.035) 0.02 ng/uL

response 124

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	72.02
34.00	0.00	0.00
0.00	0.00	0.00

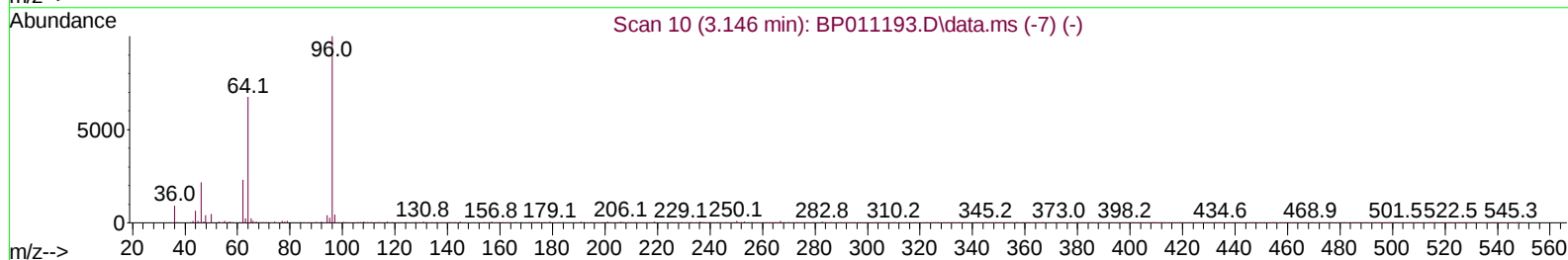
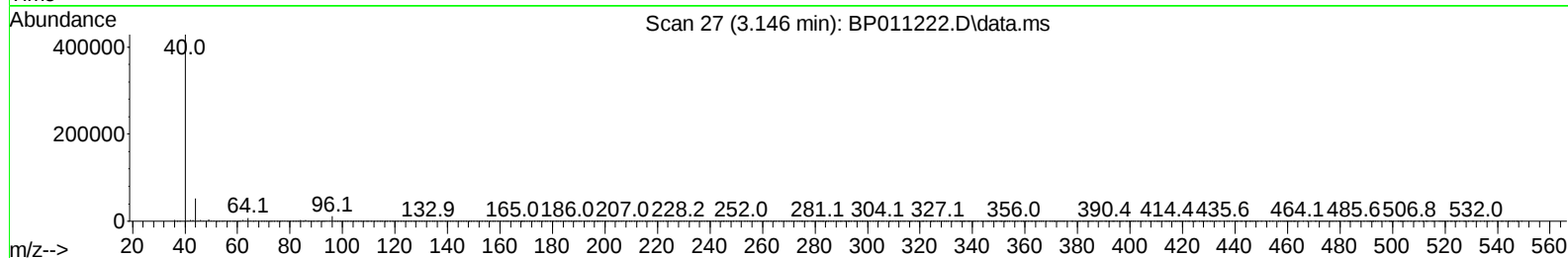
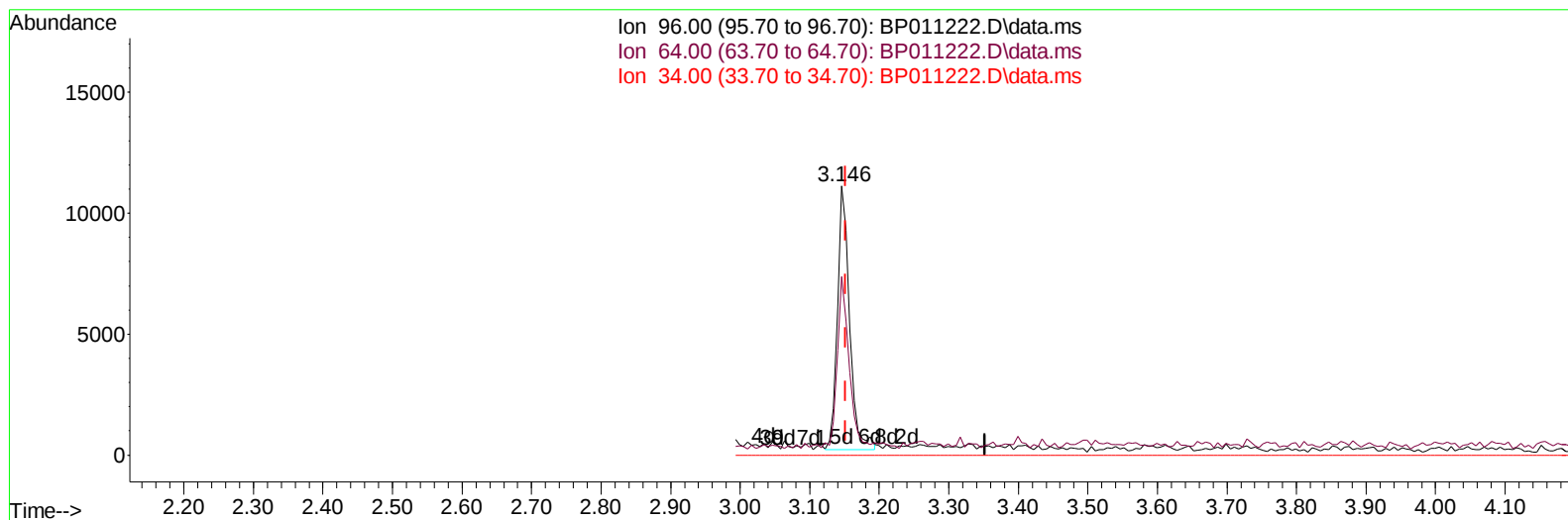
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TIC: BP011222.D\data.ms

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

3.146min (-0.006) 2.55 ng/uL m

response 13086

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	66.49#
34.00	0.00	0.00
0.00	0.00	0.00

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 Operator : CG/JU
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 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	235372	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	1003130	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.316	164	494131	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	942398	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	944200	20.000	ng/ul	# 0.00
88) Perylene-d12	23.557	264	1012812	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	13086m	2.546	ng/uL	0.00
4) Pyridine-d5	3.564	84	48746	3.353	ng/ul	0.00
7) Phenol-d5	6.828	99	96927	4.718	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	373365	25.701	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	304547	18.811	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113	183359	10.910	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	200853	25.764	ng/ul	0.00
24) 2-Nitrophenol-d4	9.528	143	176459	23.630	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	281374	21.381	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	426729	20.002	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	1022340	30.367	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1218447	28.801	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	11811	1.950	ng/ul	0.01
60) Fluorene-d10	15.316	176	833686	29.470	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	83421	17.560	ng/ul	0.00
73) Anthracene-d10	17.186	188	1458849	35.132	ng/ul	0.00
81) Pyrene-d10	19.445	212	1690533	33.320	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.410	264	1787214	35.527	ng/ul	0.00

Target Compounds Qvalue

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

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