

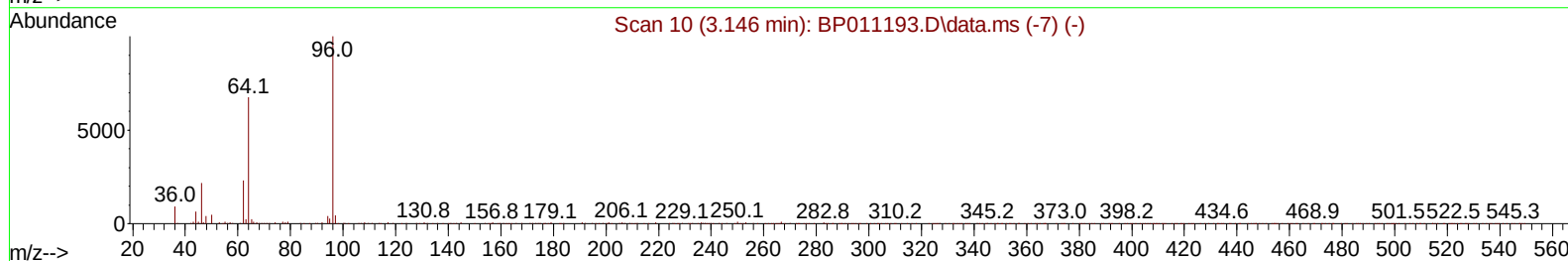
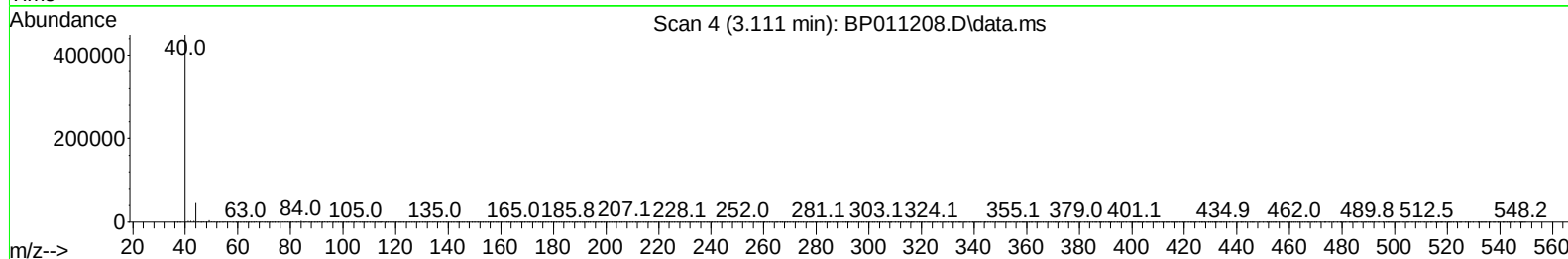
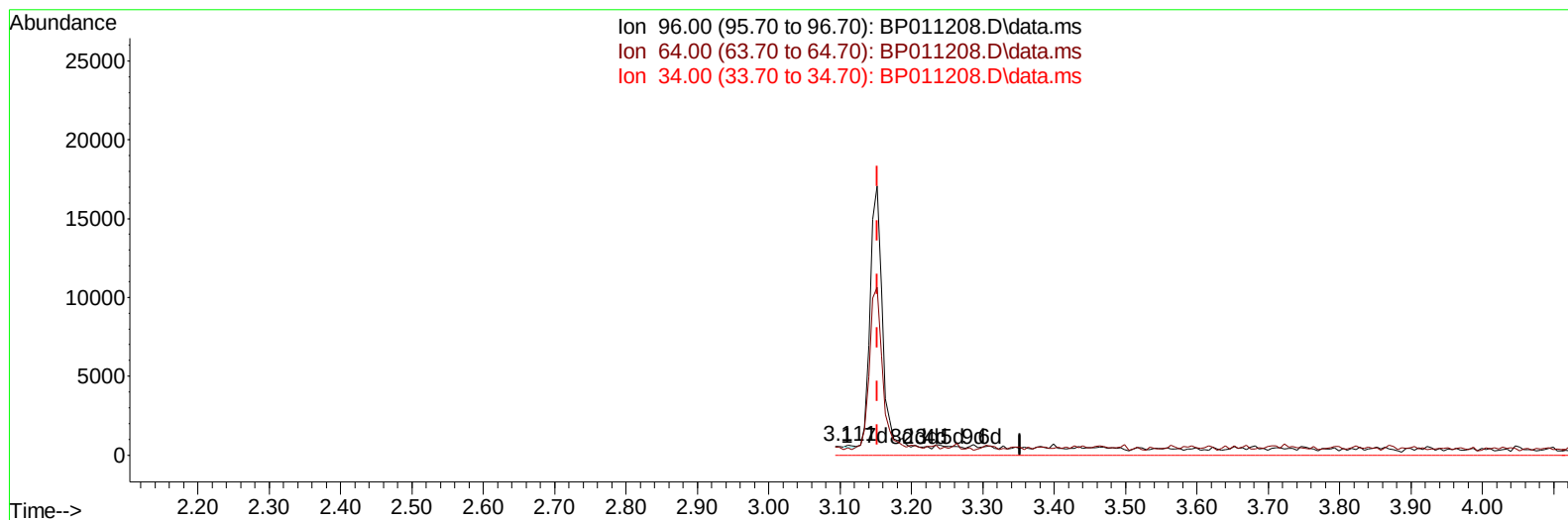
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011208.D
 Acq On : 01 Aug 2022 19:55
 Operator : CG/JU
 Sample : N3792-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBUN3

Manual IntegrationsAPPROVED

Quant Time: Aug 01 23:01:39 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: BP011208.D\data.ms

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

3.111min (-0.041) 0.02 ng/uL

response	93	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	73.37
34.00	0.00	0.00
0.00	0.00	0.00

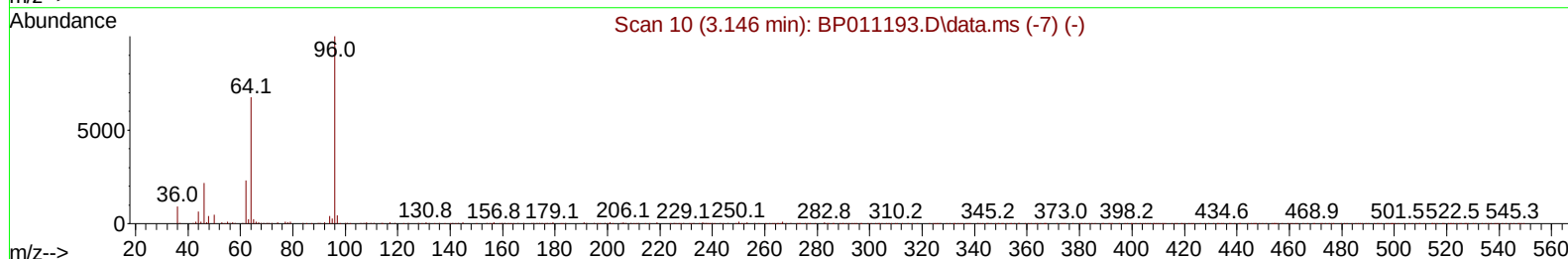
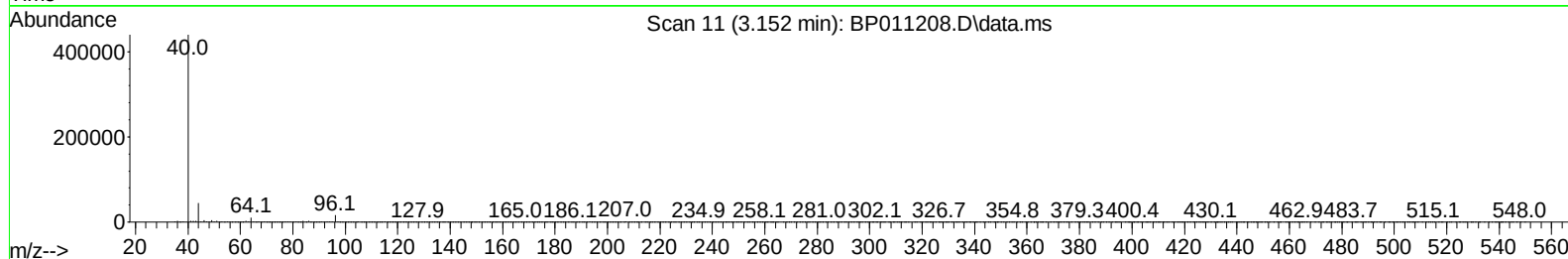
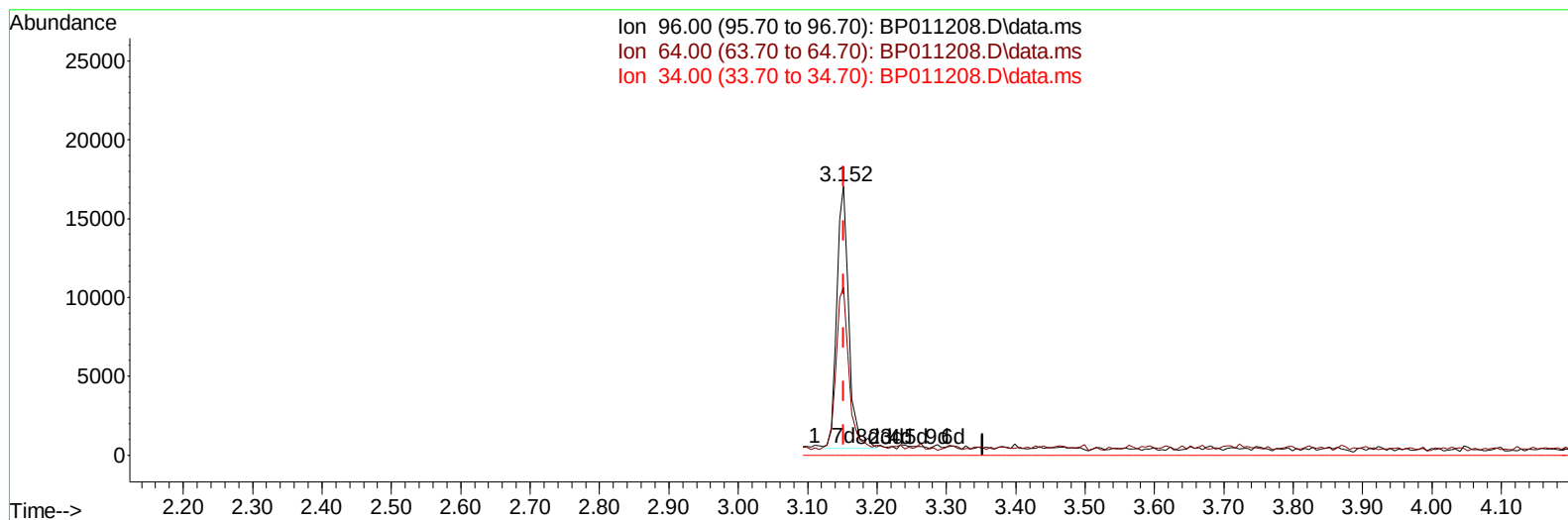
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TIC: BP011208.D\data.ms

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

3.152min (-0.000) 4.02 ng/uL m

response 19897

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	62.44#
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	7.646	152	226635	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	969876	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.316	164	503095	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	969711	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	914431	20.000	ng/ul	# 0.00
88) Perylene-d12	23.557	264	961271	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	19897m	4.020	ng/uL	0.00
4) Pyridine-d5	3.569	84	78540	5.610	ng/ul	0.01
7) Phenol-d5	6.834	99	140306	7.093	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	461542	32.996	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	407696	26.153	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	255820	15.809	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	262054	34.767	ng/ul	0.00
24) 2-Nitrophenol-d4	9.528	143	242766	33.624	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	374222	29.411	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	534125	25.894	ng/ul	0.00
46) Dimethylphthalate-d6	13.739	166	1306768	38.124	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1627642	37.788	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	25636	4.158	ng/ul	0.01
60) Fluorene-d10	15.316	176	1124759	39.051	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	126634	25.906	ng/ul	0.00
73) Anthracene-d10	17.186	188	1893309	44.311	ng/ul	0.00
81) Pyrene-d10	19.445	212	2177018	44.306	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.410	264	2137379	44.766	ng/ul	0.00

Target Compounds Qvalue

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

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