

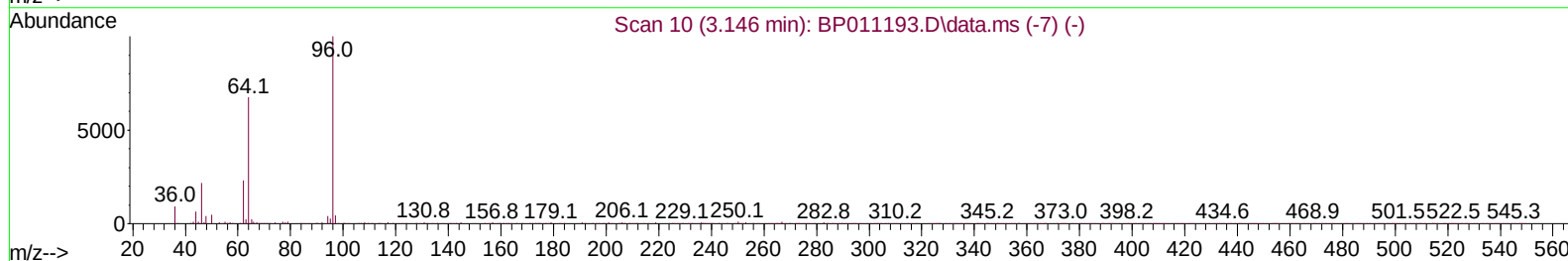
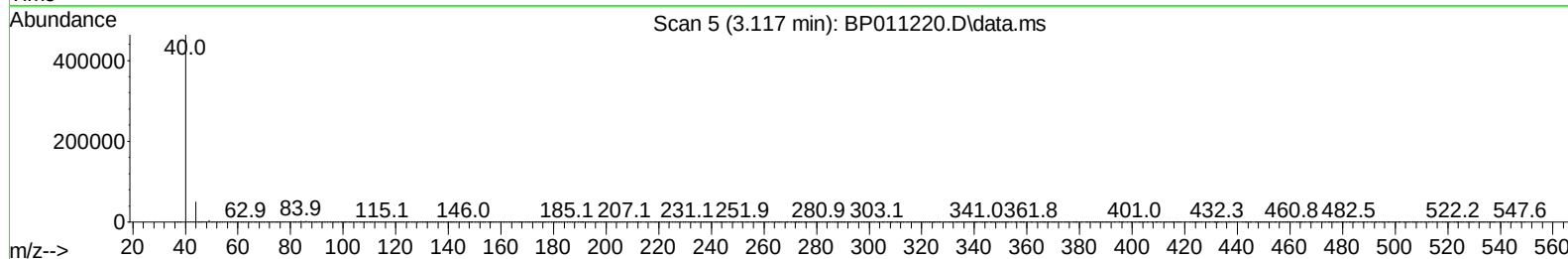
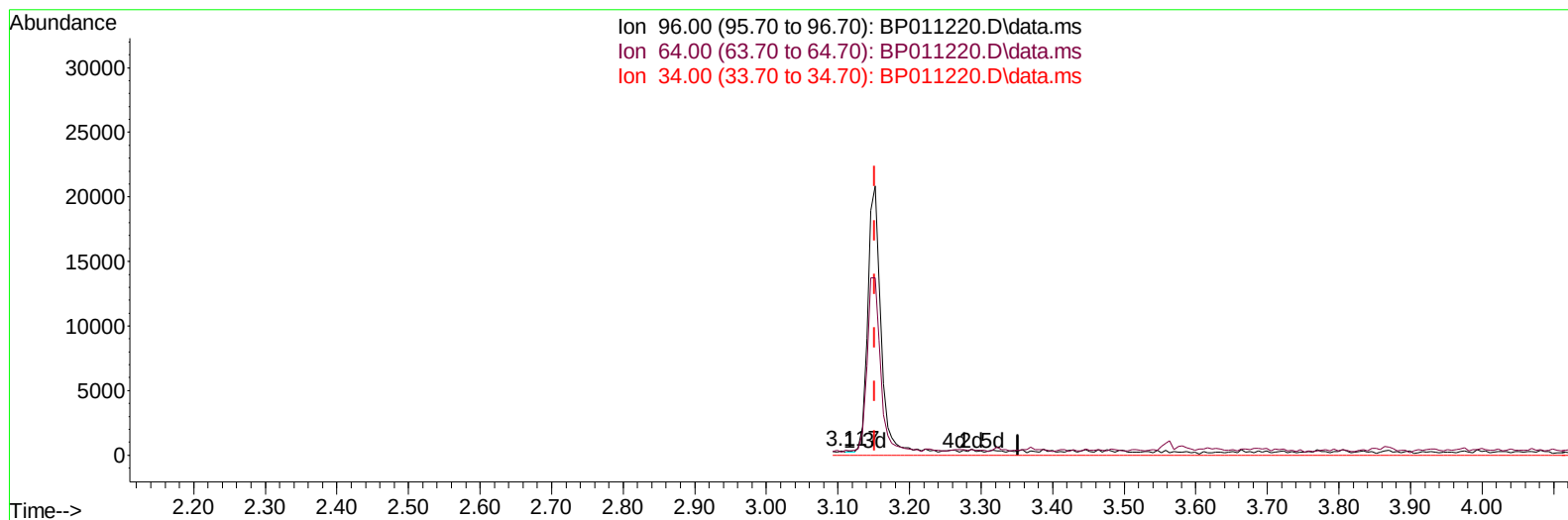
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011220.D
 Acq On : 02 Aug 2022 08:23
 Operator : CG/JU
 Sample : PB146570BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK570

Manual IntegrationsAPPROVED

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

Quant Time: Aug 02 16:00:53 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



TIC: BP011220.D\data.ms

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

3.117min (-0.035) 0.01 ng/uL

response 45

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

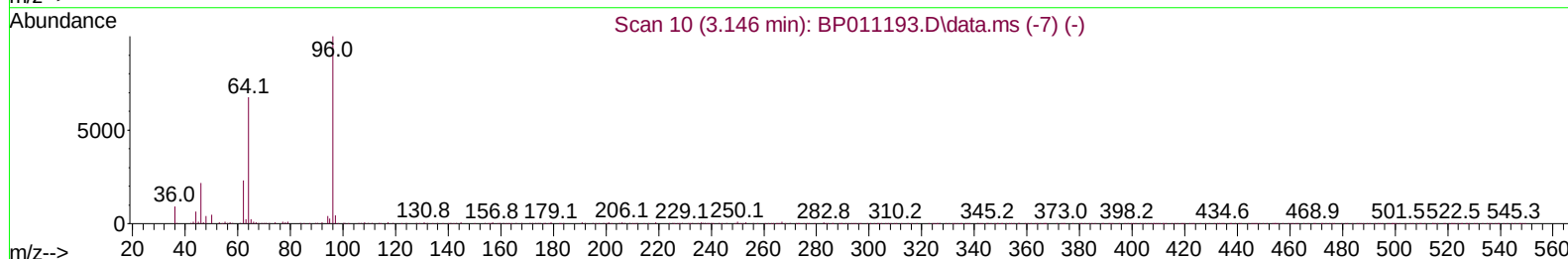
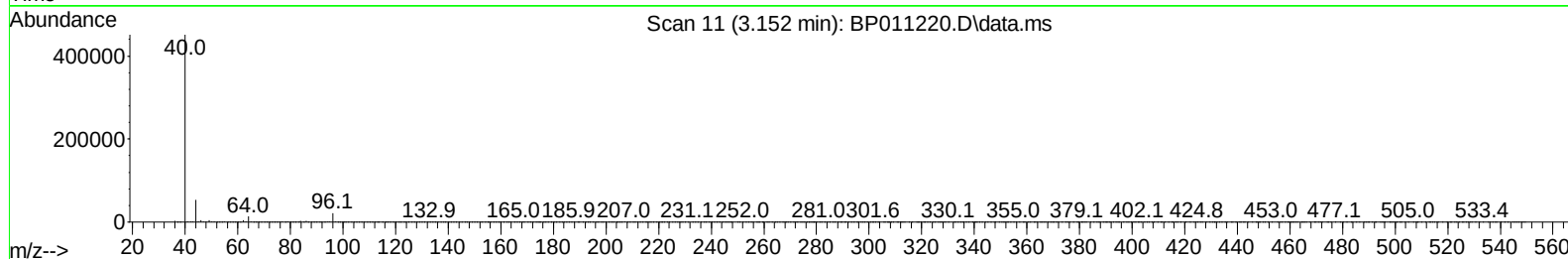
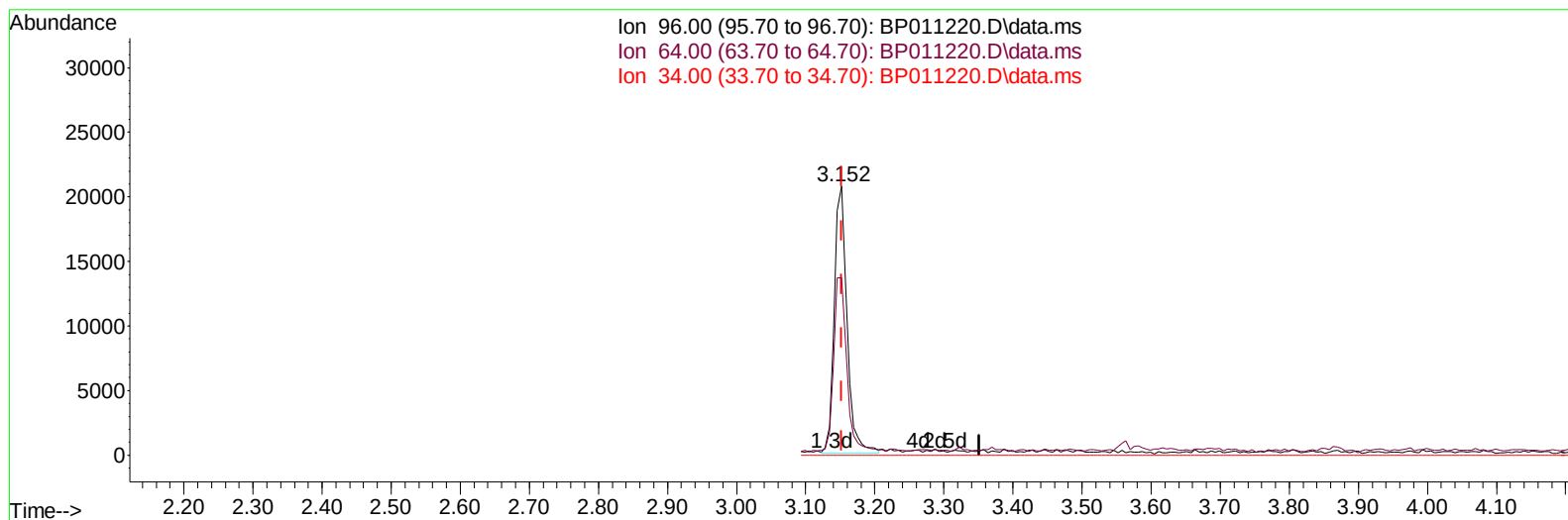
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TIC: BP011220.D\data.ms

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

3.152min (-0.000) 5.03 ng/uL m

response 26518

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	65.81#
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	7.652	152	241232	20.000	ng/ul	0.00
20) Naphthalene-d8	10.446	136	1016798	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.322	164	524962	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	1032102	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.221	240	997653	20.000	ng/ul	# 0.01
88) Perylene-d12	23.568	264	1031277	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	26518m	5.034	ng/uL	0.00
4) Pyridine-d5	3.558	84	352065	23.625	ng/ul	0.00
7) Phenol-d5	6.834	99	675111	32.063	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	468688	31.480	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	545839	32.895	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	515792	29.946	ng/ul	0.00
21) Nitrobenzene-d5	8.816	128	258270	32.683	ng/ul	0.00
24) 2-Nitrophenol-d4	9.534	143	250051	33.035	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	424197	31.800	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	637629	29.486	ng/ul	0.00
46) Dimethylphthalate-d6	13.745	166	1242364	34.735	ng/ul	0.00
49) Acenaphthylene-d8	14.010	160	1563797	34.793	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143	176078	27.369	ng/ul	0.02
60) Fluorene-d10	15.322	176	1038037	34.539	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	110326	21.205	ng/ul	0.01
73) Anthracene-d10	17.192	188	1632037	35.887	ng/ul	0.00
81) Pyrene-d10	19.451	212	1830259	34.141	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.421	264	1843919	35.998	ng/ul	0.02

Target Compounds Qvalue

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

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