

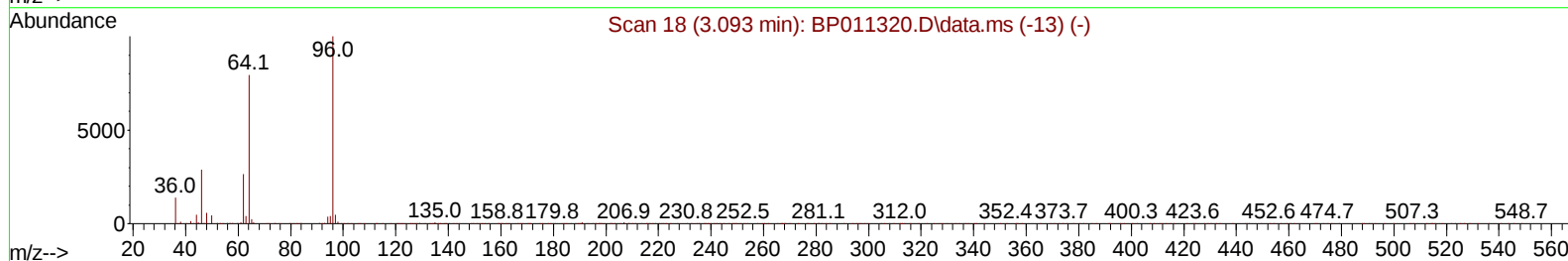
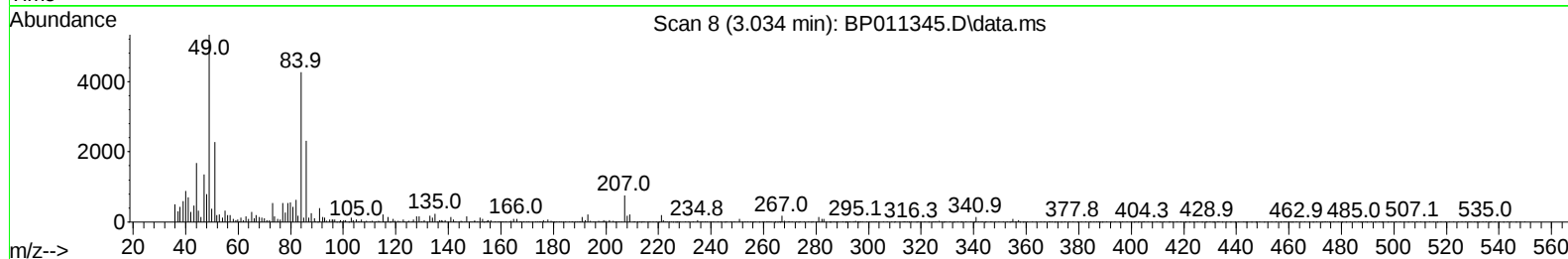
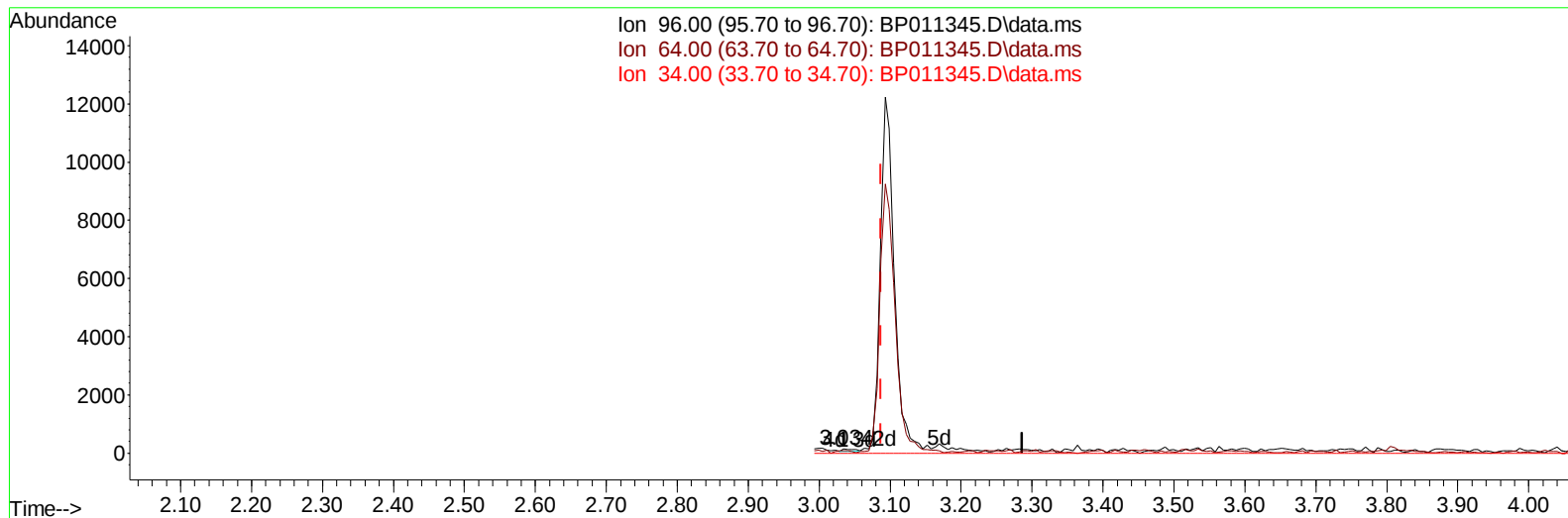
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011345.D
 Acq On : 10 Aug 2022 00:13
 Operator : CG/JU
 Sample : N3900-14
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CA2

Manual IntegrationsAPPROVED

Quant Time: Aug 10 01:28:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 18:47:19 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/10/2022
 Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011345.D\data.ms

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

3.034min (-0.053) 0.02 ng/uL

response 121

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	59.44
34.00	0.00	0.00
0.00	0.00	0.00

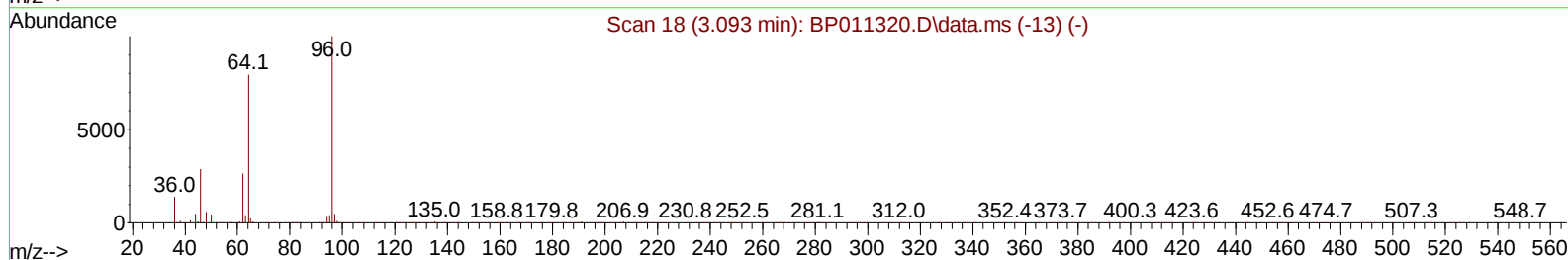
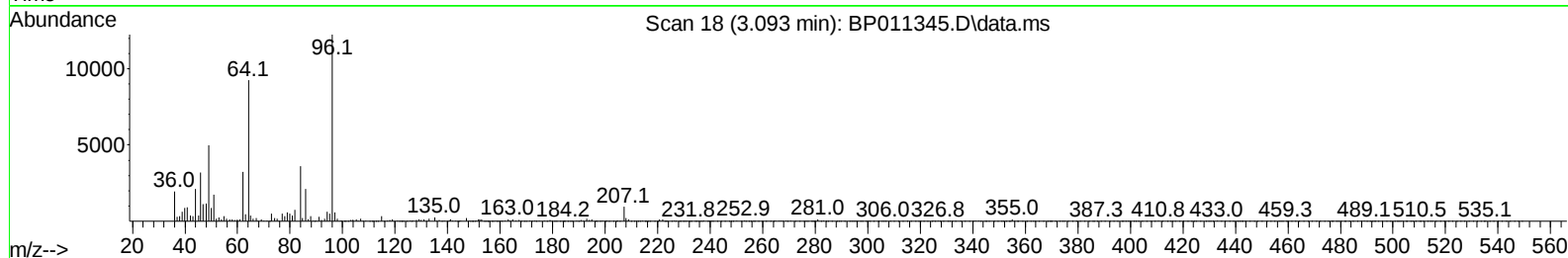
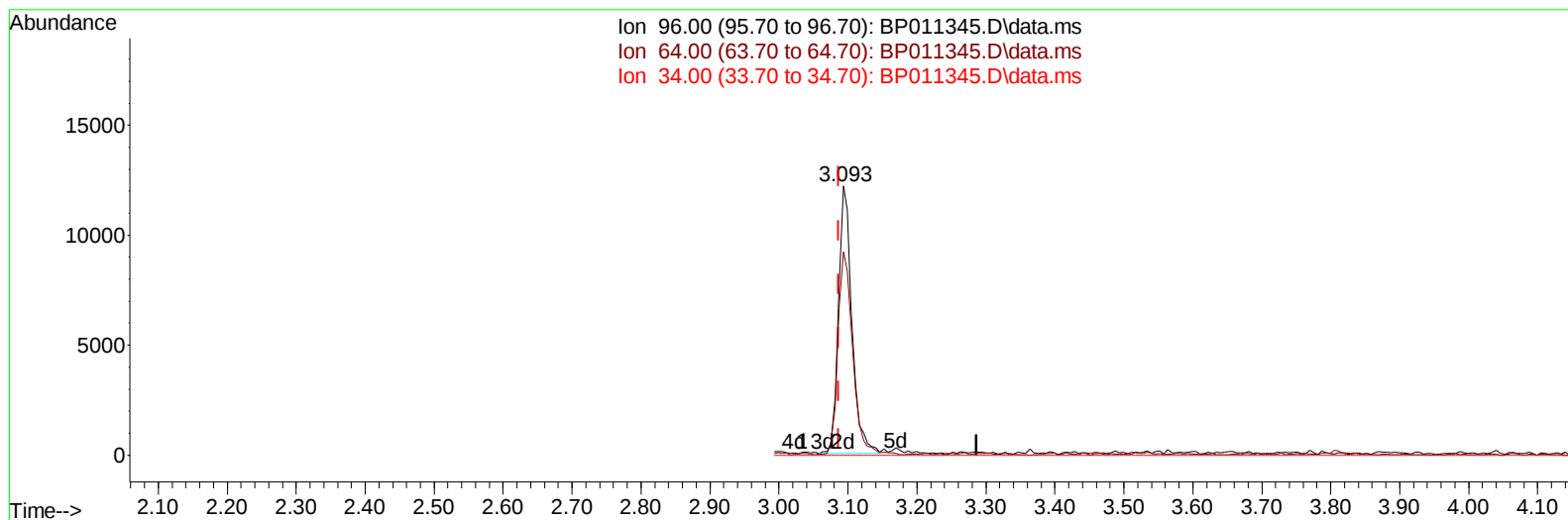
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(3) 1,4-Dioxane-d8 (S)

3.093min (+ 0.006) 3.35 ng/uL m

response 16569

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	75.57#
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	7.587	152	151613	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	638792	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	358709	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	715023	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	669375	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	654539	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	16569m	3.350	ng/uL	0.00
4) Pyridine-d5	3.517	84	58317	4.675	ng/ul	0.02
7) Phenol-d5	6.775	99	64878	4.835	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	233544	24.549	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	210161	19.963	ng/ul	0.00
15) 4-Methylphenol-d8	8.311	113	128341	12.182	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	140380	32.290	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	131801	33.113	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	203417	25.307	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	215929	15.366	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	861698	34.622	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	953616	31.829	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	16575	3.646	ng/ul	0.01
60) Fluorene-d10	15.263	176	691789	32.501	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	107900	32.488	ng/ul	0.00
73) Anthracene-d10	17.134	188	1159997	36.093	ng/ul	0.00
81) Pyrene-d10	19.392	212	1349044	36.581	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	1238693	37.526	ng/ul	-0.01

Target Compounds Qvalue

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

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