

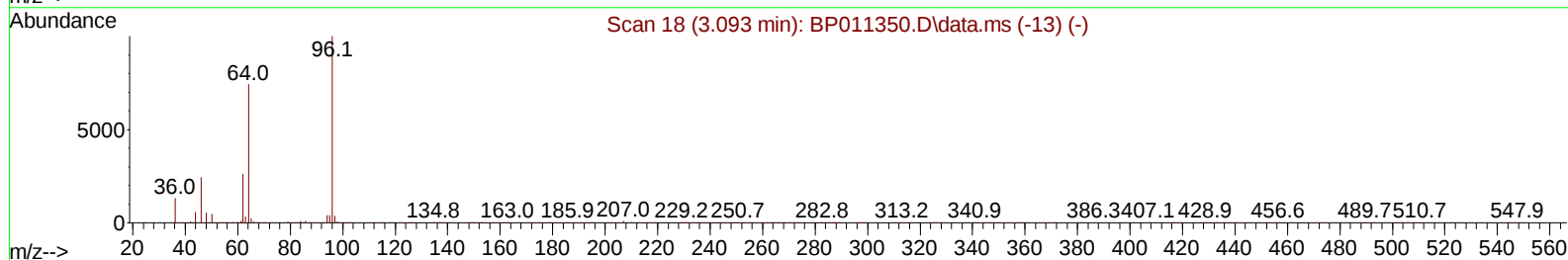
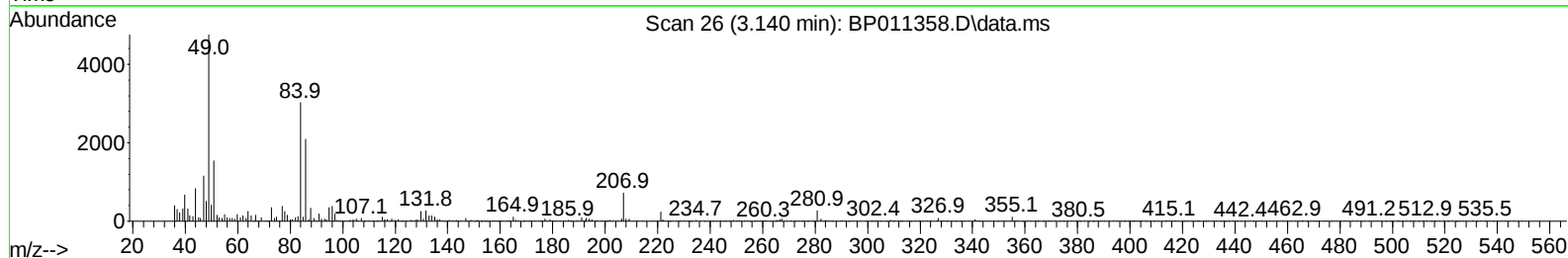
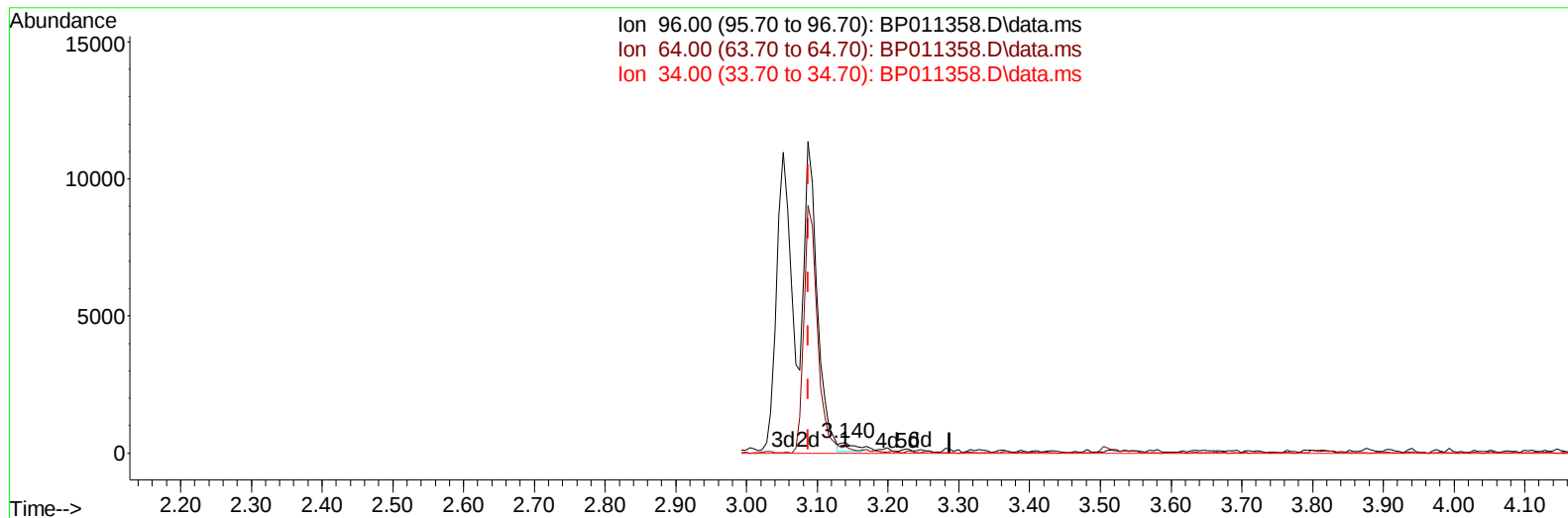
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011358.D
 Acq On : 10 Aug 2022 08:17
 Operator : CG/JU
 Sample : N4069-07
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJL3

Manual IntegrationsAPPROVED

Quant Time: Aug 10 23:53:47 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 08/11/2022
 Supervised By :Sohil Jodhani 08/12/2022



TIC: BP011358.D\data.ms

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

3.140min (+ 0.053) 0.13 ng/uL

response 482

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	64.63
34.00	0.00	0.00
0.00	0.00	0.00

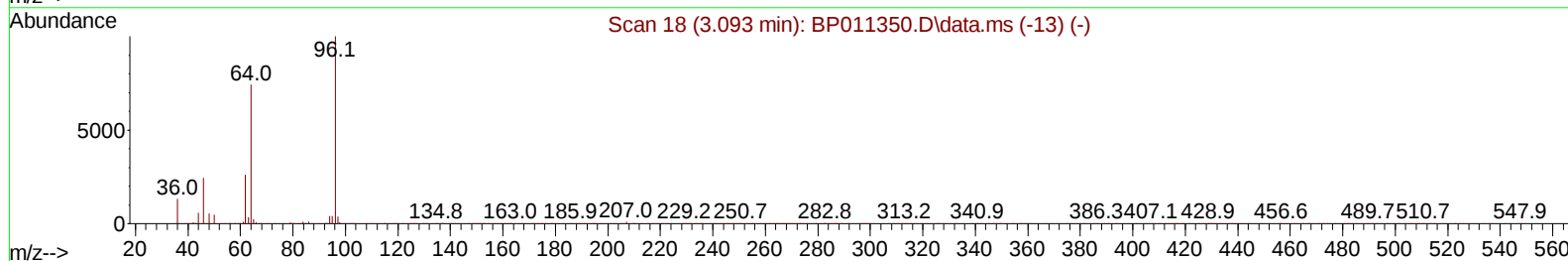
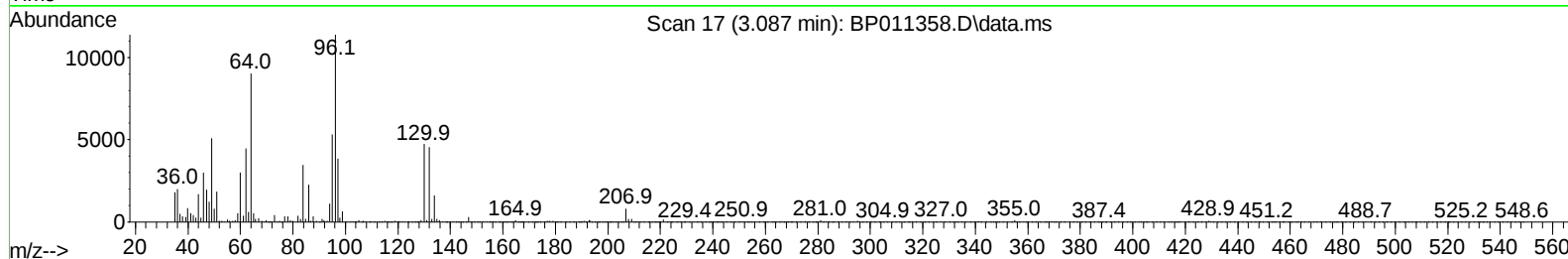
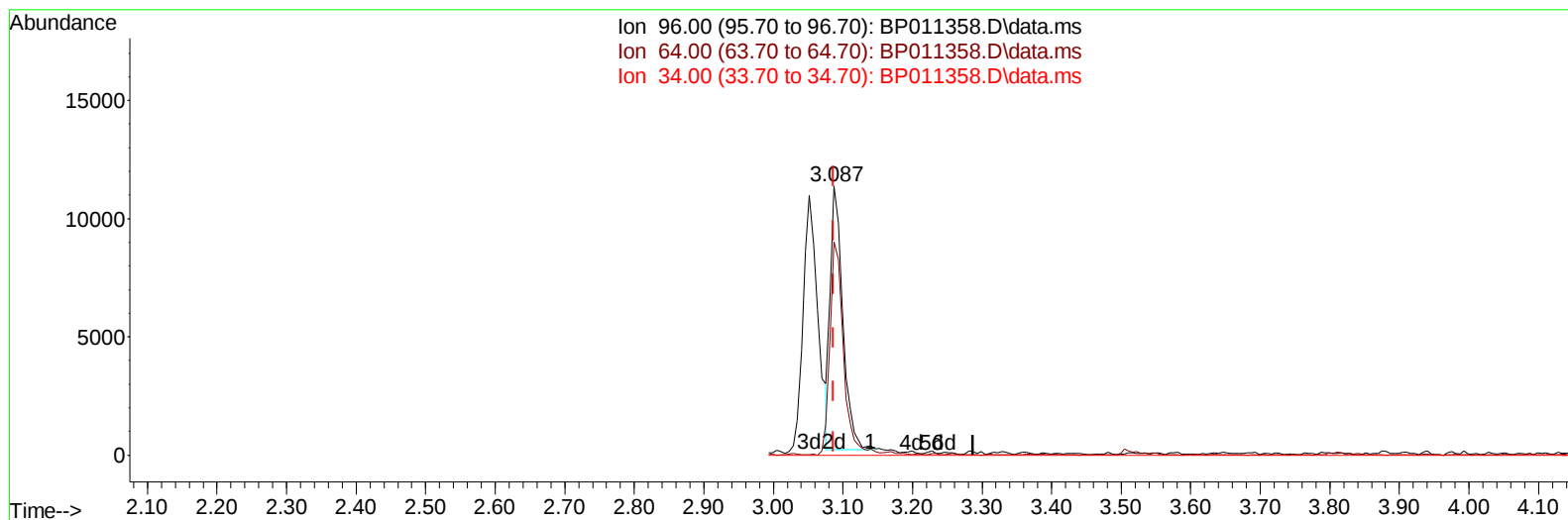
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TIC: BP011358.D\data.ms

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

3.087min (+ 0.000) 3.76 ng/uL m

response 13987

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	79.37#
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	113985	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	478864	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	276100	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	554257	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	519609	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	536122	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	13987m	3.762	ng/uL	0.00
4) Pyridine-d5	3.505	84	75491	8.049	ng/ul	0.00
7) Phenol-d5	6.769	99	66500	6.592	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	189202	26.453	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	185510	23.438	ng/ul	0.00
15) 4-Methylphenol-d8	8.304	113	124974	15.779	ng/ul	-0.01
21) Nitrobenzene-d5	8.746	128	104670	32.117	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.463	143	103931	34.832	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.004	165	175894	29.191	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	109271	10.373	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	663833	34.652	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	751629	32.594	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.492	143	12470	3.564	ng/ul	0.00
60) Fluorene-d10	15.263	176	536183	32.728	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.392	200	83819	32.558	ng/ul	0.00
73) Anthracene-d10	17.128	188	913421	36.665	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1047448	36.589	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	976756	36.127	ng/ul	0.00

Target Compounds Qvalue

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

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