

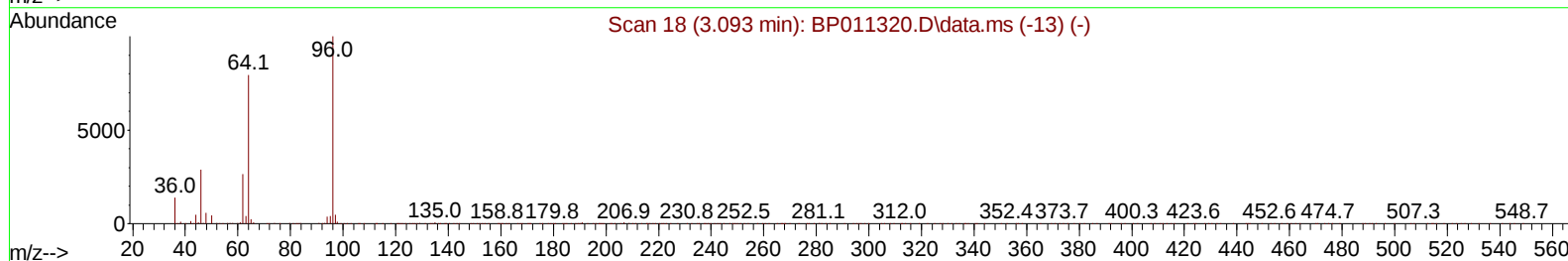
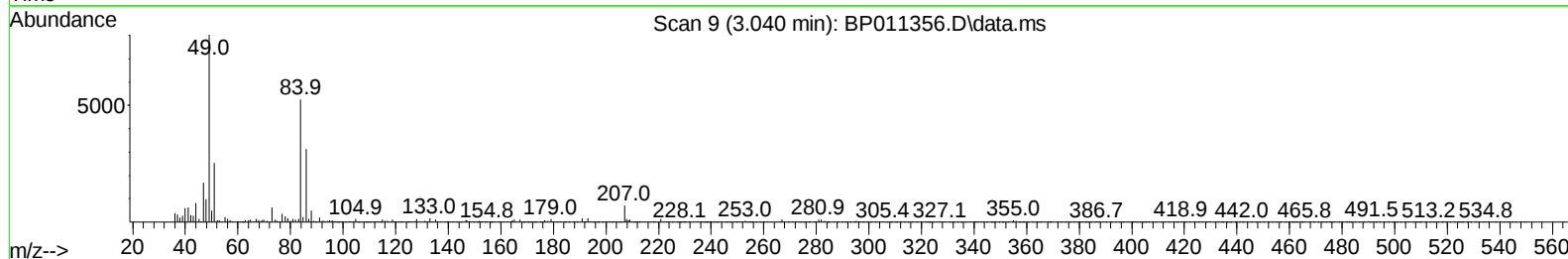
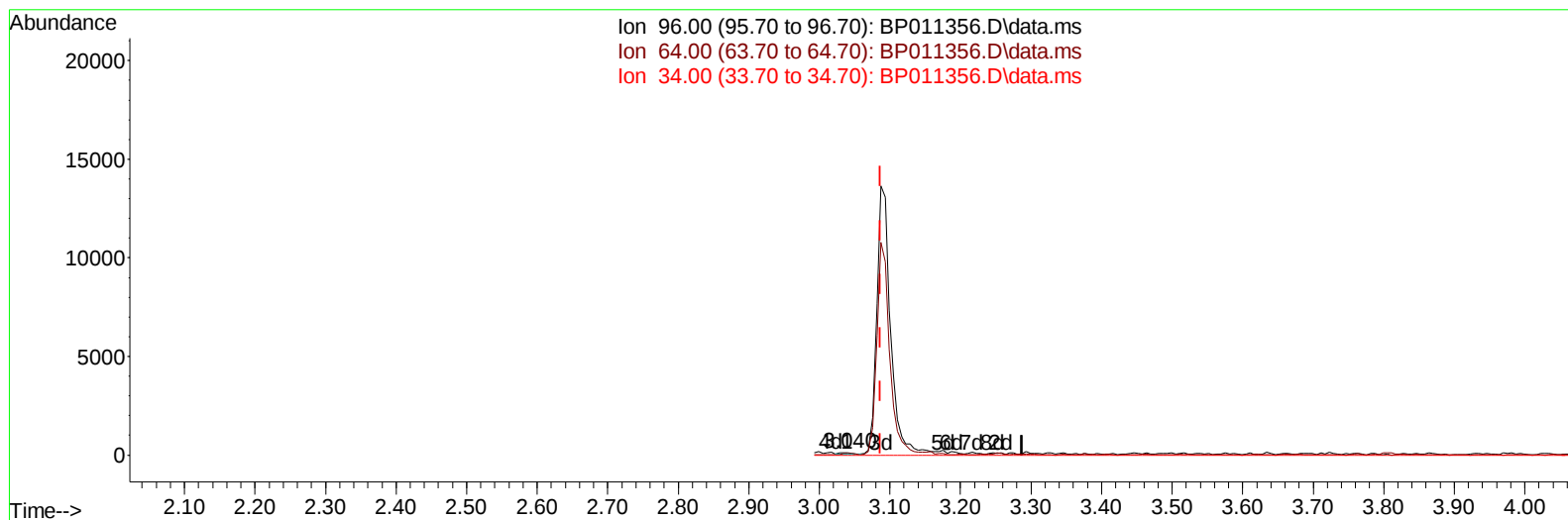
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011356.D
 Acq On : 10 Aug 2022 07:08
 Operator : CG/JU
 Sample : N4069-03
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJK1

Manual IntegrationsAPPROVED

Quant Time: Aug 10 07:46:05 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: BP011356.D\data.ms

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

3.040min (-0.047) 0.02 ng/uL

response 73

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	67.48
34.00	0.00	0.00
0.00	0.00	0.00

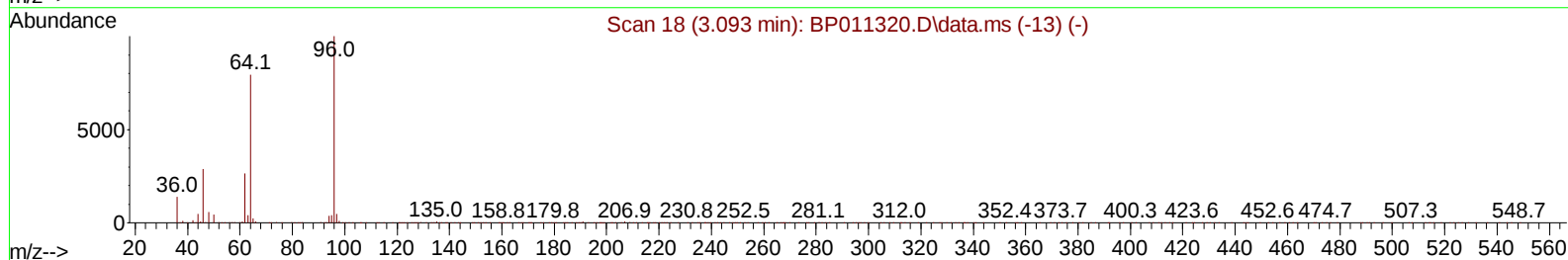
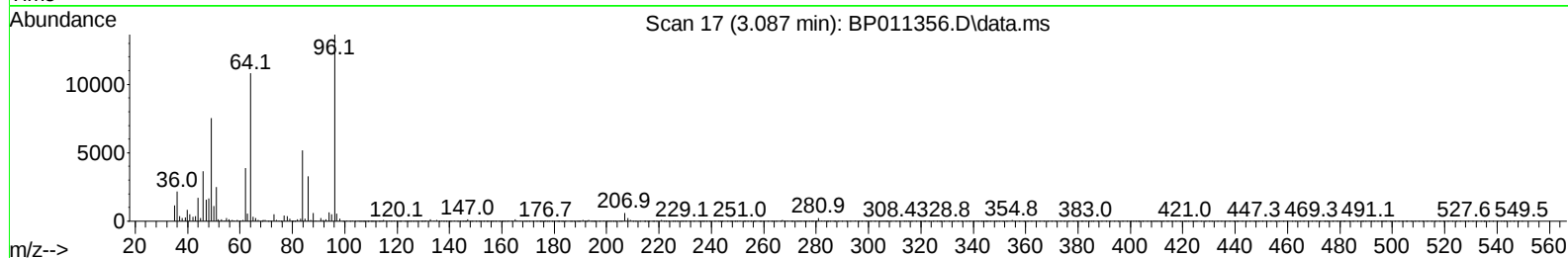
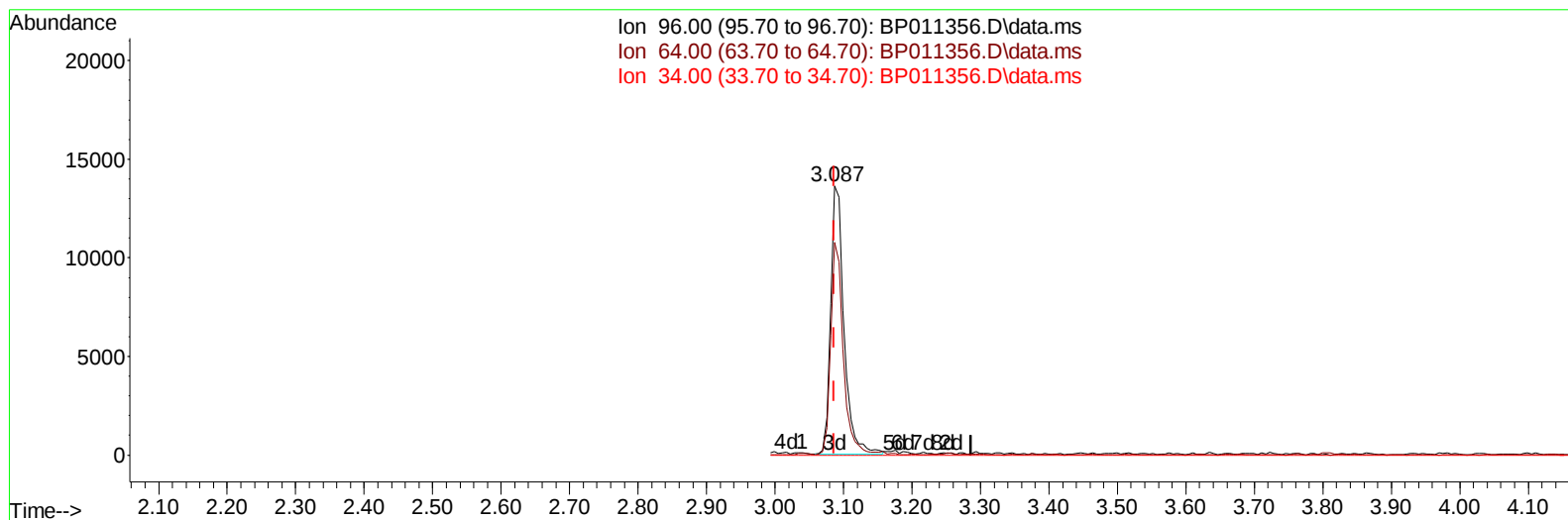
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(3) 1,4-Dioxane-d8 (S)

3.087min (+ 0.000) 5.01 ng/uL m

response 18479

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	79.21#
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	113020	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	483193	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	274649	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	557798	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	522315	20.000	ng/ul	-0.01
88) Perylene-d12	23.469	264	527084	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	18479m	5.012	ng/uL	0.00
4) Pyridine-d5	3.517	84	23806	2.560	ng/ul	0.02
7) Phenol-d5	6.769	99	79722	7.971	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	211423	29.812	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	207572	26.450	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	142027	18.085	ng/ul	-0.01
21) Nitrobenzene-d5	8.752	128	119253	36.264	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	118563	39.380	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	195169	32.100	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	156336	14.708	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	766296	40.212	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	846083	36.883	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	18461	5.304	ng/ul	0.00
60) Fluorene-d10	15.263	176	614753	37.722	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.393	200	91905	35.472	ng/ul	0.00
73) Anthracene-d10	17.128	188	1021946	40.761	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1182412	41.090	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.322	264	1120912	42.169	ng/ul	0.00

Target Compounds Qvalue

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

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