

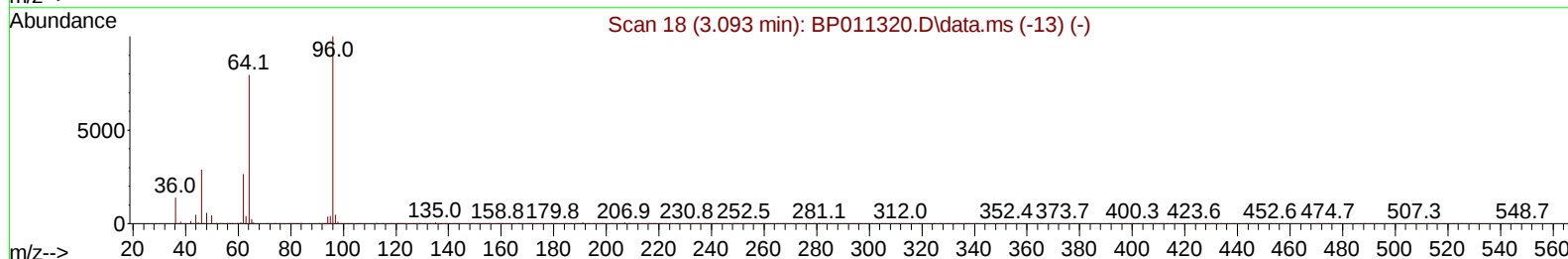
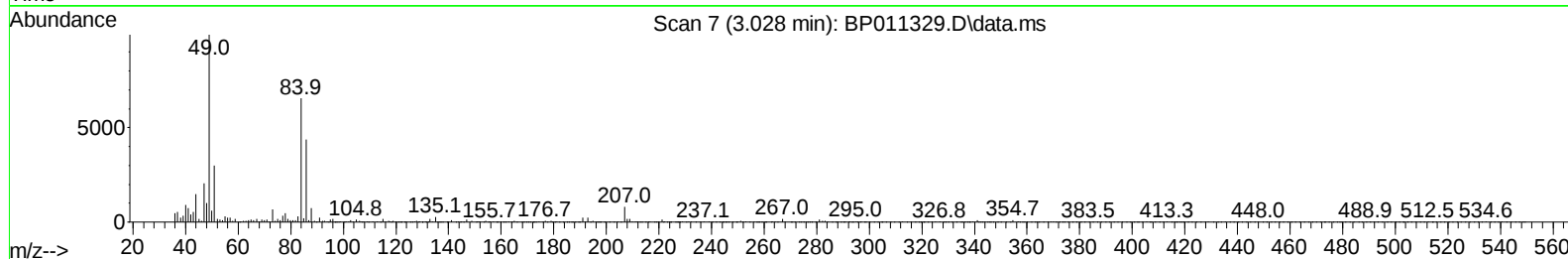
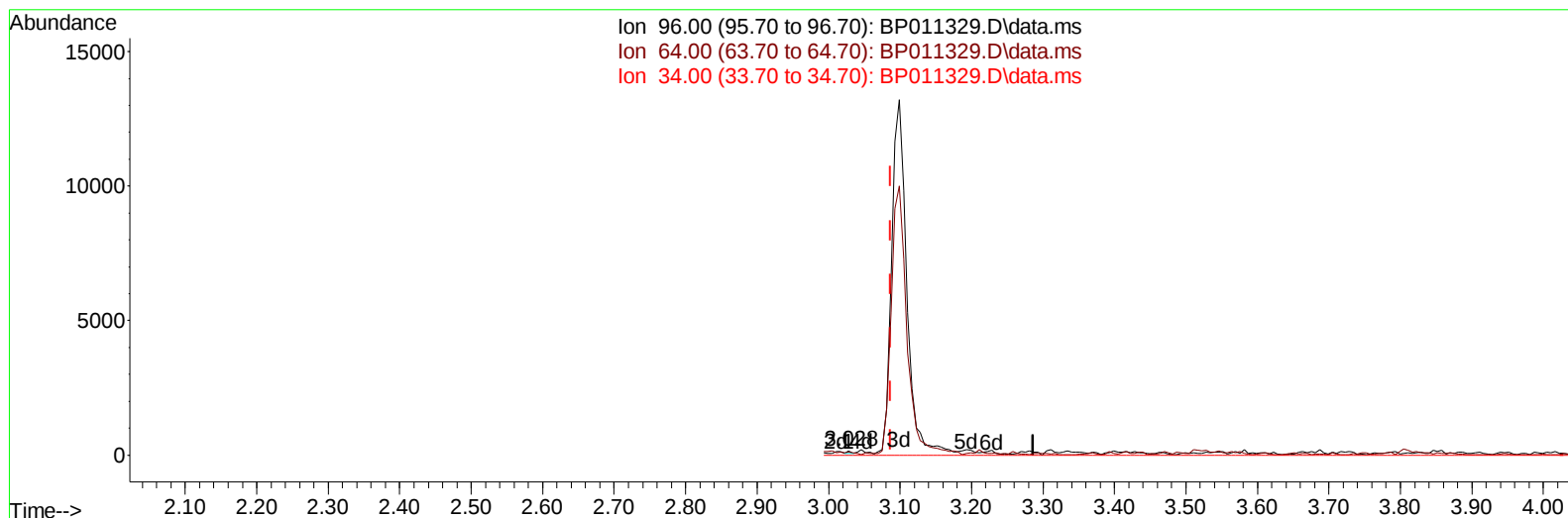
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011329.D
 Acq On : 09 Aug 2022 14:55
 Operator : CG/JU
 Sample : N4026-20
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJJ4

Manual IntegrationsAPPROVED

Quant Time: Aug 09 23:11:37 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: BP011329.D\data.ms

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

3.028min (-0.059) 0.01 ng/uL

response	39	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	68.90
34.00	0.00	0.00
0.00	0.00	0.00

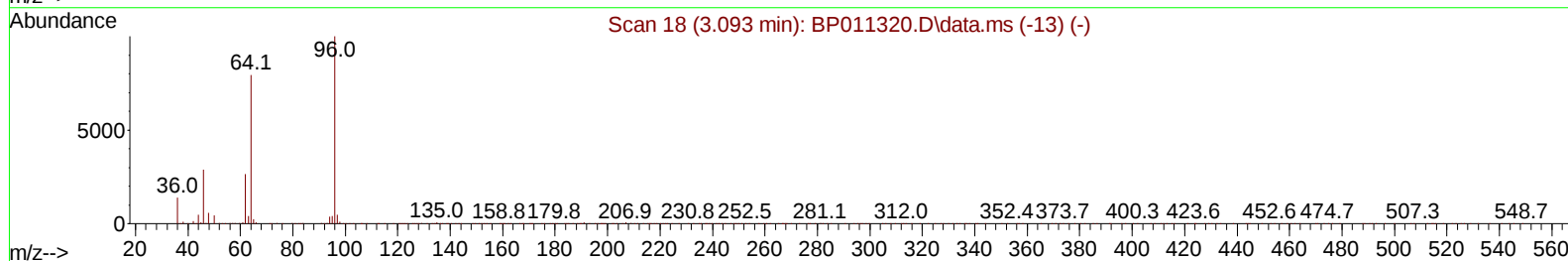
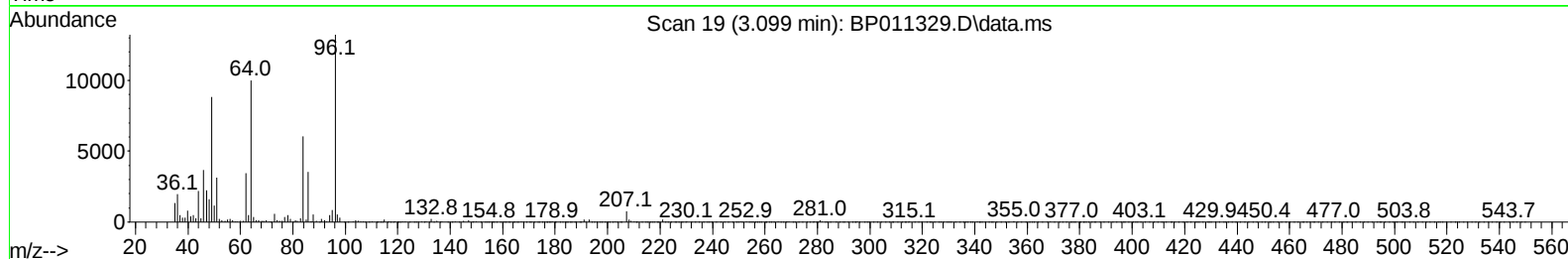
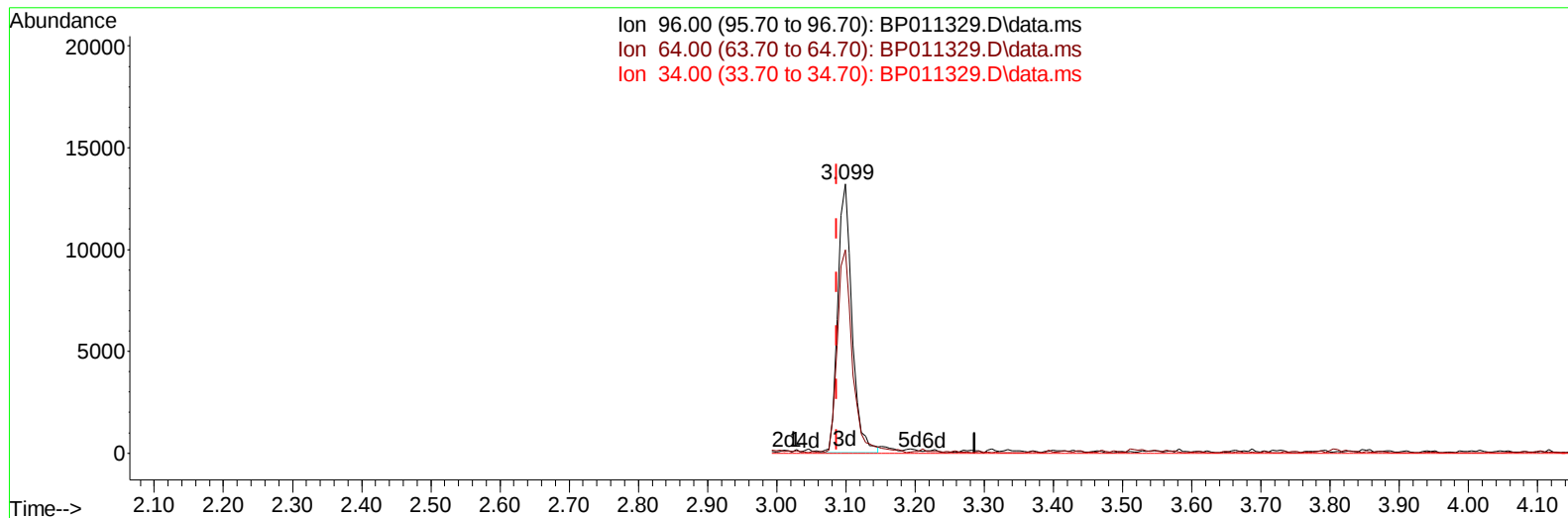
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TIC: BP011329.D\data.ms

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

3.099min (+ 0.012) 4.24 ng/uL m

response 18926

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	75.67#
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	136730	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	589091	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	336999	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	679457	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	637642	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	636646	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	18926m	4.243	ng/uL	0.01
4) Pyridine-d5	3.517	84	84857	7.542	ng/ul	0.02
7) Phenol-d5	6.775	99	77631	6.416	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	256674	29.917	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	224735	23.671	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	149368	15.721	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	139992	34.918	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	125356	34.151	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	228182	30.783	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	169569	13.085	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	947147	40.506	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	1027048	36.489	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	18171	4.255	ng/ul	0.00
60) Fluorene-d10	15.263	176	775867	38.799	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	107922	34.196	ng/ul	0.00
73) Anthracene-d10	17.133	188	1267886	41.515	ng/ul	0.00
81) Pyrene-d10	19.398	212	1483013	42.215	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.327	264	1368478	42.623	ng/ul	0.00

Target Compounds Qvalue

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

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