

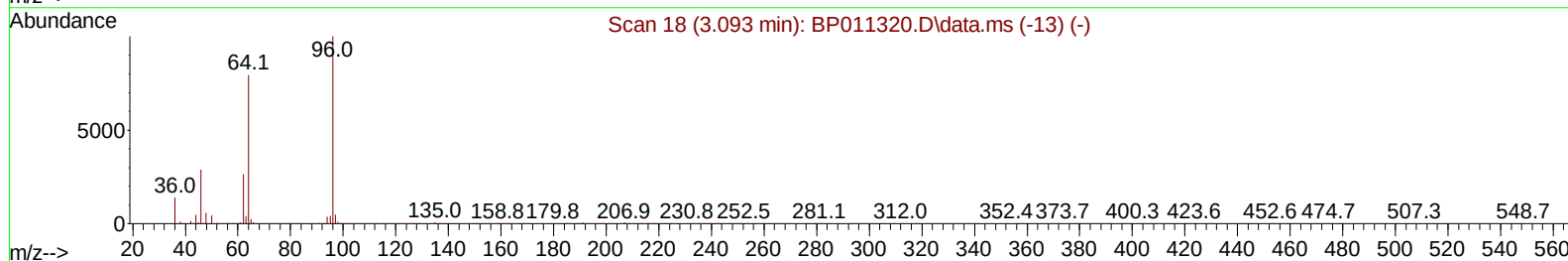
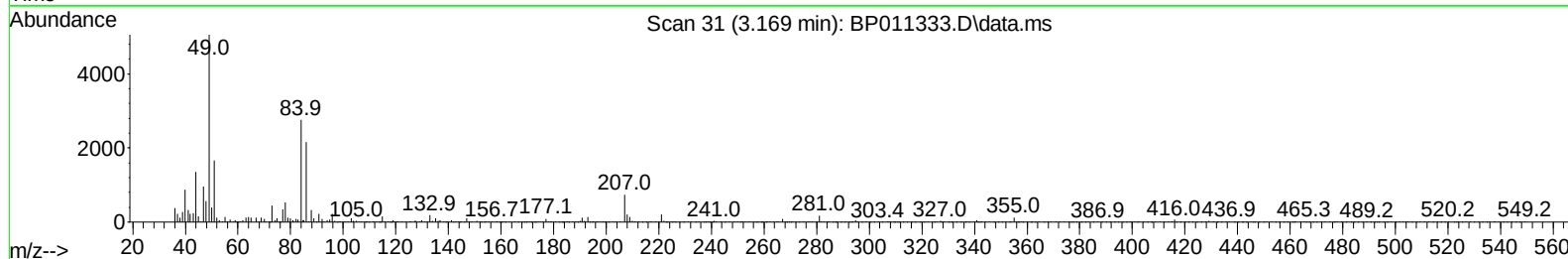
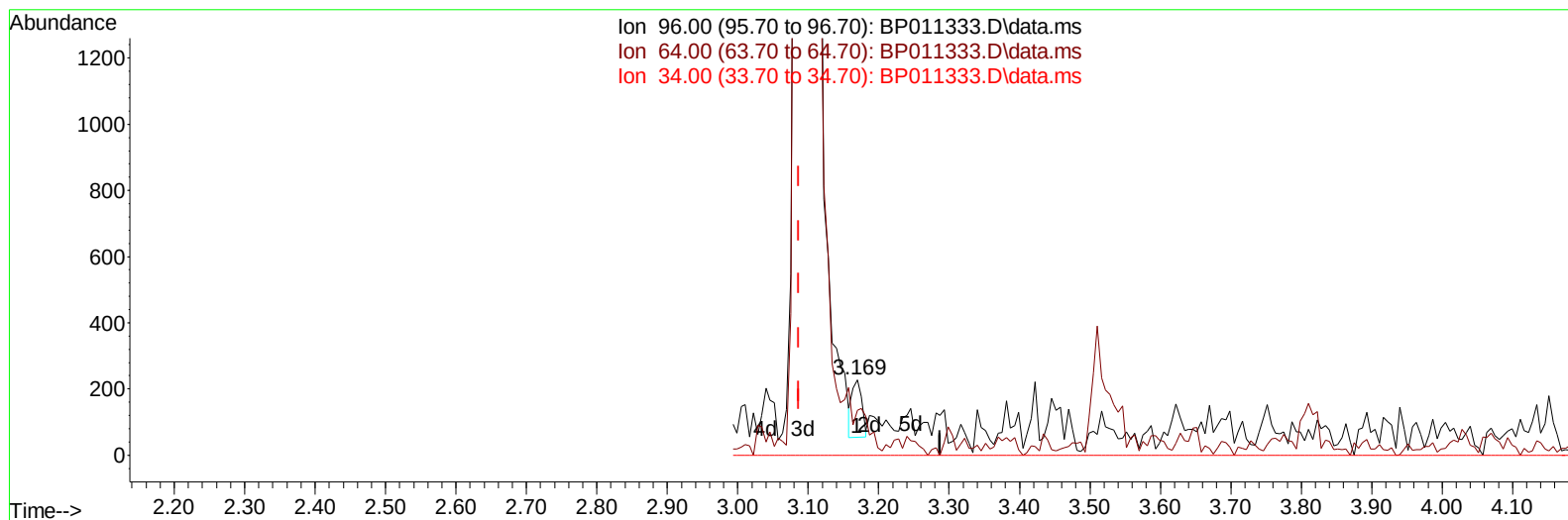
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011333.D
 Acq On : 09 Aug 2022 17:14
 Operator : CG/JU
 Sample : PB146591BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK591

Manual IntegrationsAPPROVED

Quant Time: Aug 09 23:13:53 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: BP011333.D\data.ms

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

3.169min (+ 0.082) 0.04 ng/uL

response 165

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	59.21
34.00	0.00	0.00
0.00	0.00	0.00

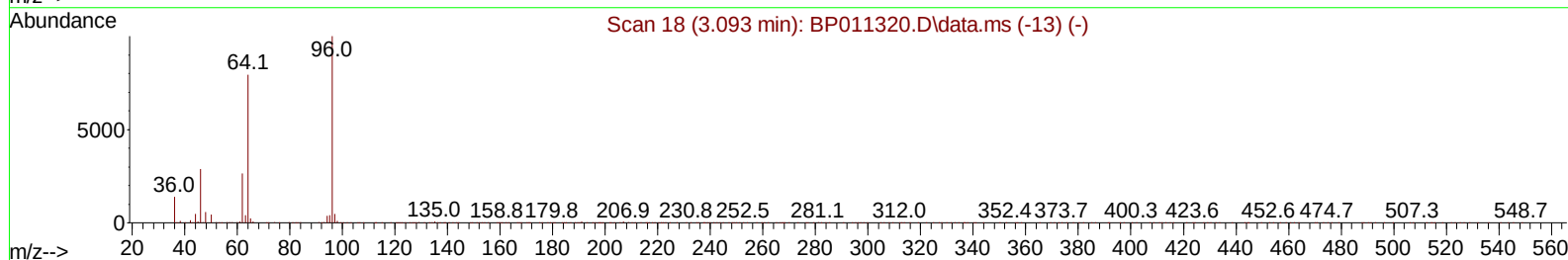
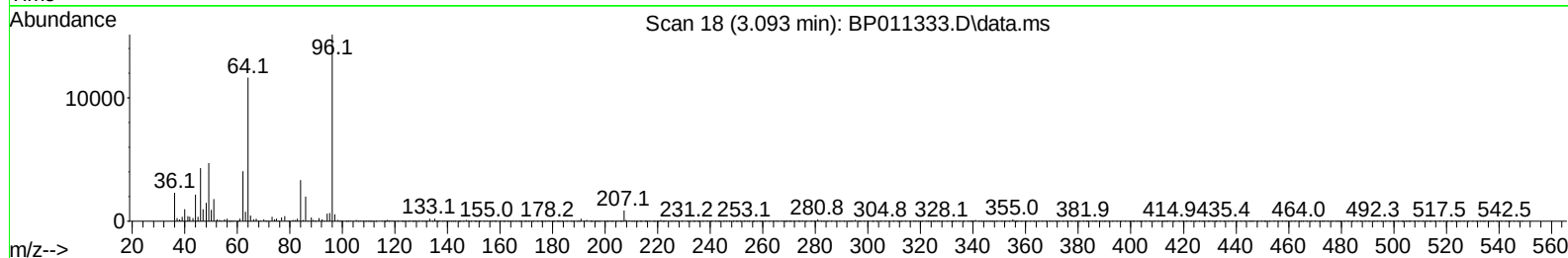
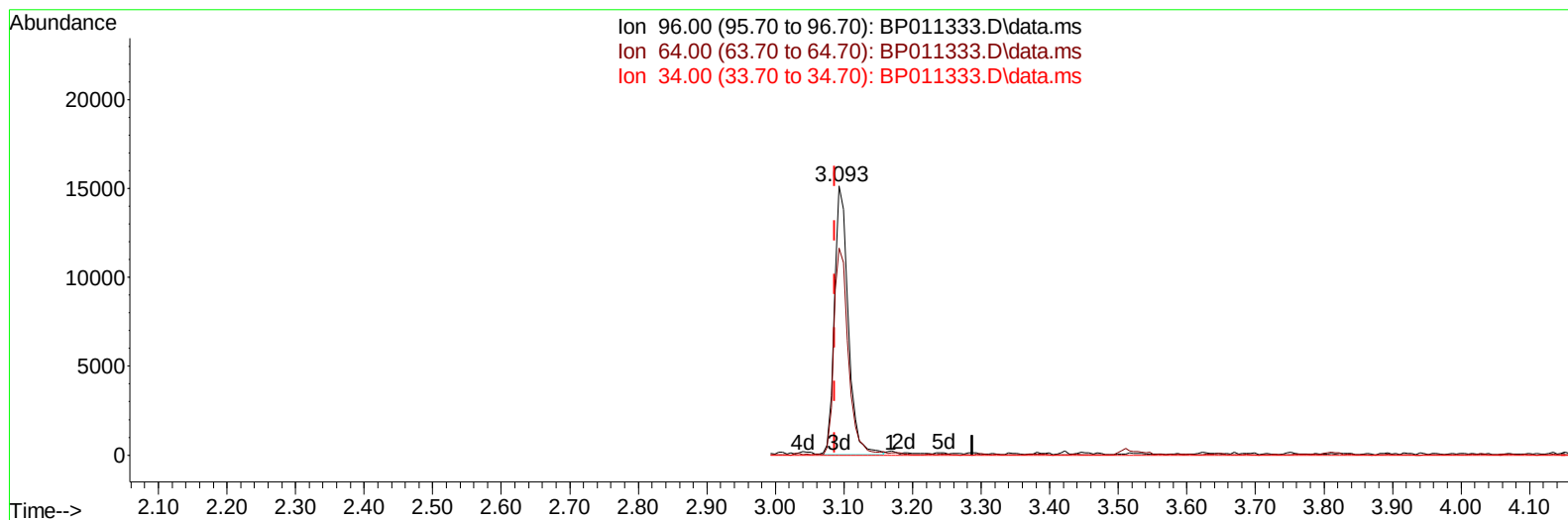
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(3) 1,4-Dioxane-d8 (S)**3.093min (+ 0.006) 5.61 ng/uL m****response 21221**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	76.91#
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	115881	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	485200	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	274116	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	542520	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	490428	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	488356	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	21221m	5.614	ng/uL	0.00
4) Pyridine-d5	3.511	84	148534	15.578	ng/ul	0.01
7) Phenol-d5	6.775	99	284216	27.714	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	196493	27.023	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	229685	28.545	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	217381	26.997	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	103210	31.255	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	98781	32.674	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	171938	28.162	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	175870	16.477	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	611744	32.164	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	696304	30.413	ng/ul	0.00
54) 4-Nitrophenol-d4	14.486	143	75284	21.672	ng/ul	0.00
60) Fluorene-d10	15.263	176	492788	30.297	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	63881	25.350	ng/ul	0.00
73) Anthracene-d10	17.133	188	778621	31.930	ng/ul	0.00
81) Pyrene-d10	19.392	212	861920	31.900	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	777117	31.554	ng/ul	0.00

Target Compounds Qvalue

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

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