

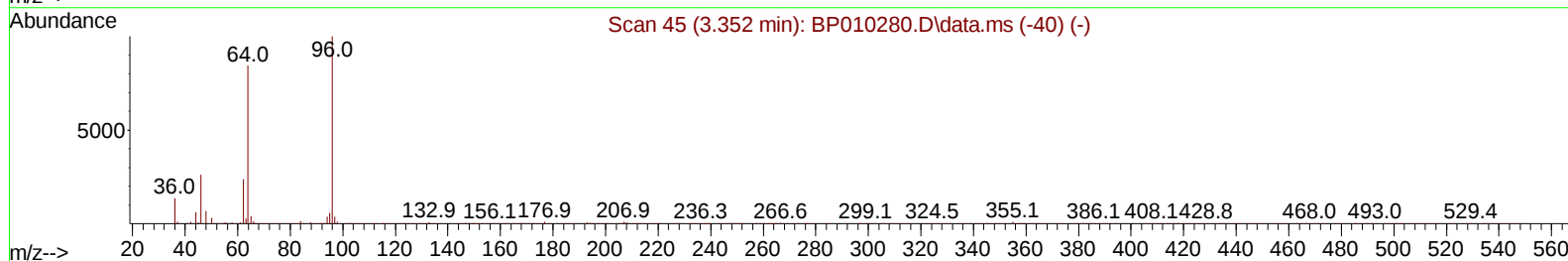
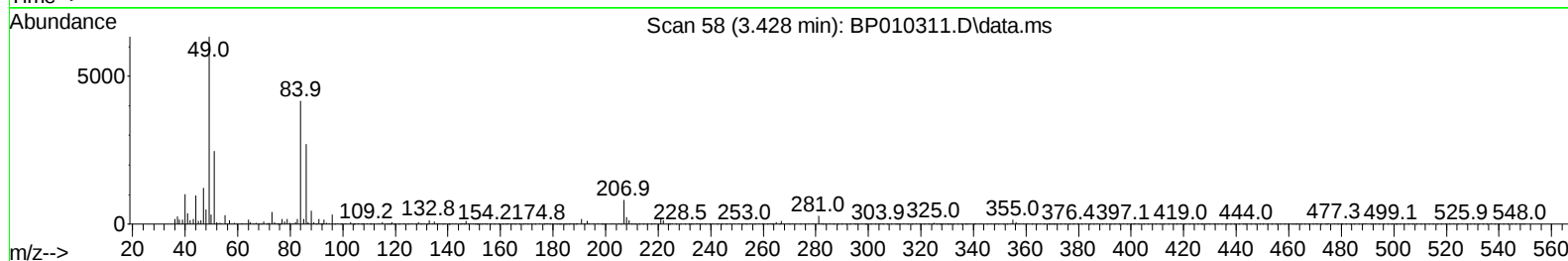
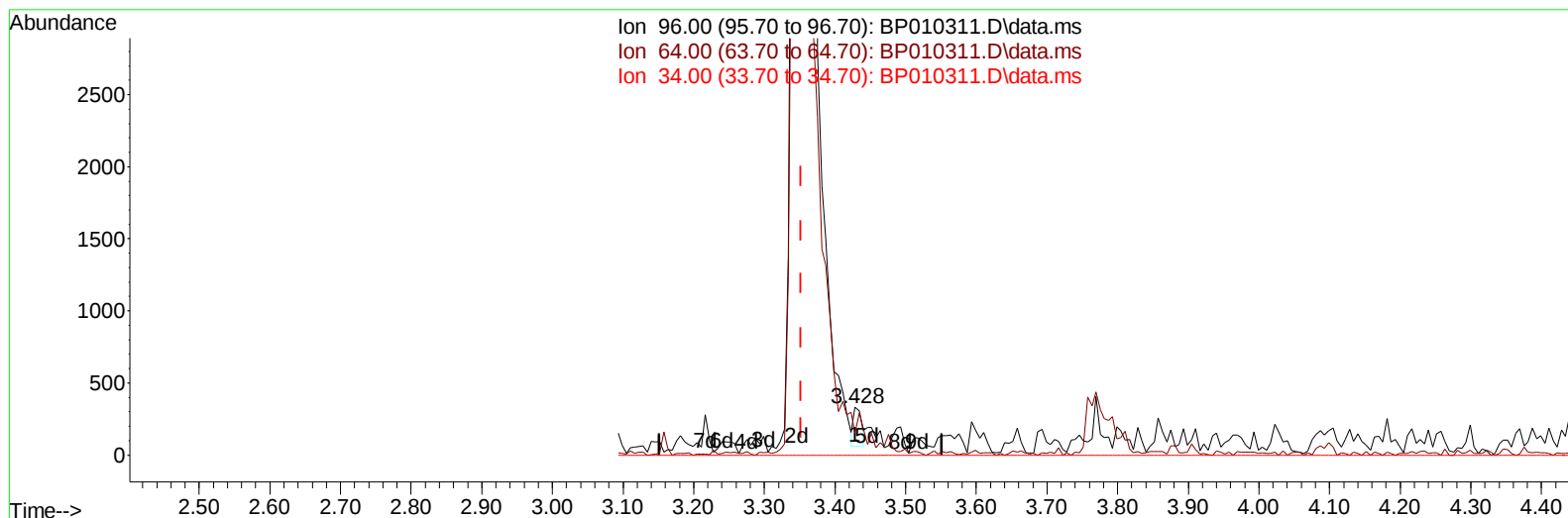
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010311.D
 Acq On : 12 May 2022 06:39
 Operator : CG/JU
 Sample : PB144620BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK620

Manual IntegrationsAPPROVED

Quant Time: May 12 07:39:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010311.D\data.ms

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

3.428min (+ 0.076) 0.06 ng/uL

response 223

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.92
34.00	0.00	0.00
0.00	0.00	0.00

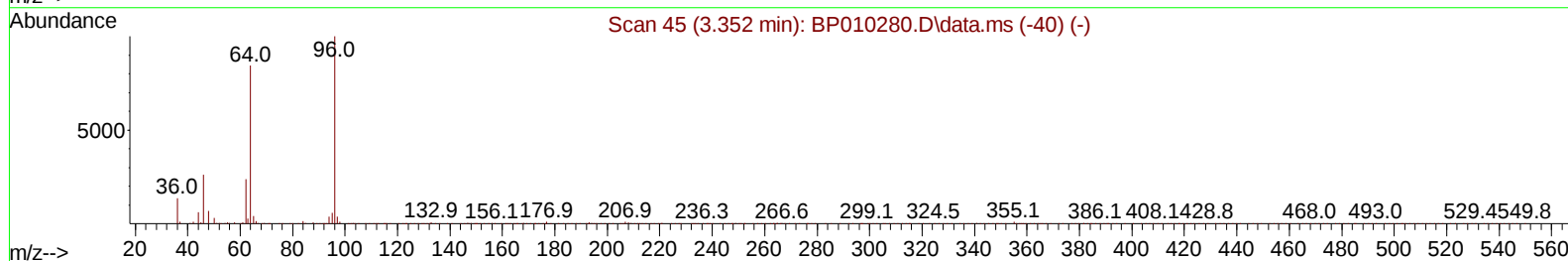
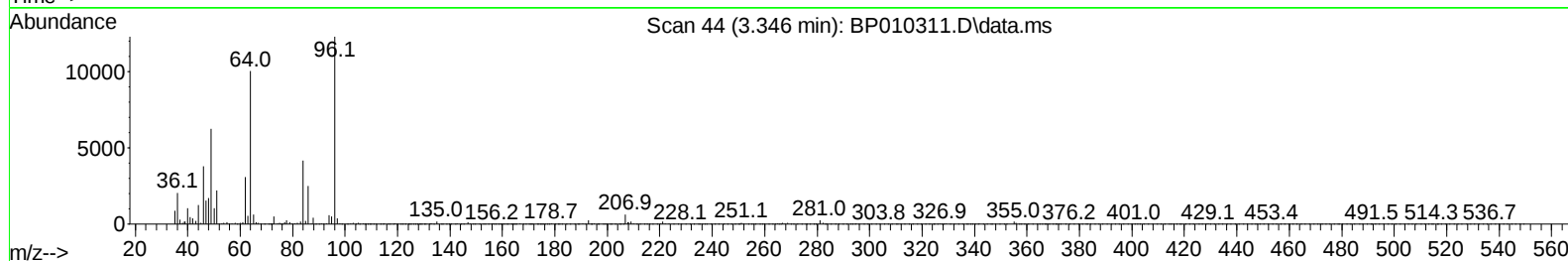
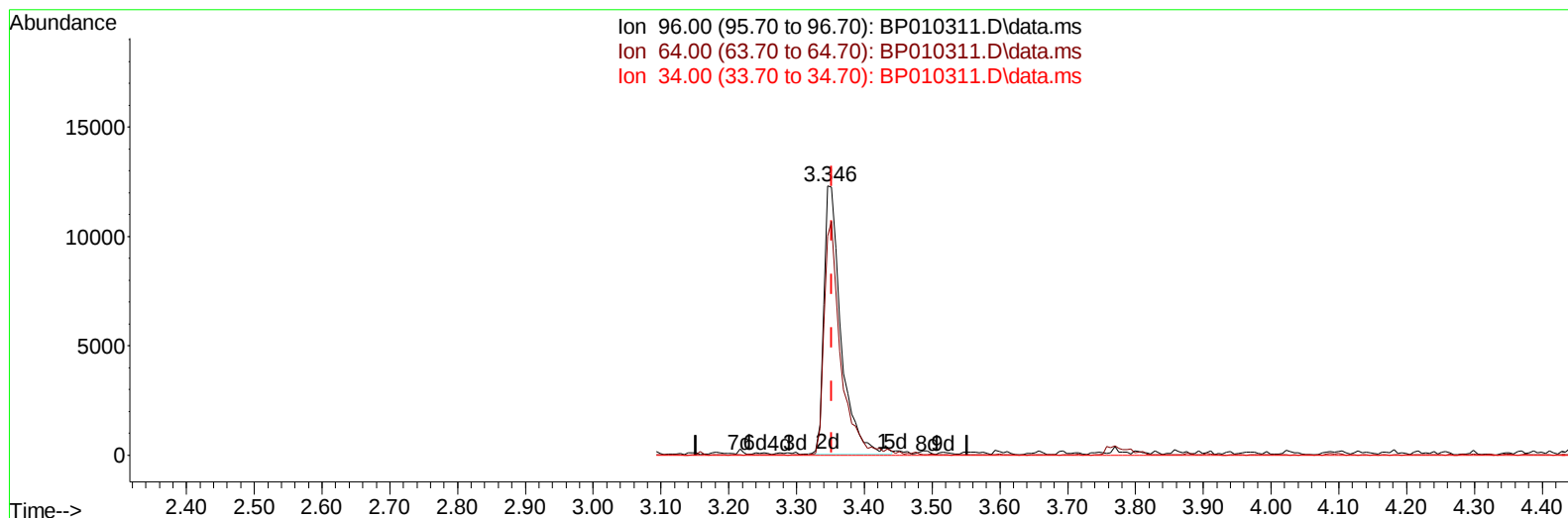
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TIC: BP010311.D\data.ms

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

3.346min (-0.006) 5.85 ng/uL m

response 21782

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.55#
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	7.905	152	158223	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	698718	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.534	164	500169	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1088008	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	1084078	20.000	ng/ul	# 0.00
88) Perylene-d12	23.798	264	1044425	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	21782m	5.848	ng/uL	0.00
4) Pyridine-d5	3.764	84	298408	28.792	ng/ul	0.00
7) Phenol-d5	7.069	99	367602	25.015	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	261833	28.801	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	298215	27.554	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	312965	25.904	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	156618	27.955	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	174585	27.331	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	297250	25.567	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	442102	25.757	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1104812	27.726	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1228640	25.862	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	157656	20.627	ng/ul	-0.01
60) Fluorene-d10	15.528	176	941954	28.005	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	161619	21.692	ng/ul	0.00
73) Anthracene-d10	17.386	188	1498881	28.132	ng/ul	0.00
81) Pyrene-d10	19.616	212	1847413	30.838	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1669966	29.982	ng/ul	-0.02

Target Compounds Qvalue

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

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