

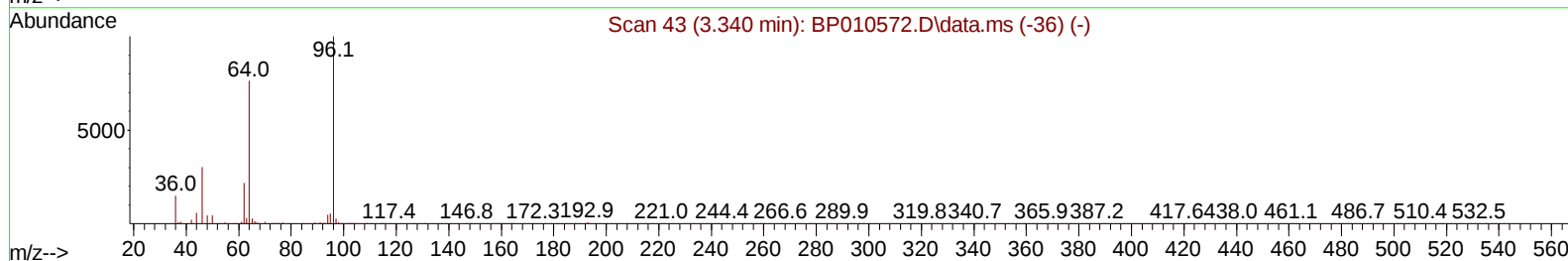
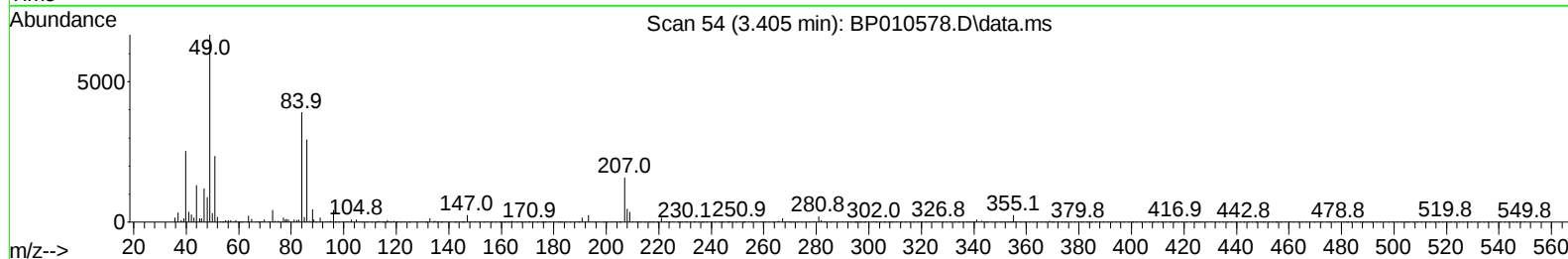
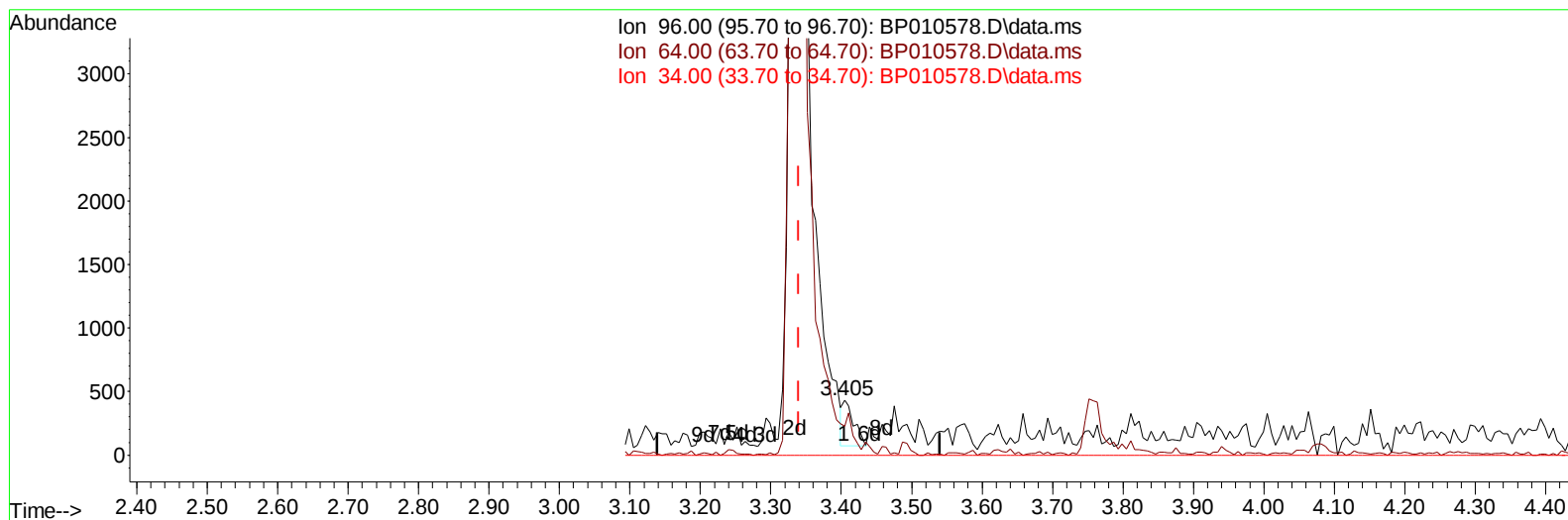
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052722\
 Data File : BP010578.D
 Acq On : 27 May 2022 13:31
 Operator : CG/JU
 Sample : PB145142BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK142

Manual IntegrationsAPPROVED

Quant Time: May 28 00:15:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 24 05:50:40 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/31/2022
 Supervised By :mohammad ahmed 05/31/2022



TIC: BP010578.D\data.ms

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

3.405min (+ 0.065) 0.14 ng/uL

response 382

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	52.76
34.00	0.00	0.00
0.00	0.00	0.00

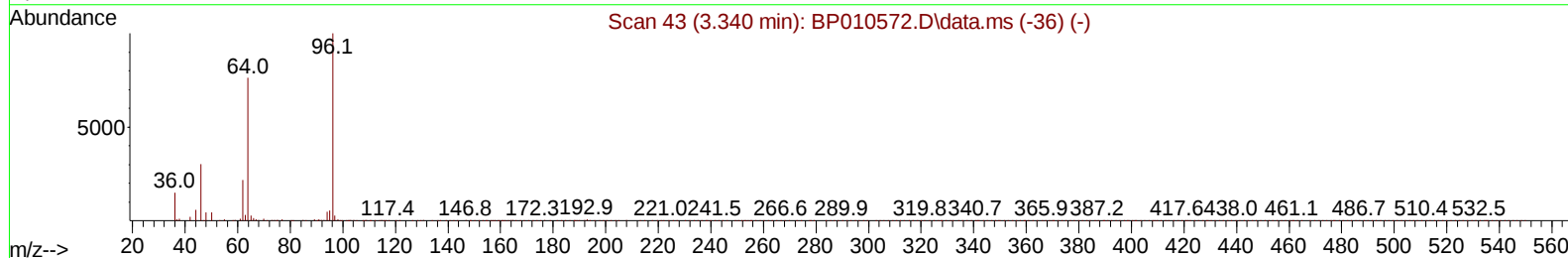
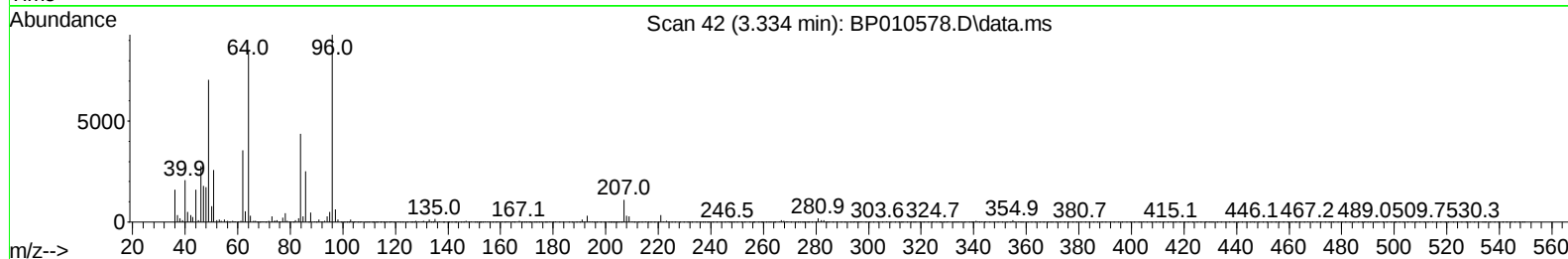
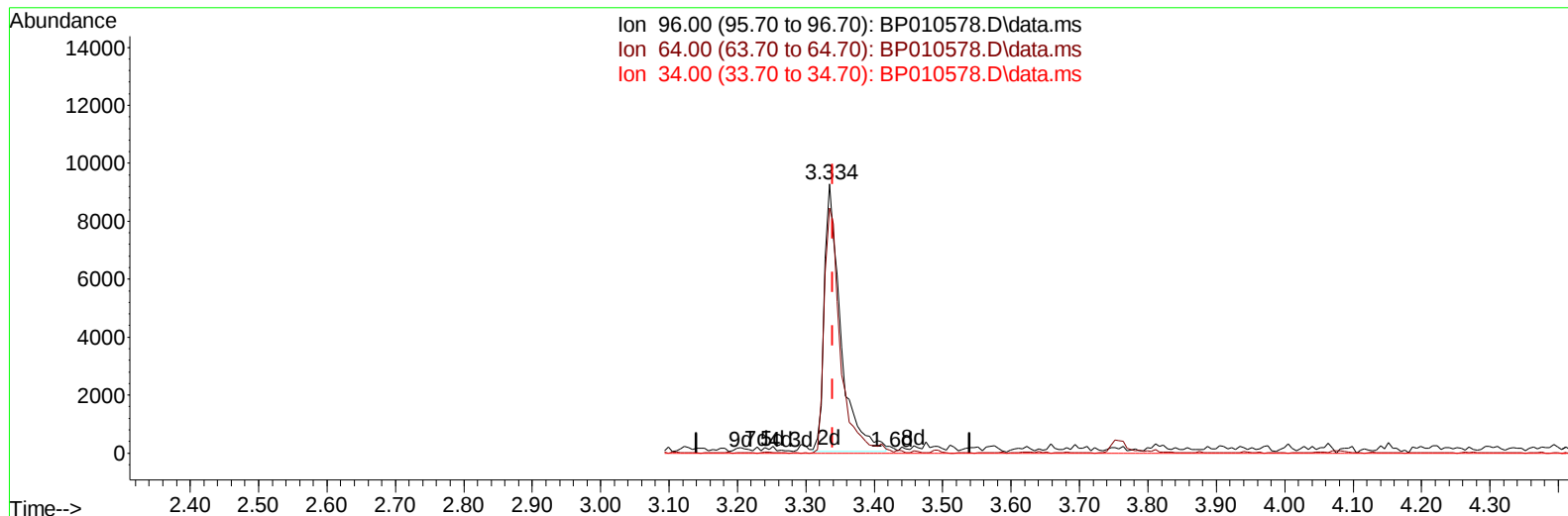
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TIC: BP010578.D\data.ms

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

3.334min (-0.006) 5.52 ng/uL m

response 15490

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	91.23#
34.00	0.00	0.00
0.00	0.00	0.00

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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	119713	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	523053	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.510	164	367825	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.263	188	835318	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.351	240	766009	20.000	ng/ul	# 0.00
88) Perylene-d12	23.768	264	653017	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	15490m	5.515	ng/uL	0.00
4) Pyridine-d5	3.752	84	217748	27.305	ng/ul	0.00
7) Phenol-d5	7.046	99	270550	26.926	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	199175	29.504	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	215695	28.054	ng/ul	-0.01
15) 4-Methylphenol-d8	8.587	113	231092	27.063	ng/ul	-0.01
21) Nitrobenzene-d5	9.040	128	120781	29.837	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	125442	29.240	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	213550	26.326	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	263959	21.885	ng/ul	-0.01
46) Dimethylphthalate-d6	13.922	166	854929	31.547	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	945255	30.239	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	107130	22.209	ng/ul	0.00
60) Fluorene-d10	15.504	176	725623	30.777	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	112765	23.083	ng/ul	0.00
73) Anthracene-d10	17.363	188	1197392	31.657	ng/ul	0.00
81) Pyrene-d10	19.598	212	1444276	35.503	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	1062545	32.404	ng/ul	0.00

Target Compounds Qvalue

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

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