

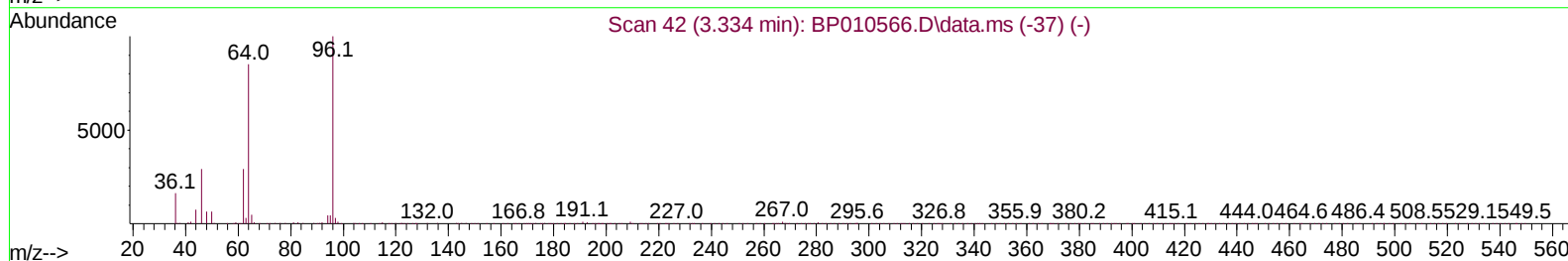
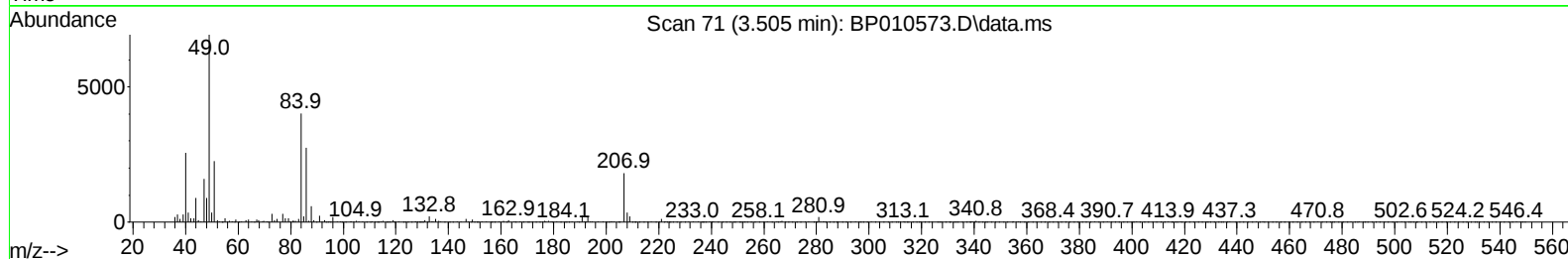
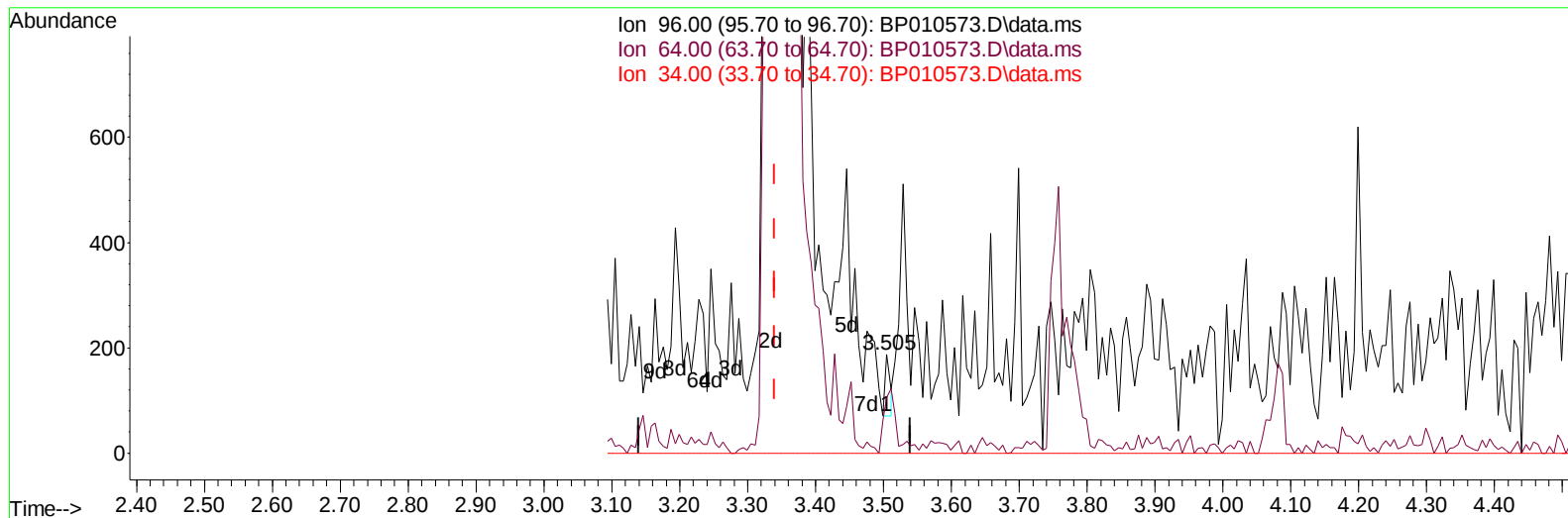
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052722\
 Data File : BP010573.D
 Acq On : 27 May 2022 09:17
 Operator : CG/JU
 Sample : PB145094BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK094

Manual IntegrationsAPPROVED

Quant Time: May 27 12:41:24 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 05/27/2022
 Supervised By :mohammad ahmed 05/31/2022



TIC: BP010573.D\data.ms

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

3.505min (+ 0.165) 0.02 ng/uL

response 59

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.68
34.00	0.00	0.00
0.00	0.00	0.00

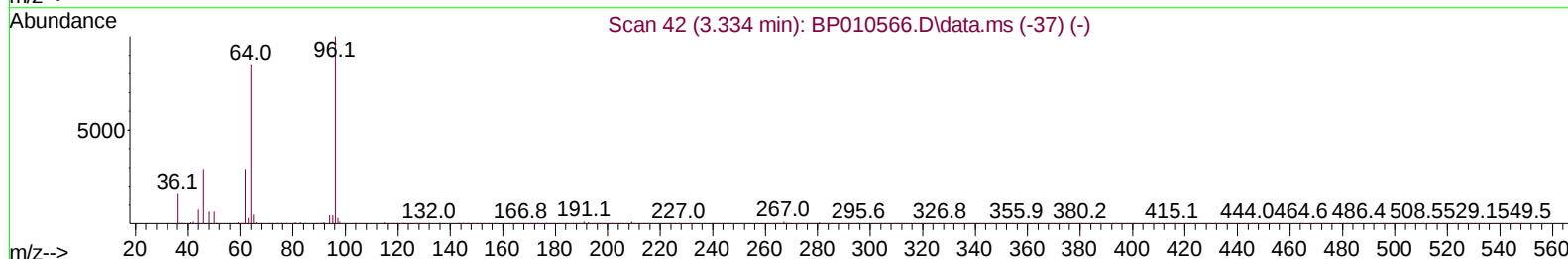
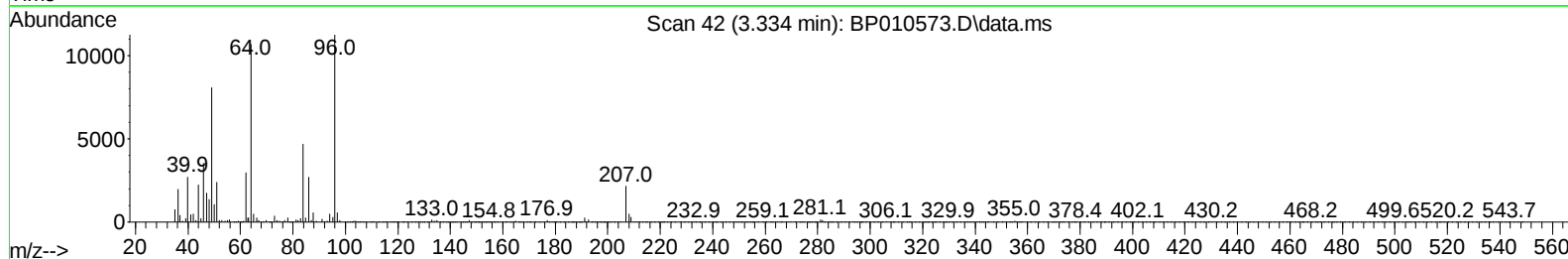
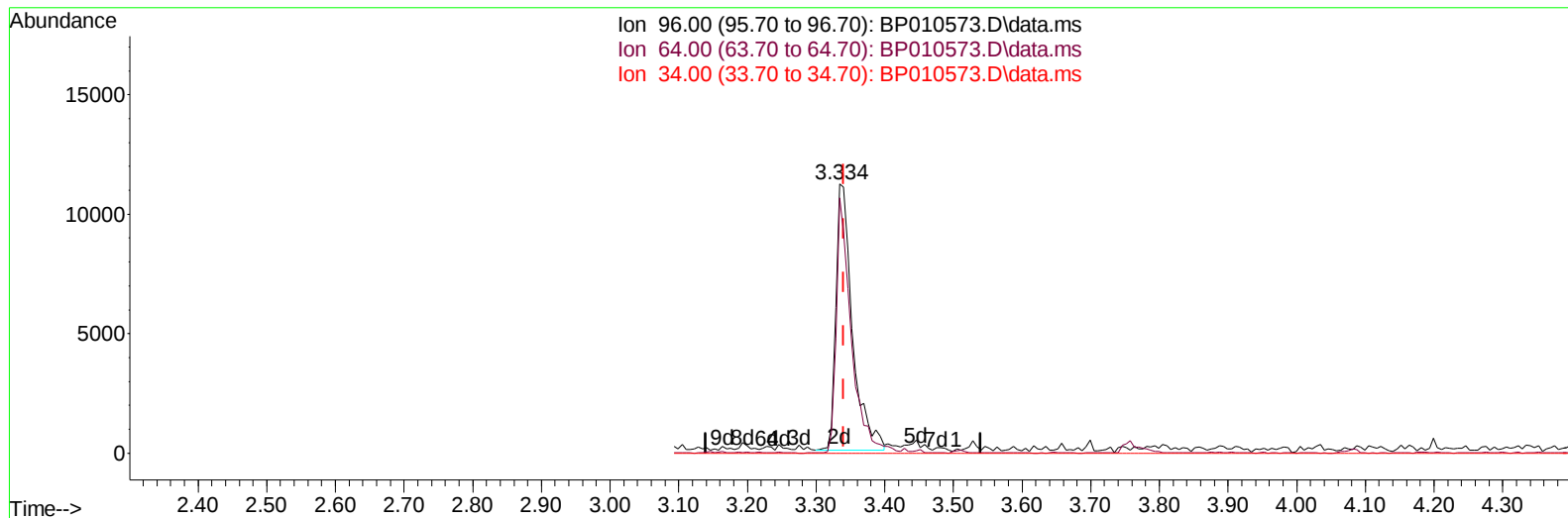
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TIC: BP010573.D\data.ms

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

3.334min (-0.006) 5.83 ng/uL m

response 18818

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	94.79#
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	7.881	152	137633	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	596398	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.516	164	419315	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	921611	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	773782	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	620461	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	18818m	5.828	ng/uL	0.00
4) Pyridine-d5	3.752	84	255142	27.829	ng/ul	0.00
7) Phenol-d5	7.052	99	311935	27.003	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	230456	29.693	ng/ul	0.00
11) 2-Chlorophenol-d4	7.417	132	254563	28.799	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	269478	27.449	ng/ul	0.00
21) Nitrobenzene-d5	9.040	128	138510	30.008	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	147286	30.110	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	247680	26.778	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	303892	22.097	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	962814	31.165	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	1070629	30.044	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	108077	19.654	ng/ul	0.00
60) Fluorene-d10	15.510	176	808248	30.072	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	121511	22.544	ng/ul	0.00
73) Anthracene-d10	17.369	188	1321330	31.662	ng/ul	0.00
81) Pyrene-d10	19.604	212	1548916	37.693	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.616	264	1004869	32.253	ng/ul	0.00

Target Compounds Qvalue

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

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