

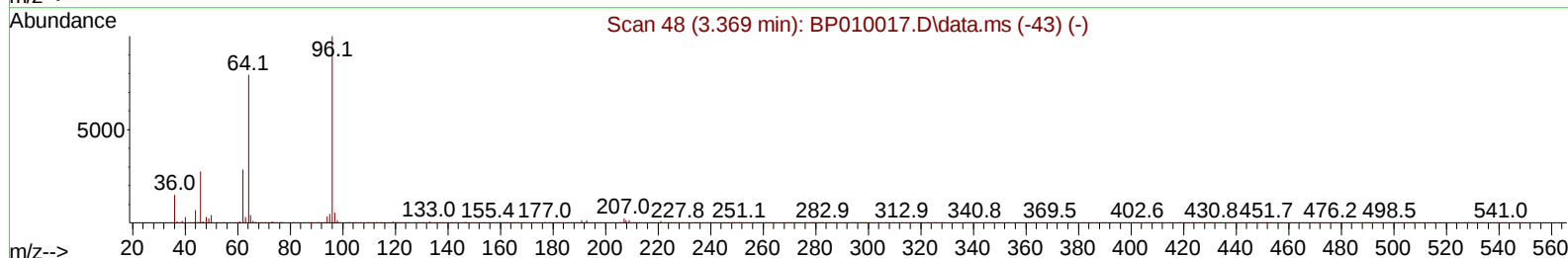
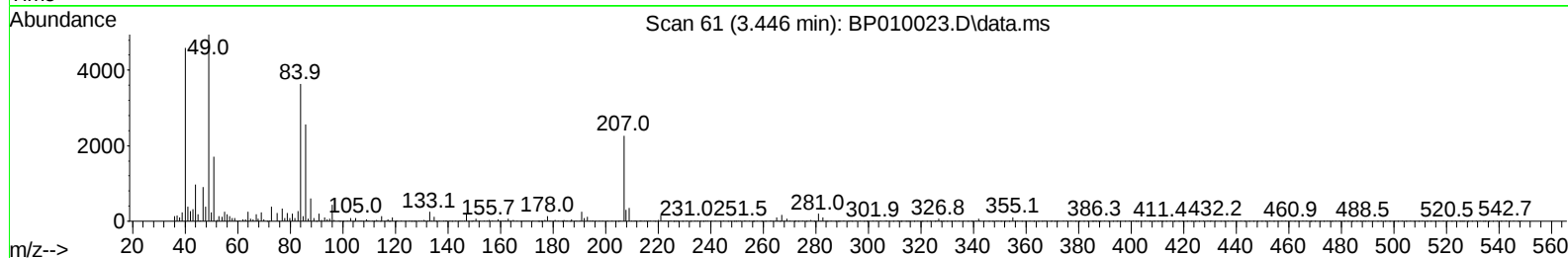
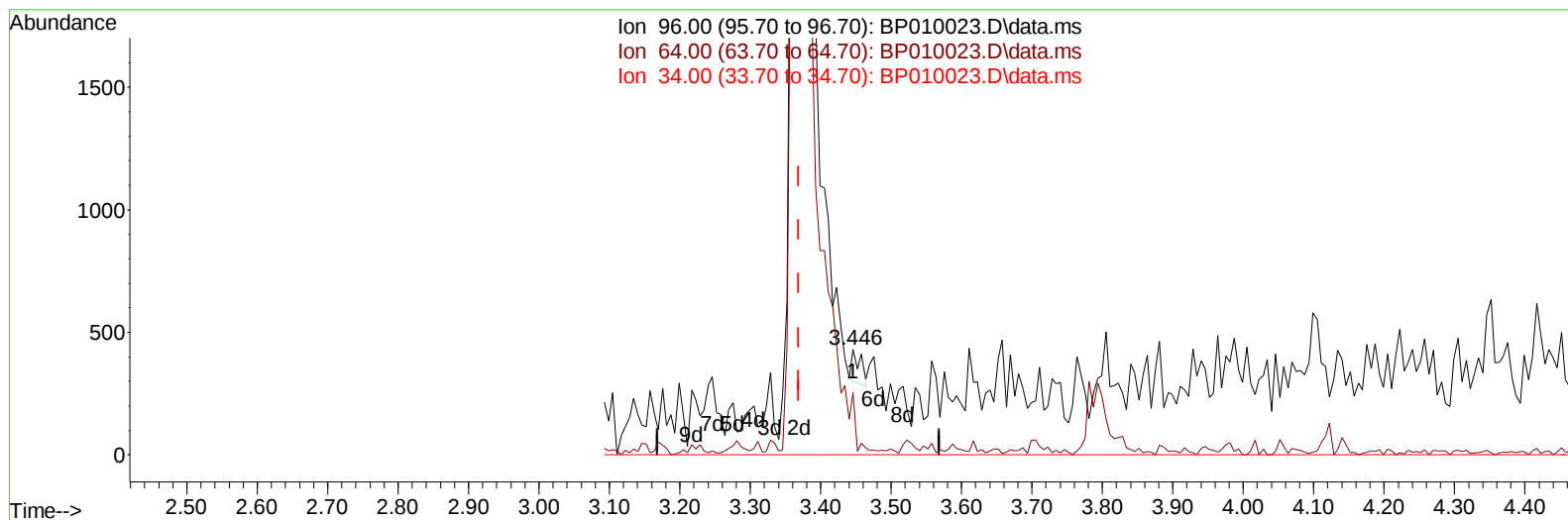
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010023.D
 Acq On : 21 Apr 2022 20:07
 Operator : CG/JU
 Sample : N2373-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFD3

Manual IntegrationsAPPROVED

Quant Time: Apr 22 04:19:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010023.D\data.ms

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

3.446min (+ 0.077) 0.05 ng/uL

response 116

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	59.07
34.00	0.00	0.00
0.00	0.00	0.00

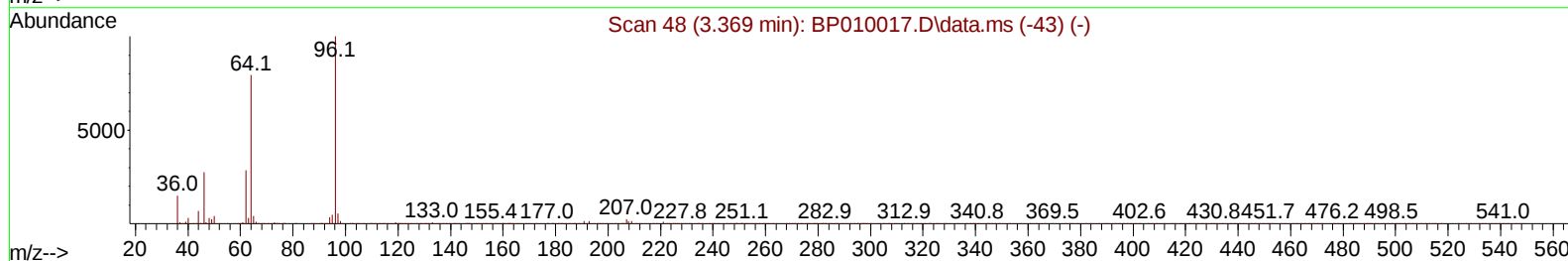
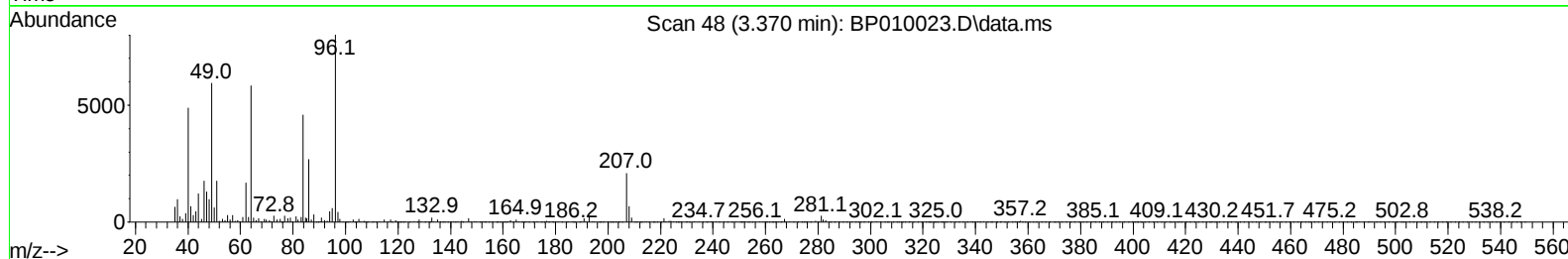
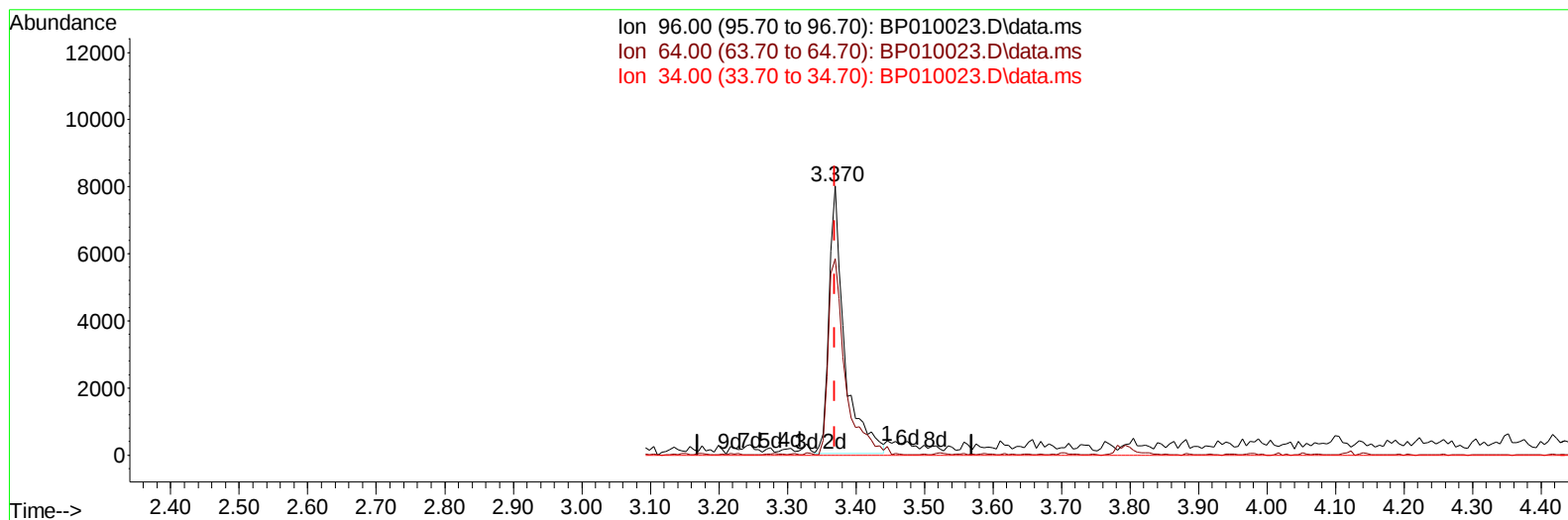
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TIC: BP010023.D\data.ms

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

3.370min (+ 0.001) 5.65 ng/uL m

response 12476

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.02#
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	7.940	152	96211	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	443766	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	322692	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	739402	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	777254	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	736584	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	12476m	5.655	ng/uL	0.00
4) Pyridine-d5	3.787	84	165425	26.652	ng/ul	0.00
7) Phenol-d5	7.099	99	252227	28.822	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	166432	31.694	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	206618	31.880	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	226975	31.304	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	112035	31.743	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	123481	32.034	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	220085	29.329	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	316719	29.436	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	818256	32.597	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	926933	30.511	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	113664	24.596	ng/ul	0.00
60) Fluorene-d10	15.563	176	710032	32.883	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	95188	20.360	ng/ul	0.00
73) Anthracene-d10	17.422	188	1139592	31.858	ng/ul	0.00
81) Pyrene-d10	19.651	212	1438217	32.537	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1298567	33.628	ng/ul	0.00

Target Compounds Qvalue

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

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