

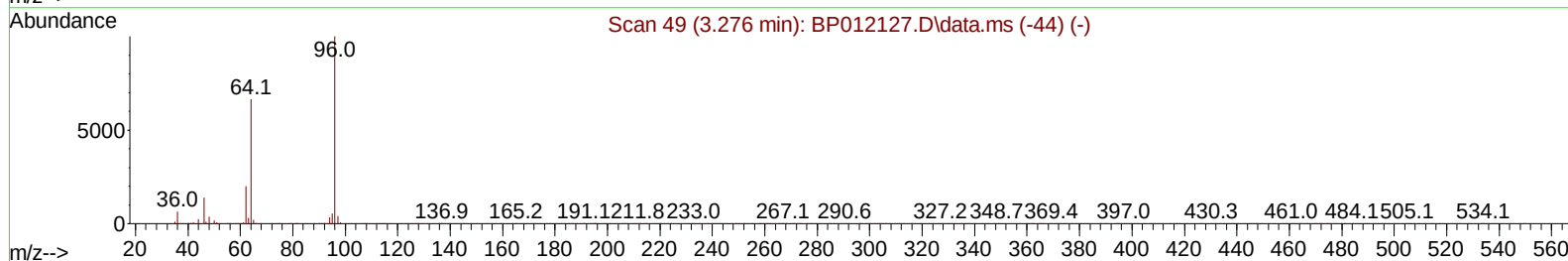
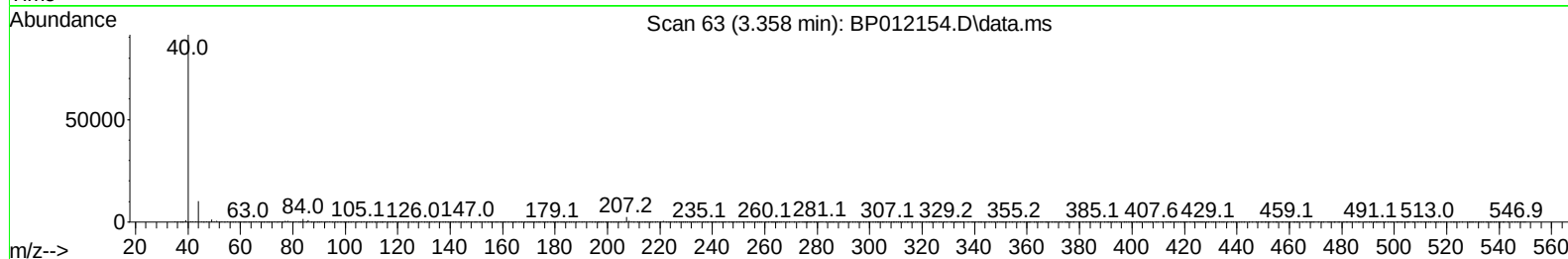
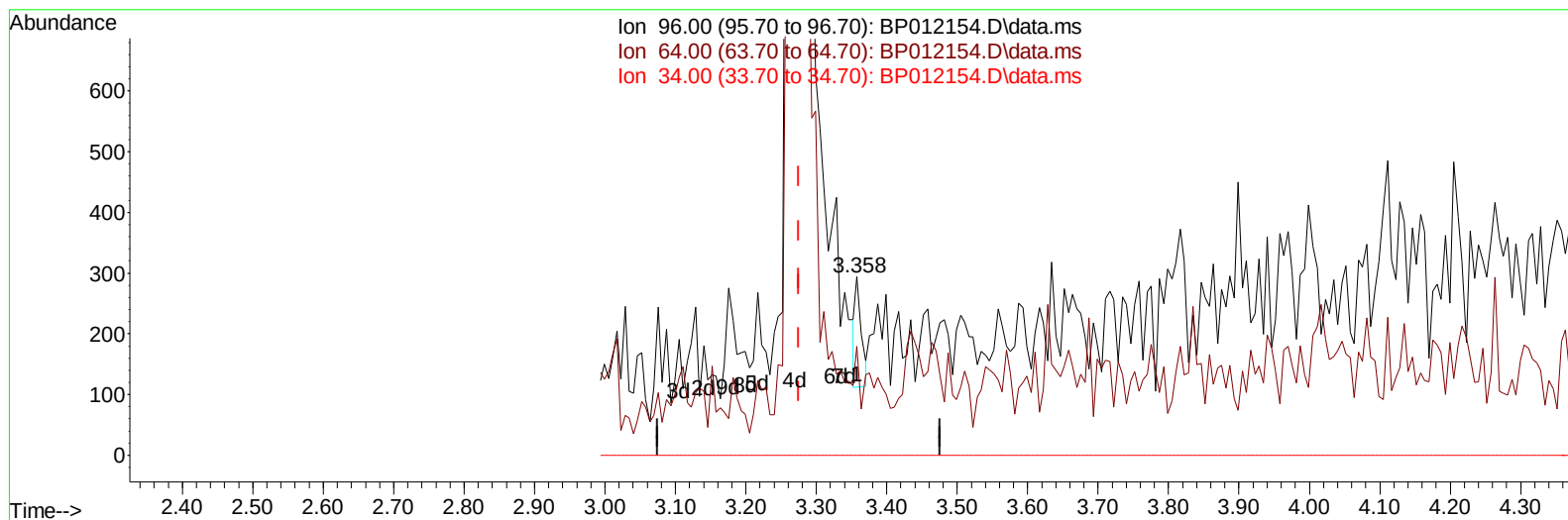
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012154.D
 Acq On : 15 Oct 2022 01:01
 Operator : CG/JU
 Sample : N4984-09
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGNB6

Manual IntegrationsAPPROVED

Quant Time: Oct 15 02:25:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/15/2022
 Supervised By :mohammad ahmed 10/16/2022



TIC: BP012154.D\data.ms

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

3.358min (+ 0.082) 0.02 ng/uL

response 109

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.10	60.88
34.00	0.00	0.00
0.00	0.00	0.00

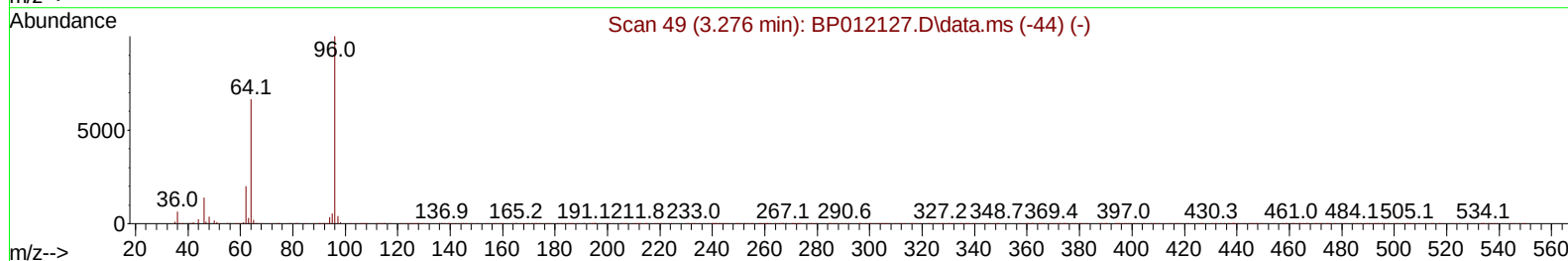
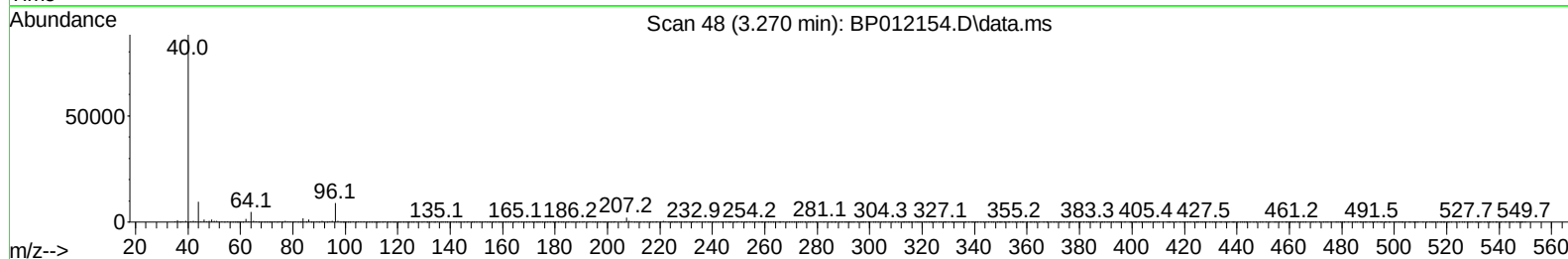
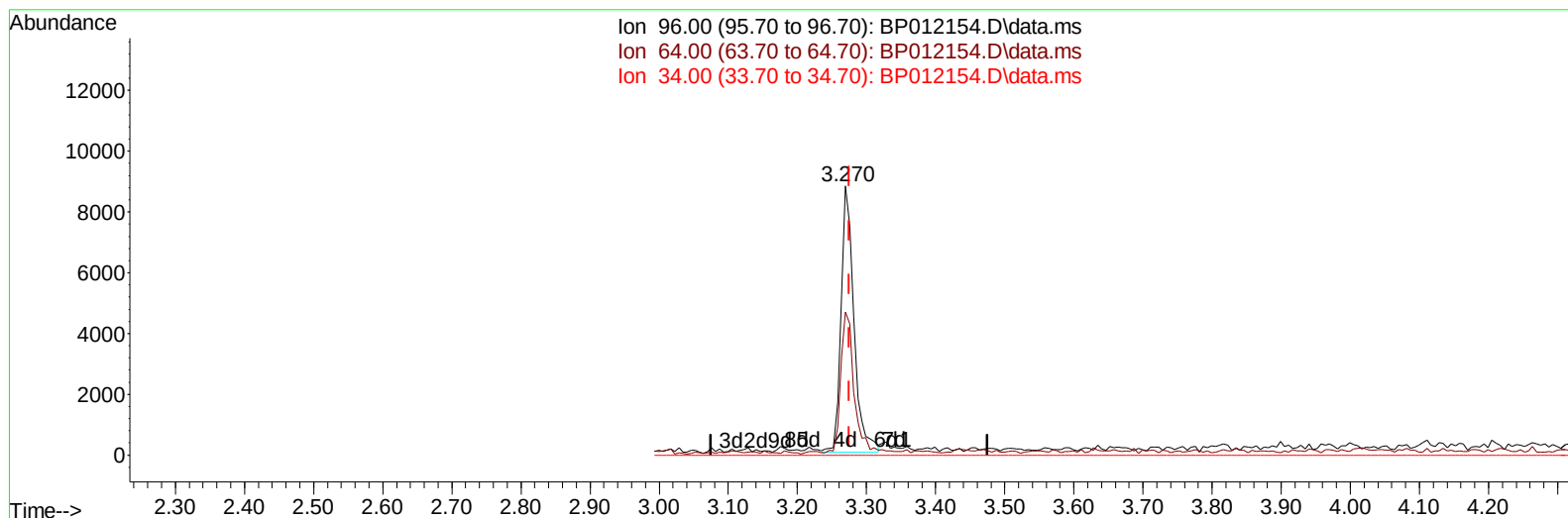
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(3) 1,4-Dioxane-d8 (S)

3.270min (-0.006) 1.88 ng/uL m

response 11378

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.10	53.21#
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	7.840	152	238595	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.646	136	967937	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.493	164	552609	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1192276	20.000	ng/ul	0.00
79) Chrysene-d12	21.339	240	1231537	20.000	ng/ul	# 0.00
88) Perylene-d12	23.757	264	1315655	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	11378m	1.880	ng/uL	0.00
4) Pyridine-d5	3.705	84	26568	1.651	ng/ul	0.00
7) Phenol-d5	7.011	99	488248	25.444	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	285328	26.196	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.369	132	388229	25.488	ng/ul	-0.01
15) 4-Methylphenol-d8	8.558	113	362254	24.578	ng/ul	0.00
21) Nitrobenzene-d5	9.011	128	176166	25.770	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	178735	25.785	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.269	165	320453	23.854	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.793	131	377769	19.083	ng/ul	-0.01
46) Dimethylphthalate-d6	13.904	166	893165	25.521	ng/ul	-0.01
49) Acenaphthylene-d8	14.181	160	1110675	25.664	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.704	143	120000	17.465	ng/ul	0.00
60) Fluorene-d10	15.487	176	863032	27.256	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.610	200	62580	9.940	ng/ul	-0.01
73) Anthracene-d10	17.351	188	1370834	26.478	ng/ul	0.00
81) Pyrene-d10	19.586	212	1757877	27.862	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.598	264	1805296	28.227	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.034	94	26443	1.313	ng/ul#	89

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

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