

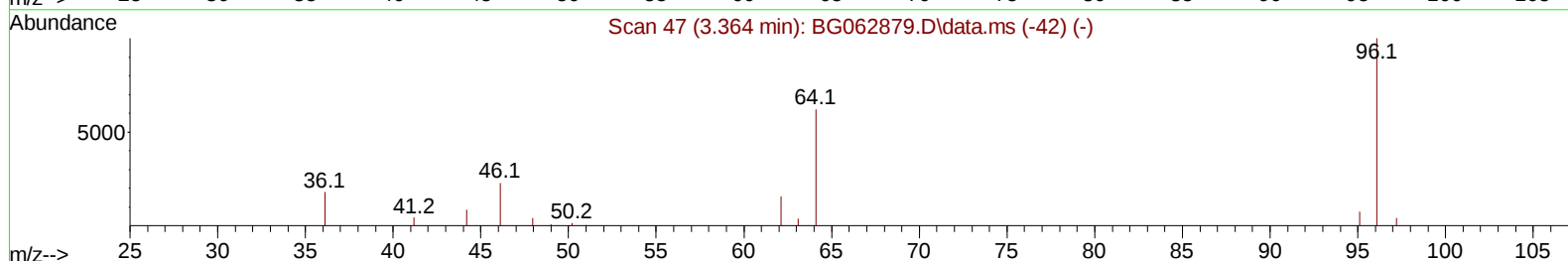
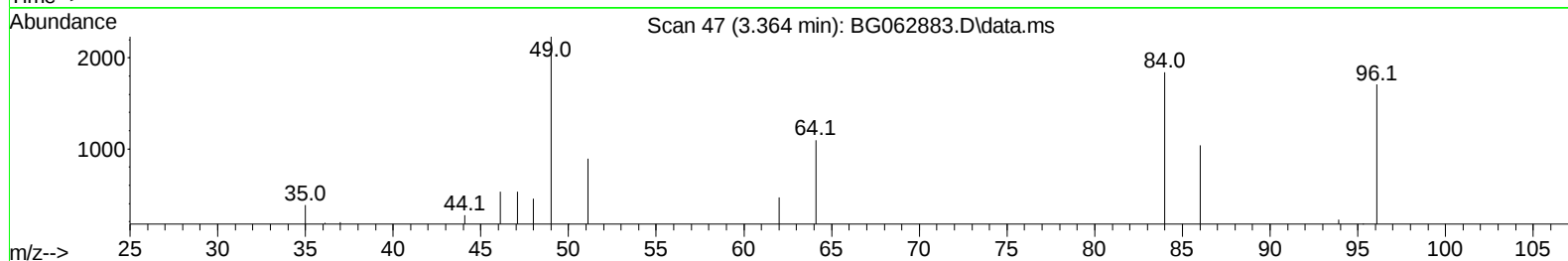
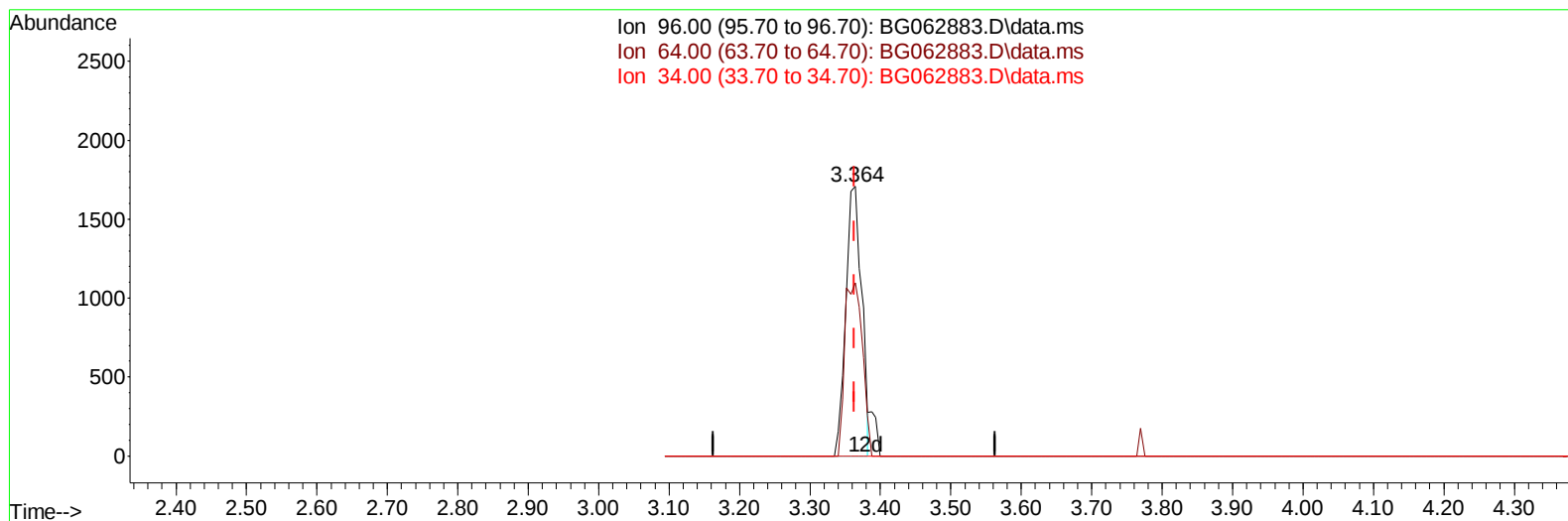
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062883.D
Acq On : 16 Sep 2024 17:32
Operator : JU/RC
Sample : P3899-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC73

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/17/2024
Supervised By :mohammad ahmed 09/17/2024

Quant Time: Sep 16 18:26:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 01:28:17 2024
Response via : Initial Calibration



TIC: BG062883.D\data.ms

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

3.364min (+ 0.001) 3.09 ng/uL

response 2638

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	62.00	64.32
34.00	0.00	0.00
0.00	0.00	0.00

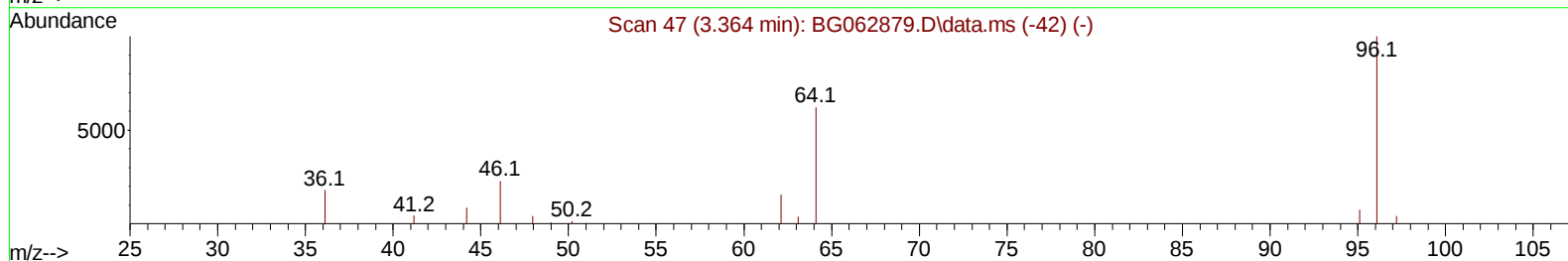
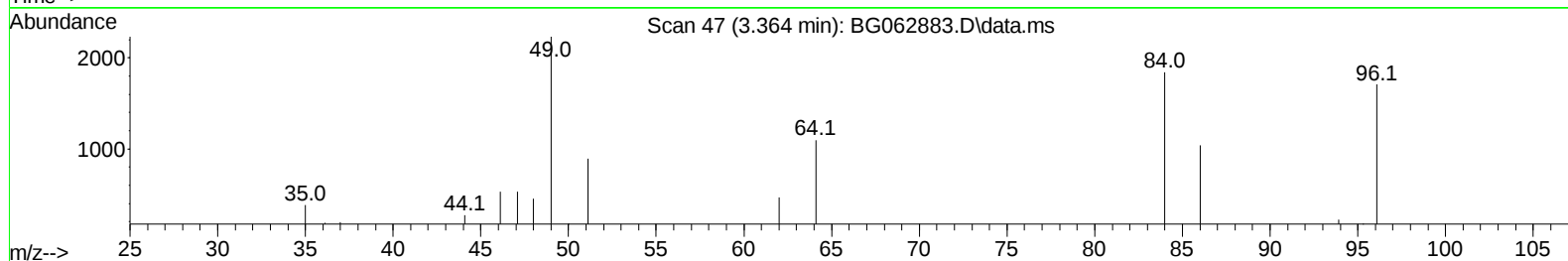
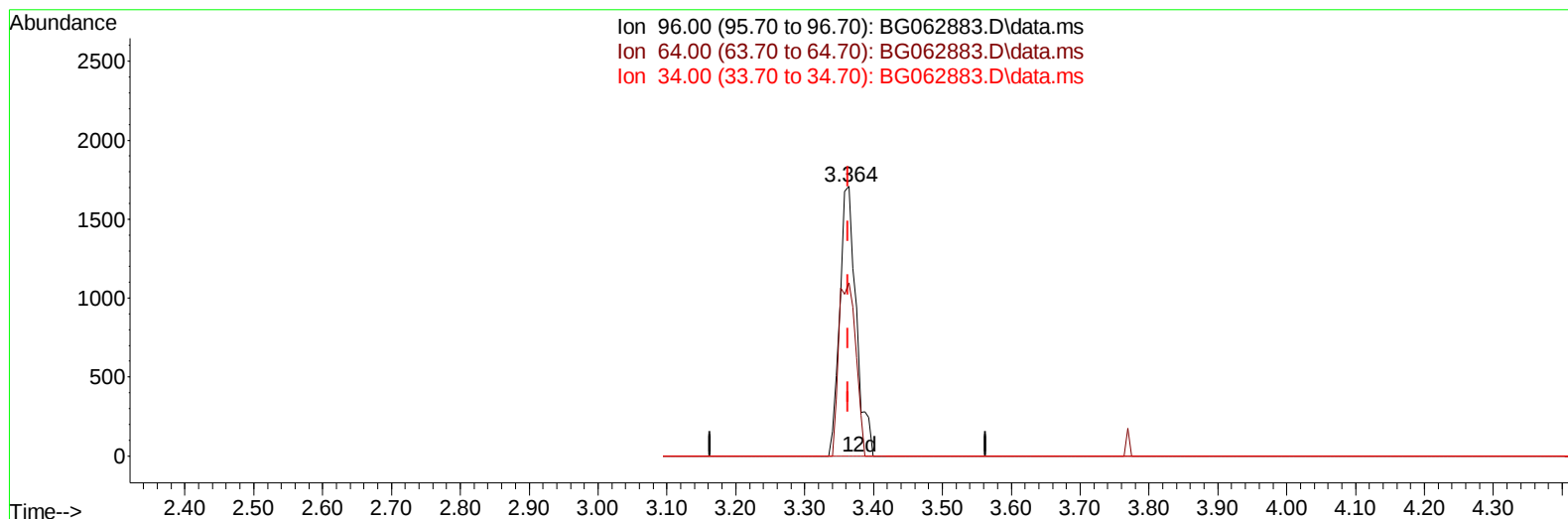
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TIC: BG062883.D\data.ms

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

3.364min (+ 0.001) 3.31 ng/uL m

response 2823

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	62.00	64.32
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	7.900	152	38548	20.000	ng/ul	0.00
20) Naphthalene-d8	10.691	136	152049	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.527	164	131195	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.277	188	385517	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.519	240	453647	20.000	ng/ul	# 0.00
88) Perylene-d12	24.586	264	518831	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	2823m	3.310	ng/uL	0.00
4) Pyridine-d5	3.769	84	28554	10.519	ng/ul	0.00
7) Phenol-d5	7.071	99	58214	17.222	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.230	67	32377	15.976	ng/ul	0.00
11) 2-Chlorophenol-d4	7.436	132	47836	19.429	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	52889	18.776	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	23937	22.027	ng/ul	0.00
24) 2-Nitrophenol-d4	9.774	143	31037	26.115	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.315	165	58421	22.572	ng/ul	0.00
31) 4-Chloroaniline-d4	10.826	131	51764	14.550	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	217396	21.518	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	238929	21.294	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	32858	19.145	ng/ul	0.00
60) Fluorene-d10	15.520	176	200853	22.062	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.638	200	42793	19.297	ng/ul	0.00
73) Anthracene-d10	17.371	188	381214	20.953	ng/ul	0.00
81) Pyrene-d10	19.663	212	524046	20.528	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.375	264	597042	21.793	ng/ul	0.00
Target Compounds						
					Qvalue	
58) 2,3,4,6-Tetrachlorophenol	15.156	232	60524	20.361	ng/ul#	97
71) Pentachlorophenol	16.930	266	497458	164.199	ng/ul	97

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

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