

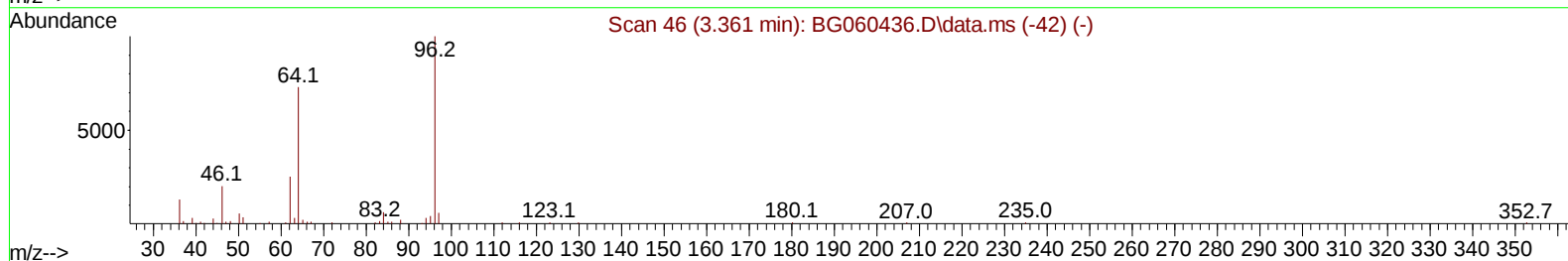
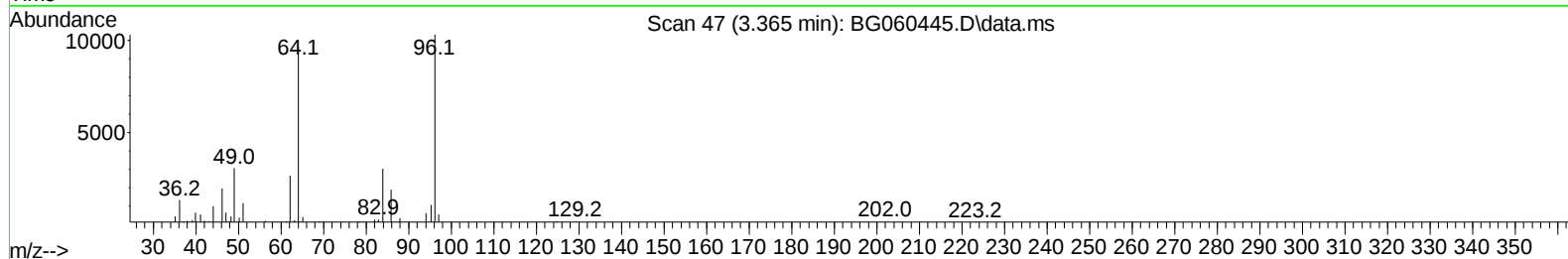
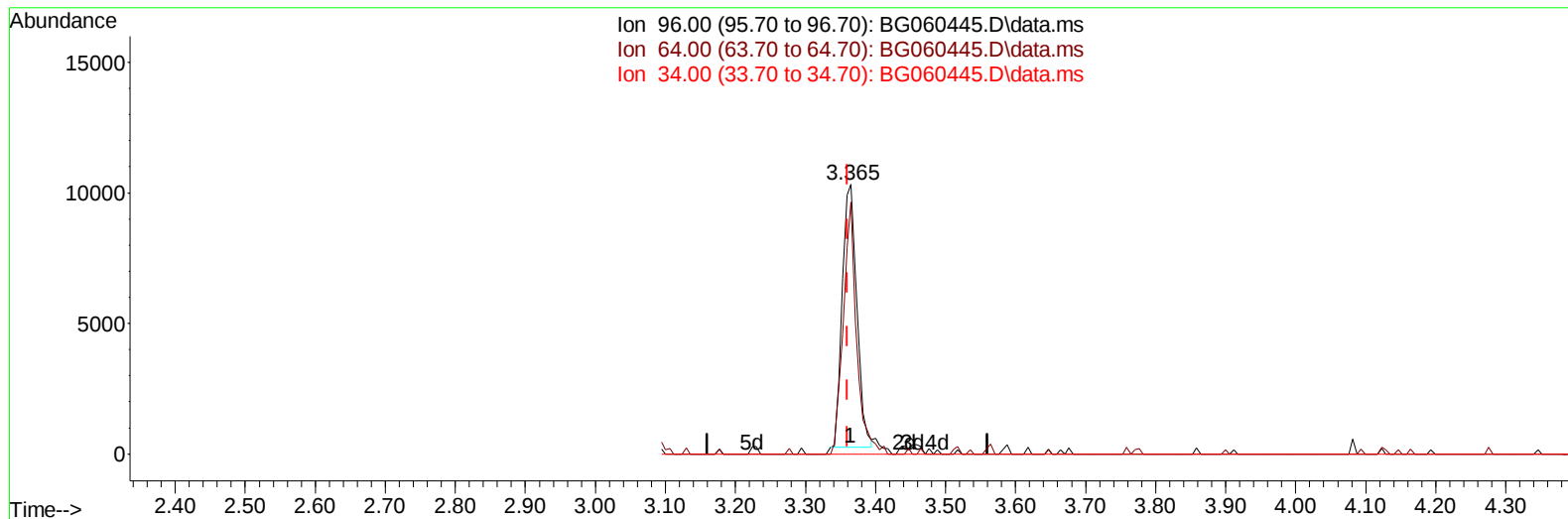
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012624\
Data File : BG060445.D
Acq On : 26 Jan 2024 17:37
Operator : MA/JU
Sample : P1238-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH703

Manual IntegrationsAPPROVED

Quant Time: Jan 26 22:04:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 26 06:06:34 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/29/2024
Supervised By :mohammad ahmed 01/30/2024



TIC: BG060445.D\data.ms

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

3.365min (+ 0.005) 4.34 ng/uL

response 14525

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.30	93.50#
34.00	0.00	0.00
0.00	0.00	0.00

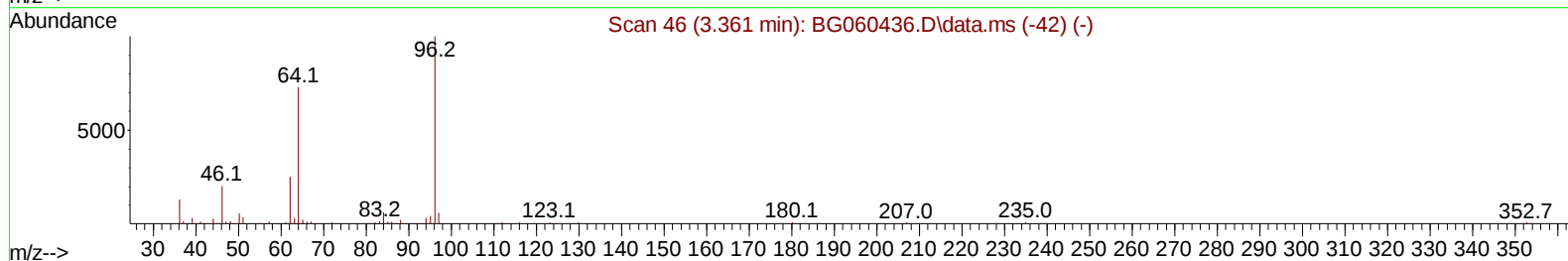
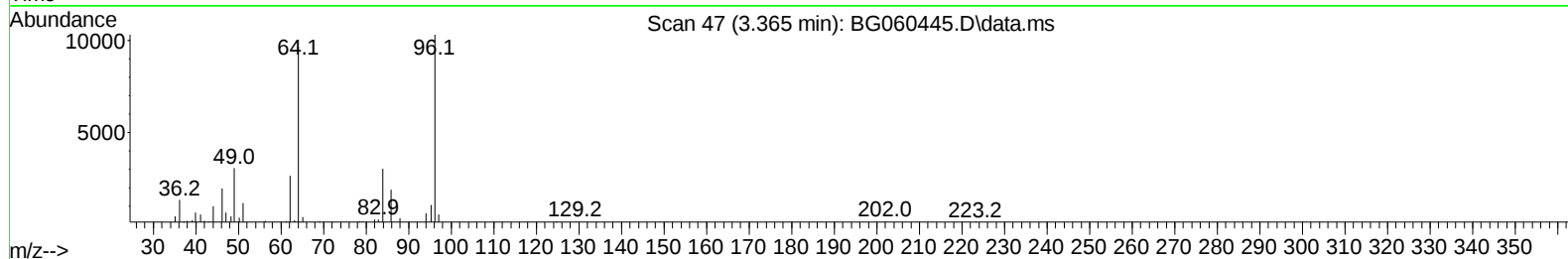
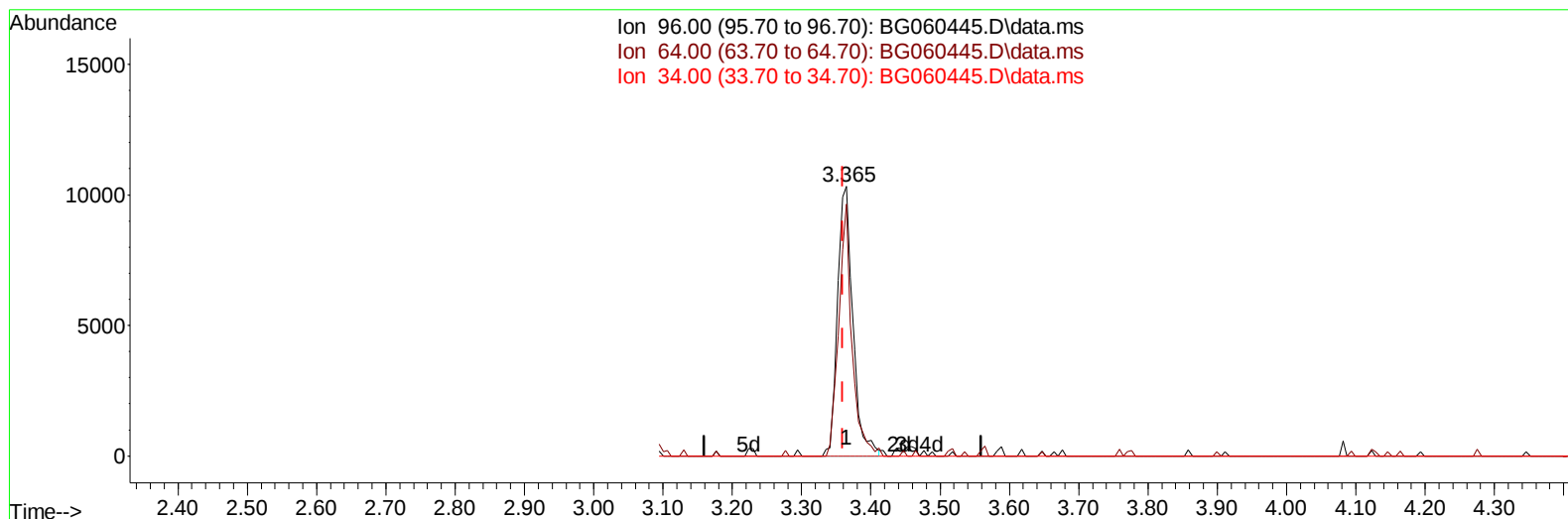
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TIC: BG060445.D\data.ms

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

3.365min (+ 0.005) 4.77 ng/uL m

response 15958

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.30	93.50#
34.00	0.00	0.00
0.00	0.00	0.00

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 QLast Update : Fri Jan 26 06:06:34 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.924	152	167697	20.000	ng/ul	0.00
20) Naphthalene-d8	10.727	136	754089	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	603104	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.307	188	1499004	20.000	ng/ul	0.00
79) Chrysene-d12	21.573	240	1436292	20.000	ng/ul	0.00
88) Perylene-d12	24.734	264	1604659	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.365	96	15958m	4.768	ng/uL	0.00
4) Pyridine-d5	3.776	84	96501	9.116	ng/ul	0.00
7) Phenol-d5	7.090	99	99450	6.244	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.248	67	237441	23.724	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	210192	19.763	ng/ul	0.00
15) 4-Methylphenol-d8	8.629	113	153404	11.693	ng/ul	0.00
21) Nitrobenzene-d5	9.088	128	144412	25.960	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	167636	25.794	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	343676	25.427	ng/ul	0.00
31) 4-Chloroaniline-d4	10.862	131	313799	17.698	ng/ul	0.00
46) Dimethylphthalate-d6	13.964	166	1314281	26.631	ng/ul	0.00
49) Acenaphthylene-d8	14.258	160	1461181	26.395	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	37401	5.291	ng/ul	0.00
60) Fluorene-d10	15.556	176	1208038	27.937	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.668	200	233942	25.204	ng/ul	0.00
73) Anthracene-d10	17.407	188	2168633	30.519	ng/ul	0.00
81) Pyrene-d10	19.699	212	2743612	31.871	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.517	264	2654870	31.539	ng/ul	0.00

Target Compounds Qvalue

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

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