

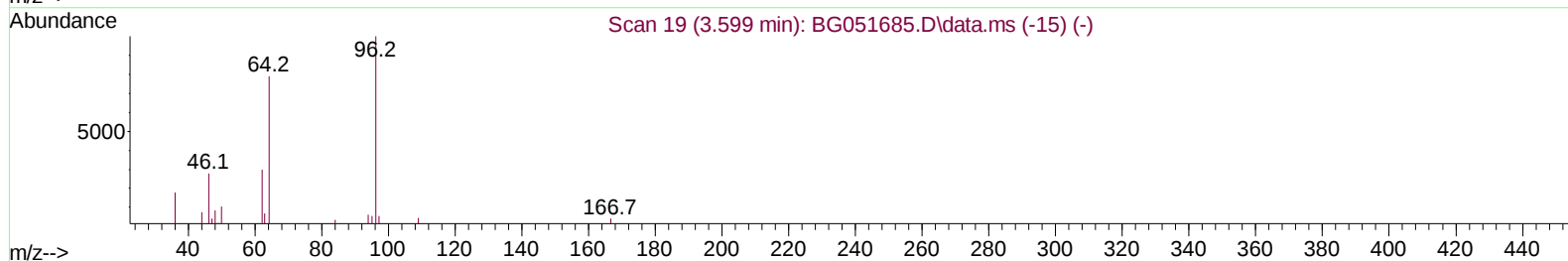
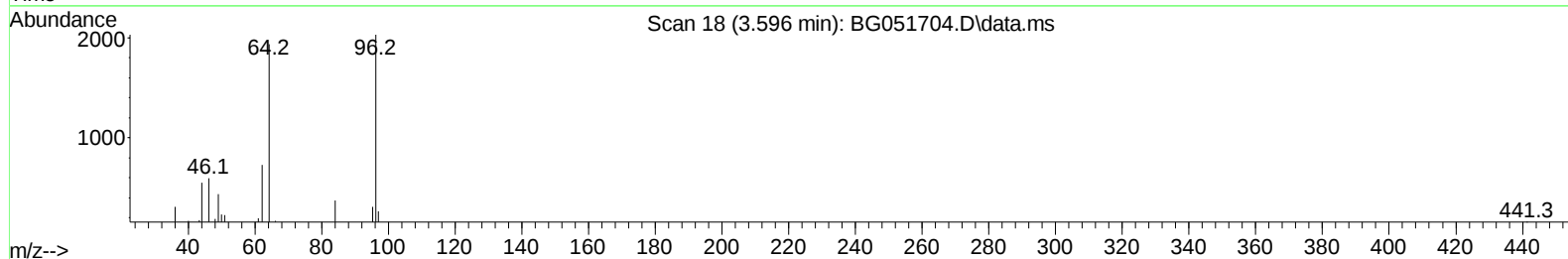
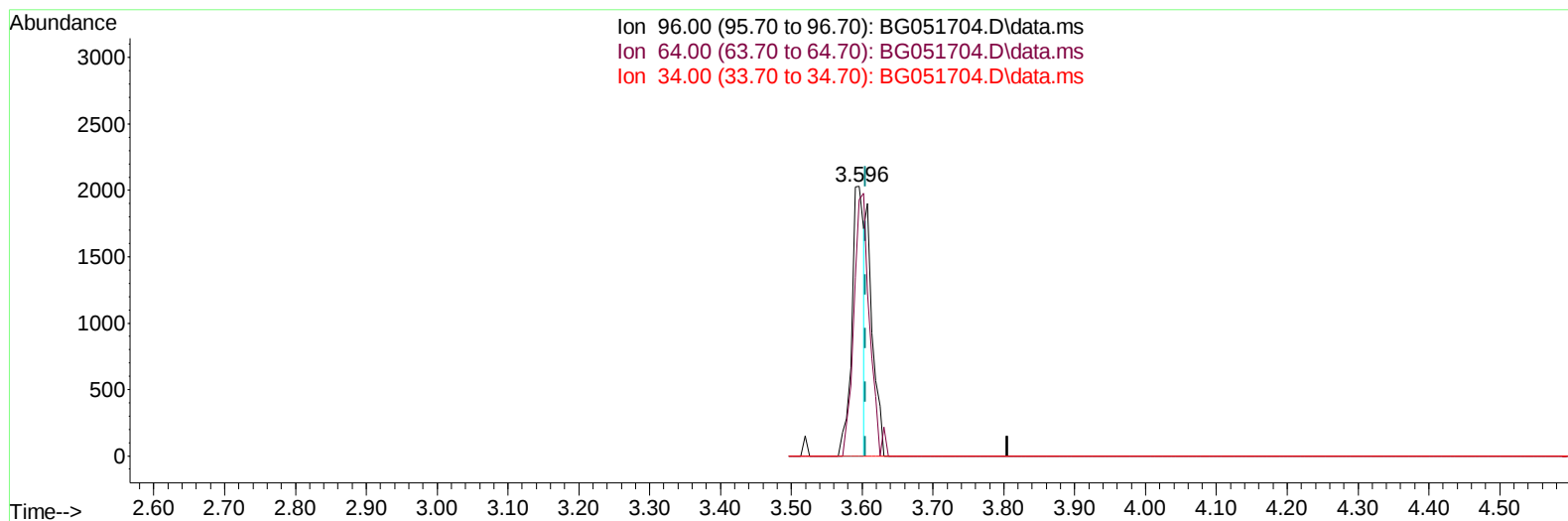
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051704.D
 Acq On : 21 Dec 2021 6:48
 Operator : CG/JU
 Sample : M5126-14
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0J06

Manual IntegrationsAPPROVED

Quant Time: Dec 21 07:33:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 14 14:19:57 2021
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/21/2021
 Supervised By :mohammad ahmed 12/28/2021



TIC: BG051704.D\data.ms

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

3.596min (-0.009) 2.97 ng/uL

response 2428

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	95.17#
34.00	0.00	0.00
0.00	0.00	0.00

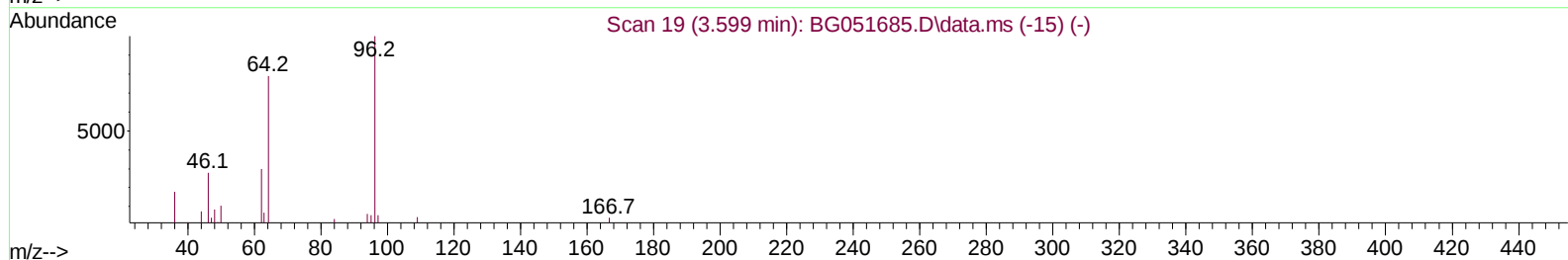
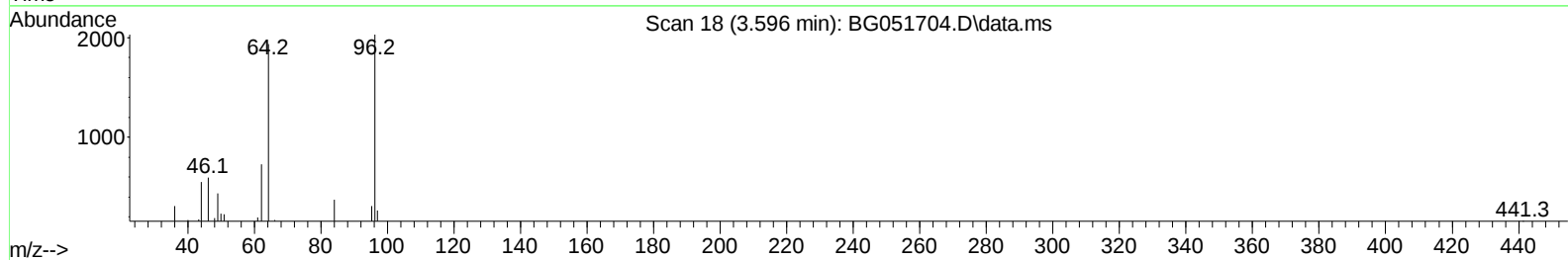
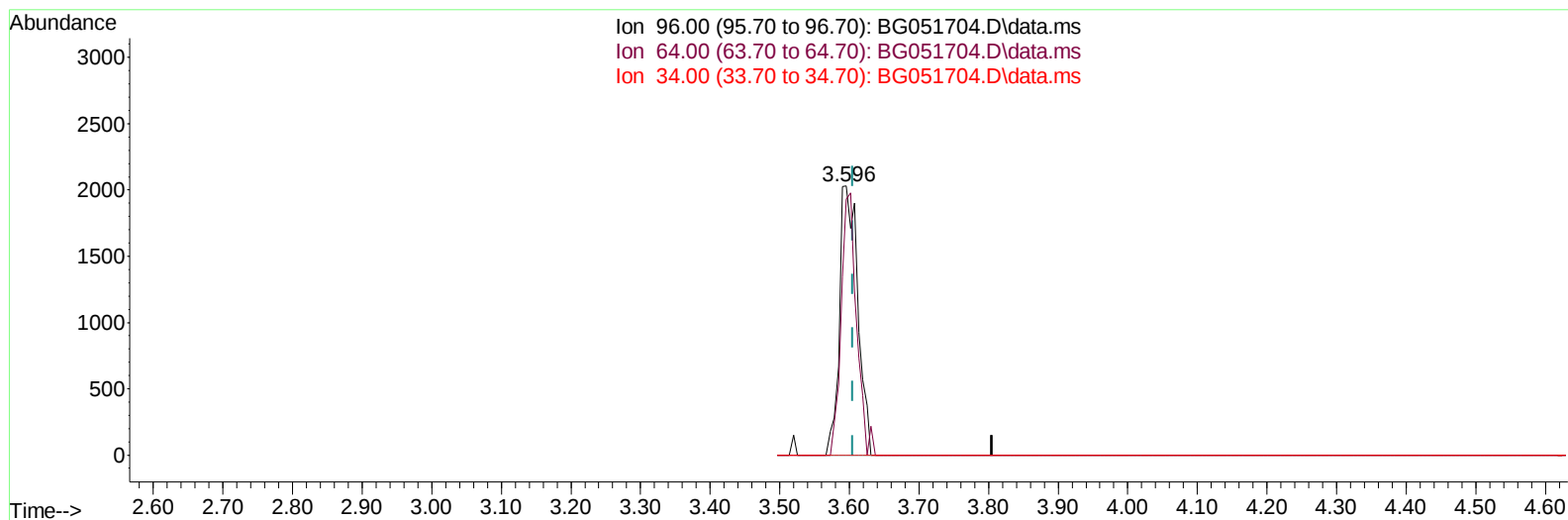
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TIC: BG051704.D\data.ms

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

3.596min (-0.009) 4.59 ng/uL m

response 3756

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	95.17#
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	8.272	152	33953	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.109	136	148473	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.910	164	102494	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.653	188	238284	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.971	240	232994	20.000	ng/ul	#-0.01
88) Perylene-d12	25.472	264	232439	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.596	96	3756m	4.594	ng/uL	0.00
4) Pyridine-d5	4.025	84	18550	7.510	ng/ul	-0.02
7) Phenol-d5	7.402	99	21801	7.191	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.573	67	66178	36.721	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.796	132	58095	27.554	ng/ul	-0.01
15) 4-Methylphenol-d8	8.965	113	40462	17.267	ng/ul	-0.01
21) Nitrobenzene-d5	9.435	128	40783	37.466	ng/ul	-0.02
24) 2-Nitrophenol-d4	10.169	143	43448	36.512	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.716	165	77565	30.058	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.227	131	69221	20.374	ng/ul	-0.01
46) Dimethylphthalate-d6	14.293	166	283475	34.555	ng/ul	-0.02
49) Acenaphthylene-d8	14.604	160	350293	34.381	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143	8237	6.174	ng/ul	0.00
60) Fluorene-d10	15.897	176	254231	34.590	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.997	200	40007	32.225	ng/ul	-0.01
73) Anthracene-d10	17.753	188	439354	38.652	ng/ul	0.00
81) Pyrene-d10	20.032	212	514207	36.669	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.225	264	493784	40.375	ng/ul	-0.03

Target Compounds Qvalue

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

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