

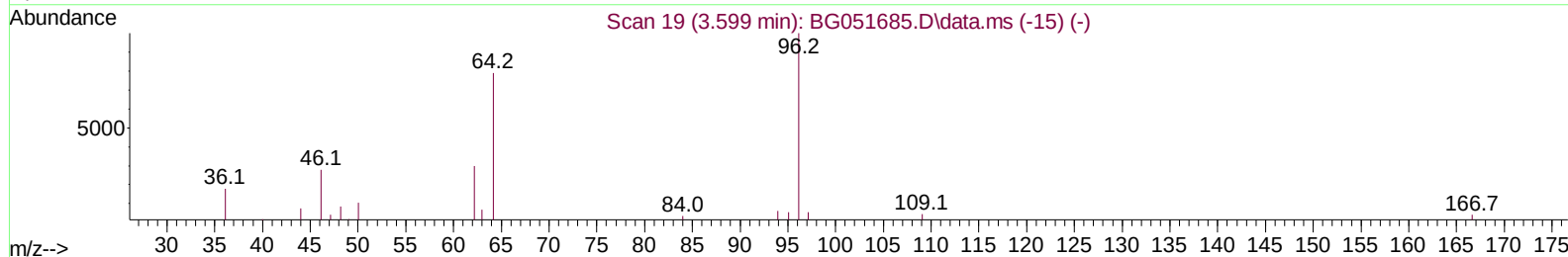
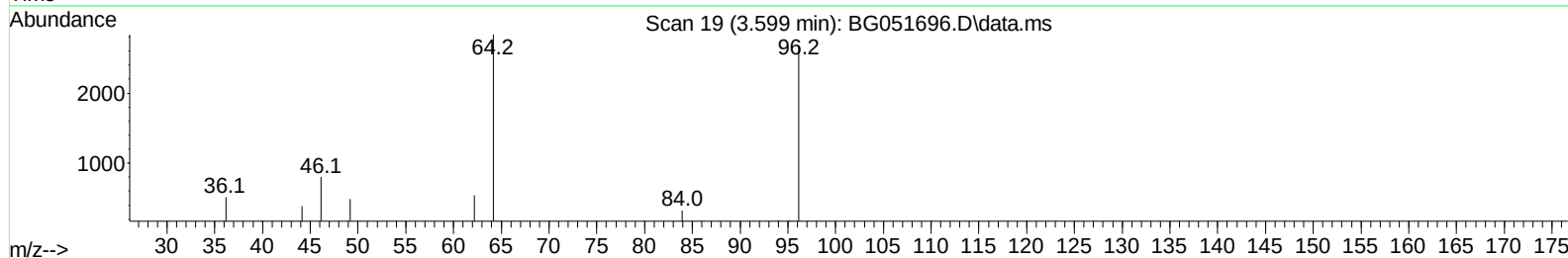
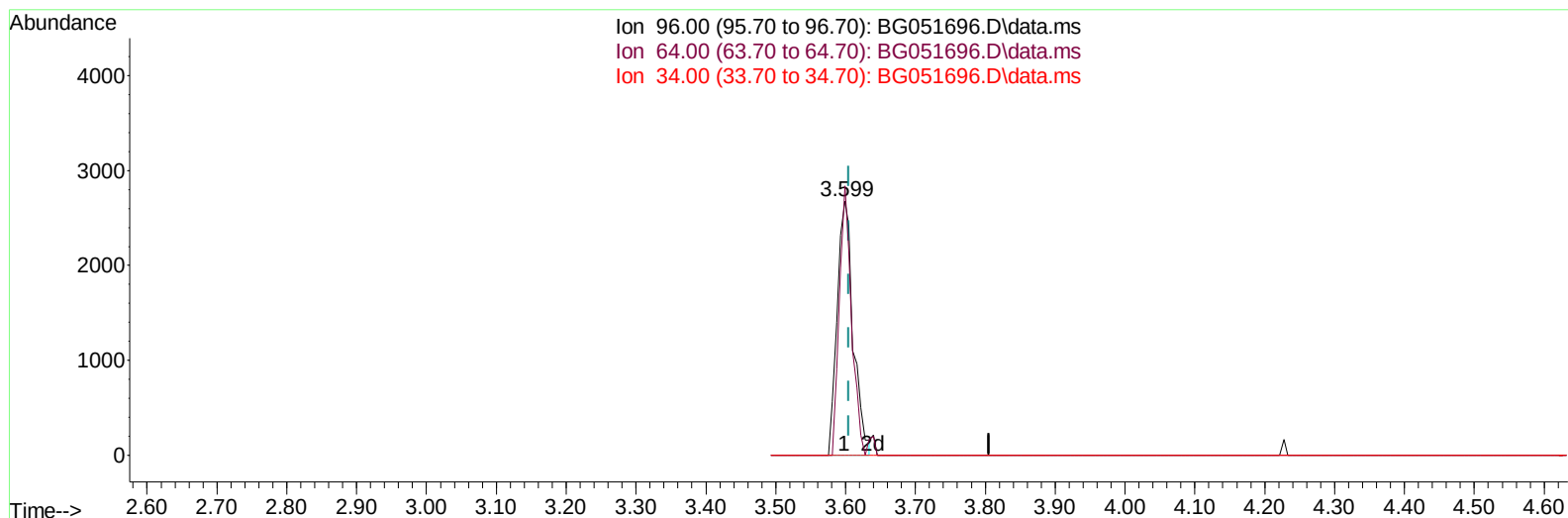
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
Data File : BG051696.D
Acq On : 21 Dec 2021 00:49
Operator : CG/JU
Sample : M5171-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLA2

Manual IntegrationsAPPROVED

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

Quant Time: Dec 21 02:29:59 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



TIC: BG051696.D\data.ms

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

3.599min (-0.006) 4.49 ng/uL

response 4321

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	105.82#
34.00	0.00	0.00
0.00	0.00	0.00

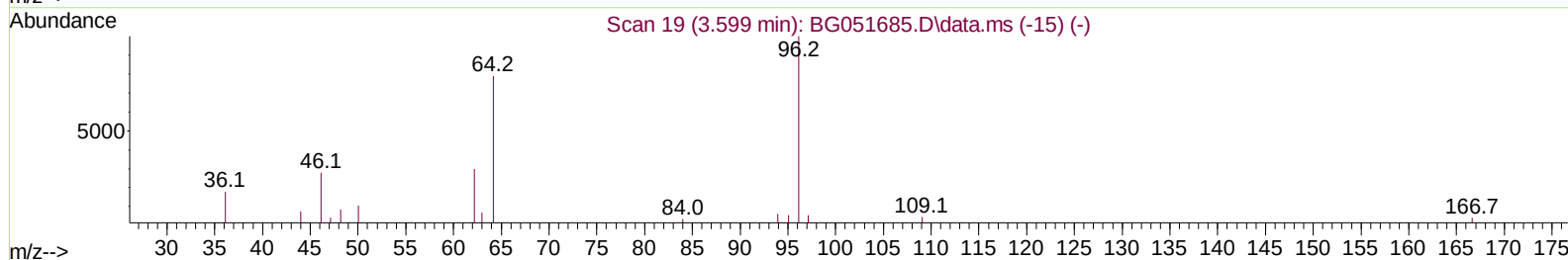
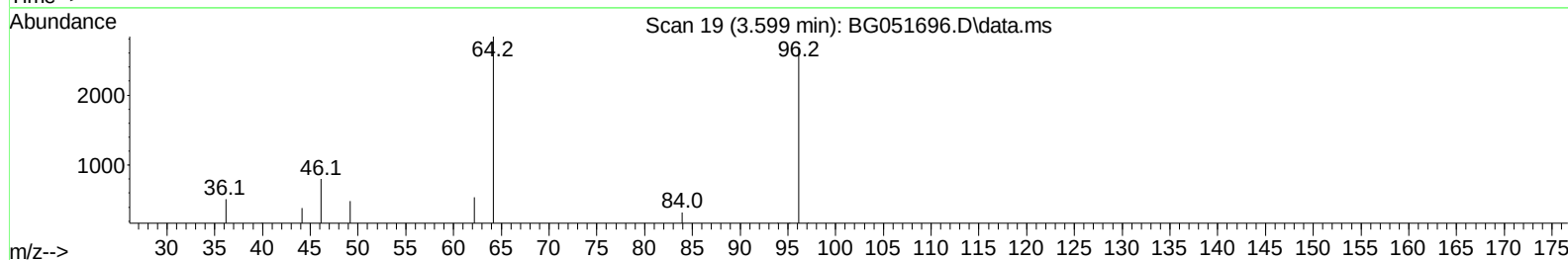
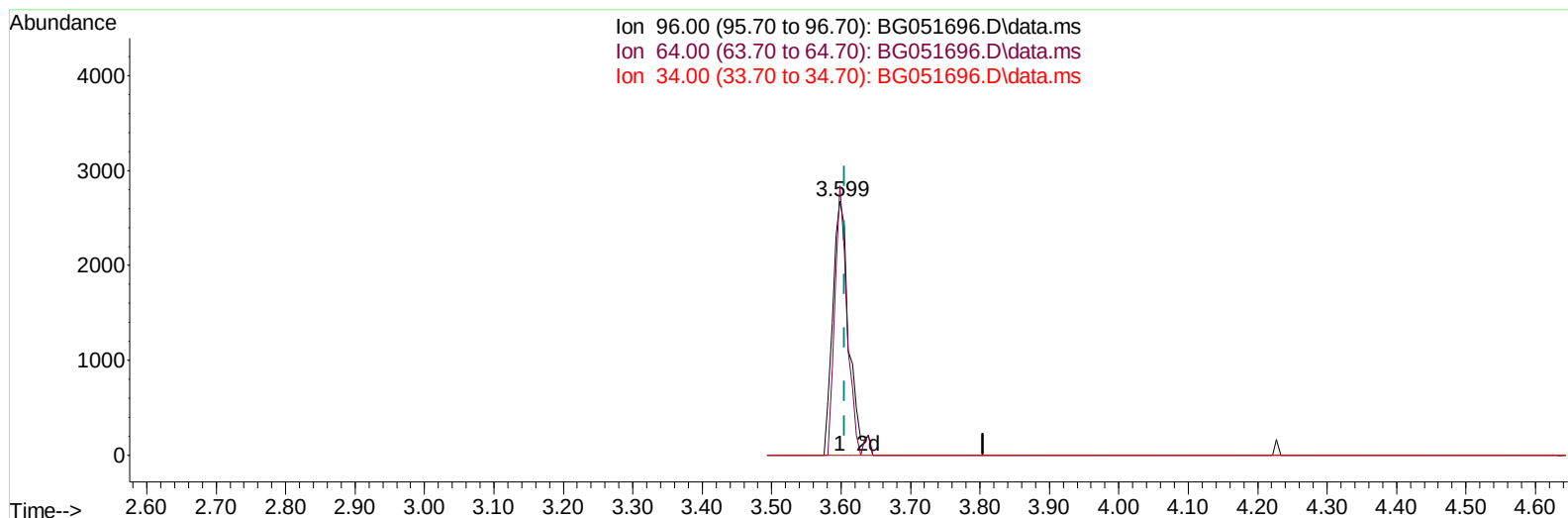
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TIC: BG051696.D\data.ms

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

3.599min (-0.006) 4.57 ng/uL m

response 4396

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	105.82#
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	8.269	152	39981	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.112	136	168905	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.907	164	122488	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.656	188	273914	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.974	240	259808	20.000	ng/ul	#-0.01
88) Perylene-d12	25.469	264	255564	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.599	96	4396m	4.566	ng/uL	0.00
4) Pyridine-d5	4.033	84	9046	3.110	ng/ul	-0.01
7) Phenol-d5	7.405	99	20740	5.810	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.576	67	61501	28.981	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.799	132	56840	22.894	ng/ul	-0.01
15) 4-Methylphenol-d8	8.968	113	39236	14.219	ng/ul	-0.01
21) Nitrobenzene-d5	9.438	128	39336	31.765	ng/ul	-0.02
24) 2-Nitrophenol-d4	10.172	143	34922	25.797	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.718	165	74100	25.242	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.229	131	62439	16.155	ng/ul	-0.01
46) Dimethylphthalate-d6	14.296	166	268392	27.376	ng/ul	-0.02
49) Acenaphthylene-d8	14.601	160	328377	26.969	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.071	143	5705	3.578	ng/ul	0.00
60) Fluorene-d10	15.894	176	240303	27.358	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.994	200	11743	8.228	ng/ul	-0.02
73) Anthracene-d10	17.750	188	420202	32.159	ng/ul	-0.01
81) Pyrene-d10	20.029	212	483802	30.940	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.234	264	442626	32.917	ng/ul	-0.02

Target Compounds Qvalue

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

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