

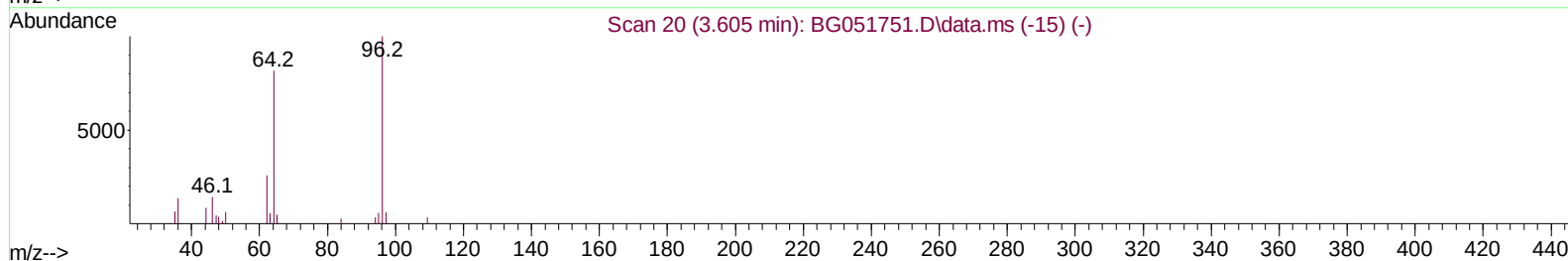
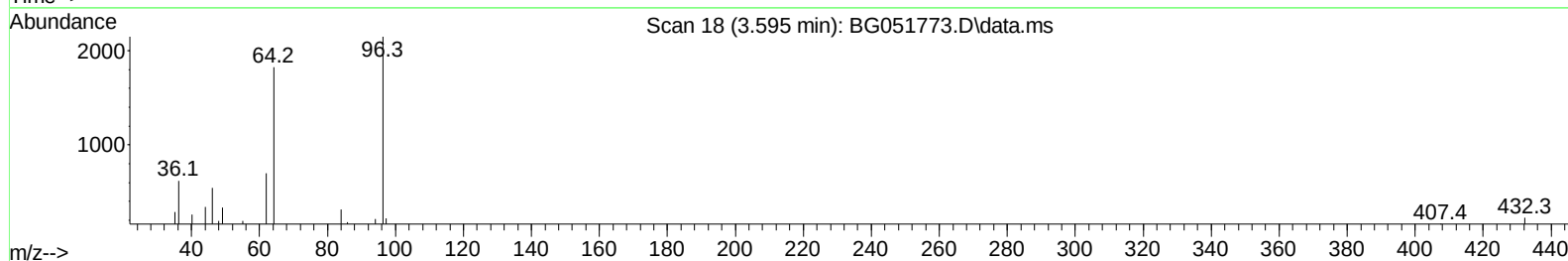
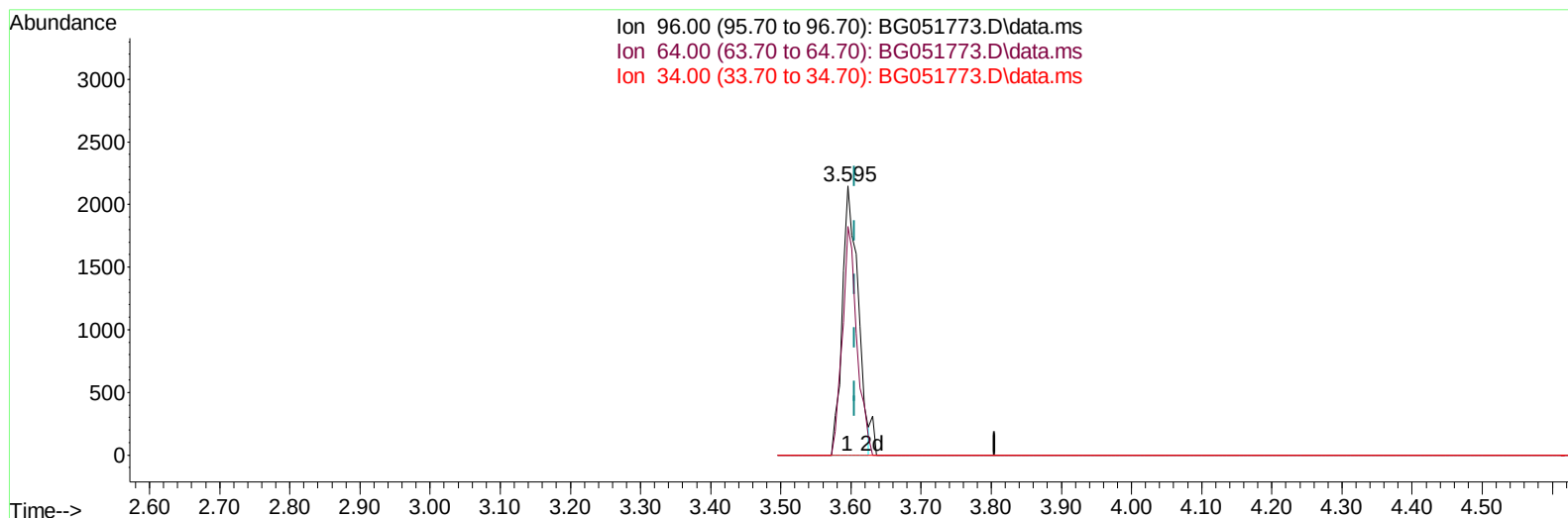
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
 Data File : BG051773.D
 Acq On : 23 Dec 2021 6:43
 Operator : CG/JU
 Sample : M5195-10
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLC2

Manual IntegrationsAPPROVED

Quant Time: Dec 23 07:32:58 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 12/23/2021
 Supervised By :mohammad ahmed 12/23/2021



TIC: BG051773.D\data.ms

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

3.595min (-0.010) 4.08 ng/uL

response 3346

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	85.10
34.00	0.00	0.00
0.00	0.00	0.00

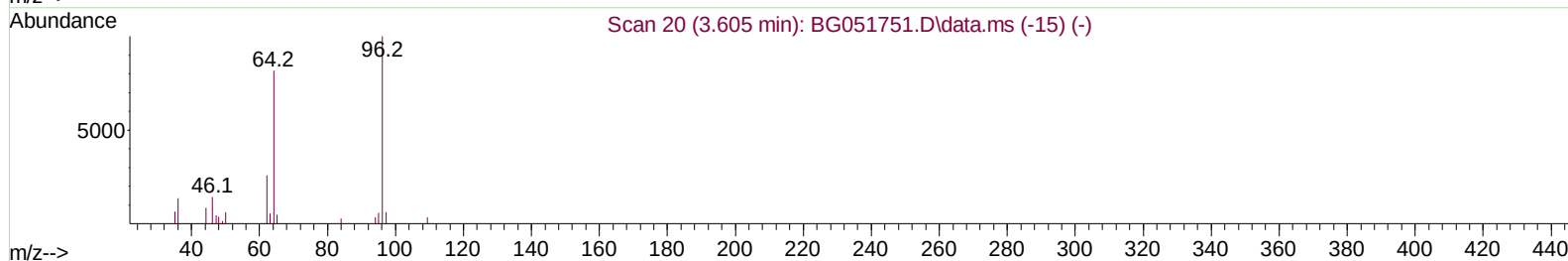
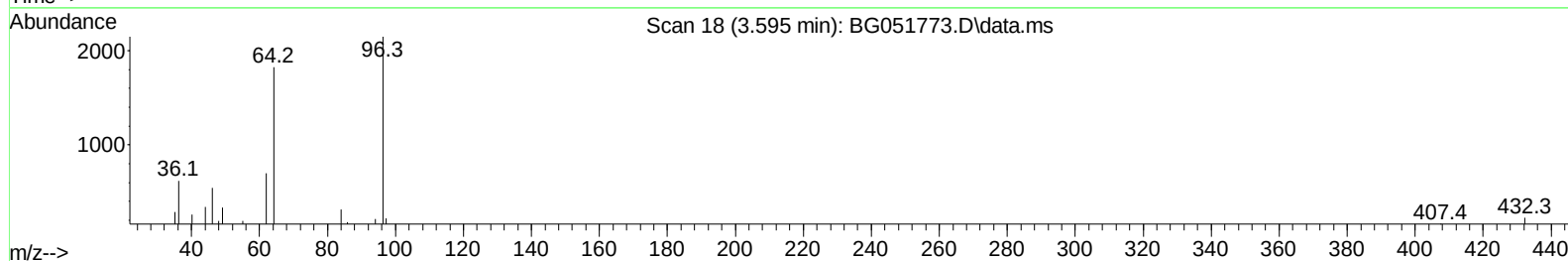
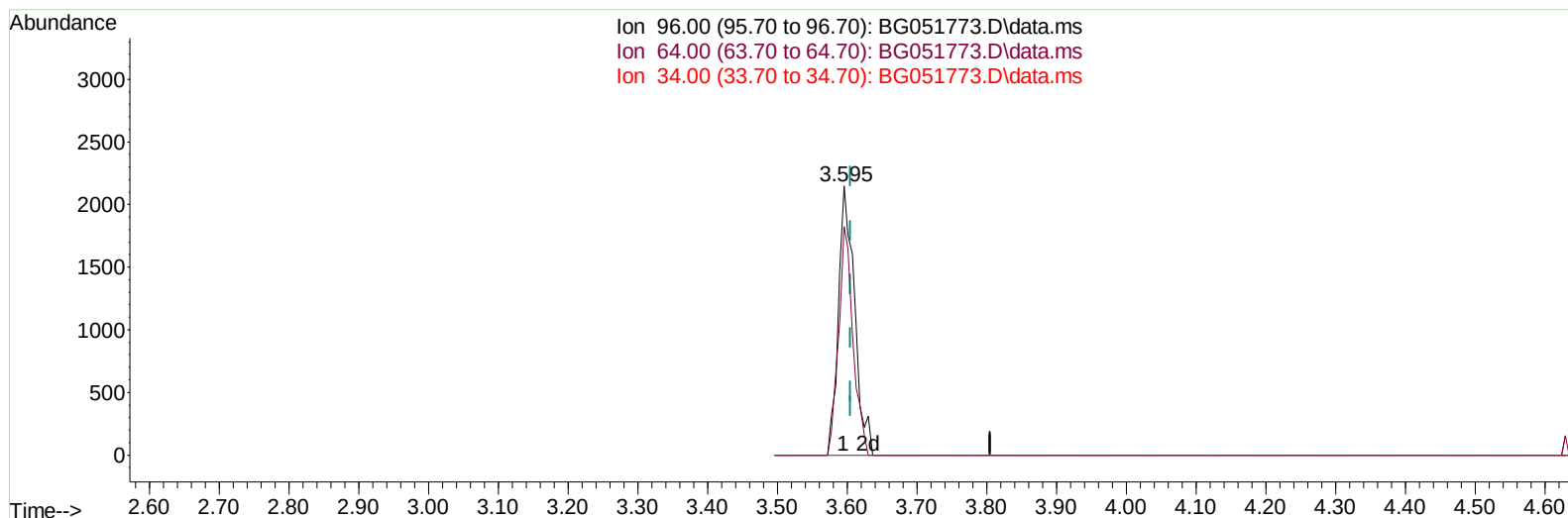
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TIC: BG051773.D\data.ms

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

3.595min (-0.010) 4.22 ng/uL m

response 3457

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	85.10
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	8.271	152	34038	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.109	136	147598	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.910	164	106179	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.653	188	238017	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.976	240	212825	20.000	ng/ul	# 0.00
88) Perylene-d12	25.472	264	215115	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.595	96	3457m	4.218	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.402	99	77084	25.364	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	7.572	67	48794	27.007	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.796	132	53546	25.333	ng/ul	-0.02
15) 4-Methylphenol-d8	8.965	113	60384	25.704	ng/ul	-0.02
21) Nitrobenzene-d5	9.435	128	30459	28.147	ng/ul	-0.02
24) 2-Nitrophenol-d4	10.175	143	26455	22.364	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.721	165	60275	23.496	ng/ul	0.00
31) 4-Chloroaniline-d4	11.226	131	46862	13.875	ng/ul	-0.02
46) Dimethylphthalate-d6	14.299	166	183870	21.636	ng/ul	-0.02
49) Acenaphthylene-d8	14.604	160	264641	25.073	ng/ul	0.00
54) 4-Nitrophenol-d4	15.074	143	12342	8.930	ng/ul	0.00
60) Fluorene-d10	15.896	176	185247	24.329	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.990	200	3852	3.106	ng/ul	-0.02
73) Anthracene-d10	17.753	188	297946	26.241	ng/ul	0.00
81) Pyrene-d10	20.026	212	330768	25.823	ng/ul	-0.02
92) Benzo(a)pyrene-d12	25.231	264	298983	26.415	ng/ul	-0.02

Target Compounds Qvalue

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

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