

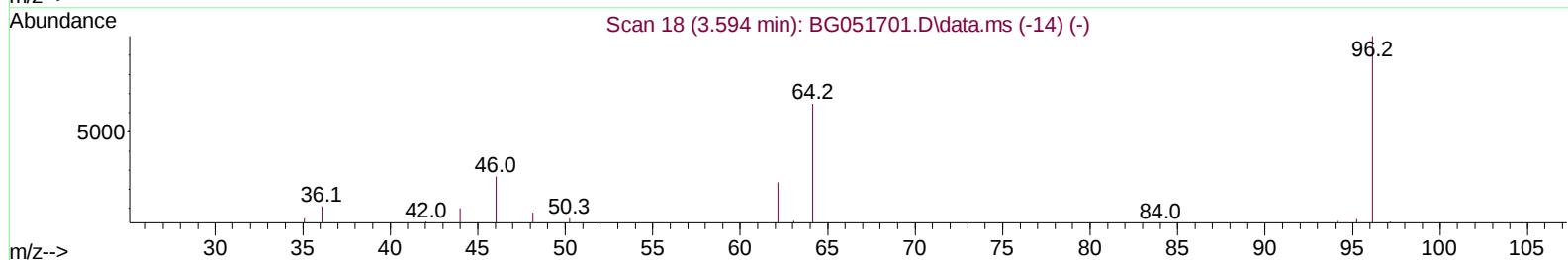
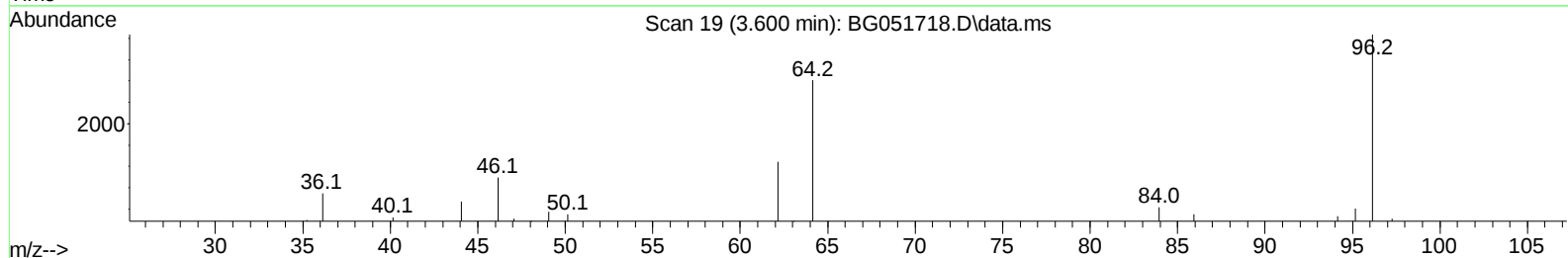
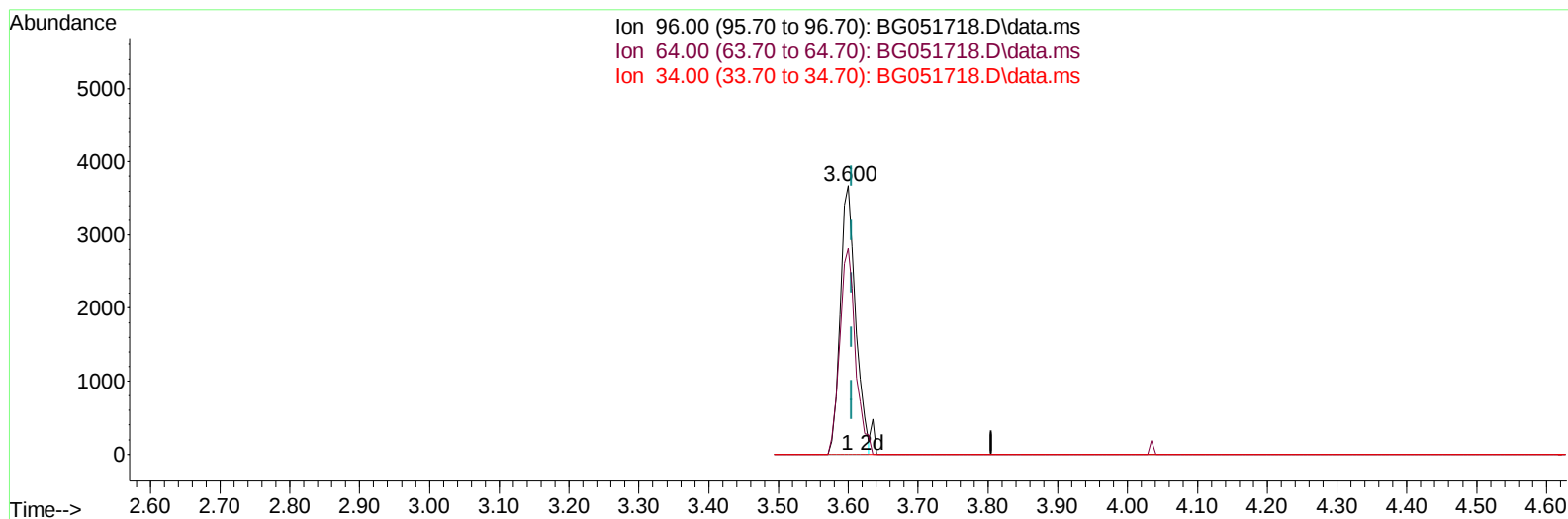
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051718.D
 Acq On : 21 Dec 2021 16:14
 Operator : CG/JU
 Sample : PB141560BL
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK560

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/22/2021
 Supervised By :mohammad ahmed 12/28/2021

Quant Time: Dec 21 23:11:08 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: BG051718.D\data.ms

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

3.600min (-0.005) 5.66 ng/uL

response 5693

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	76.71
34.00	0.00	0.00
0.00	0.00	0.00

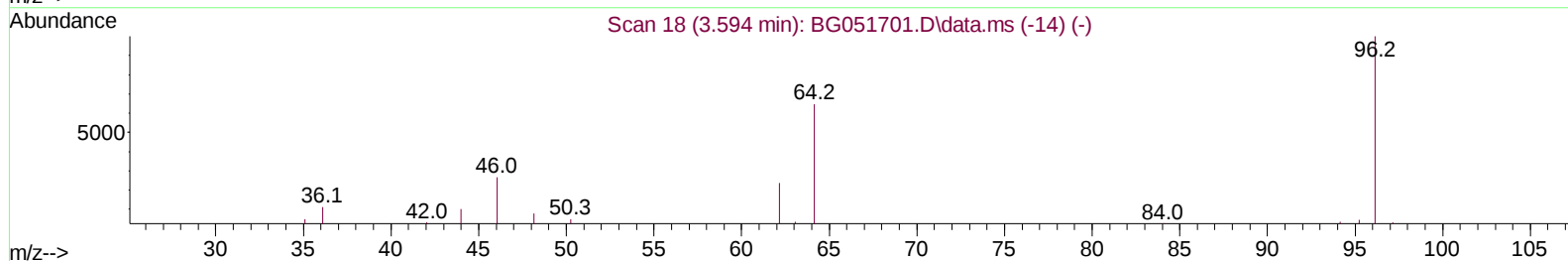
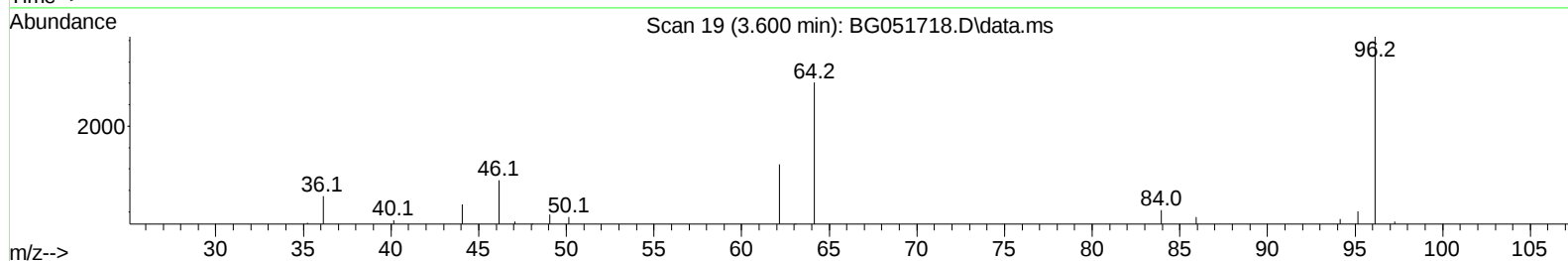
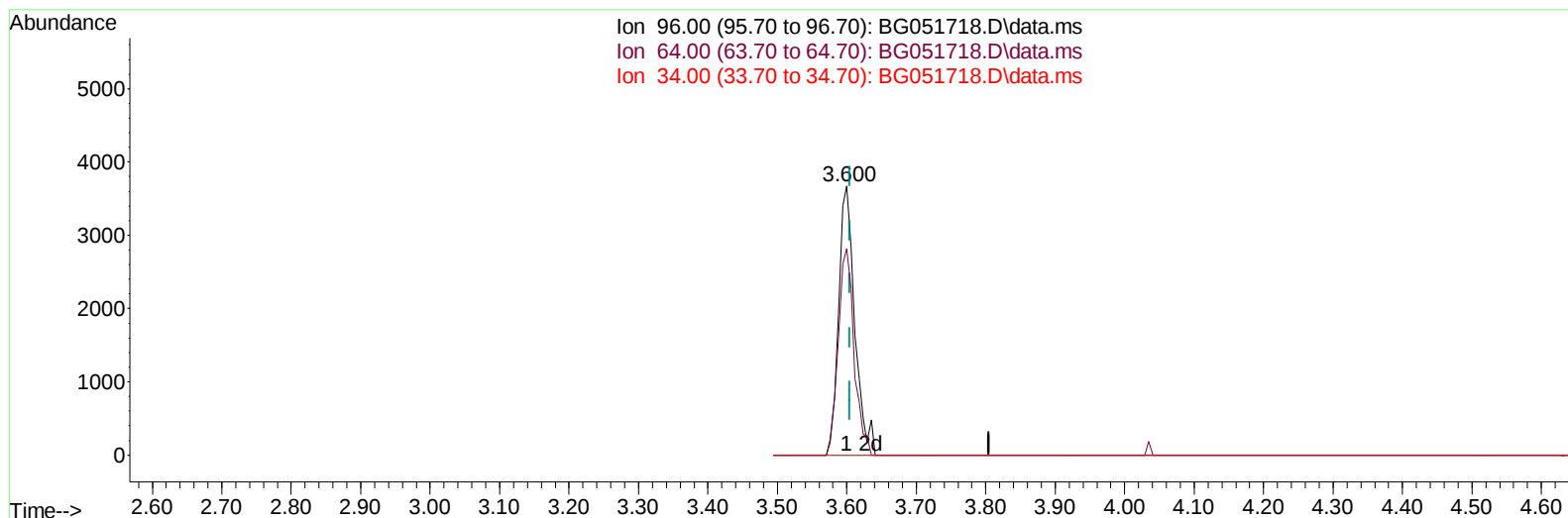
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TIC: BG051718.D\data.ms

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

3.600min (-0.005) 5.83 ng/uL m

response 5863

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	76.71
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	8.270	152	41767	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.113	136	181406	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.908	164	130782	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.651	188	304312	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.975	240	297430	20.000	ng/ul	#-0.01
88) Perylene-d12	25.470	264	293423	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.600	96	5863m	5.830	ng/uL	0.00
4) Pyridine-d5	4.023	84	76502	25.176	ng/ul	-0.02
7) Phenol-d5	7.401	99	120158	32.220	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	7.577	67	69843	31.504	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.800	132	82320	31.739	ng/ul	-0.01
15) 4-Methylphenol-d8	8.969	113	94956	32.941	ng/ul	-0.01
21) Nitrobenzene-d5	9.439	128	43765	32.906	ng/ul	-0.02
24) 2-Nitrophenol-d4	10.173	143	50152	34.495	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.714	165	94820	30.074	ng/ul	-0.02
31) 4-Chloroaniline-d4	11.231	131	108226	26.071	ng/ul	-0.01
46) Dimethylphthalate-d6	14.297	166	317031	30.286	ng/ul	-0.02
49) Acenaphthylene-d8	14.603	160	389205	29.938	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.073	143	44658	26.235	ng/ul	0.00
60) Fluorene-d10	15.895	176	275807	29.409	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.995	200	38416	24.230	ng/ul	-0.02
73) Anthracene-d10	17.751	188	448580	30.901	ng/ul	-0.01
81) Pyrene-d10	20.031	212	548733	30.654	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.229	264	495798	32.114	ng/ul	-0.02

Target Compounds Qvalue

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

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