

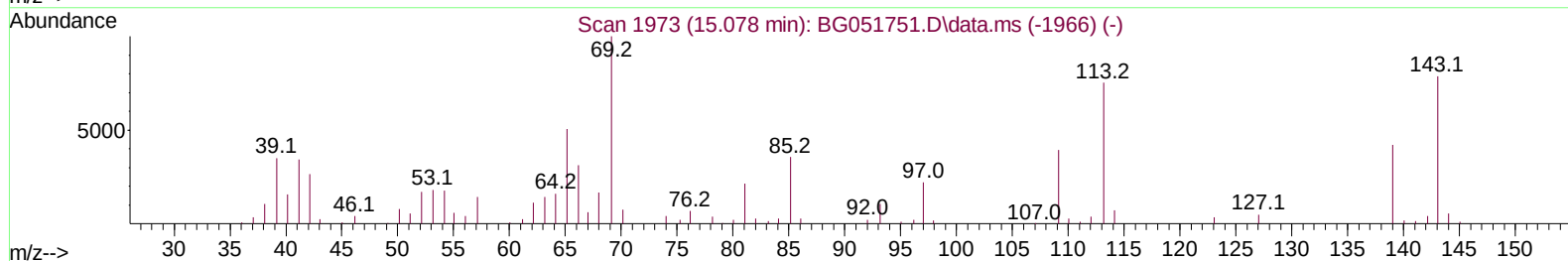
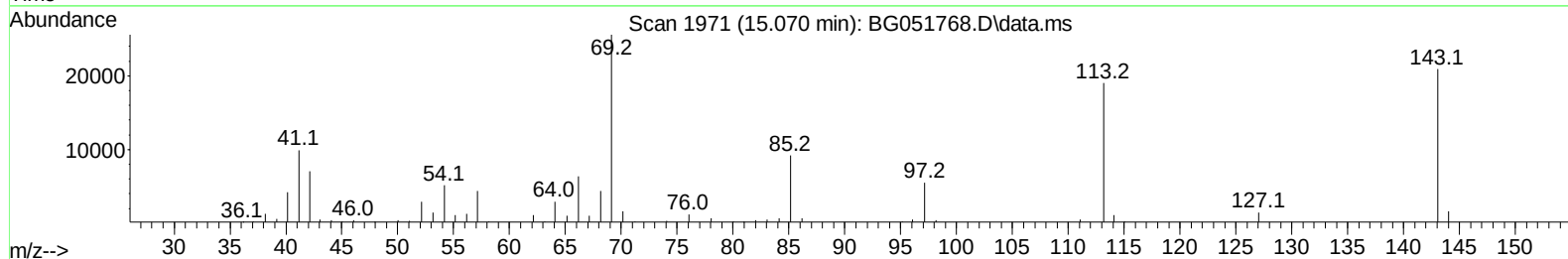
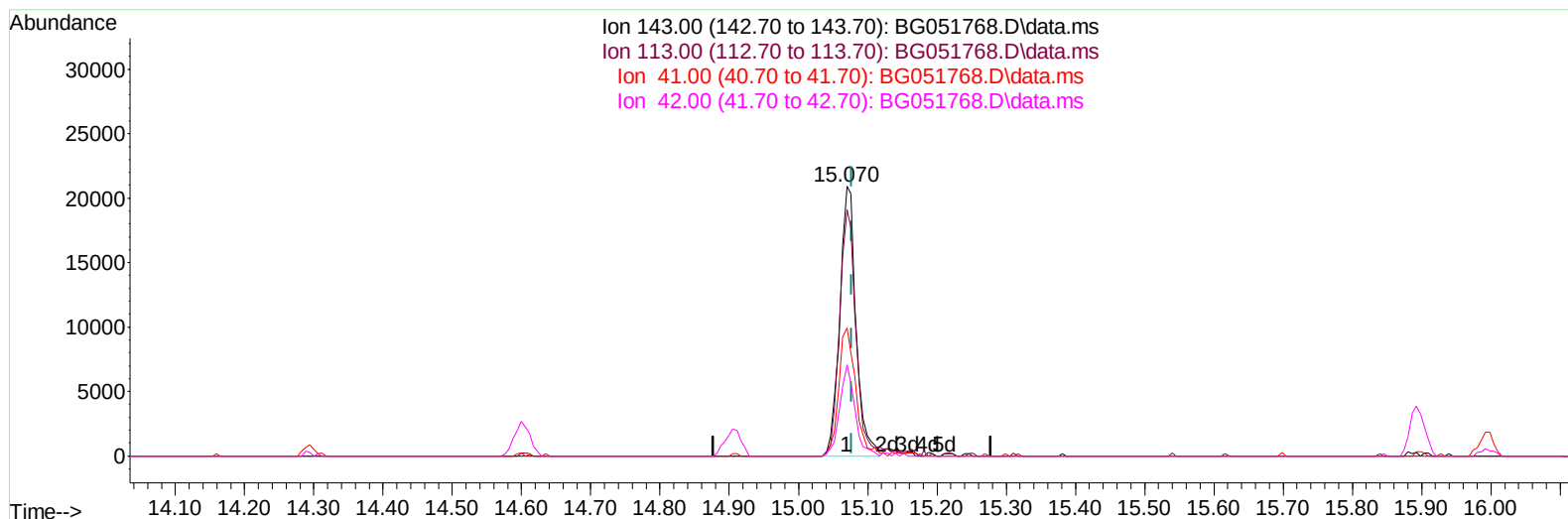
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
 Data File : BG051768.D
 Acq On : 23 Dec 2021 3:23
 Operator : CG/JU
 Sample : PB141617BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK617

Manual IntegrationsAPPROVED

Quant Time: Dec 23 05:52:27 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: BG051768.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.070min (-0.008) 24.02 ng/u1

response 34236

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	91.27
41.00	44.40	47.47
42.00	29.70	33.76

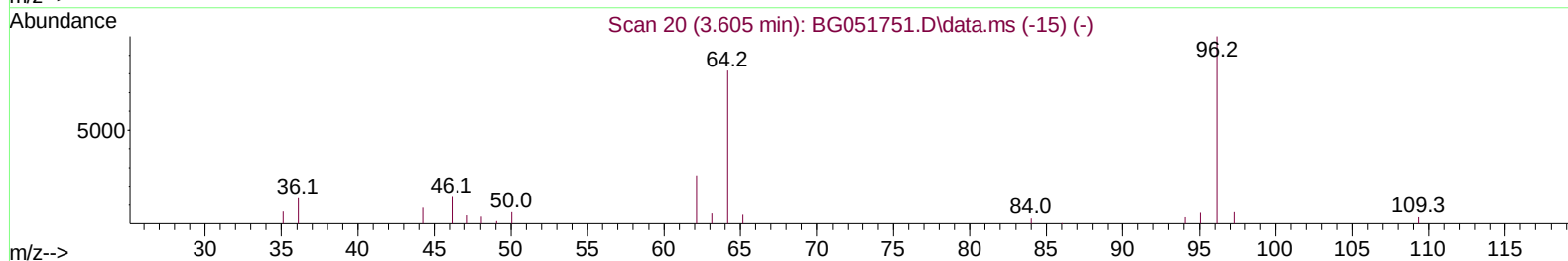
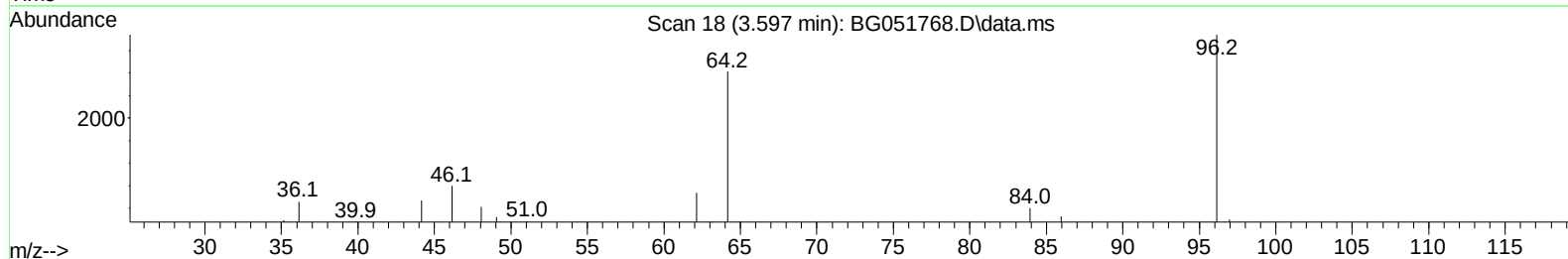
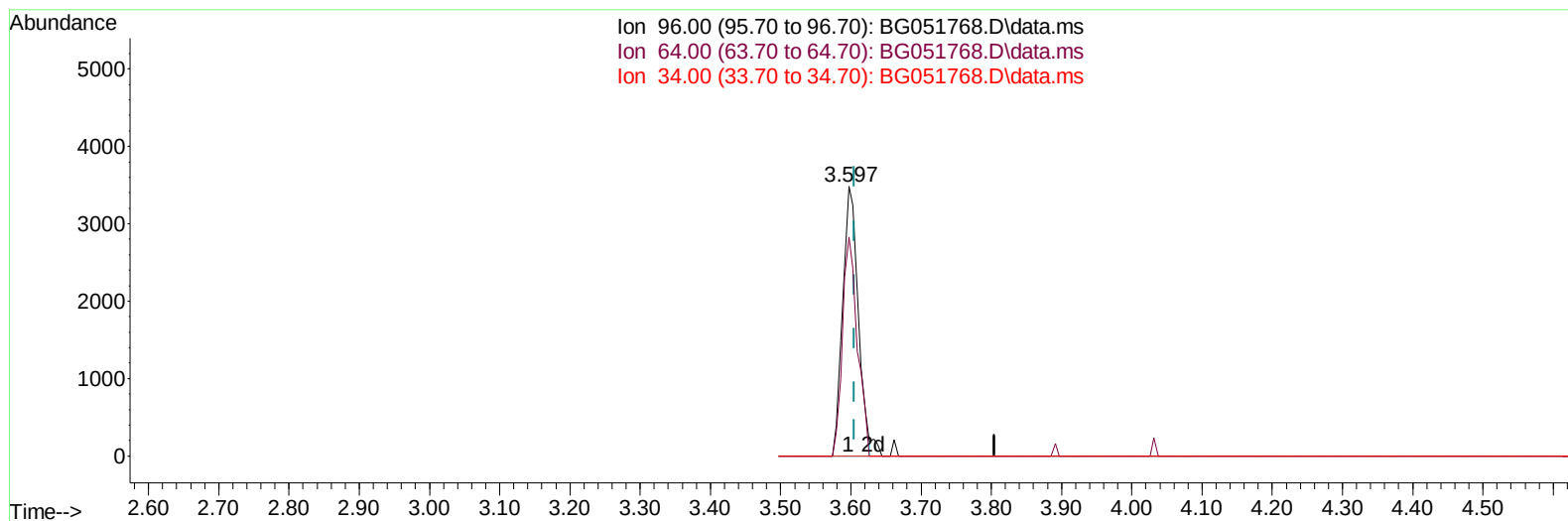
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(3) 1,4-Dioxane-d8 (S)

3.597min (-0.008) 6.05 ng/uL

response 5314

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	81.18
34.00	0.00	0.00
0.00	0.00	0.00

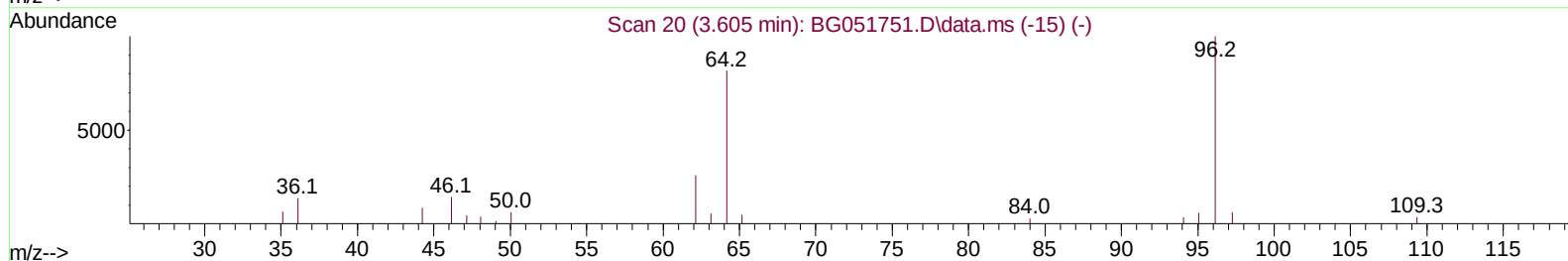
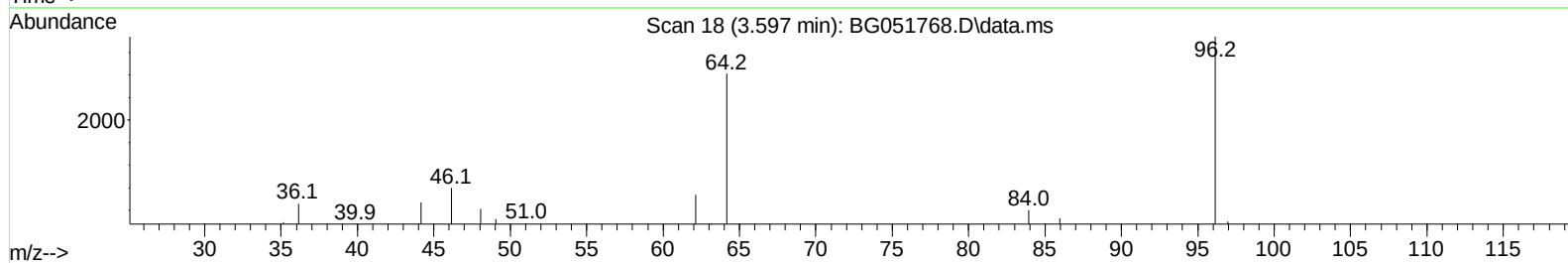
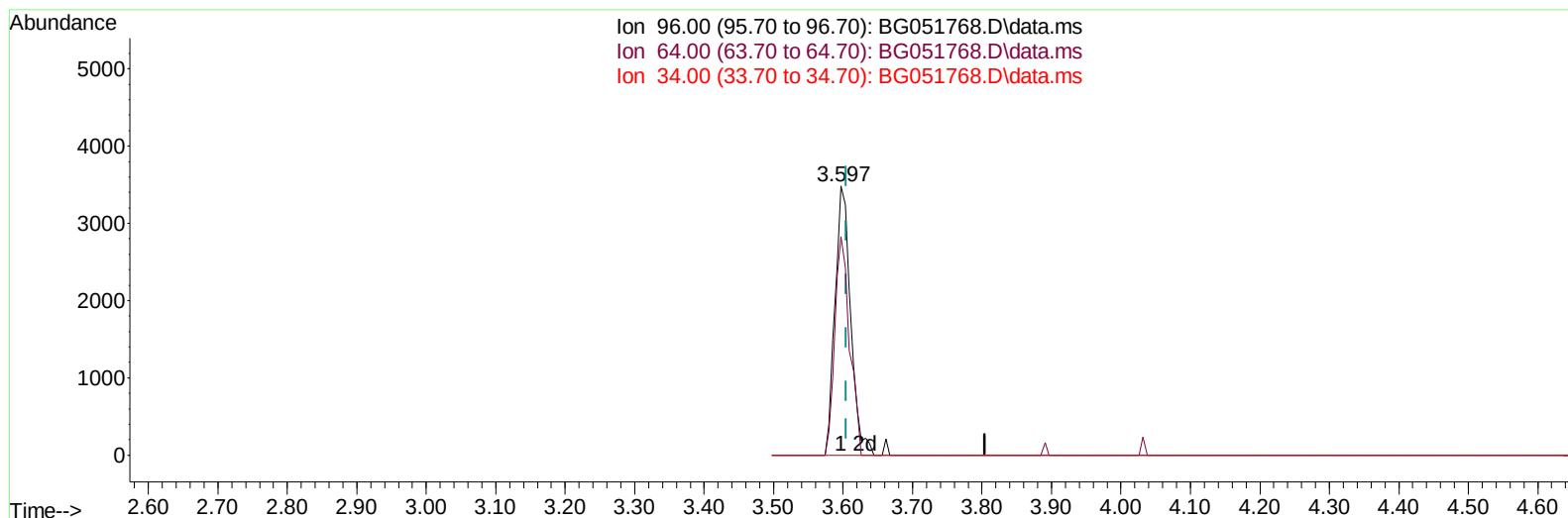
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 Data File : BG051768.D
 Acq On : 23 Dec 2021 3:23
 Operator : CG/JU
 Sample : PB141617BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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TIC: BG051768.D\data.ms

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

3.597min (-0.008) 6.20 ng/uL m

response 5444

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	81.18
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	8.273	152	36482	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.110	136	154436	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.905	164	109483	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.655	188	257229	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.972	240	256120	20.000	ng/ul	#-0.01
88) Perylene-d12	25.467	264	240225	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.597	96	5444m	6.197	ng/uL	0.00
4) Pyridine-d5	4.026	84	67477	25.423	ng/ul	-0.02
7) Phenol-d5	7.404	99	99750	30.623	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.574	67	58243	30.078	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.797	132	67315	29.713	ng/ul	-0.01
15) 4-Methylphenol-d8	8.966	113	77124	30.631	ng/ul	-0.01
21) Nitrobenzene-d5	9.442	128	34615	30.572	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.171	143	40009	32.324	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.717	165	76651	28.557	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.228	131	78329	22.164	ng/ul	-0.01
46) Dimethylphthalate-d6	14.294	166	253186	28.893	ng/ul	-0.02
49) Acenaphthylene-d8	14.600	160	320832	29.480	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.070	143	34236	24.025	ng/ul	0.00
60) Fluorene-d10	15.892	176	219950	28.015	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.998	200	31576	23.561	ng/ul	-0.01
73) Anthracene-d10	17.754	188	363794	29.648	ng/ul	0.00
81) Pyrene-d10	20.028	212	440206	28.558	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.233	264	392325	31.039	ng/ul	-0.02

Target Compounds Qvalue

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

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