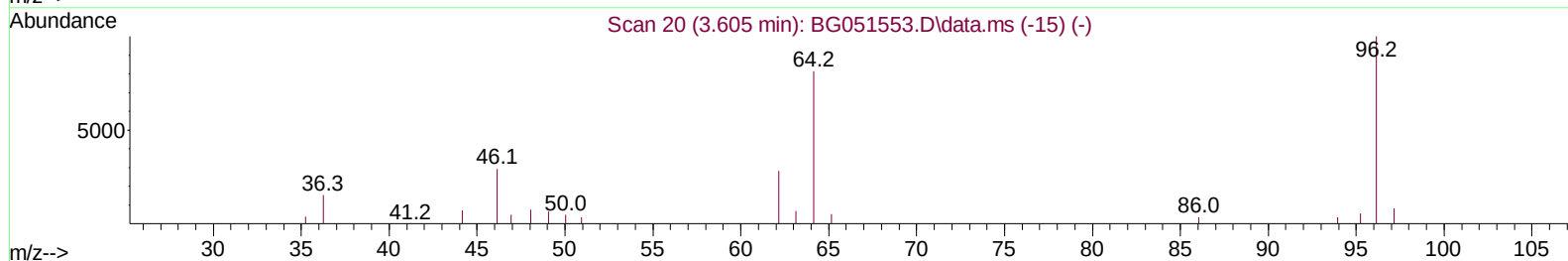
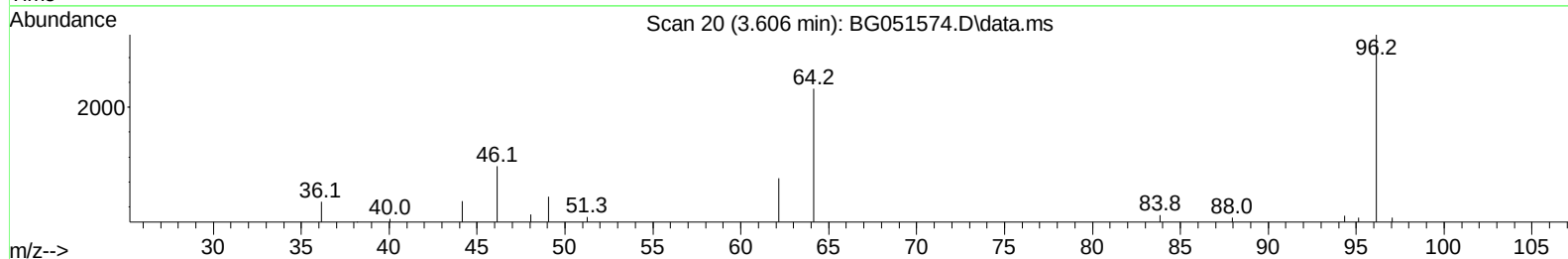
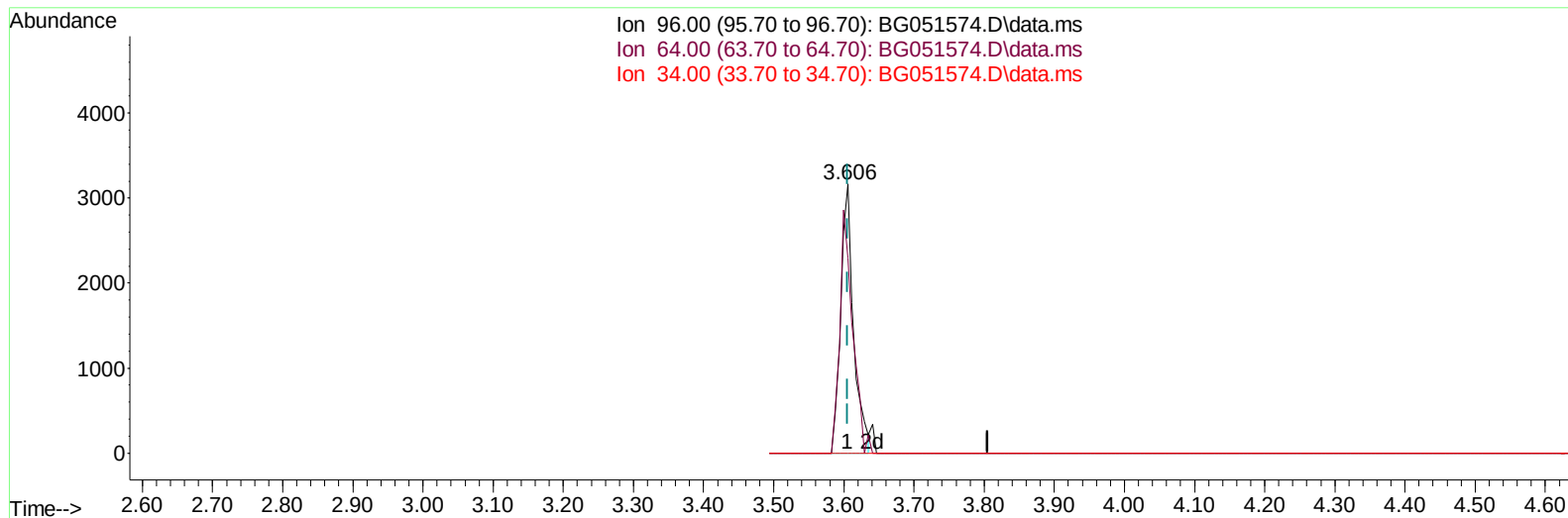


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051574.D
Acq On : 16 Dec 2021 17:29
Operator : CG/JU
Sample : M5037-07
Misc :
ALS Vial : 74 Sample Multiplier: 1

Quant Time: Dec 17 00:12:03 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



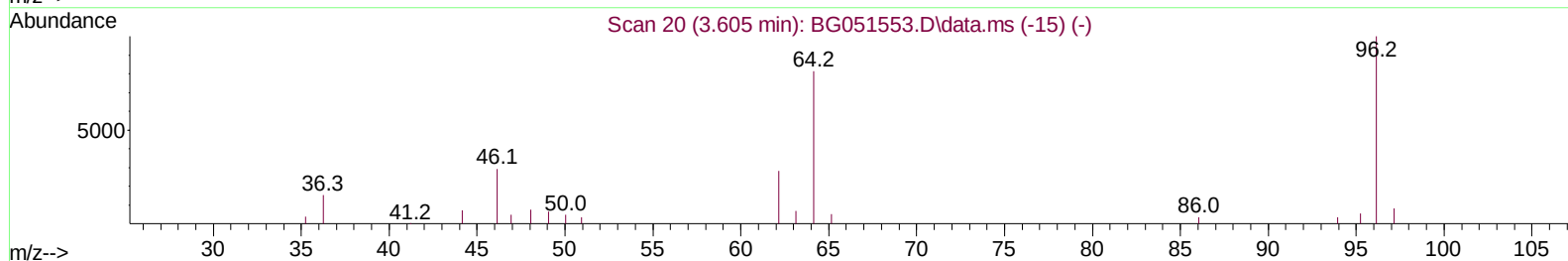
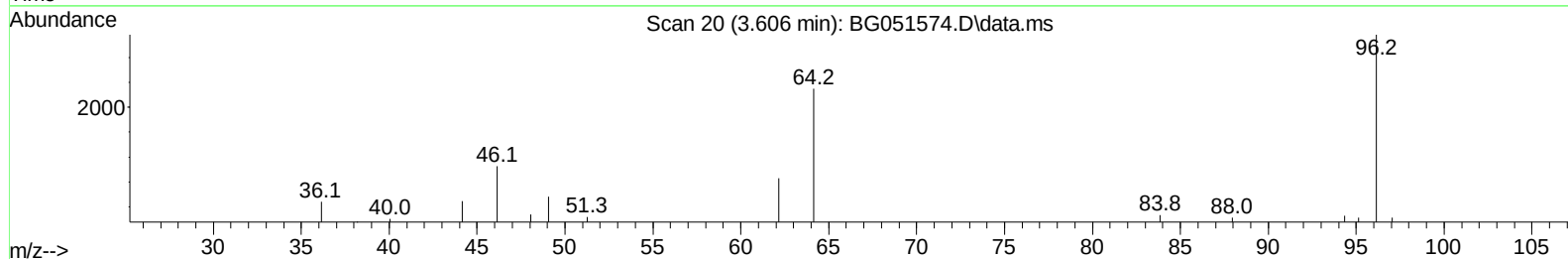
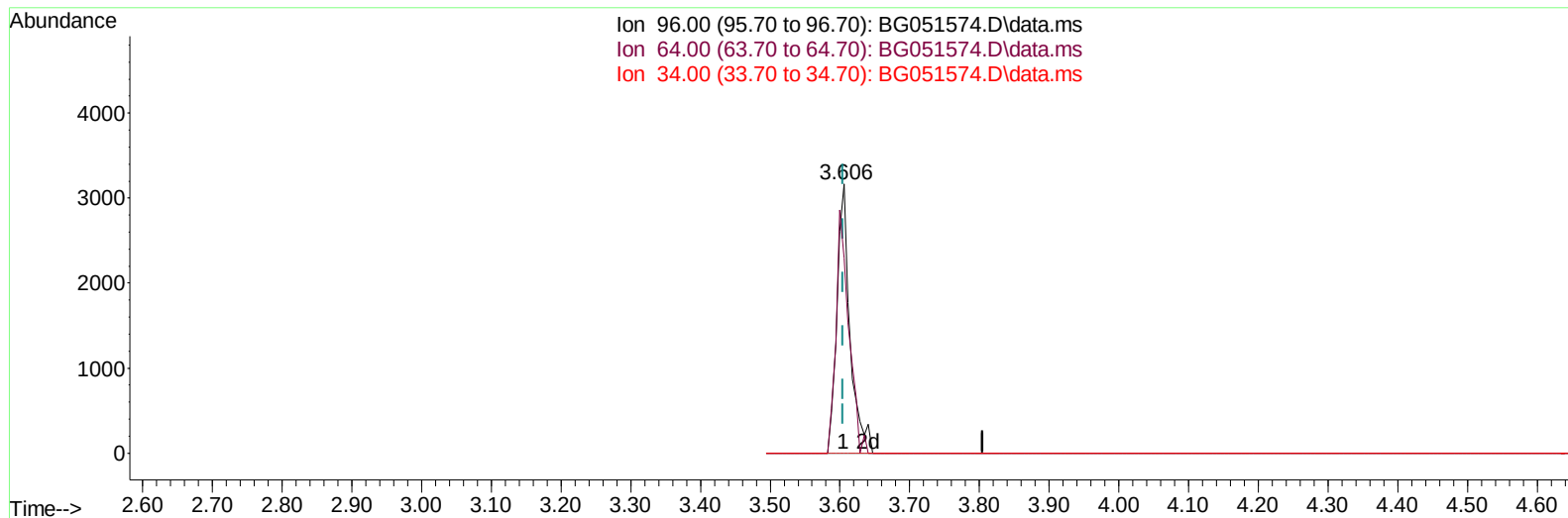
TIC: BG051574.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.606min (+ 0.001) 4.55 ng/uL****response 4006**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	72.53
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
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Acq On : 16 Dec 2021 17:29
Operator : CG/JU
Sample : M5037-07
Misc :
ALS Vial : 74 Sample Multiplier: 1

Quant Time: Dec 17 00:12:03 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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Response via : Initial Calibration



TIC: BG051574.D\data.ms

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

3.606min (+ 0.001) 4.69 ng/uL m

response 4126

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	72.53
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051574.D
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 Operator : CG/JU
 Sample : M5037-07
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Quant Time: Dec 17 00:12:03 2021
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.282	152	36550	20.000	ng/ul	0.00
20) Naphthalene-d8	11.113	136	146709	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.908	164	106942	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.657	188	257338	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.975	240	275503	20.000	ng/ul	#-0.01
88) Perylene-d12	25.470	264	283889	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.606	96	4126m	4.688	ng/uL	0.00
4) Pyridine-d5	4.034	84	26927	10.126	ng/ul	-0.01
7) Phenol-d5	7.406	99	24087	7.381	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.583	67	59955	30.904	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.800	132	56929	25.082	ng/ul	-0.01
15) 4-Methylphenol-d8	8.969	113	42433	16.821	ng/ul	-0.01
21) Nitrobenzene-d5	9.445	128	36277	33.727	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.179	143	39176	33.318	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.719	165	74910	29.378	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.231	131	88760	26.439	ng/ul	-0.01
46) Dimethylphthalate-d6	14.303	166	281975	32.943	ng/ul	-0.01
49) Acenaphthylene-d8	14.608	160	327095	30.769	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	10426	7.490	ng/ul	-0.01
60) Fluorene-d10	15.901	176	244188	31.842	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.995	200	47779	35.636	ng/ul	-0.02
73) Anthracene-d10	17.757	188	424920	34.615	ng/ul	0.00
81) Pyrene-d10	20.030	212	541092	32.633	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.235	264	532697	35.663	ng/ul	-0.02

Target Compounds	Qvalue

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

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