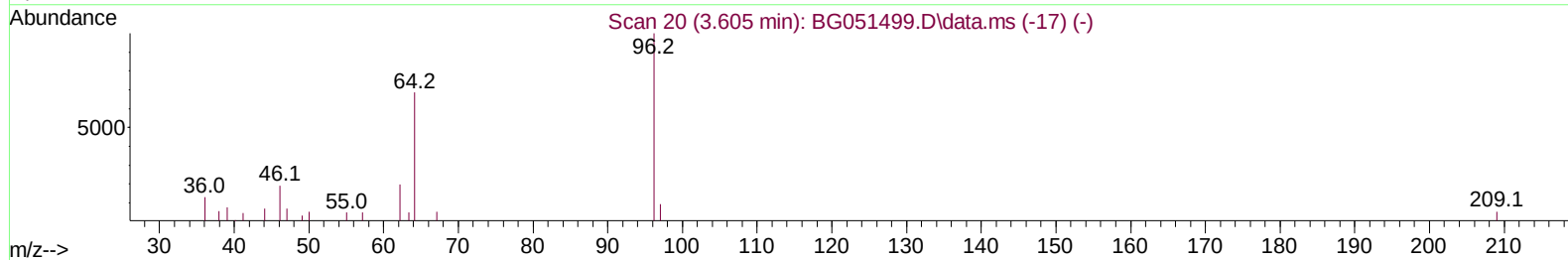
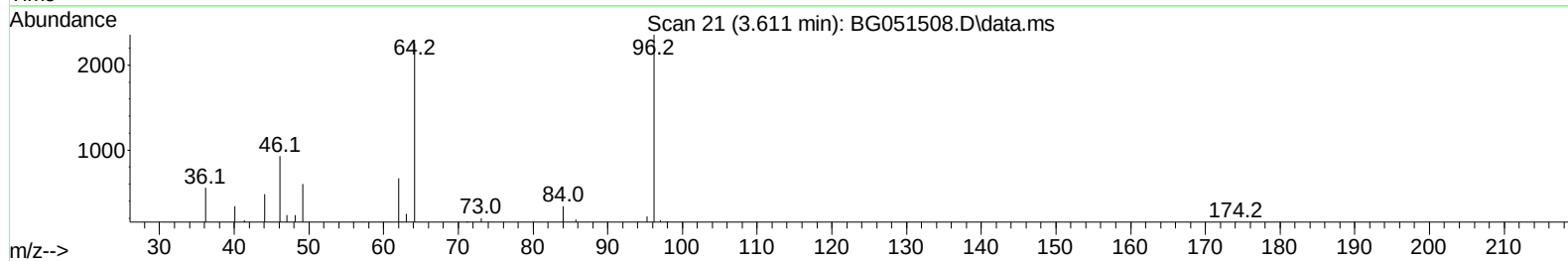
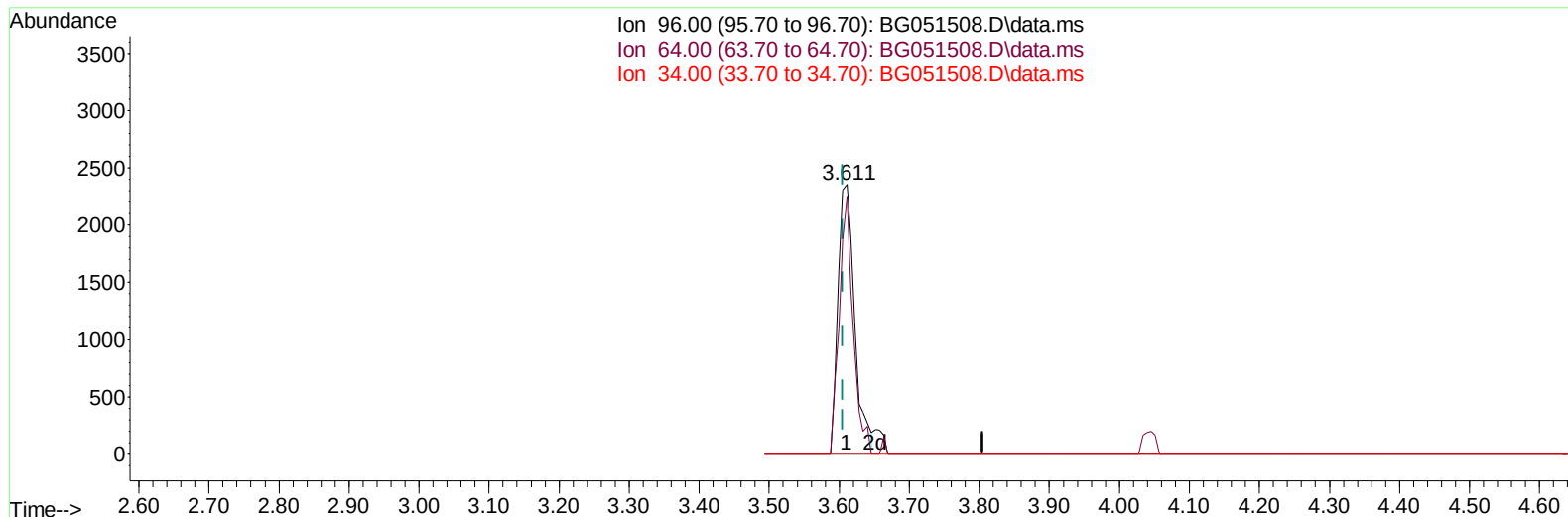


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051508.D
Acq On : 14 Dec 2021 18:01
Operator : CG/JU
Sample : PB141304BL
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Dec 14 23:48:37 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



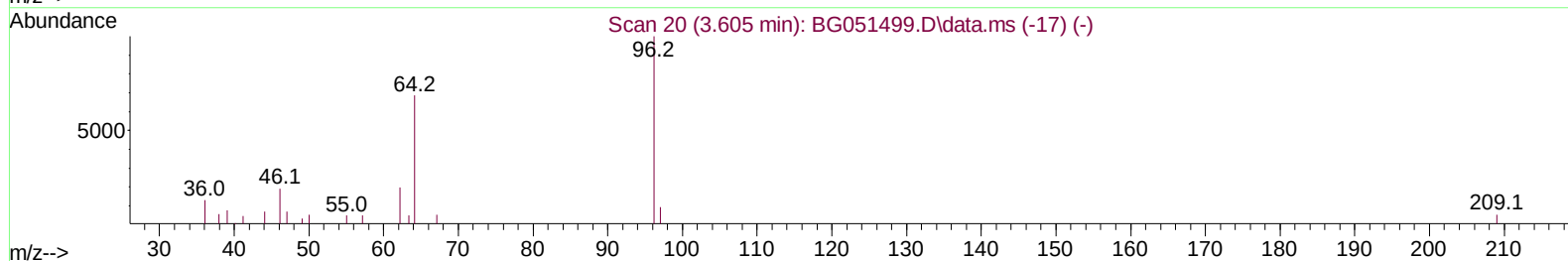
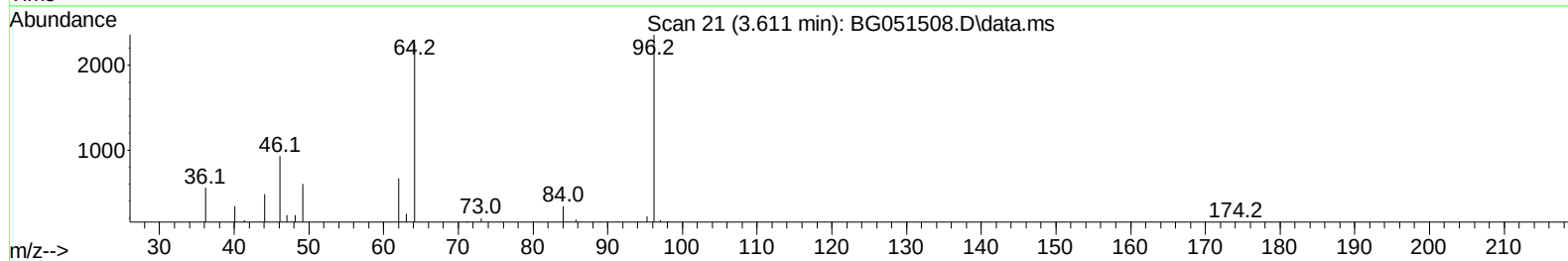
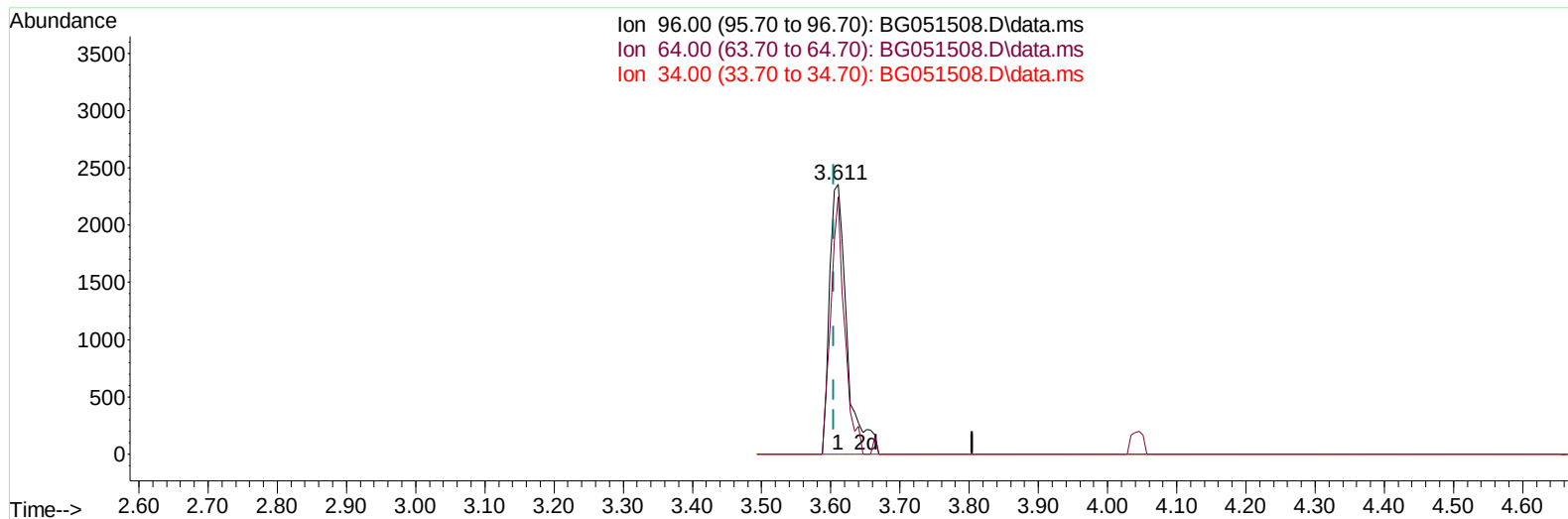
TIC: BG051508.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.611min (+ 0.006) 5.09 ng/uL****response 3916**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	95.42#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051508.D
Acq On : 14 Dec 2021 18:01
Operator : CG/JU
Sample : PB141304BL
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Dec 14 23:48:37 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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QLast Update : Tue Dec 14 14:19:57 2021
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TIC: BG051508.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.611min (+ 0.006) 5.36 ng/uL m****response 4123**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	95.42#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051508.D
 Acq On : 14 Dec 2021 18:01
 Operator : CG/JU
 Sample : PB141304BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Dec 14 23:48:37 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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.281	152	31939	20.000	ng/ul	0.00
20) Naphthalene-d8	11.118	136	127765	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.913	164	95343	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.662	188	242463	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.980	240	247636	20.000	ng/ul	# 0.00
88) Perylene-d12	25.481	264	232822	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.611	96	4123m	5.361	ng/uL	0.00
4) Pyridine-d5	4.040	84	61647	26.530	ng/ul	0.00
7) Phenol-d5	7.412	99	74603	26.161	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.588	67	48296	28.488	ng/ul	0.00
11) 2-Chlorophenol-d4	7.811	132	54789	27.624	ng/ul	0.00
15) 4-Methylphenol-d8	8.974	113	57224	25.960	ng/ul	0.00
21) Nitrobenzene-d5	9.450	128	28755	30.698	ng/ul	0.00
24) 2-Nitrophenol-d4	10.184	143	30142	29.436	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.725	165	58319	26.263	ng/ul	0.00
31) 4-Chloroaniline-d4	11.242	131	79348	27.140	ng/ul	0.00
46) Dimethylphthalate-d6	14.308	166	223170	29.244	ng/ul	0.00
49) Acenaphthylene-d8	14.614	160	253467	26.744	ng/ul	0.00
54) 4-Nitrophenol-d4	15.072	143	35634	28.714	ng/ul	0.00
60) Fluorene-d10	15.906	176	189710	27.747	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.000	200	35738	28.290	ng/ul	-0.01
73) Anthracene-d10	17.762	188	335258	28.986	ng/ul	0.00
81) Pyrene-d10	20.036	212	432398	29.012	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.240	264	375714	30.670	ng/ul	-0.01

Target Compounds	Qvalue

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

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