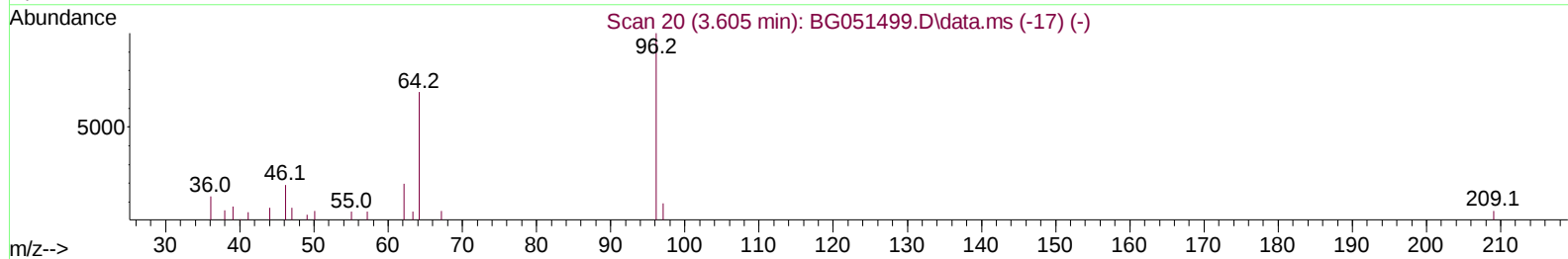
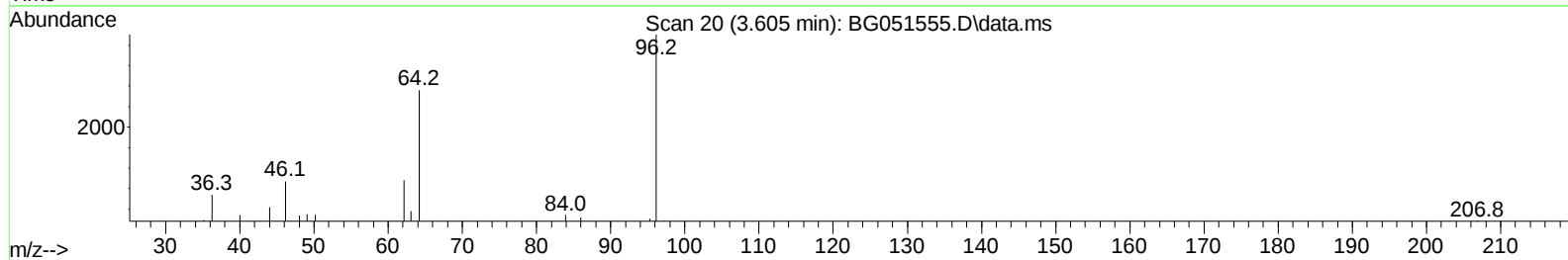
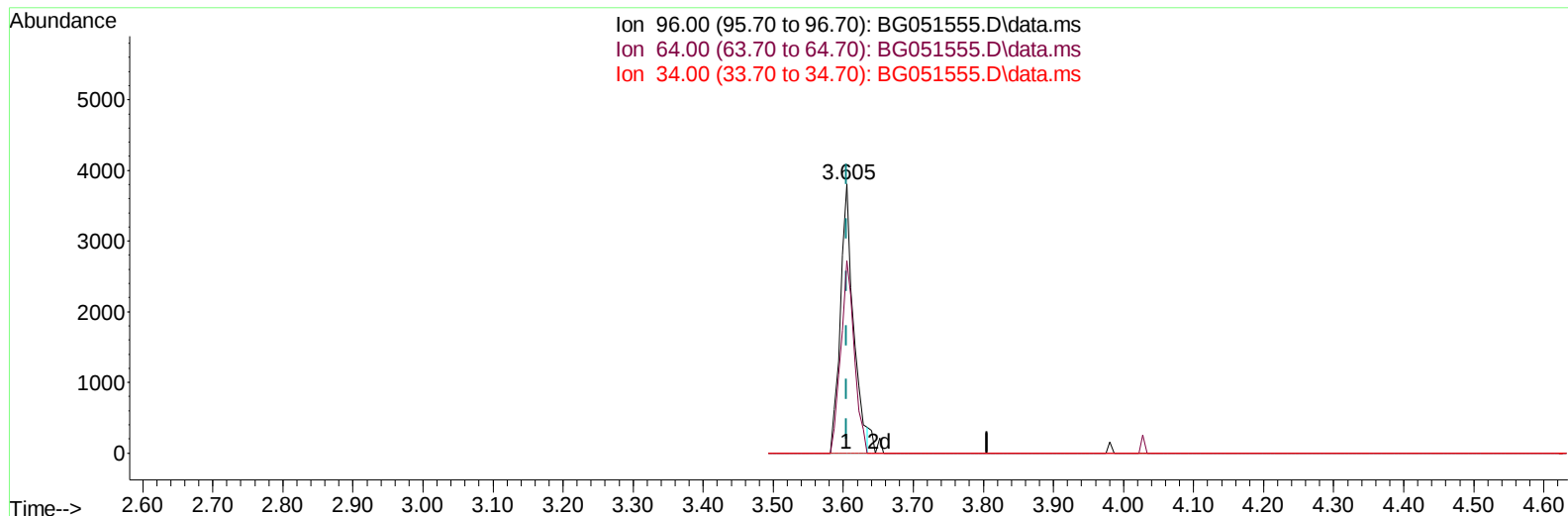


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051555.D
Acq On : 16 Dec 2021 3:52
Operator : CG/JU
Sample : PB141413BL
Misc :
ALS Vial : 54 Sample Multiplier: 1

Quant Time: Dec 16 05:35:22 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



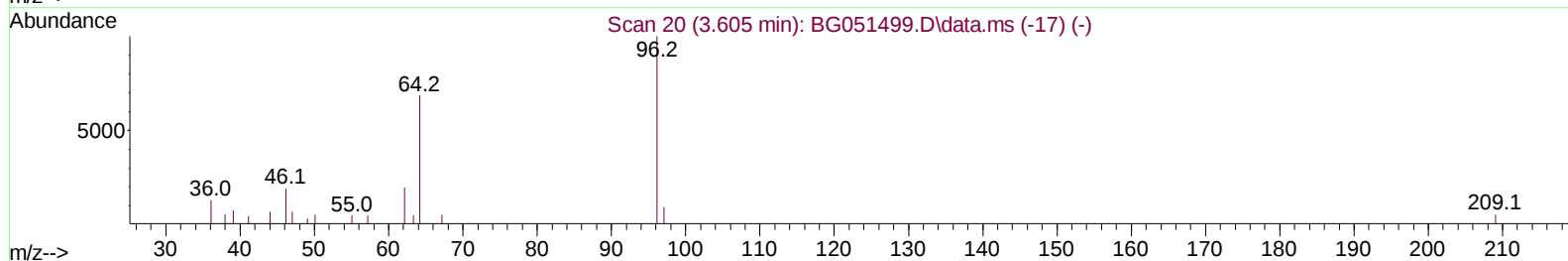
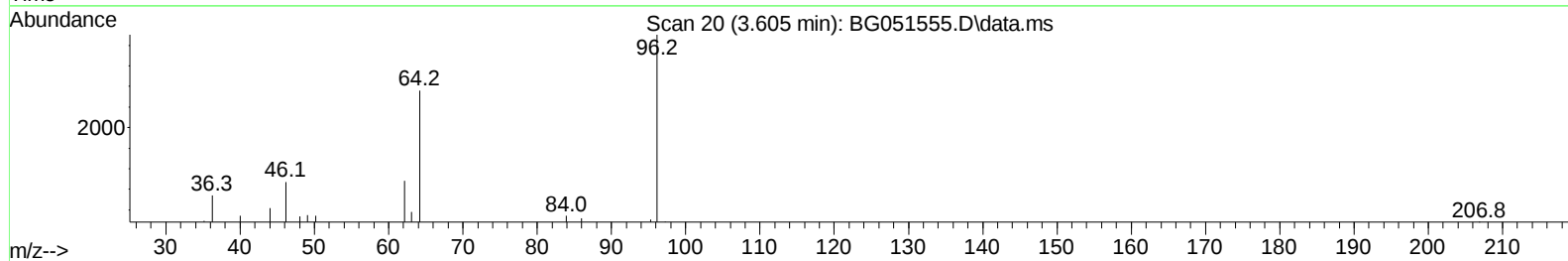
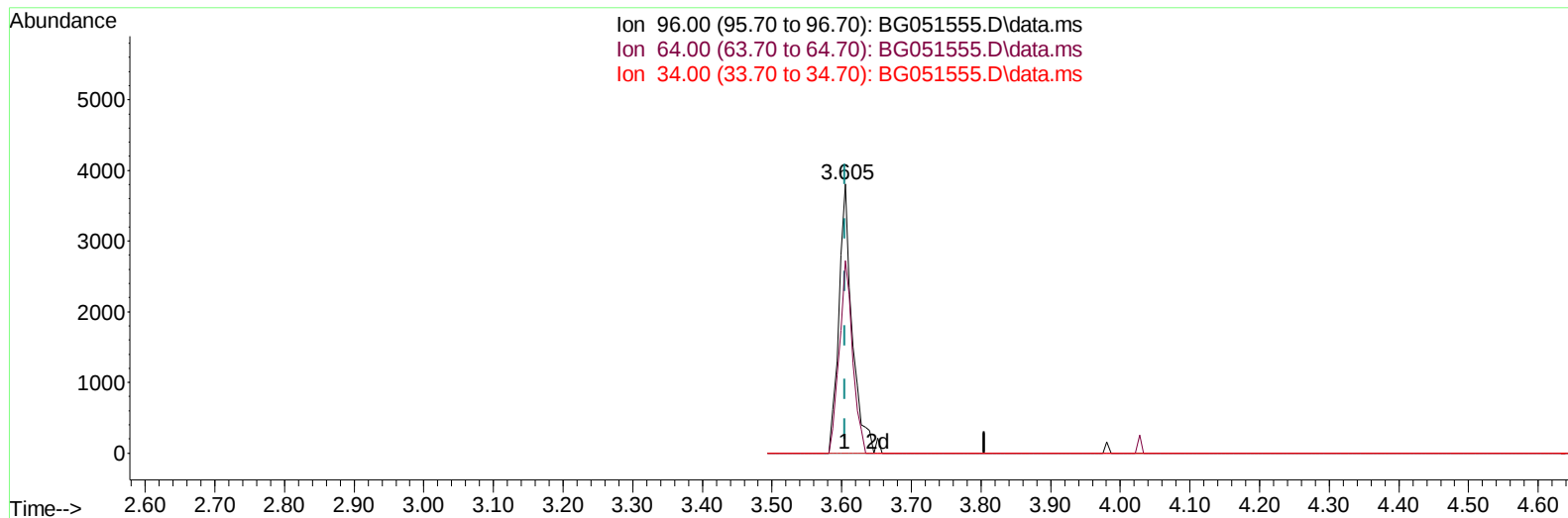
TIC: BG051555.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.605min (-0.000) 7.11 ng/uL****response 4966**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	71.49
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051555.D
Acq On : 16 Dec 2021 3:52
Operator : CG/JU
Sample : PB141413BL
Misc :
ALS Vial : 54 Sample Multiplier: 1

Quant Time: Dec 16 05:35:22 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: BG051555.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.605min (-0.000) 7.27 ng/uL m****response 5078**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	71.49
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051555.D
 Acq On : 16 Dec 2021 3:52
 Operator : CG/JU
 Sample : PB141413BL
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Quant Time: Dec 16 05:35:22 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.275	152	29023	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.118	136	118139	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.913	164	86710	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.656	188	223100	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.974	240	228603	20.000	ng/ul	#-0.01
88) Perylene-d12	25.469	264	225970	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.605	96	5078m	7.266	ng/uL	0.00
4) Pyridine-d5	4.034	84	72629	34.397	ng/ul	-0.01
7) Phenol-d5	7.406	99	90243	34.824	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.582	67	54360	35.287	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.799	132	64528	35.803	ng/ul	-0.01
15) 4-Methylphenol-d8	8.968	113	69885	34.889	ng/ul	-0.01
21) Nitrobenzene-d5	9.450	128	30967	35.753	ng/ul	0.00
24) 2-Nitrophenol-d4	10.178	143	36700	38.760	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.719	165	72352	35.237	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.236	131	91241	33.750	ng/ul	0.00
46) Dimethylphthalate-d6	14.302	166	250569	36.104	ng/ul	-0.01
49) Acenaphthylene-d8	14.608	160	293301	34.028	ng/ul	0.00
54) 4-Nitrophenol-d4	15.066	143	39065	34.613	ng/ul	-0.01
60) Fluorene-d10	15.900	176	215781	34.703	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.994	200	37174	31.981	ng/ul	-0.02
73) Anthracene-d10	17.756	188	381520	35.849	ng/ul	0.00
81) Pyrene-d10	20.030	212	476787	34.654	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.228	264	437602	36.805	ng/ul	-0.02

Target Compounds	Qvalue

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

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