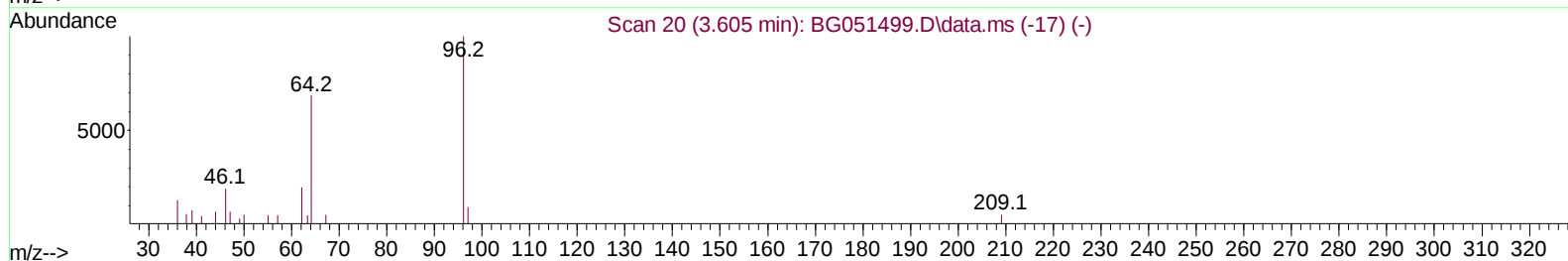
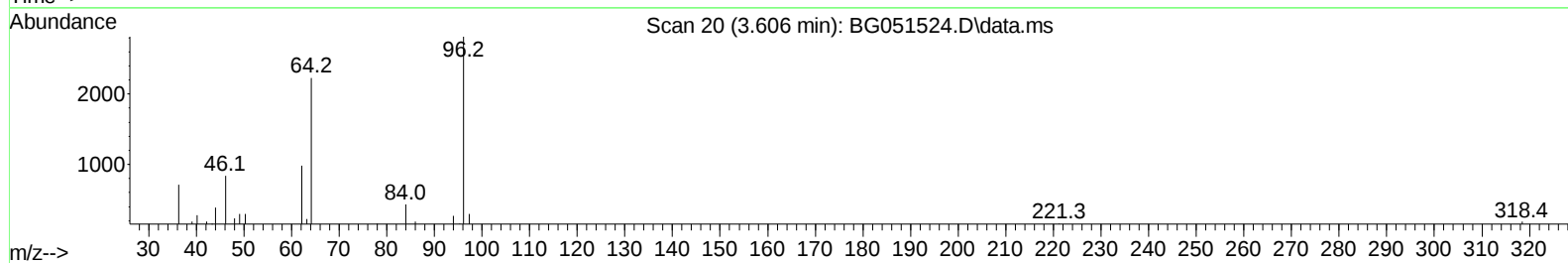
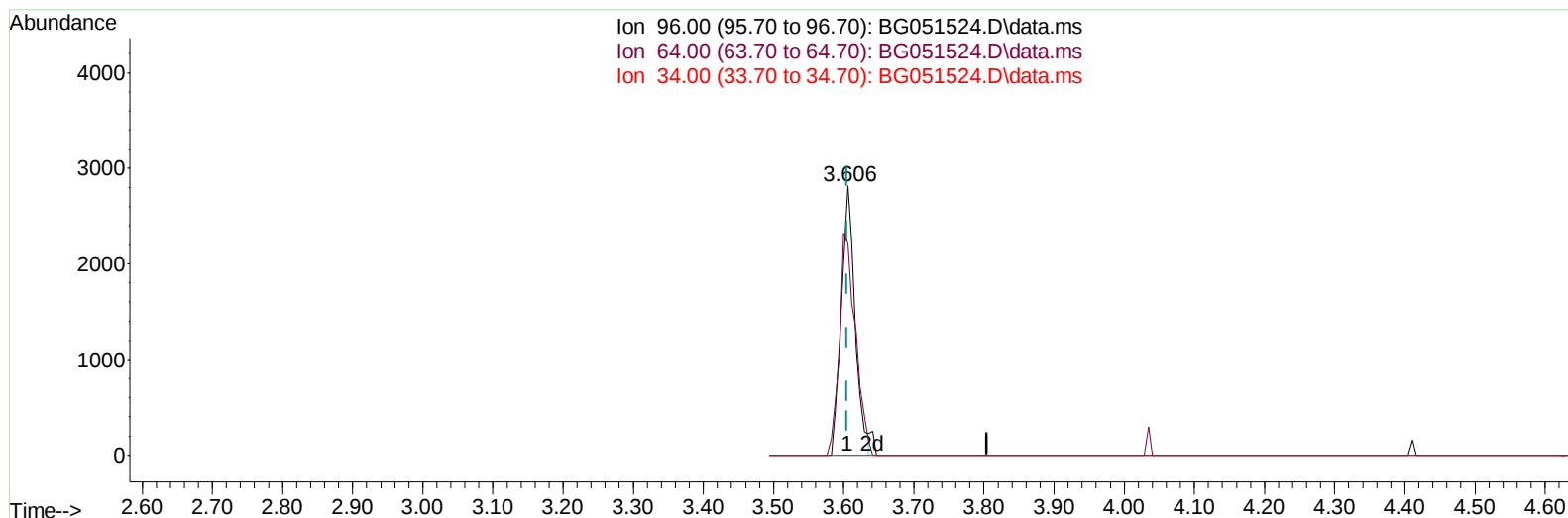


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051524.D
Acq On : 15 Dec 2021 5:27
Operator : CG/JU
Sample : PB141354BL
Misc :
ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 15 06:58:56 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



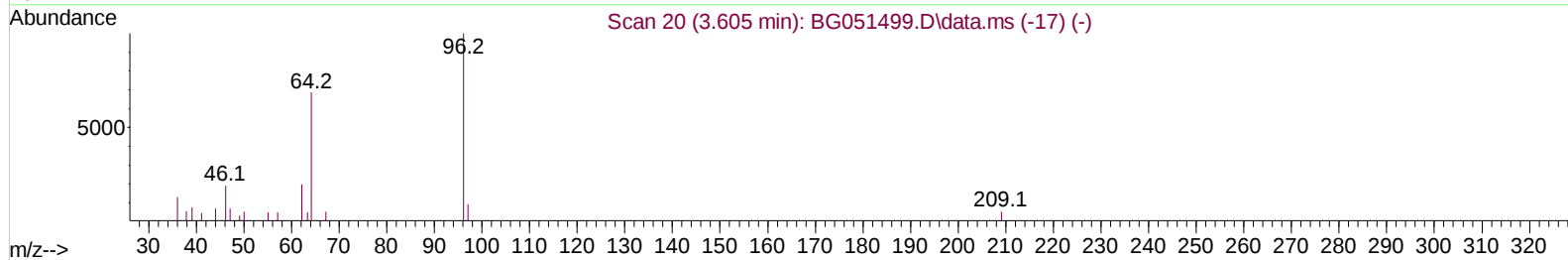
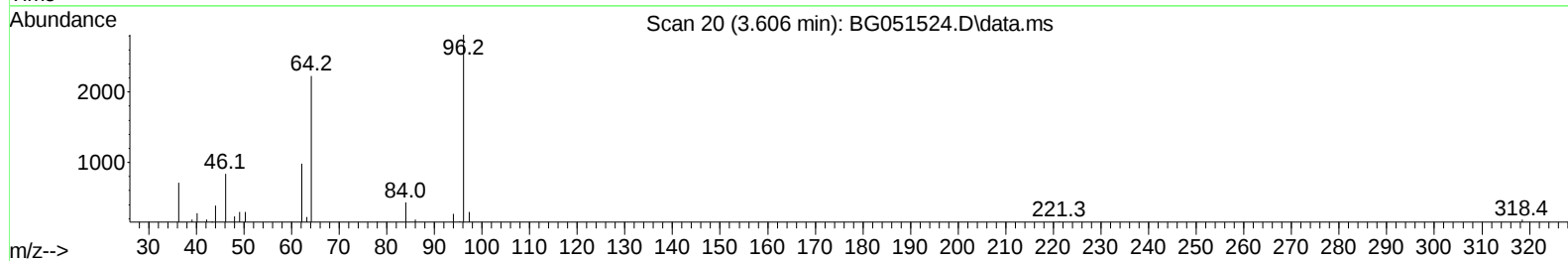
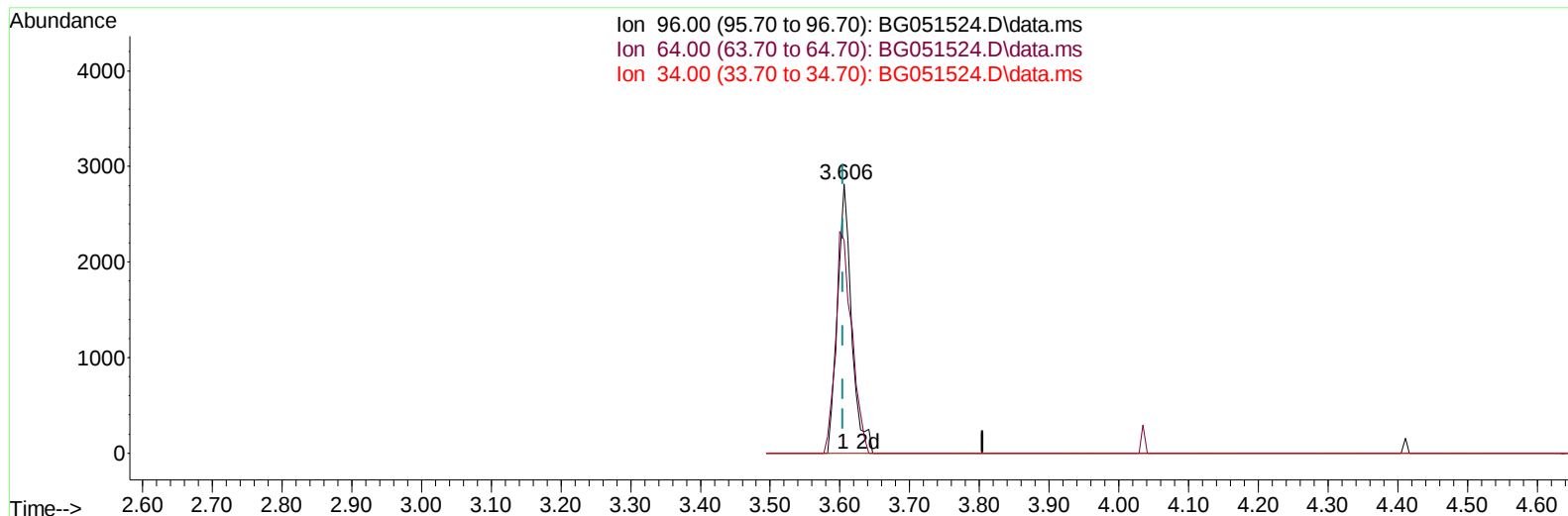
TIC: BG051524.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.606min (+ 0.001) 5.04 ng/uL****response 3849**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	79.21
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051524.D
Acq On : 15 Dec 2021 5:27
Operator : CG/JU
Sample : PB141354BL
Misc :
ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 15 06:58:56 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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TIC: BG051524.D\data.ms

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

3.606min (+ 0.001) 5.16 ng/uL m

response 3937

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	79.21
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051524.D
 Acq On : 15 Dec 2021 5:27
 Operator : CG/JU
 Sample : PB141354BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 15 06:58:56 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
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.276	152	31713	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.113	136	126829	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.914	164	92773	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.657	188	232475	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.975	240	235719	20.000	ng/ul	#-0.01
88) Perylene-d12	25.470	264	226354	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.606	96	3937m	5.156	ng/uL	0.00
4) Pyridine-d5	4.035	84	72131	31.263	ng/ul	-0.01
7) Phenol-d5	7.407	99	87876	31.034	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.583	67	52392	31.125	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.800	132	64368	32.685	ng/ul	-0.01
15) 4-Methylphenol-d8	8.969	113	68169	31.145	ng/ul	-0.01
21) Nitrobenzene-d5	9.445	128	31704	34.096	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.179	143	35627	35.049	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.720	165	67196	30.484	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.237	131	89319	30.776	ng/ul	0.00
46) Dimethylphthalate-d6	14.303	166	243405	32.780	ng/ul	-0.01
49) Acenaphthylene-d8	14.609	160	288221	31.253	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	39764	32.930	ng/ul	-0.01
60) Fluorene-d10	15.901	176	206998	31.115	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.995	200	40035	33.054	ng/ul	-0.02
73) Anthracene-d10	17.757	188	364079	32.830	ng/ul	0.00
81) Pyrene-d10	20.031	212	451632	31.835	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.235	264	395186	33.181	ng/ul	-0.02

Target Compounds	Qvalue

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

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