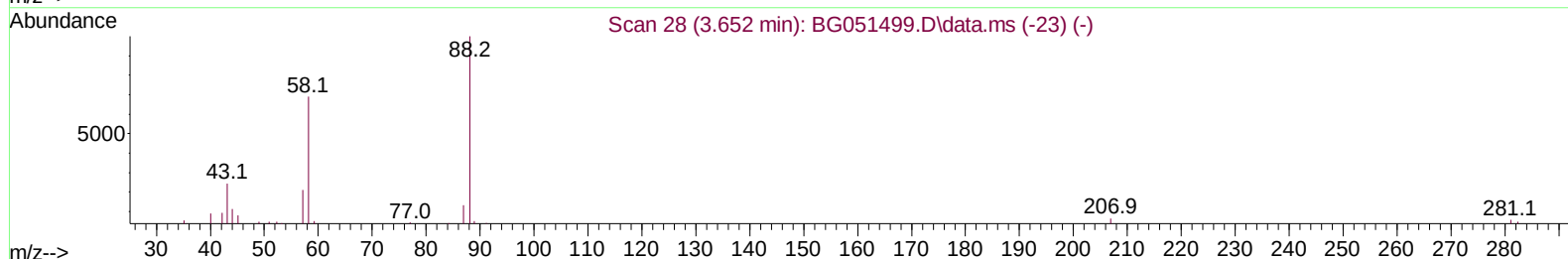
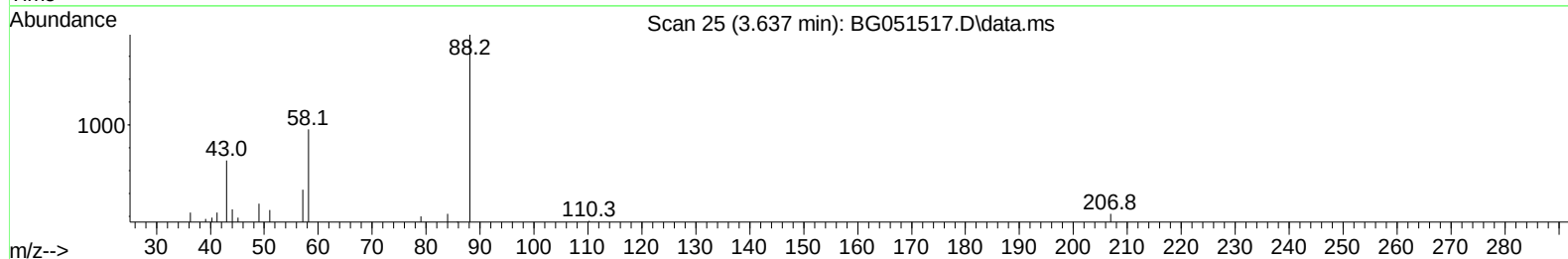
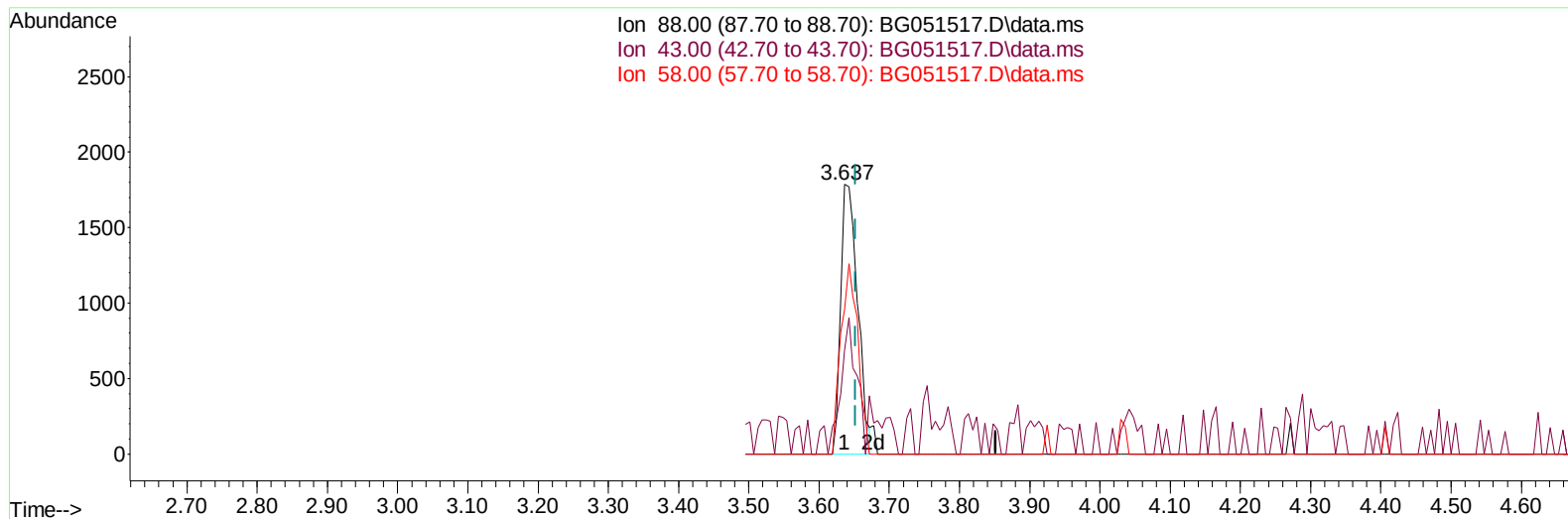


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051517.D
Acq On : 15 Dec 2021 00:05
Operator : CG/JU
Sample : M4995-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 15 01:35:50 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: BG051517.D\data.ms

(2) 1,4-Dioxane

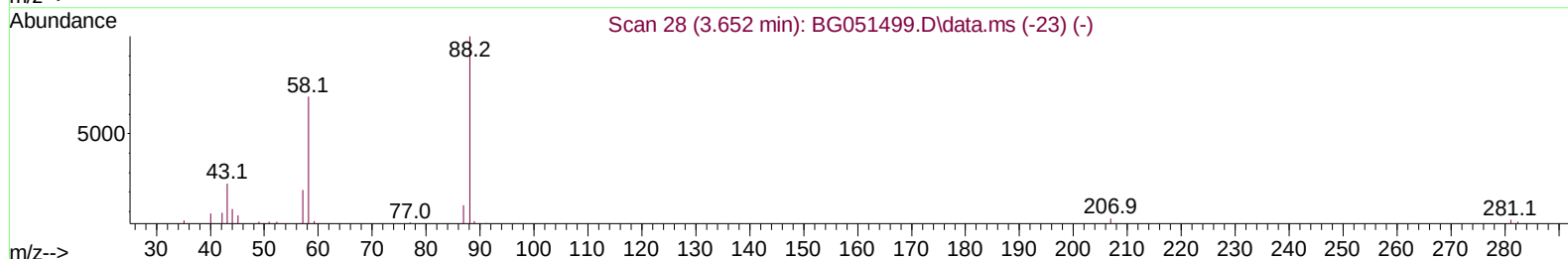
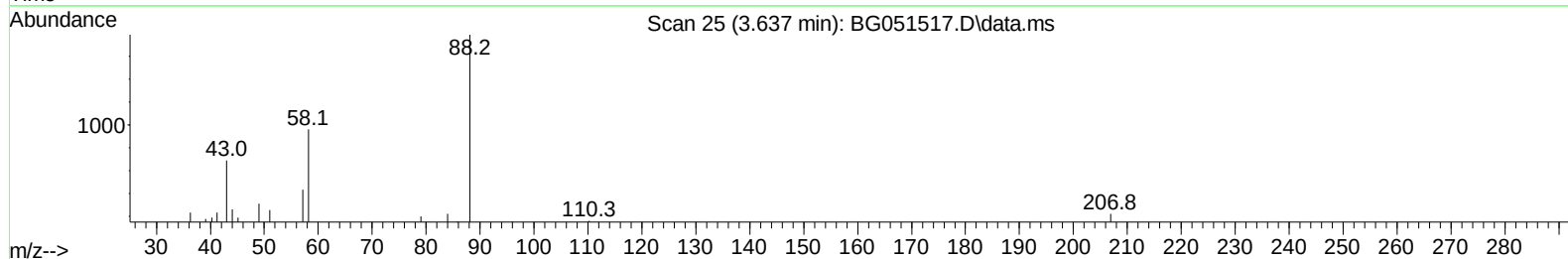
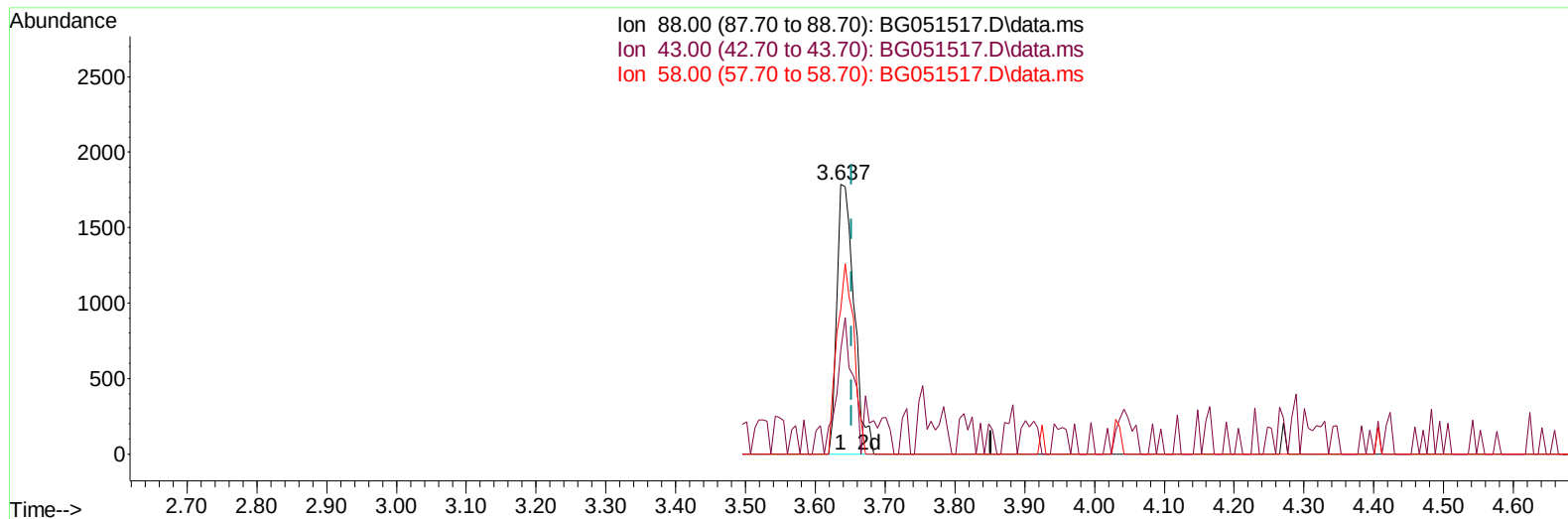
3.637min (-0.015) 4.11 ng/uL

response 2997

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	28.70	38.63#
58.00	78.00	53.70#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051517.D
Acq On : 15 Dec 2021 00:05
Operator : CG/JU
Sample : M4995-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 15 01:35:50 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: BG051517.D\data.ms

(2) 1,4-Dioxane

3.637min (-0.015) 4.20 ng/uL m

response 3063

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	28.70	38.63#
58.00	78.00	53.70#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051517.D
 Acq On : 15 Dec 2021 00:05
 Operator : CG/JU
 Sample : M4995-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 15 01:35:50 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
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.283	152	29007	20.000	ng/ul	# 0.00
20) Naphthalene-d8	11.121	136	117365	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.915	164	86624	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.659	188	226150	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.976	240	230317	20.000	ng/ul	# 0.00
88) Perylene-d12	25.478	264	224550	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.601	96	3735	5.348	ng/uL	0.00
4) Pyridine-d5	4.036	84	17536	8.310	ng/ul	0.00
7) Phenol-d5	7.408	99	16235	6.268	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.584	67	46069	29.922	ng/ul	0.00
11) 2-Chlorophenol-d4	7.801	132	42329	23.499	ng/ul	0.00
15) 4-Methylphenol-d8	8.970	113	29815	14.893	ng/ul	0.00
21) Nitrobenzene-d5	9.446	128	28935	33.627	ng/ul	0.00
24) 2-Nitrophenol-d4	10.181	143	31003	32.959	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.721	165	55862	27.385	ng/ul	0.00
31) 4-Chloroaniline-d4	11.238	131	72099	26.845	ng/ul	0.00
46) Dimethylphthalate-d6	14.304	166	232480	33.531	ng/ul	0.00
49) Acenaphthylene-d8	14.610	160	271602	31.542	ng/ul	0.00
54) 4-Nitrophenol-d4	15.068	143	7548	6.694	ng/ul	0.00
60) Fluorene-d10	15.902	176	209039	33.652	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.996	200	41934	35.590	ng/ul	-0.02
73) Anthracene-d10	17.759	188	373860	34.655	ng/ul	0.00
81) Pyrene-d10	20.032	212	461422	33.288	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.237	264	415090	35.133	ng/ul	-0.02
Target Compounds						
2) 1,4-Dioxane	3.637	88	3063m	4.202	ng/uL	
30) Naphthalene	11.168	128	9960	1.575	ng/ul	99

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

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