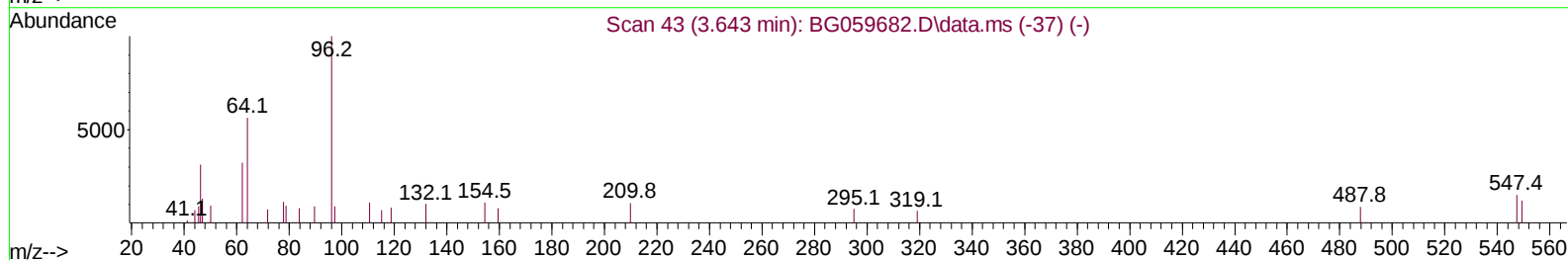
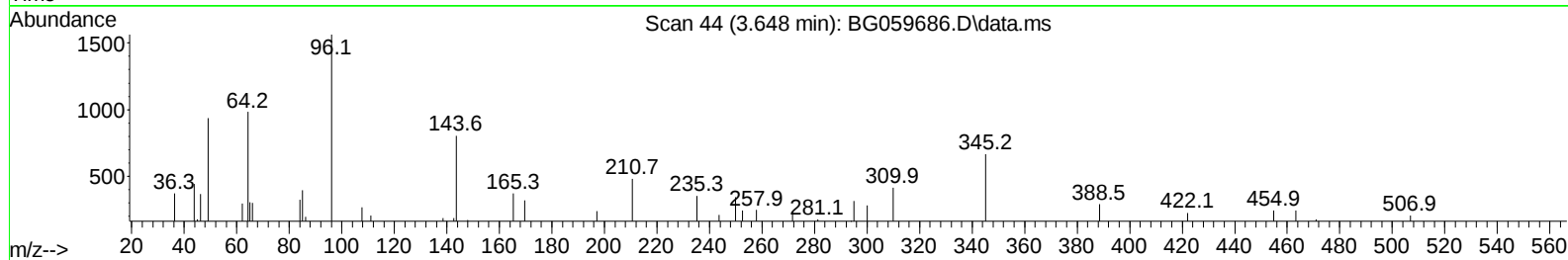
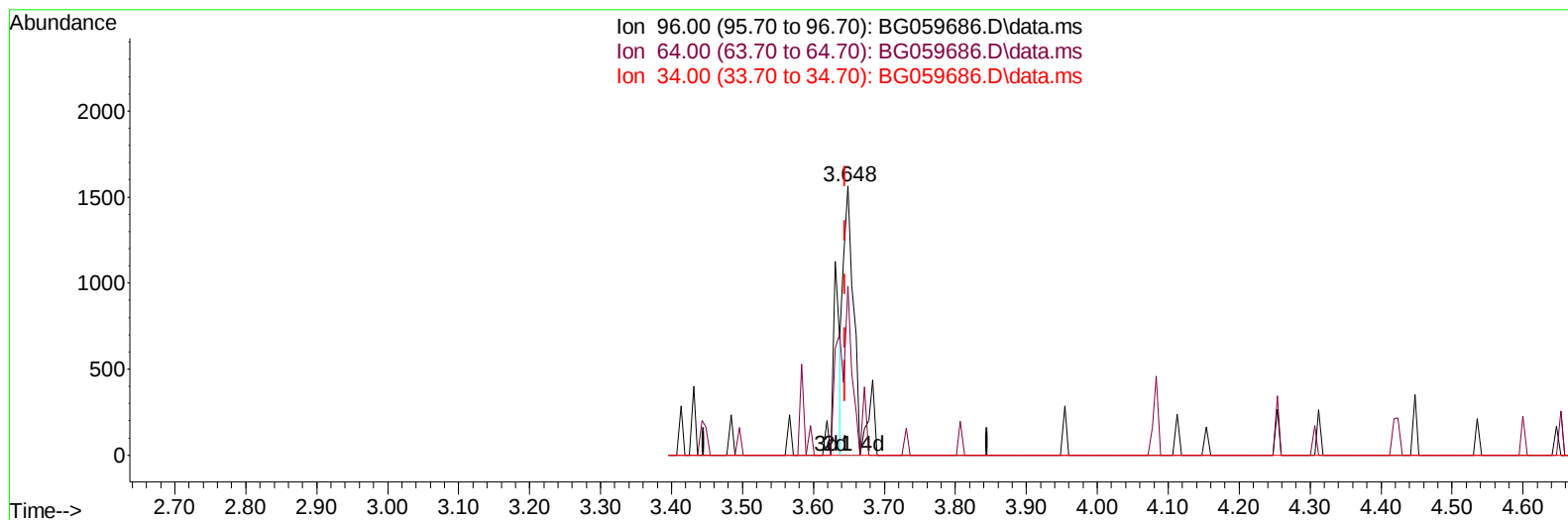


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
Data File : BG059686.D
Acq On : 6 Nov 2023 12:29
Operator : MA/JU
Sample : 05013-12
Misc :
ALS Vial : 79 Sample Multiplier: 1

Quant Time: Nov 06 13:02:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Nov 04 02:46:50 2023
Response via : Initial Calibration



TIC: BG059686.D\data.ms

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

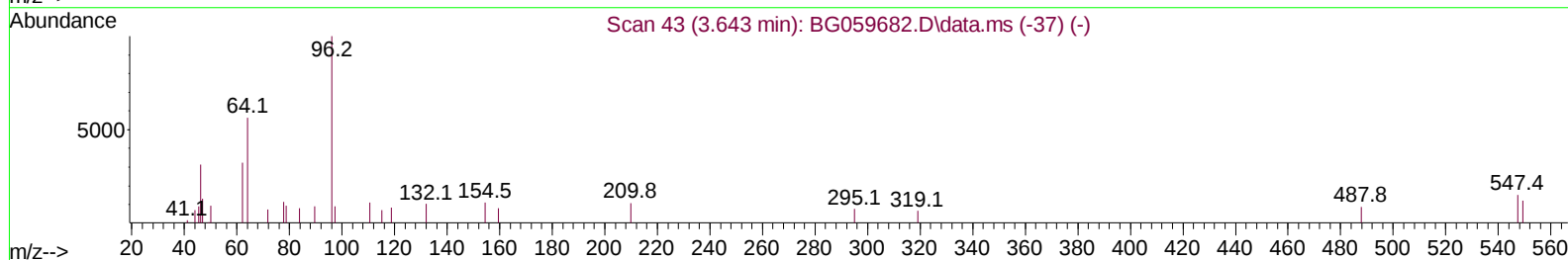
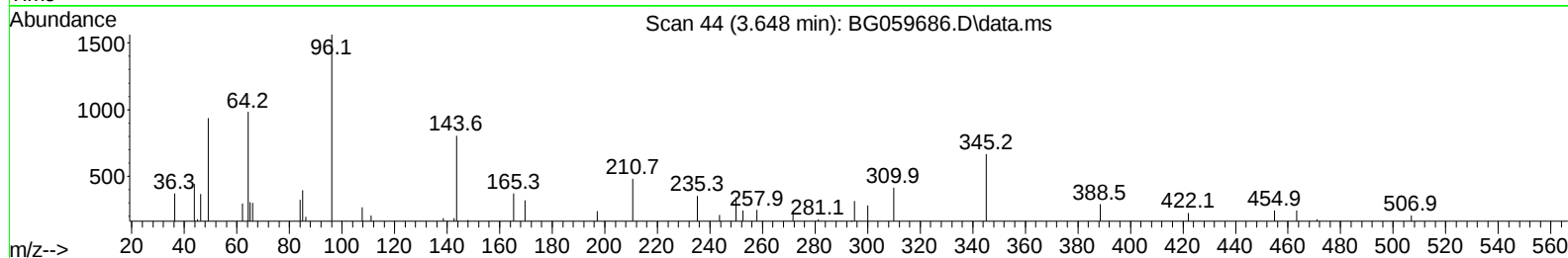
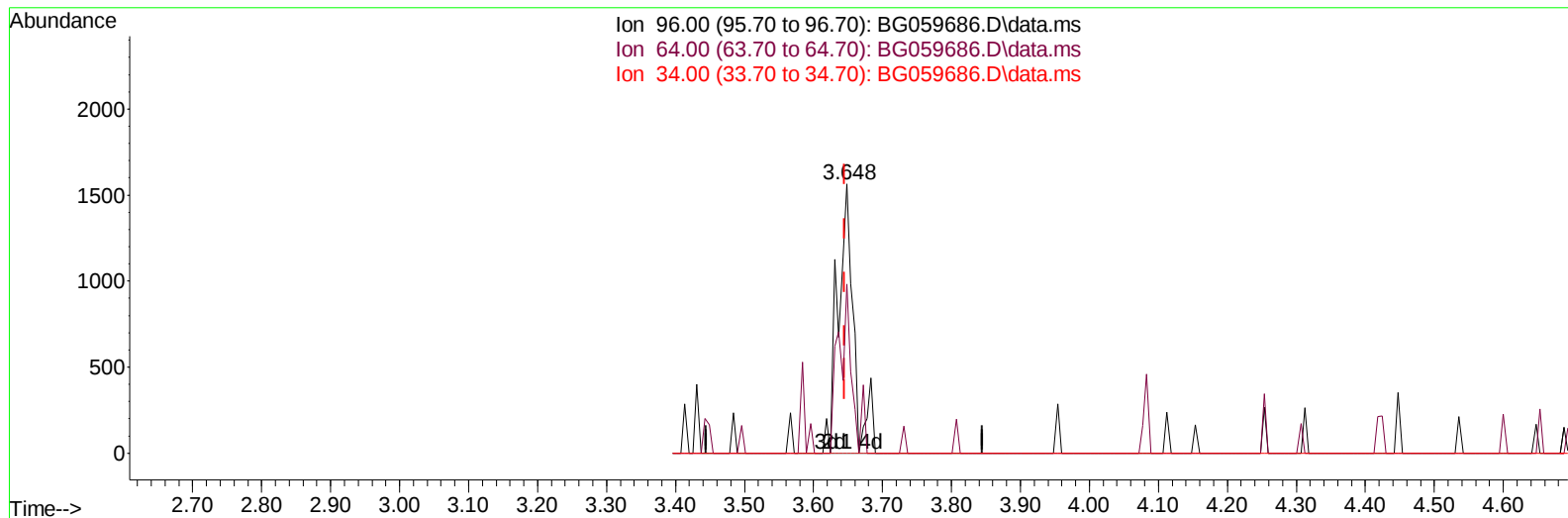
3.648min (+ 0.003) 2.21 ng/uL

response 1548

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.30	62.83#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
Data File : BG059686.D
Acq On : 6 Nov 2023 12:29
Operator : MA/JU
Sample : 05013-12
Misc :
ALS Vial : 79 Sample Multiplier: 1

Quant Time: Nov 06 13:02:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Nov 04 02:46:50 2023
Response via : Initial Calibration



TIC: BG059686.D\data.ms

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

3.648min (+ 0.003) 3.61 ng/uL m

response 2532

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.30	62.83#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
 Data File : BG059686.D
 Acq On : 6 Nov 2023 12:29
 Operator : MA/JU
 Sample : 05013-12
 Misc :
 ALS Vial : 79 Sample Multiplier: 1

Quant Time: Nov 06 13:03:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 04 02:46:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.349	152	26258	20.000	ng/ul	0.00
20) Naphthalene-d8	11.198	136	137672	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.976	164	95734	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.732	188	228271	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.062	240	262562	20.000	ng/ul	0.00
88) Perylene-d12	25.658	264	317993	20.000	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.648	96	2532m	3.614	ng/uL	0.00
4) Pyridine-d5	4.077	84	17885	8.681	ng/ul	0.00
7) Phenol-d5	7.456	99	13087	4.875	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.667	67	32845	22.883	ng/ul	0.00
11) 2-Chlorophenol-d4	7.867	132	37793	20.728	ng/ul	0.00
15) 4-Methylphenol-d8	9.036	113	30307	13.794	ng/ul	0.00
21) Nitrobenzene-d5	9.547	128	29371	29.172	ng/ul	0.00
24) 2-Nitrophenol-d4	10.270	143	34463	28.364	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.787	165	68109	31.551	ng/ul	0.00
31) 4-Chloroaniline-d4	11.339	131	88725	26.119	ng/ul	0.00
46) Dimethylphthalate-d6	14.371	166	270519	35.209	ng/ul	0.00
49) Acenaphthylene-d8	14.682	160	270824	29.560	ng/ul	0.00
54) 4-Nitrophenol-d4	15.153	143	8463	6.239	ng/ul	0.00
60) Fluorene-d10	15.963	176	236153	32.648	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.063	200	43223	29.915	ng/ul	0.00
73) Anthracene-d10	17.826	188	384769	32.206	ng/ul	0.00
81) Pyrene-d10	20.100	212	587016	34.937	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.411	264	677040	36.370	ng/ul	0.00

Target Compounds	Qvalue

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

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