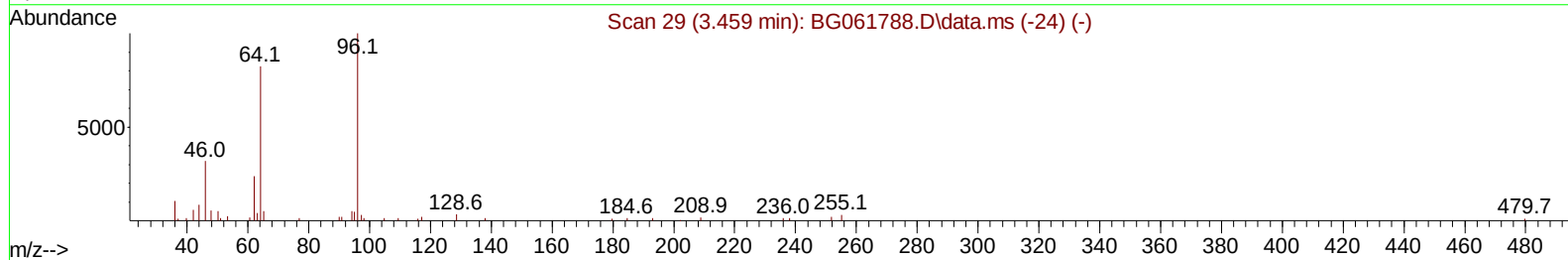
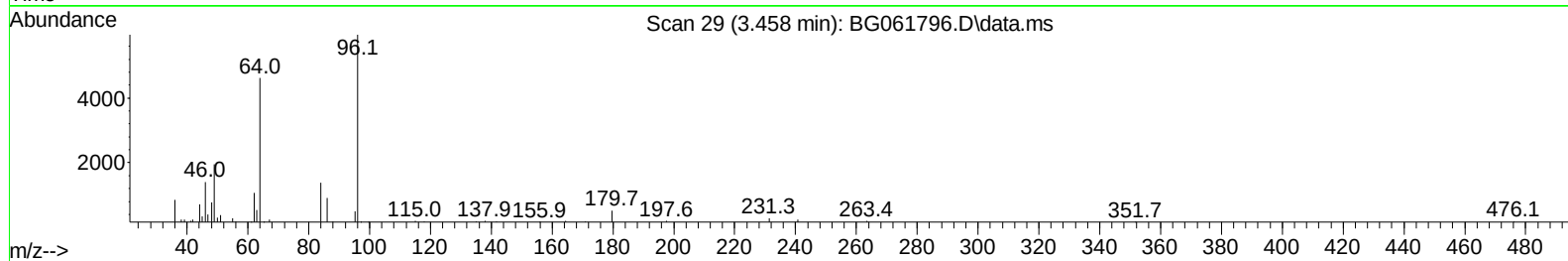
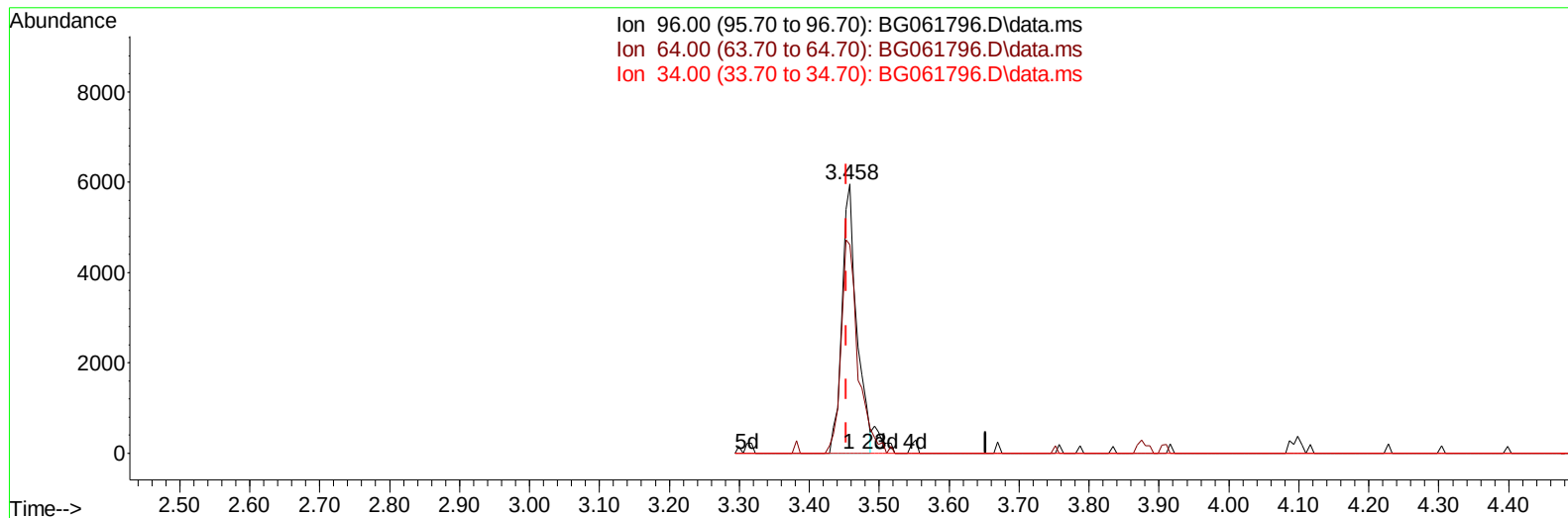


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061796.D
Acq On : 29 Jun 2024 7:55
Operator : MA/JU
Sample : P2897-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 01 00:28:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 12:23:53 2024
Response via : Initial Calibration



TIC: BG061796.D\data.ms

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

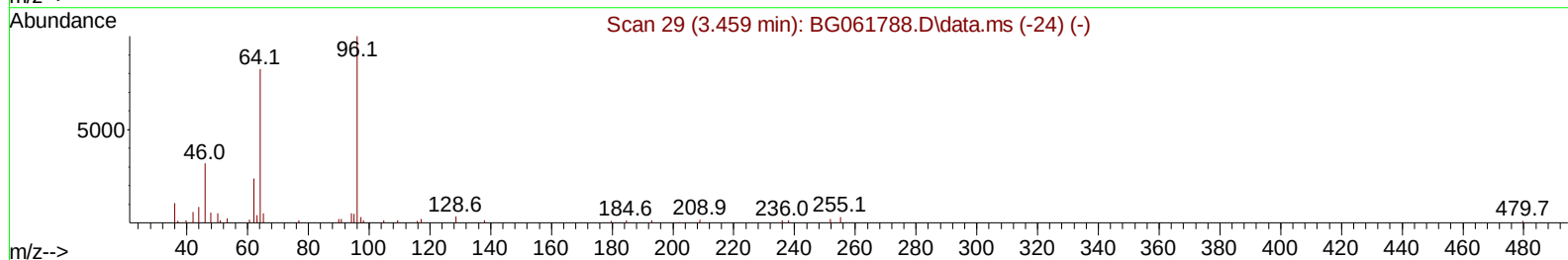
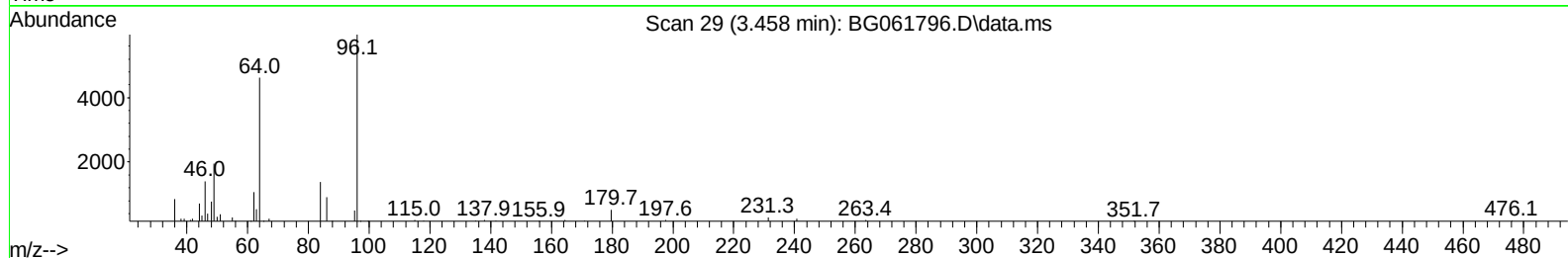
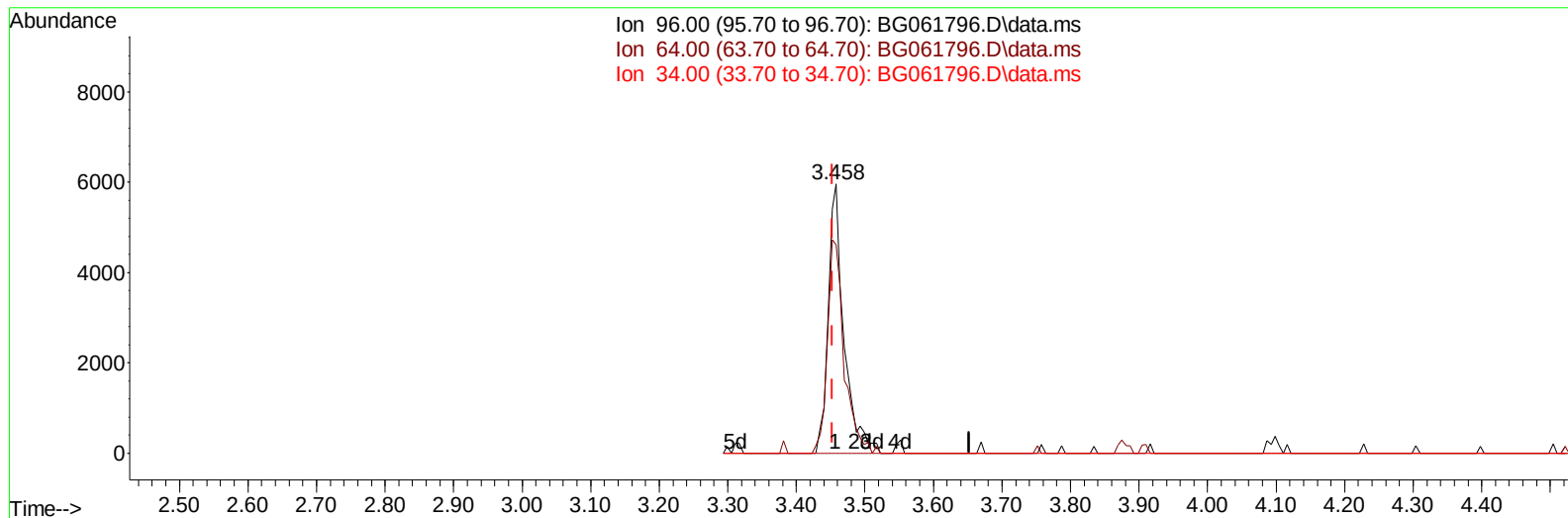
3.458min (+ 0.005) 4.05 ng/uL

response 8898

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	77.54
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061796.D
Acq On : 29 Jun 2024 7:55
Operator : MA/JU
Sample : P2897-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 01 00:28:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 12:23:53 2024
Response via : Initial Calibration



TIC: BG061796.D\data.ms

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

3.458min (+ 0.005) 4.33 ng/uL m

response 9508

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	77.54
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
 Data File : BG061796.D
 Acq On : 29 Jun 2024 7:55
 Operator : MA/JU
 Sample : P2897-16
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 01 00:34:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 27 12:23:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.058	152	79337	20.000	ng/ul	0.00
20) Naphthalene-d8	10.867	136	380430	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.686	164	250046	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.424	188	551324	20.000	ng/ul	0.00
79) Chrysene-d12	21.684	240	440191	20.000	ng/ul	0.00
88) Perylene-d12	24.898	264	504107	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.458	96	9508m	4.329	ng/uL	0.00
4) Pyridine-d5	3.875	84	100376	15.117	ng/ul	0.00
7) Phenol-d5	7.201	99	204894	23.789	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.377	67	127523	23.752	ng/ul	0.00
11) 2-Chlorophenol-d4	7.582	132	146468	24.782	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	162979	24.478	ng/ul	0.00
21) Nitrobenzene-d5	9.222	128	73885	28.603	ng/ul	0.00
24) 2-Nitrophenol-d4	9.939	143	83572	31.944	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.479	165	162661	26.258	ng/ul	0.00
31) 4-Chloroaniline-d4	10.996	131	225159	24.007	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	504046	26.195	ng/ul	0.00
49) Acenaphthylene-d8	14.380	160	580229	27.515	ng/ul	0.00
54) 4-Nitrophenol-d4	14.868	143	73038	22.101	ng/ul	0.00
60) Fluorene-d10	15.673	176	454499	26.981	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.779	200	42105	17.860	ng/ul	0.00
73) Anthracene-d10	17.524	188	704865	28.009	ng/ul	0.00
81) Pyrene-d10	19.809	212	736282	29.945	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.680	264	695351	28.881	ng/ul	0.00

Target Compounds	Qvalue

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

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