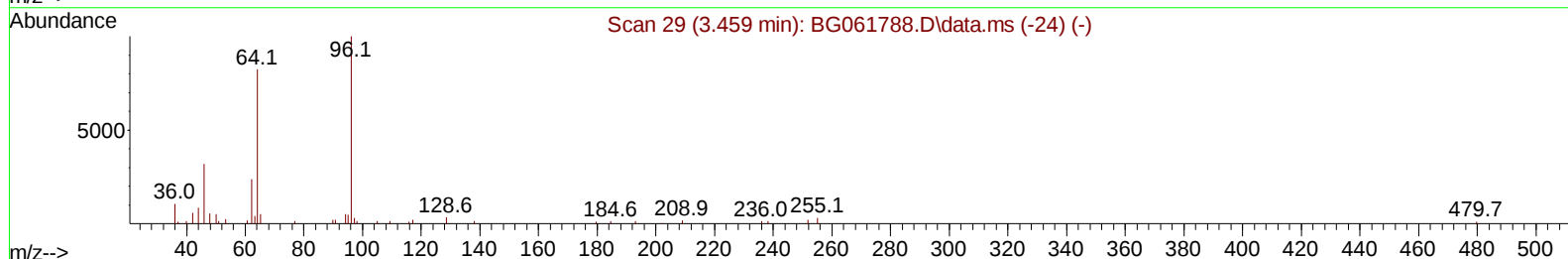
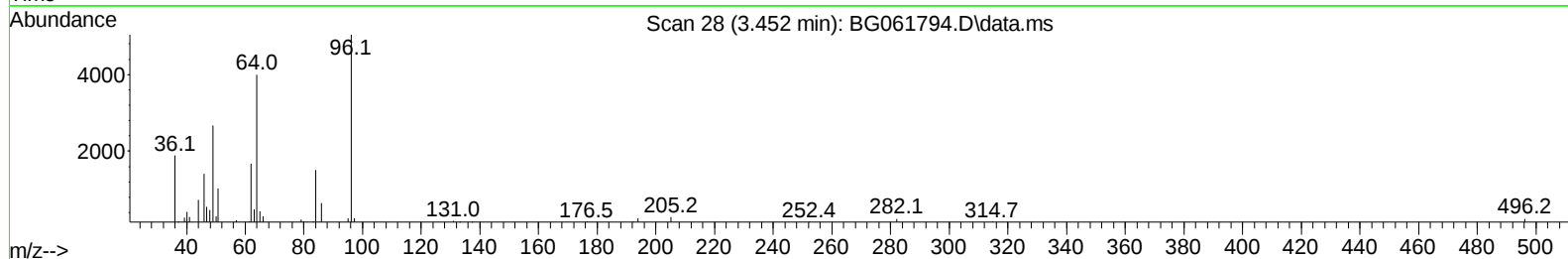
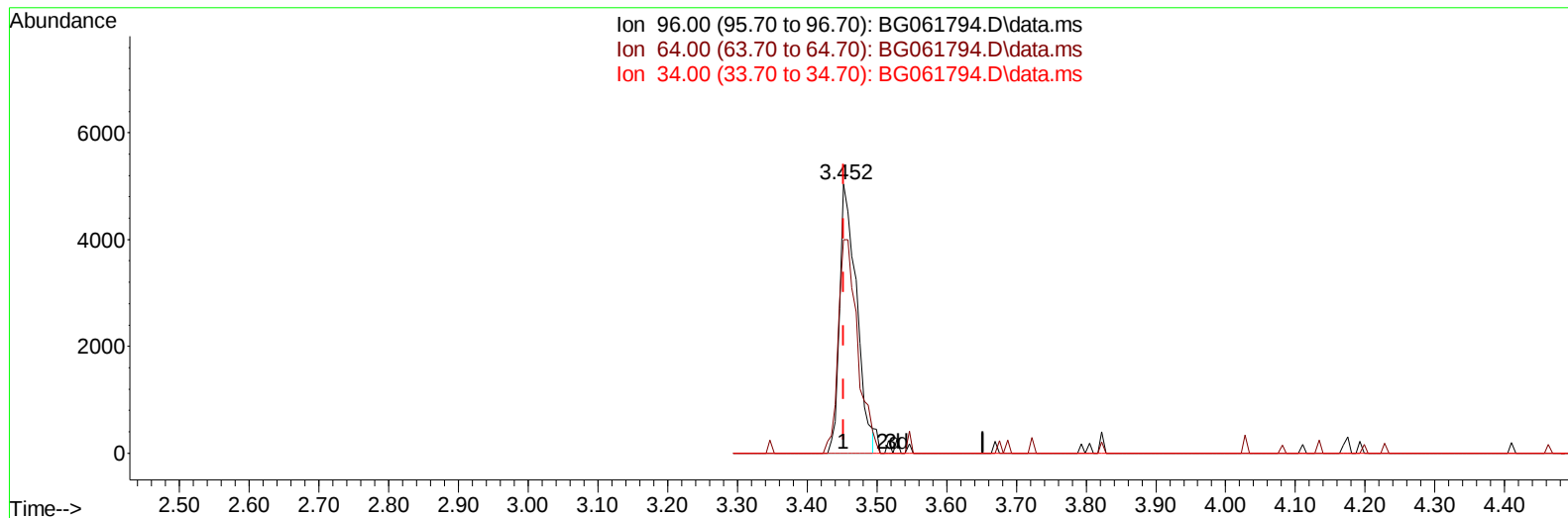


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061794.D
Acq On : 29 Jun 2024 6:34
Operator : MA/JU
Sample : P2897-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Jul 01 00:28:18 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: BG061794.D\data.ms

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

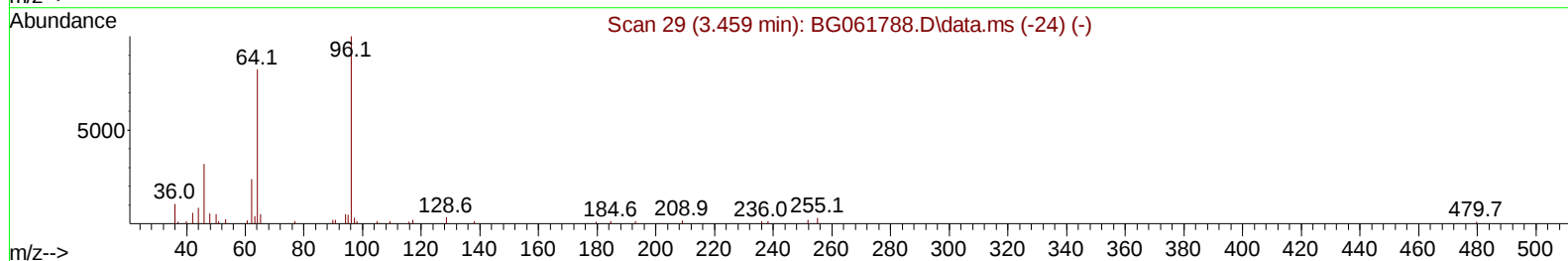
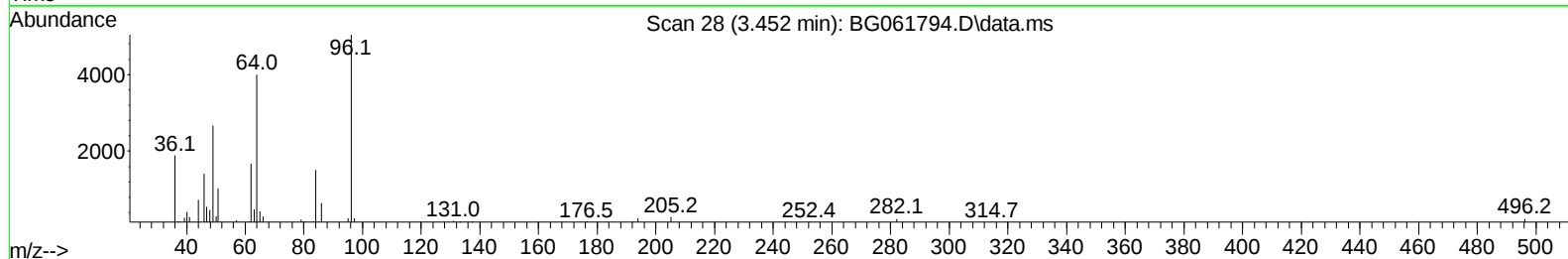
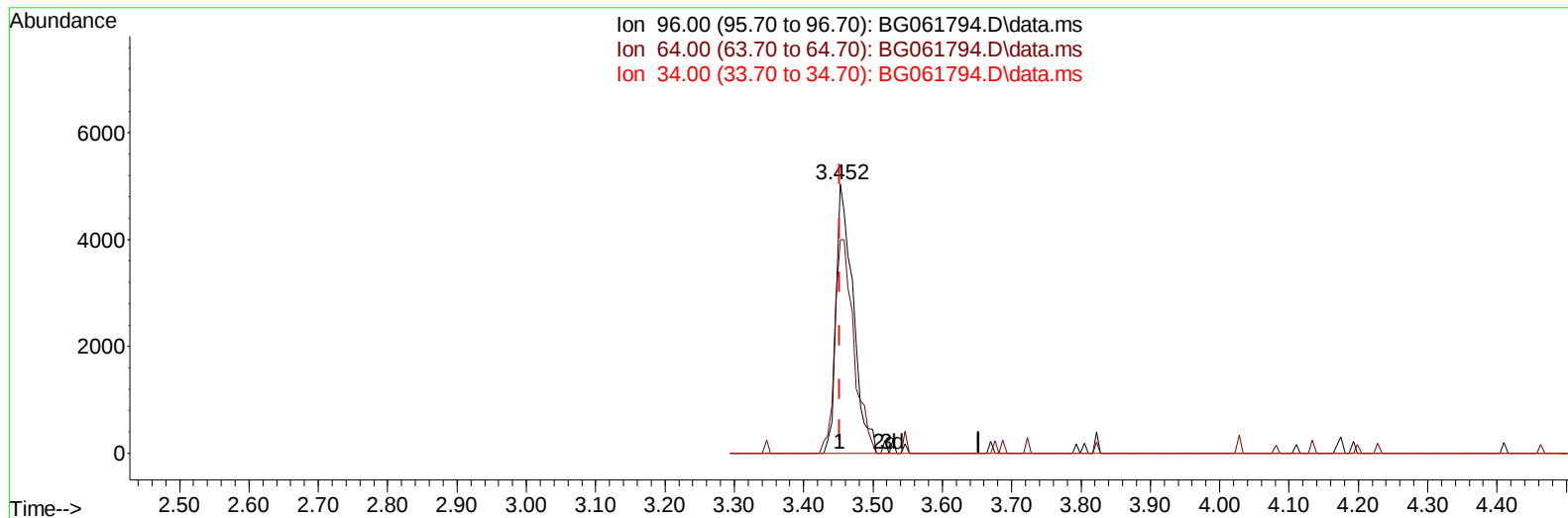
3.452min (-0.000) 4.57 ng/uL

response 8444

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	79.40#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
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Sample : P2897-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Jul 01 00:28:18 2024
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Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 12:23:53 2024
Response via : Initial Calibration



TIC: BG061794.D\data.ms

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

3.452min (-0.000) 4.66 ng/uL m

response 8600

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	79.40#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
 Data File : BG061794.D
 Acq On : 29 Jun 2024 6:34
 Operator : MA/JU
 Sample : P2897-12
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Jul 01 00:30:55 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	8.053	152	66693	20.000	ng/ul	0.00
20) Naphthalene-d8	10.867	136	326589	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.680	164	214394	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.424	188	487216	20.000	ng/ul	0.00
79) Chrysene-d12	21.684	240	380254	20.000	ng/ul	0.00
88) Perylene-d12	24.904	264	438459	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.452	96	8600m	4.658	ng/uL	0.00
4) Pyridine-d5	3.881	84	25609	4.588	ng/ul	0.01
7) Phenol-d5	7.201	99	185439	25.612	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.377	67	116367	25.783	ng/ul	0.00
11) 2-Chlorophenol-d4	7.583	132	131331	26.433	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	143791	25.690	ng/ul	0.00
21) Nitrobenzene-d5	9.222	128	67957	30.645	ng/ul	0.00
24) 2-Nitrophenol-d4	9.945	143	78183	34.811	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.479	165	145472	27.354	ng/ul	0.00
31) 4-Chloroaniline-d4	10.996	131	182807	22.705	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	462798	28.051	ng/ul	0.00
49) Acenaphthylene-d8	14.381	160	528324	29.220	ng/ul	0.00
54) 4-Nitrophenol-d4	14.862	143	66324	23.407	ng/ul	0.00
60) Fluorene-d10	15.673	176	410413	28.415	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.785	200	37730	18.110	ng/ul	0.00
73) Anthracene-d10	17.524	188	638058	28.690	ng/ul	0.00
81) Pyrene-d10	19.804	212	668000	31.450	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.680	264	636320	30.386	ng/ul	0.00

Target Compounds	Qvalue

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

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