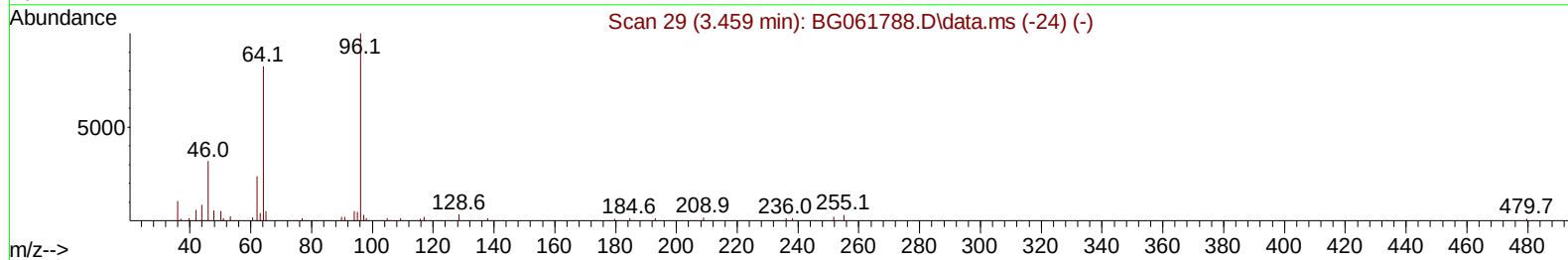
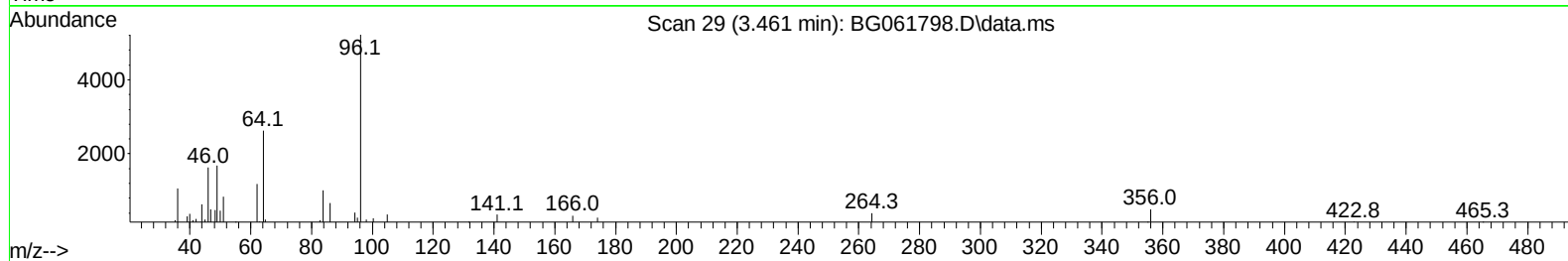
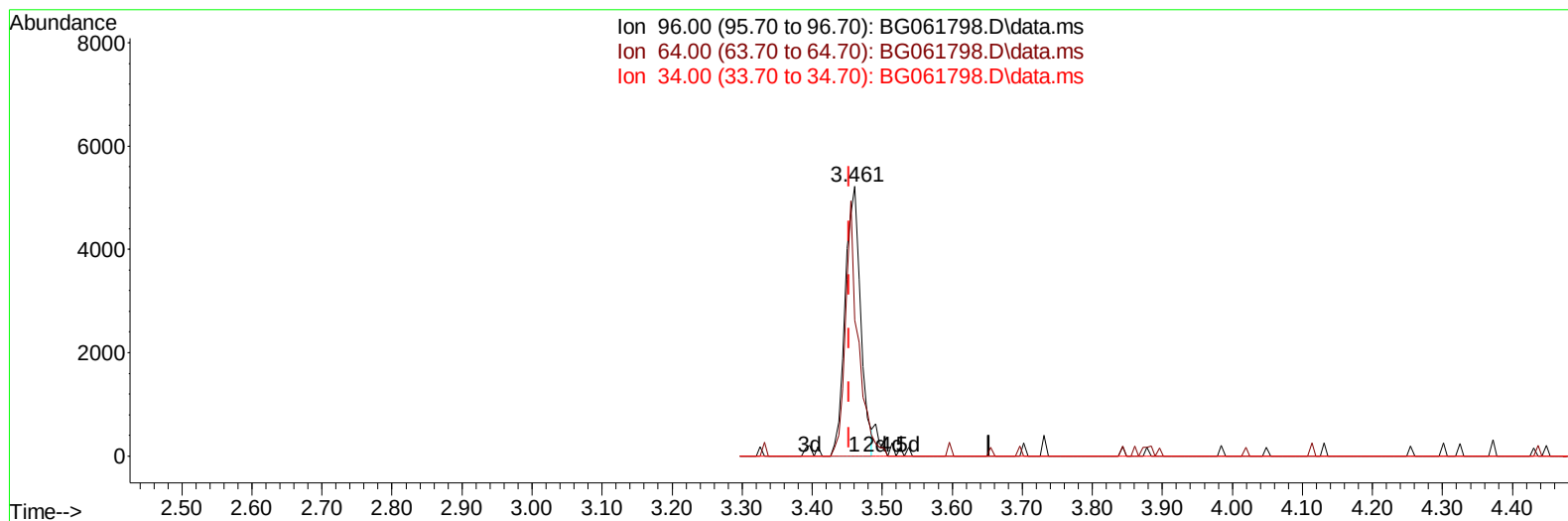


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061798.D
Acq On : 29 Jun 2024 9:16
Operator : MA/JU
Sample : P2897-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jul 01 00:28:50 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: BG061798.D\data.ms

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

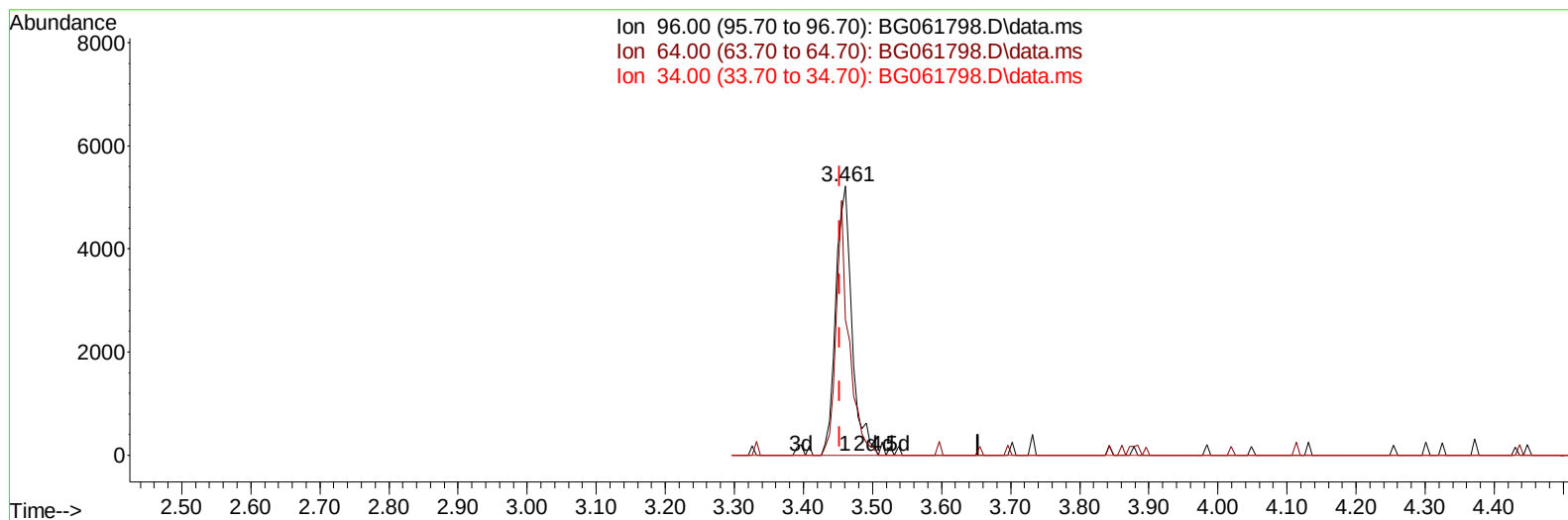
3.461min (+ 0.008) 4.03 ng/uL

response 8135

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	50.32#
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Jul 01 00:45:29 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	8.056	152	72920	20.000	ng/ul	0.00
20) Naphthalene-d8	10.864	136	346038	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.683	164	234499	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.427	188	511755	20.000	ng/ul	0.00
79) Chrysene-d12	21.687	240	405455	20.000	ng/ul	0.00
88) Perylene-d12	24.901	264	467797	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.461	96	8503m	4.212	ng/uL	0.00
4) Pyridine-d5	3.872	84	107510	17.616	ng/ul	0.00
7) Phenol-d5	7.204	99	188170	23.770	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.380	67	110912	22.476	ng/ul	0.00
11) 2-Chlorophenol-d4	7.586	132	126864	23.354	ng/ul	0.00
15) 4-Methylphenol-d8	8.749	113	145623	23.796	ng/ul	0.00
21) Nitrobenzene-d5	9.219	128	68202	29.027	ng/ul	0.00
24) 2-Nitrophenol-d4	9.942	143	78402	32.946	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.476	165	145011	25.735	ng/ul	0.00
31) 4-Chloroaniline-d4	10.999	131	206480	24.203	ng/ul	0.00
46) Dimethylphthalate-d6	14.090	166	478727	26.529	ng/ul	0.00
49) Acenaphthylene-d8	14.378	160	538669	27.238	ng/ul	0.00
54) 4-Nitrophenol-d4	14.865	143	71412	23.042	ng/ul	0.00
60) Fluorene-d10	15.670	176	431524	27.315	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.782	200	41589	19.005	ng/ul	0.00
73) Anthracene-d10	17.527	188	684076	29.284	ng/ul	0.00
81) Pyrene-d10	19.807	212	716266	31.626	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.683	264	671088	30.036	ng/ul	0.00

Target Compounds	Qvalue

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

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