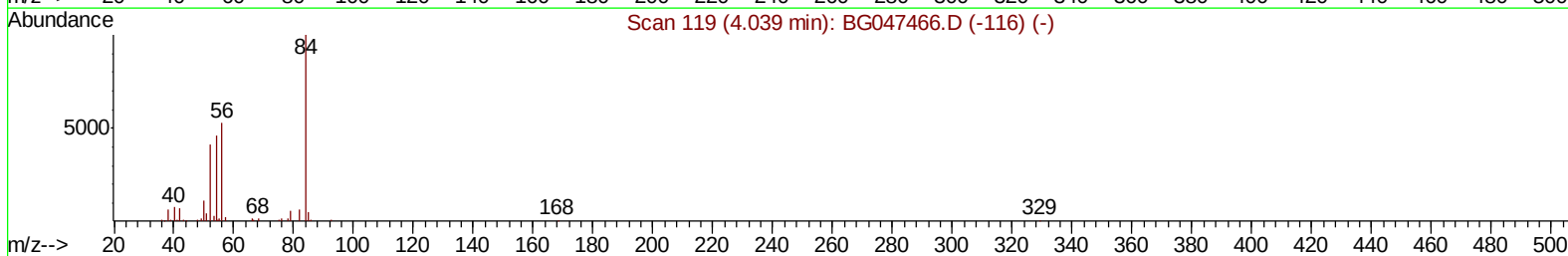
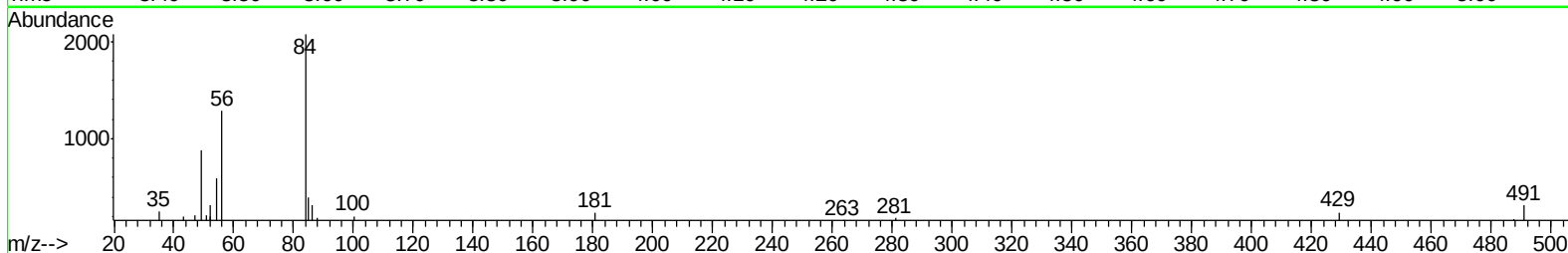
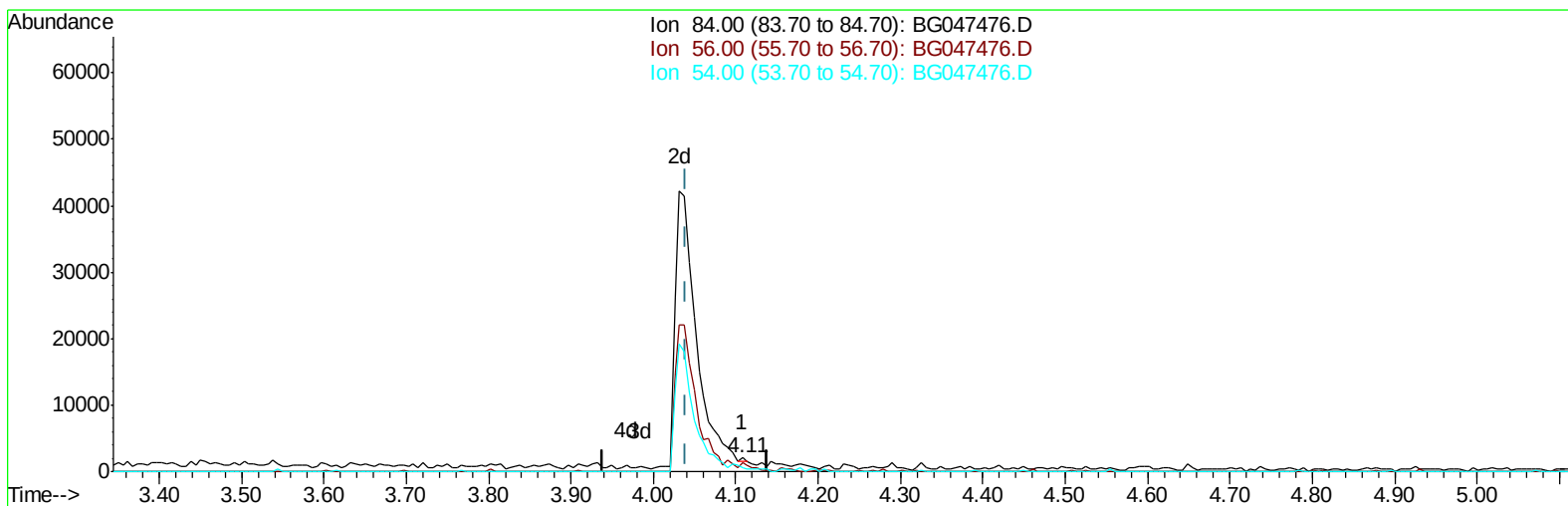


Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG112720\
Data File : BG047476.D
Acq On : 27 Nov 2020 11:08
Operator : CG/JU
Sample : PB133234BL
Misc : MDL-WATER-BL-04
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 27 11:50:58 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG111820.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 19 04:10:57 2020
Response via : Initial Calibration



TIC: BG047476.D

(4) Pyridine-d5 (S)

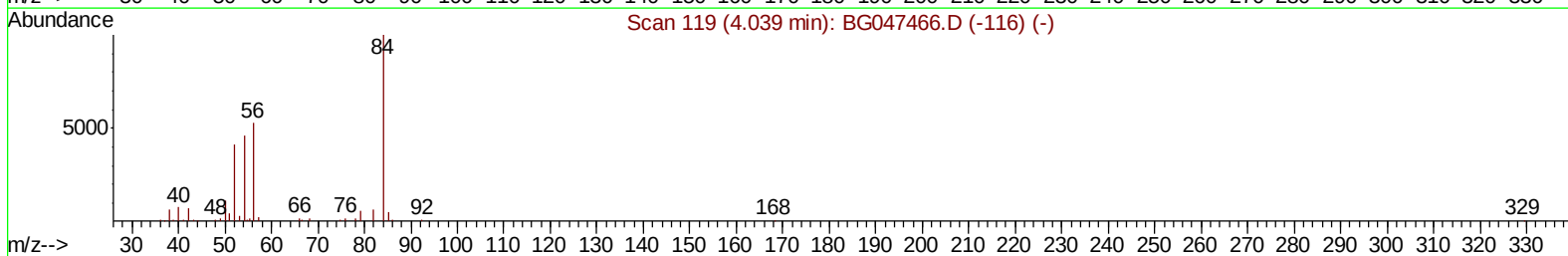
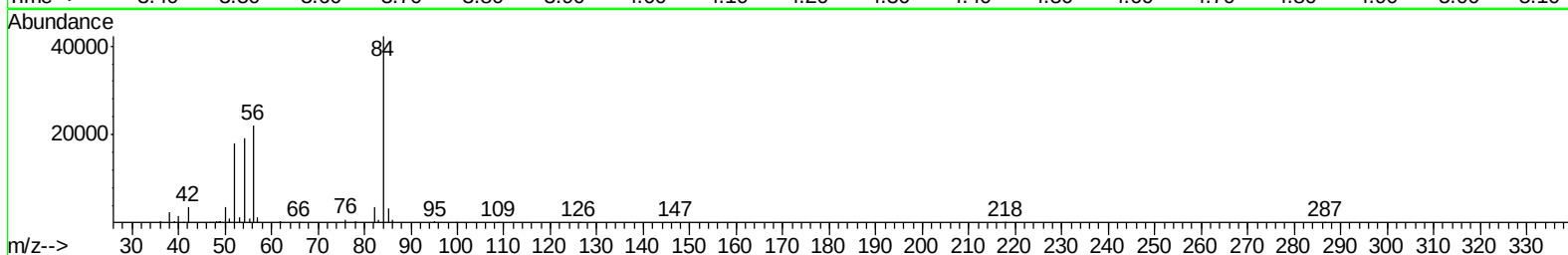
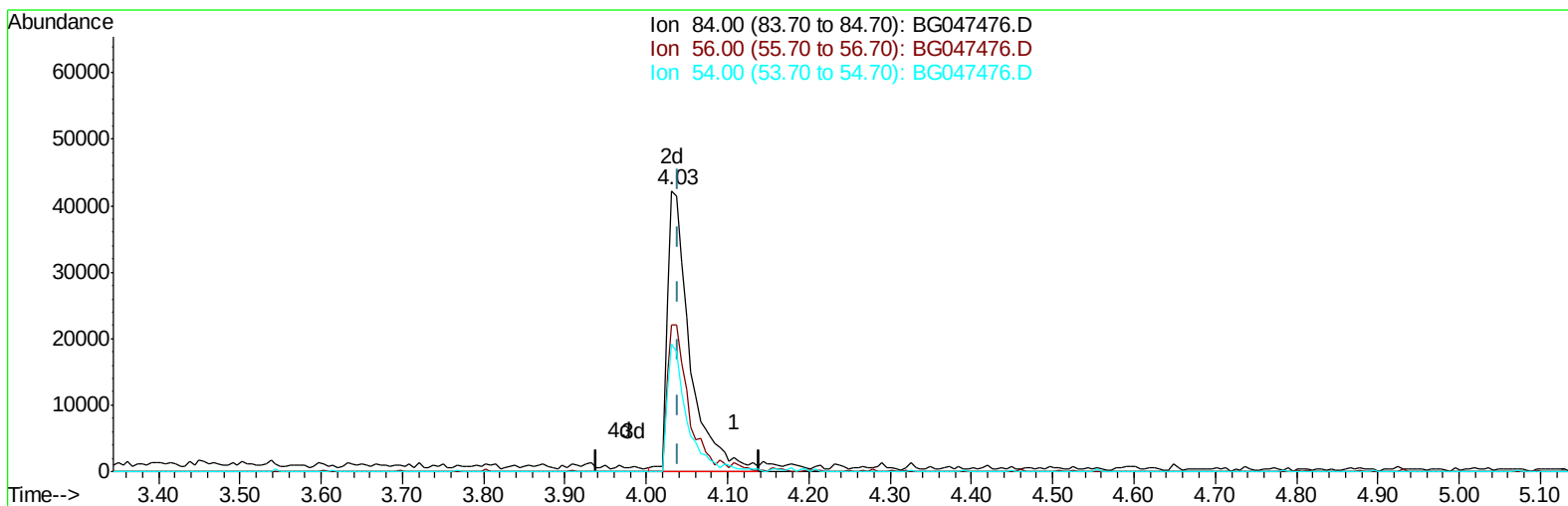
4.108min (+0.069) 0.11ng/ul

response 244

Ion	Exp%	Act%
84.00	100	100
56.00	54.00	62.01
54.00	28.60	28.78
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG112720\
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TIC: BG047476.D

(4) Pyridine-d5 (S)

4.032min (-0.008) 37.25ng/ul m

response 81933

Ion	Exp%	Act%
84.00	100	100
56.00	54.00	52.25
54.00	28.60	45.63#
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : PB133234BL
 Misc : MDL-WATER-BL-04
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 27 11:54:56 2020

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG111820.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Nov 19 04:10:57 2020

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	28383	20.00	ng/ul	0.00
20) Naphthalene-d8	11.09	136	113292	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.88	164	90344	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	224407	20.00	ng/ul	0.00
79) Chrysene-d12	21.89	240	231326	20.00	ng/ul	0.00
88) Perylene-d12	25.29	264	217680	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.61	96	5690	7.94	ng/uL	0.00
4) Pyridine-d5	4.03	84	81933m	37.25	ng/ul	0.00
7) Phenol-d5	7.40	99	98795	36.32	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.58	67	53114	33.37	ng/ul	0.00
11) 2-Chlorophenol-d4	7.79	132	75978	40.93	ng/ul	0.00
15) 4-Methylphenol-d8	8.96	113	76958	36.44	ng/ul	0.00
21) Nitrobenzene-d5	9.44	128	36704	39.03	ng/ul	0.00
24) 2-Nitrophenol-d4	10.16	143	45077	45.81	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.70	165	88741	39.20	ng/ul	0.00
31) 4-Chloroaniline-d4	11.22	131	103249	36.92	ng/ul	0.00
46) Dimethylphthalate-d6	14.27	166	308689	40.86	ng/ul	0.00
49) Acenaphthylene-d8	14.58	160	323419	37.52	ng/ul	0.00
54) 4-Nitrophenol-d4	15.04	143	41896	37.98	ng/ul	0.00
60) Fluorene-d10	15.86	176	264135	38.69	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.96	200	43391	36.50	ng/ul	0.00
73) Anthracene-d10	17.72	188	444741	40.94	ng/ul	0.00
81) Pyrene-d10	19.98	212	517406	39.64	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.07	264	519028	43.25	ng/ul	0.00

Target Compounds Ovalue

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

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