

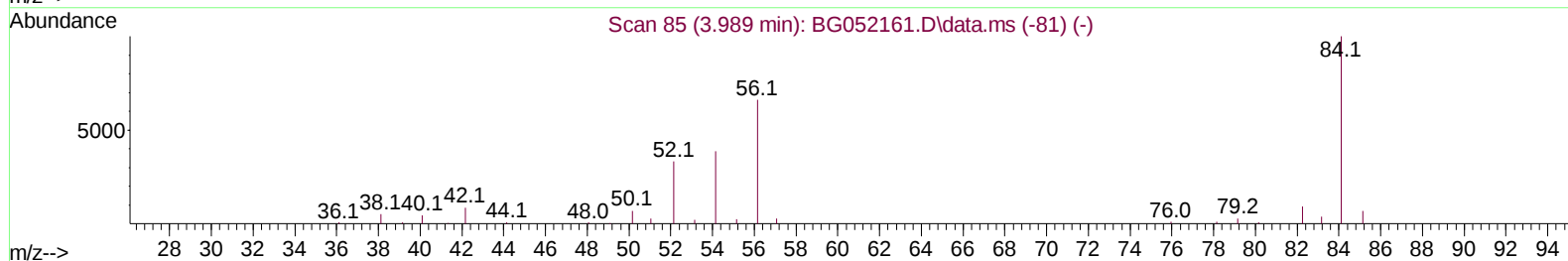
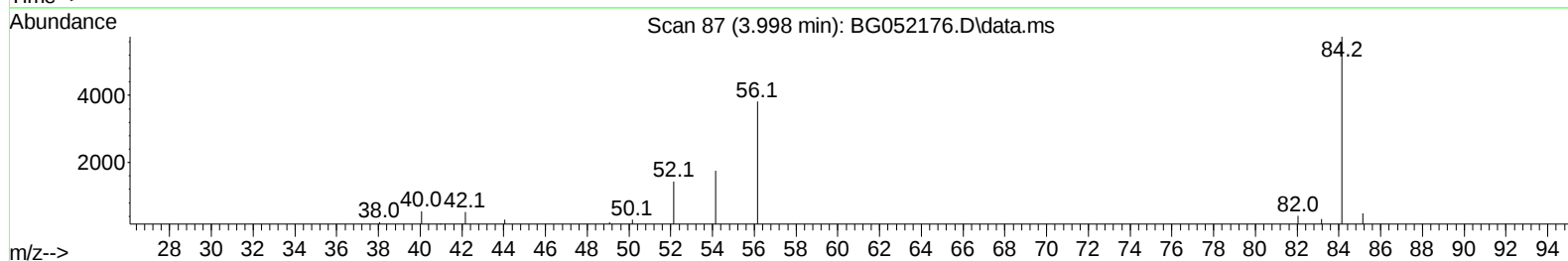
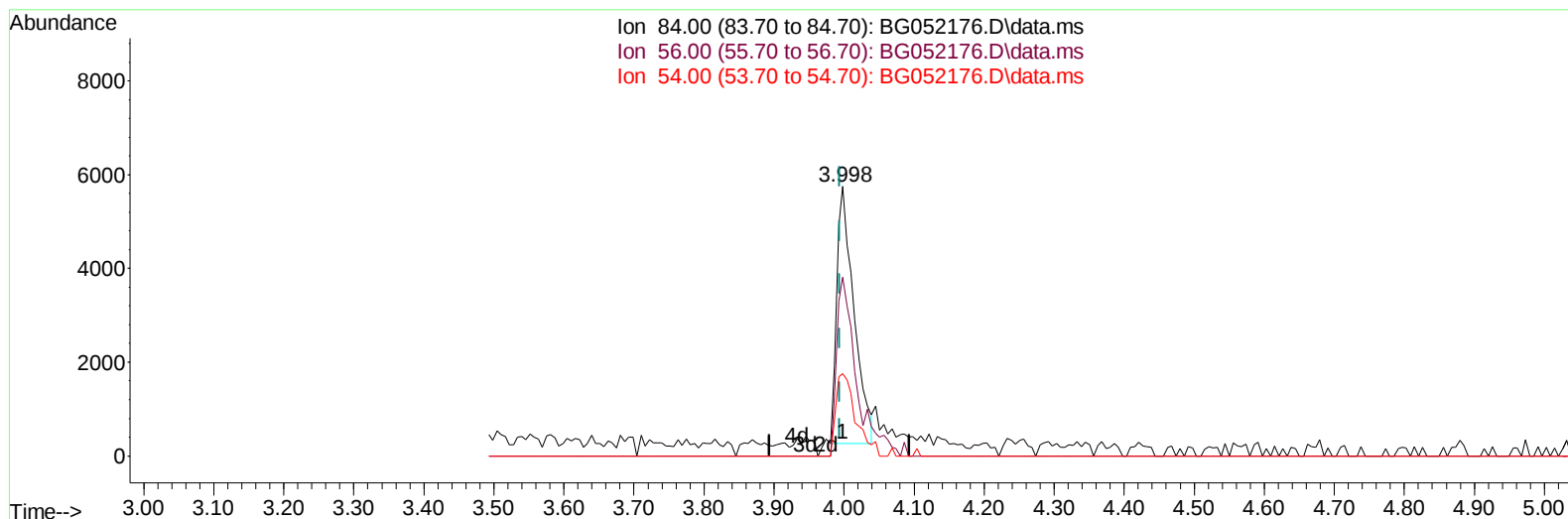
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052176.D
 Acq On : 27 Jan 2022 8:34
 Operator : CG/JU
 Sample : N1230-16
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q60

Manual IntegrationsAPPROVED

Quant Time: Jan 27 09:07:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/27/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052176.D\data.ms

(4) Pyridine-d5 (S)

3.998min (+ 0.004) 3.97 ng/u1

response 9406

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	66.37
54.00	31.50	30.48
0.00	0.00	0.00

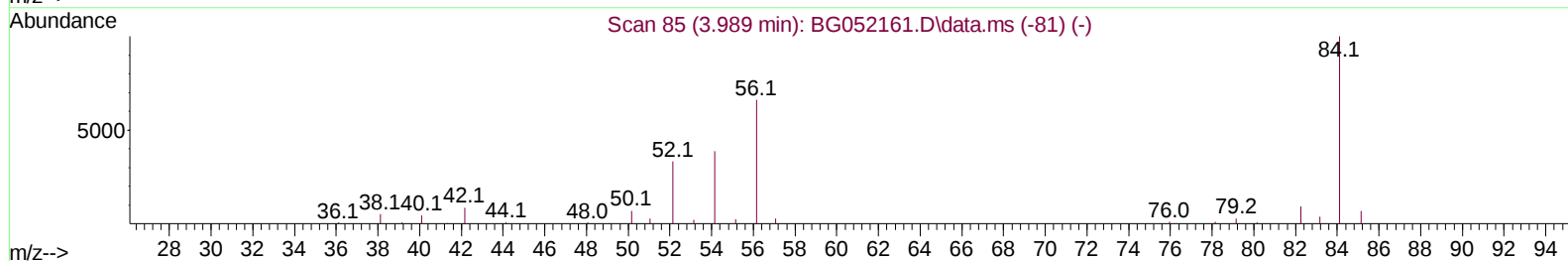
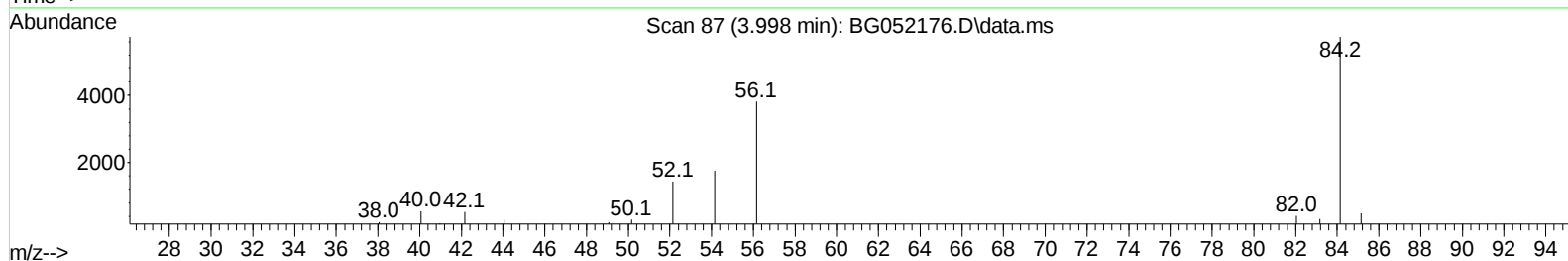
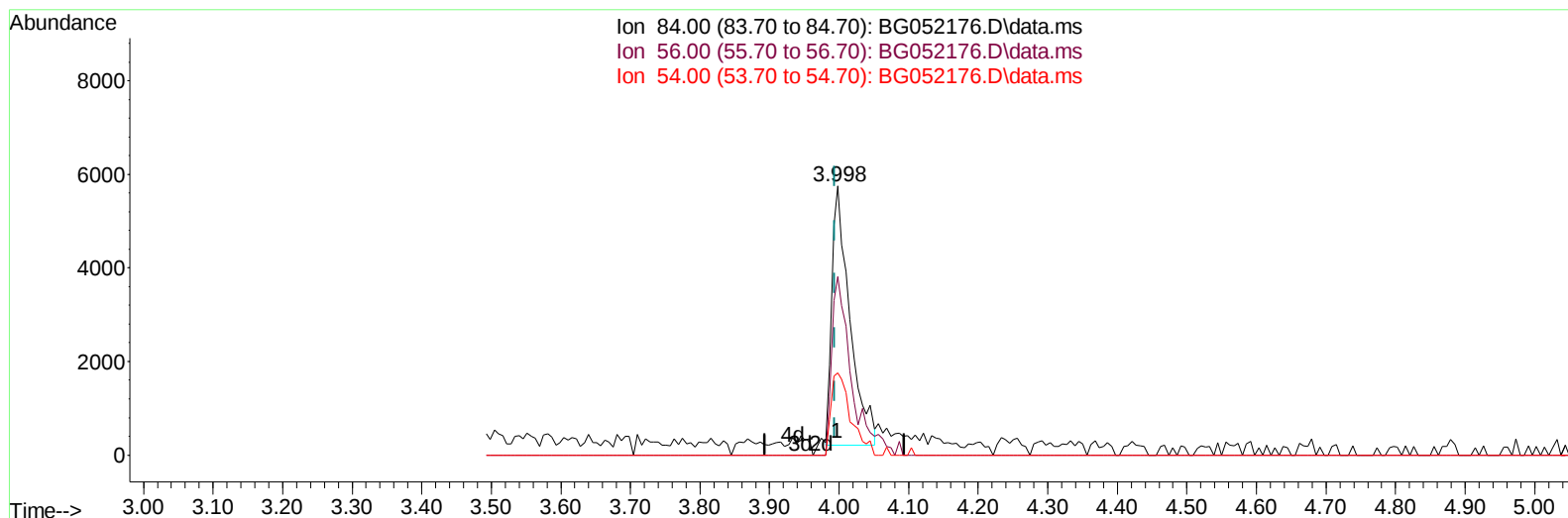
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TIC: BG052176.D\data.ms

(4) Pyridine-d5 (S)

3.998min (+ 0.004) 4.22 ng/ul m

response 9993

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	66.37
54.00	31.50	30.48
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.233	152		27243	20.000	ng/ul	0.00
20) Naphthalene-d8	11.077	136		114481	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.883	164		87731	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188		218112	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.944	240		222801	20.000	ng/ul	# 0.00
88) Perylene-d12	25.428	264		218805	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.563	96		3357	4.403	ng/uL	0.00
4) Pyridine-d5	3.998	84		9993m	4.216	ng/ul	0.00
7) Phenol-d5	7.382	99		17592	6.598	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.540	67		45622	26.953	ng/ul	0.00
11) 2-Chlorophenol-d4	7.763	132		40441	22.908	ng/ul	0.00
15) 4-Methylphenol-d8	8.938	113		30997	14.955	ng/ul	0.00
21) Nitrobenzene-d5	9.408	128		27880	30.820	ng/ul	0.00
24) 2-Nitrophenol-d4	10.143	143		30022	30.008	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.689	165		55432	28.346	ng/ul	0.00
31) 4-Chloroaniline-d4	11.200	131		57839	21.238	ng/ul	0.00
46) Dimethylphthalate-d6	14.266	166		225590	31.109	ng/ul	0.00
49) Acenaphthylene-d8	14.578	160		266798	30.458	ng/ul	0.00
54) 4-Nitrophenol-d4	15.065	143		7776	6.414	ng/ul	0.00
60) Fluorene-d10	15.870	176		199639	31.576	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200		32390	23.540	ng/ul	0.00
73) Anthracene-d10	17.726	188		366180	35.391	ng/ul	0.00
81) Pyrene-d10	20.006	212		445844	34.770	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.187	264		414313	35.968	ng/ul	0.00

Target Compounds Qvalue

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

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