

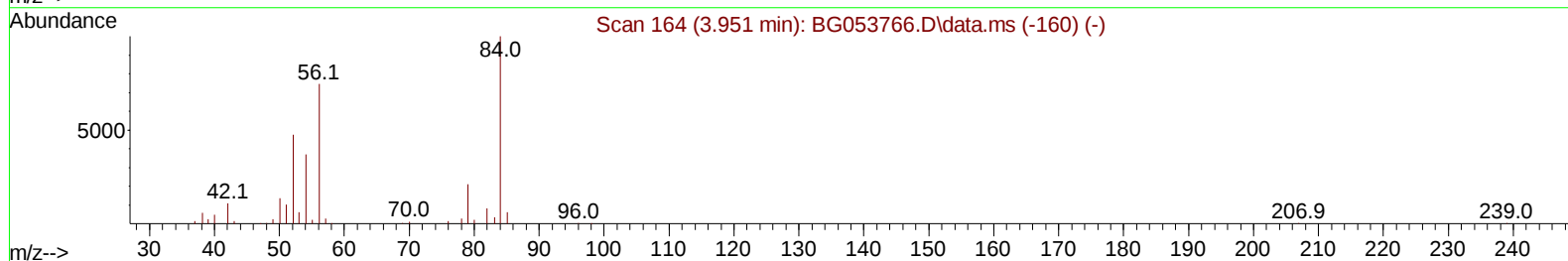
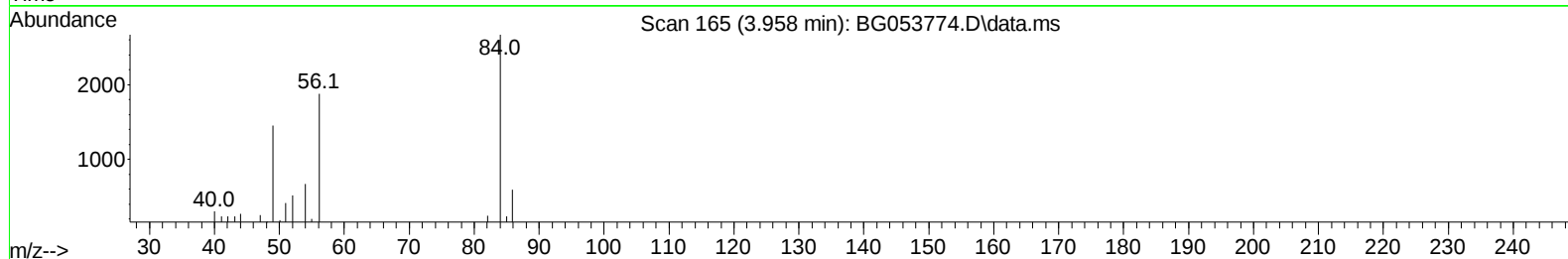
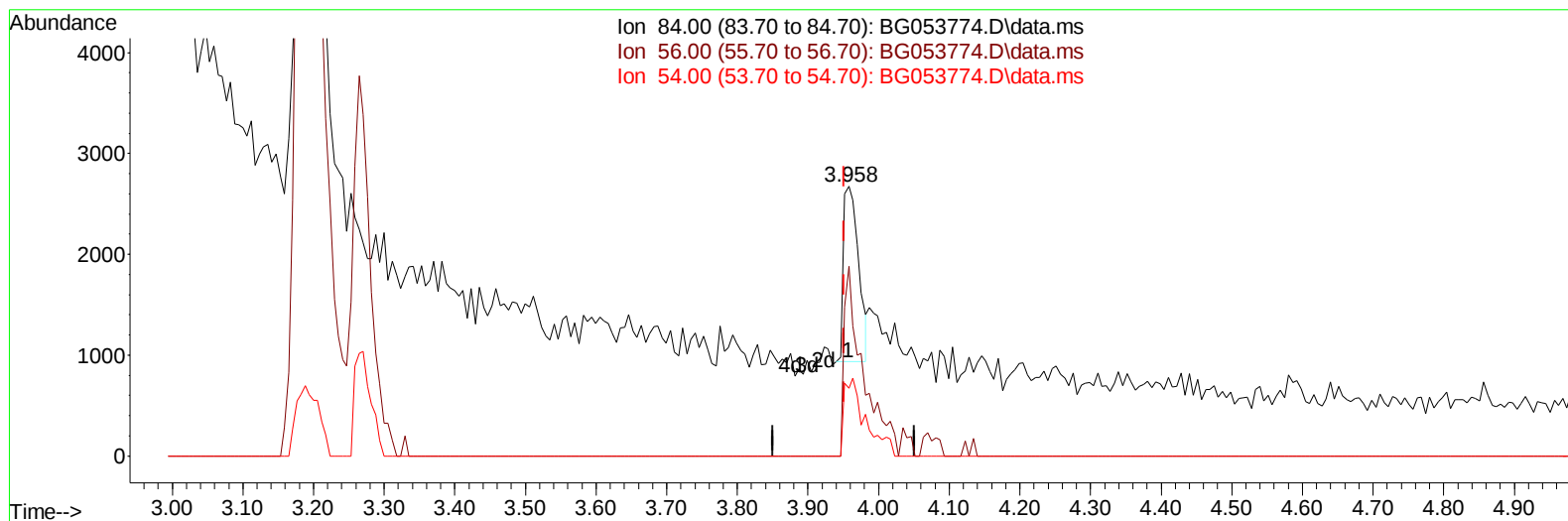
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
 Data File : BG053774.D
 Acq On : 1 Jun 2022 00:36
 Operator : CG/JU
 Sample : N3039-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0JQ9

Manual IntegrationsAPPROVED

Quant Time: Jun 01 02:20:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 01 00:25:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/01/2022
 Supervised By :mohammad ahmed 06/06/2022



TIC: BG053774.D\data.ms

(4) Pyridine-d5 (S)

3.958min (+ 0.007) 2.13 ng/u1

response 2594

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	71.00	70.47
54.00	35.20	25.19#
0.00	0.00	0.00

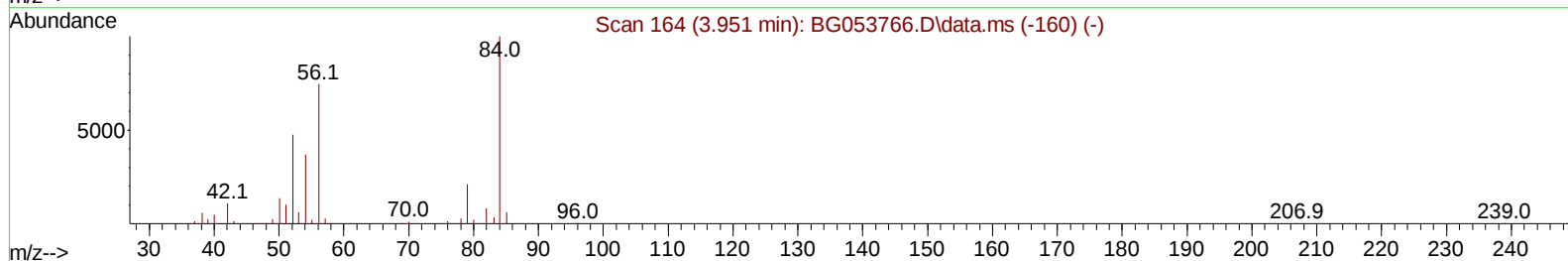
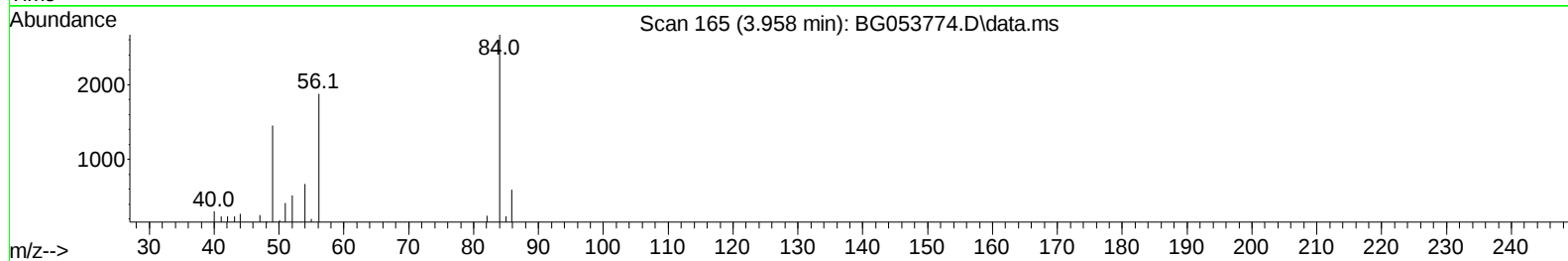
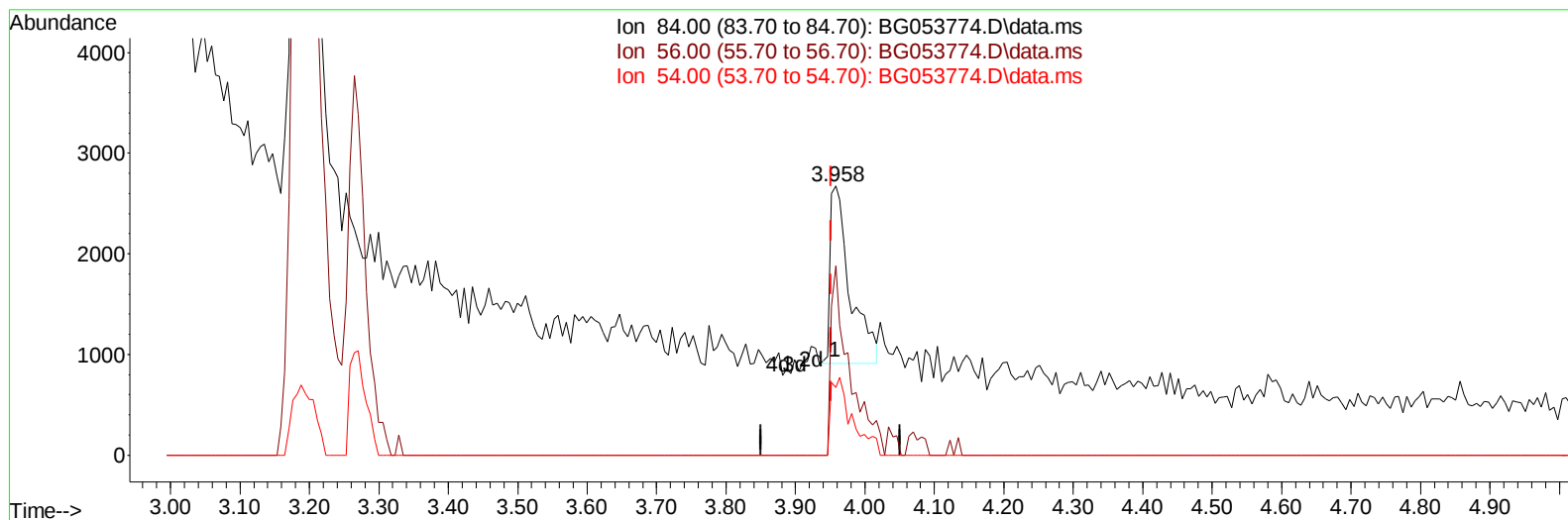
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TIC: BG053774.D\data.ms

(4) Pyridine-d5 (S)

3.958min (+ 0.007) 2.86 ng/u1 m

response 3476

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	71.00	70.47
54.00	35.20	25.19#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.100	152	20869	20.000	ng/ul	0.00
20) Naphthalene-d8	10.908	136	84196	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.722	164	57247	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.465	188	141116	20.000	ng/ul	0.00
79) Chrysene-d12	21.743	240	167614	20.000	ng/ul	0.00
88) Perylene-d12	25.015	264	167790	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	2401	4.352	ng/uL	0.00
4) Pyridine-d5	3.958	84	3476m	2.857	ng/ul	0.00
7) Phenol-d5	7.242	99	10889	6.861	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.418	67	32628	31.806	ng/ul	0.00
11) 2-Chlorophenol-d4	7.630	132	31054	24.609	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	20640	15.701	ng/ul	0.00
21) Nitrobenzene-d5	9.257	128	17694	35.156	ng/ul	0.00
24) 2-Nitrophenol-d4	9.986	143	16017	32.063	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.521	165	36847	27.159	ng/ul	0.00
31) 4-Chloroaniline-d4	11.032	131	40301	21.653	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	148339	33.721	ng/ul	0.00
49) Acenaphthylene-d8	14.416	160	158669	32.279	ng/ul	0.00
54) 4-Nitrophenol-d4	14.880	143	3614	5.975	ng/ul	-0.01
60) Fluorene-d10	15.709	176	131087	34.025	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.814	200	13945	24.716	ng/ul	0.00
73) Anthracene-d10	17.565	188	242289	38.013	ng/ul	0.00
81) Pyrene-d10	19.845	212	316893	35.521	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.792	264	313881	37.405	ng/ul	-0.01

Target Compounds Qvalue

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

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