

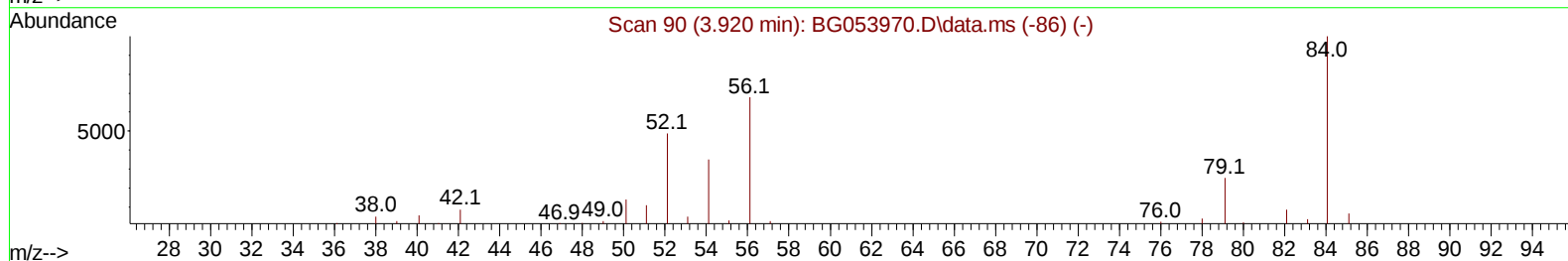
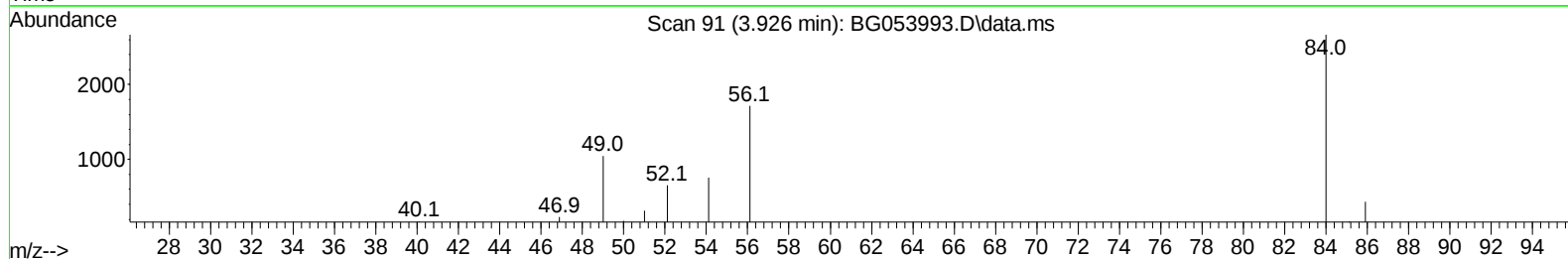
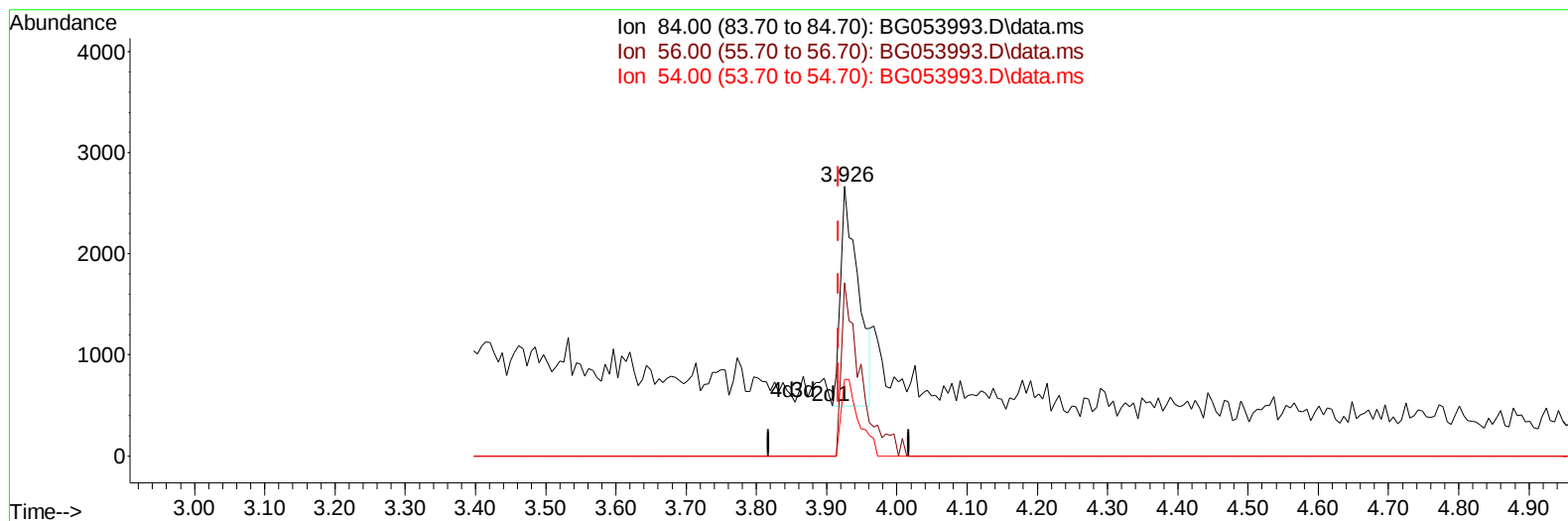
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062122\
 Data File : BG053993.D
 Acq On : 22 Jun 2022 1:12
 Operator : CG/JU
 Sample : N3355-17
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EY0D3

Manual IntegrationsAPPROVED

Quant Time: Jun 22 03:10:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 20 15:29:13 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/22/2022
 Supervised By :mohammad ahmed 06/28/2022



TIC: BG053993.D\data.ms

(4) Pyridine-d5 (S)

3.926min (+ 0.008) 2.99 ng/u1

response 3820

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	71.00	64.22
54.00	35.20	28.51
0.00	0.00	0.00

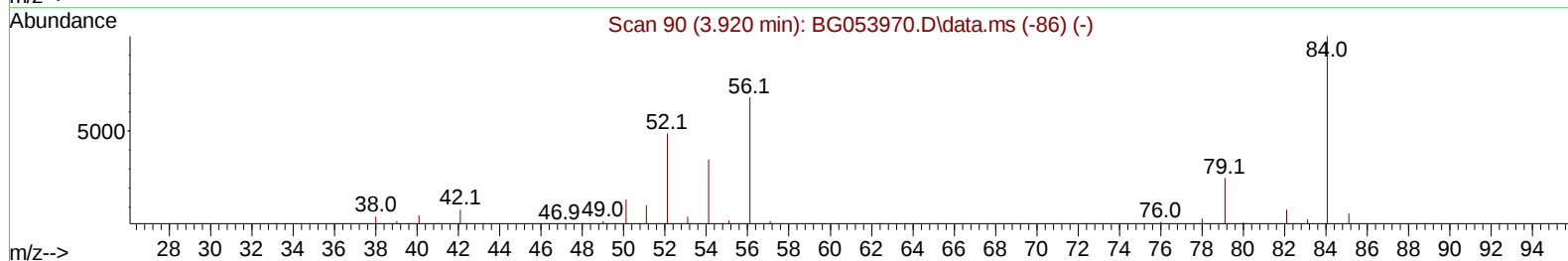
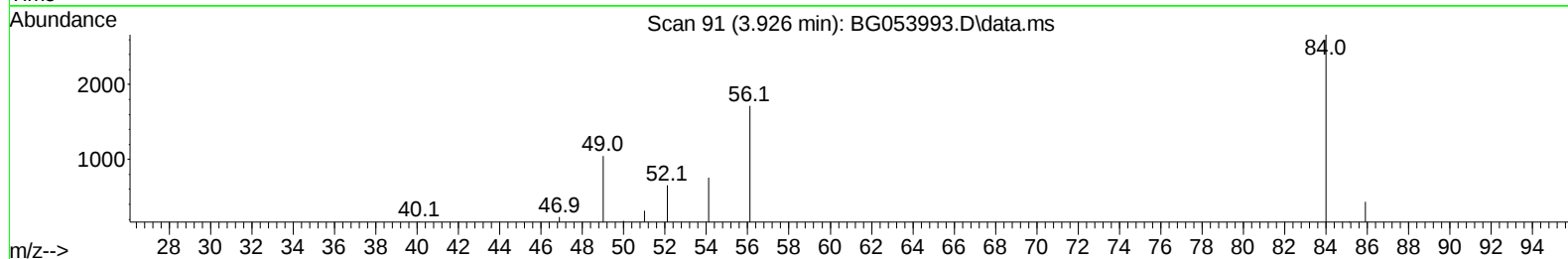
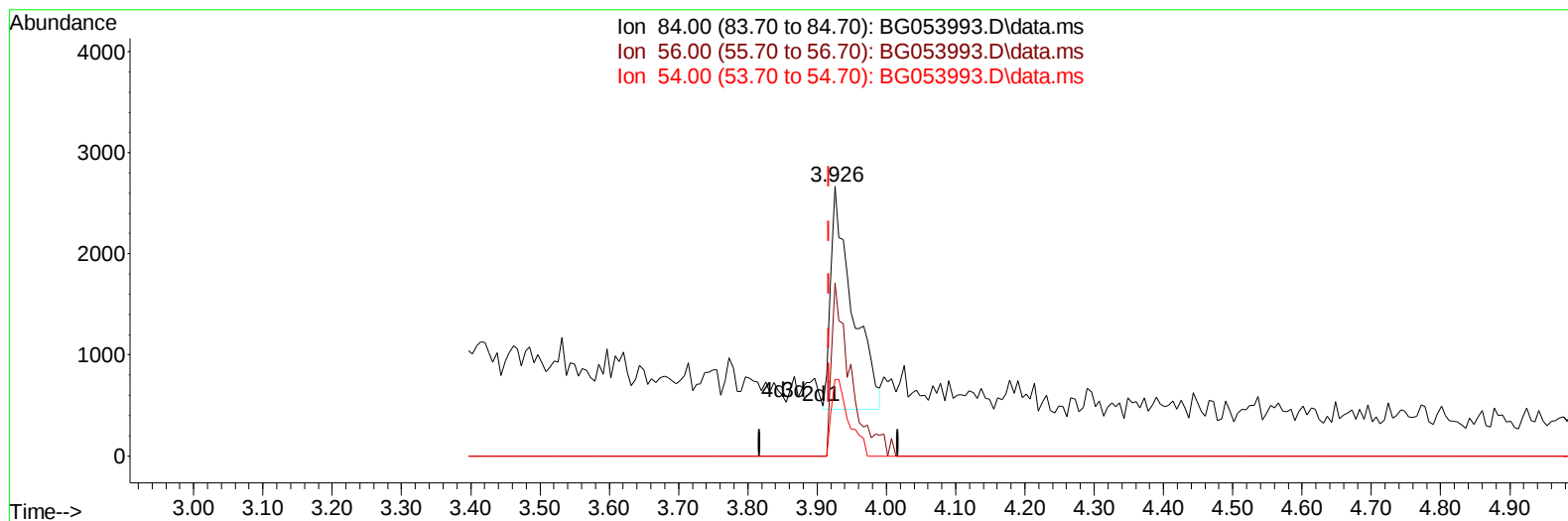
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TIC: BG053993.D\data.ms

(4) Pyridine-d5 (S)

3.926min (+ 0.008) 3.74 ng/u1 m

response 4772

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	71.00	64.22
54.00	35.20	28.51
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	8.080	152	22330	20.000	ng/ul	0.00
20) Naphthalene-d8	10.888	136	88244	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.701	164	67108	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	182977	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.717	240	219407	20.000	ng/ul	# 0.00
88) Perylene-d12	24.983	264	219842	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.503	96	2946	6.277	ng/uL	0.00
4) Pyridine-d5	3.926	84	4772m	3.738	ng/ul	0.00
7) Phenol-d5	7.234	99	11628	7.269	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.398	67	30695	31.808	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	32240	25.272	ng/ul	0.00
15) 4-Methylphenol-d8	8.779	113	22662	16.614	ng/ul	0.00
21) Nitrobenzene-d5	9.237	128	20398	31.142	ng/ul	0.00
24) 2-Nitrophenol-d4	9.966	143	23511	33.783	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.500	165	45374	29.469	ng/ul	0.00
31) 4-Chloroaniline-d4	11.017	131	34033	17.496	ng/ul	0.00
46) Dimethylphthalate-d6	14.102	166	176568	33.445	ng/ul	0.00
49) Acenaphthylene-d8	14.396	160	185247	32.269	ng/ul	0.00
54) 4-Nitrophenol-d4	14.883	143	3876	5.362	ng/ul	0.00
60) Fluorene-d10	15.694	176	158982	33.526	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.800	200	29083	29.137	ng/ul	0.00
73) Anthracene-d10	17.545	188	281944	34.119	ng/ul	0.00
81) Pyrene-d10	19.825	212	381035	33.308	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.760	264	387680	34.935	ng/ul	0.00

Target Compounds Qvalue

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

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