

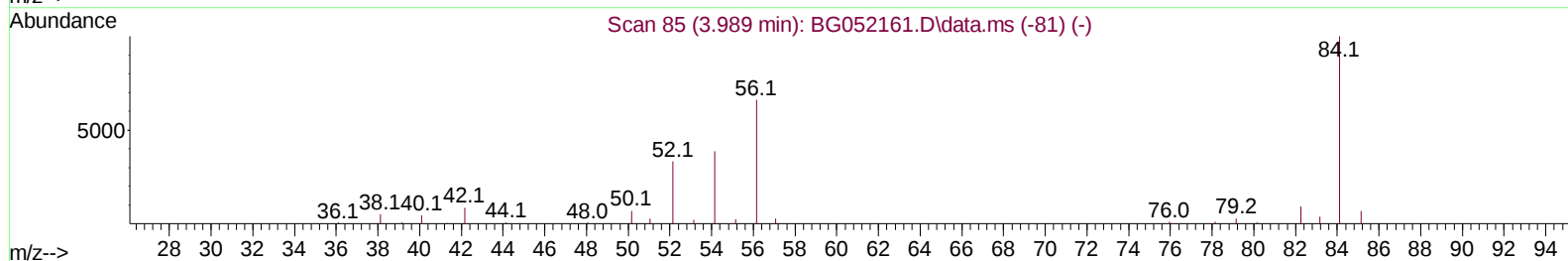
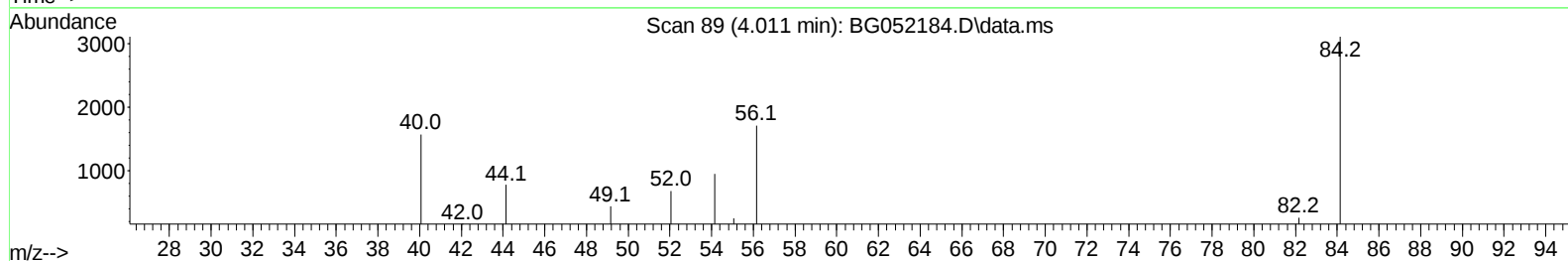
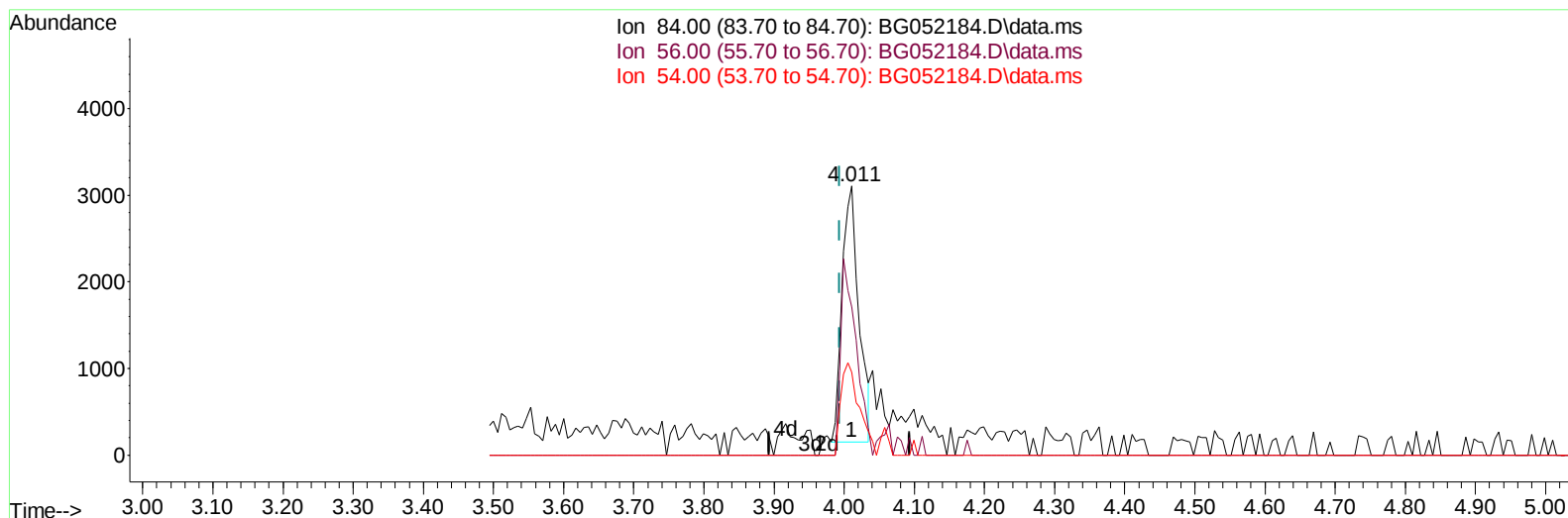
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052184.D
 Acq On : 27 Jan 2022 14:12
 Operator : CG/JU
 Sample : N1279-04
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COVE1

Manual IntegrationsAPPROVED

Quant Time: Jan 27 23:59:01 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 01/28/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052184.D\data.ms

(4) Pyridine-d5 (S)

4.011min (+ 0.017) 2.56 ng/u1

response 4913

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	55.15
54.00	31.50	30.75
0.00	0.00	0.00

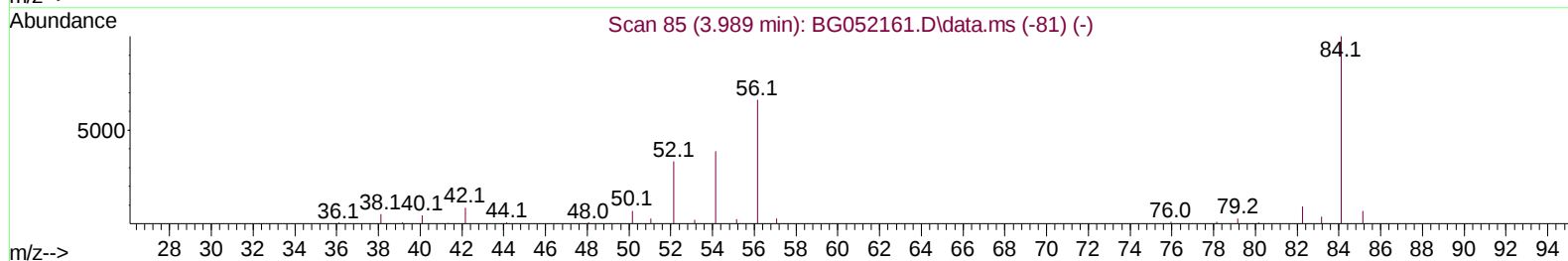
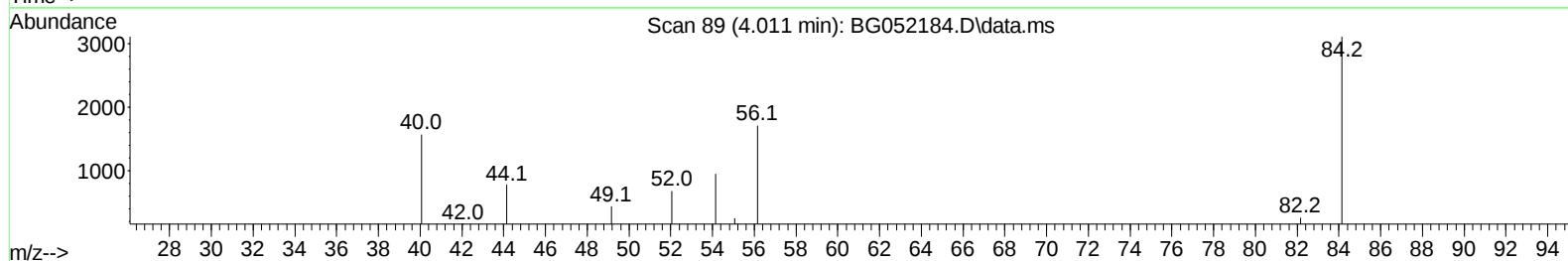
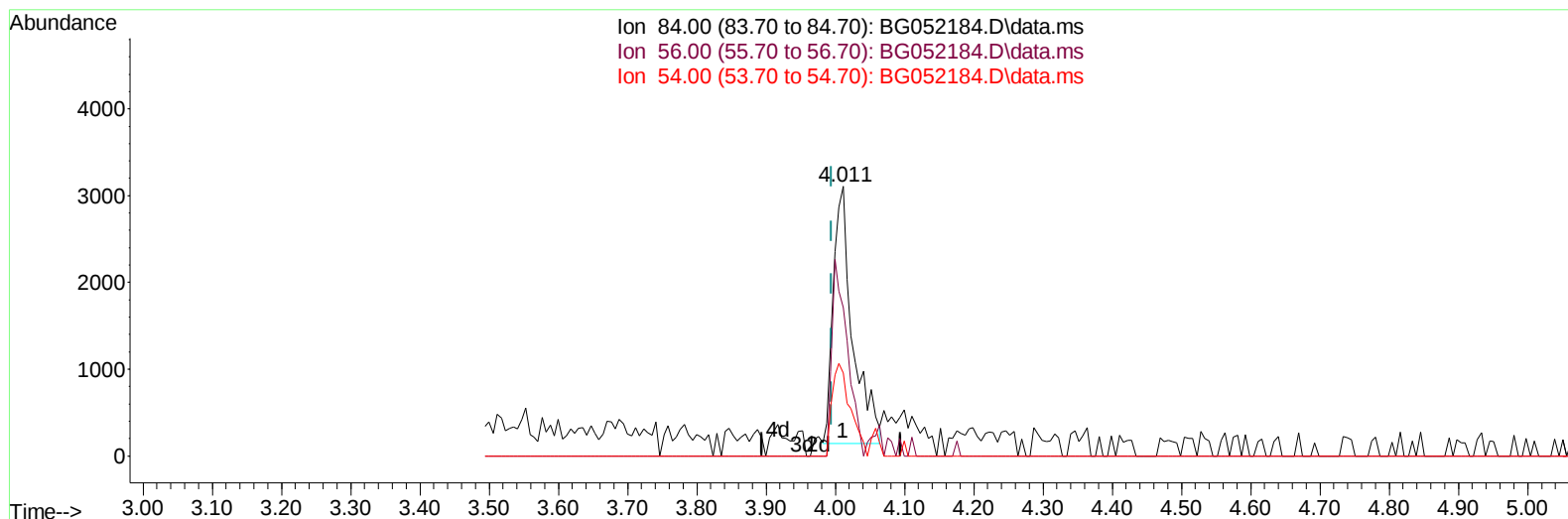
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(4) Pyridine-d5 (S)

4.011min (+ 0.017) 3.01 ng/u1 m

response 5777

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	55.15
54.00	31.50	30.75
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.235	152	22026	20.000	ng/ul	0.00
20) Naphthalene-d8	11.072	136	94197	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.879	164	72612	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.628	188	177429	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.945	240	186110	20.000	ng/ul	# 0.00
88) Perylene-d12	25.423	264	182581	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.559	96	2601	4.220	ng/uL	-0.01
4) Pyridine-d5	4.011	84	5777m	3.015	ng/ul	0.02
7) Phenol-d5	7.377	99	12903	5.985	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	40264	29.422	ng/ul	0.00
11) 2-Chlorophenol-d4	7.765	132	35466	24.848	ng/ul	0.00
15) 4-Methylphenol-d8	8.940	113	24914	14.867	ng/ul	0.00
21) Nitrobenzene-d5	9.404	128	24362	32.730	ng/ul	0.00
24) 2-Nitrophenol-d4	10.138	143	25860	31.414	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.690	165	46888	29.140	ng/ul	0.00
31) 4-Chloroaniline-d4	11.201	131	53252	23.765	ng/ul	0.00
46) Dimethylphthalate-d6	14.268	166	199860	33.299	ng/ul	0.00
49) Acenaphthylene-d8	14.579	160	237260	32.725	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	4336	4.321	ng/ul	0.00
60) Fluorene-d10	15.871	176	178342	34.080	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.977	200	24967	22.306	ng/ul	0.00
73) Anthracene-d10	17.728	188	323222	38.402	ng/ul	0.00
81) Pyrene-d10	20.007	212	394056	36.790	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.182	264	365739	38.051	ng/ul	0.00

Target Compounds Qvalue

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

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