

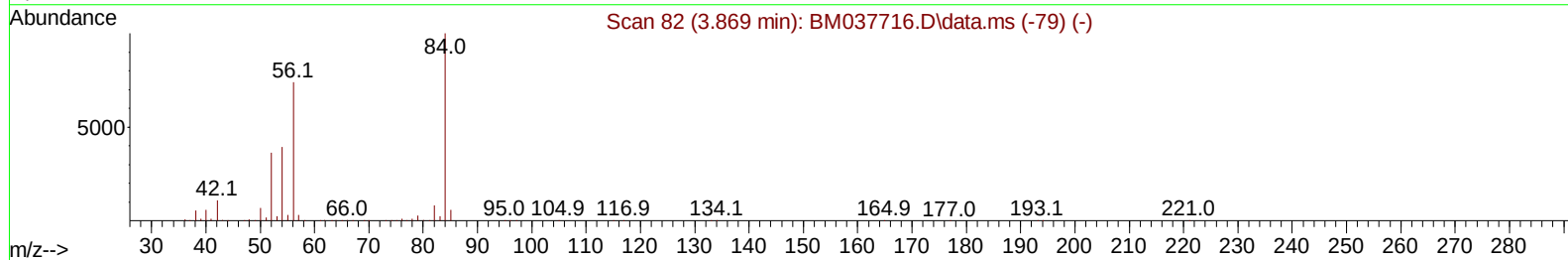
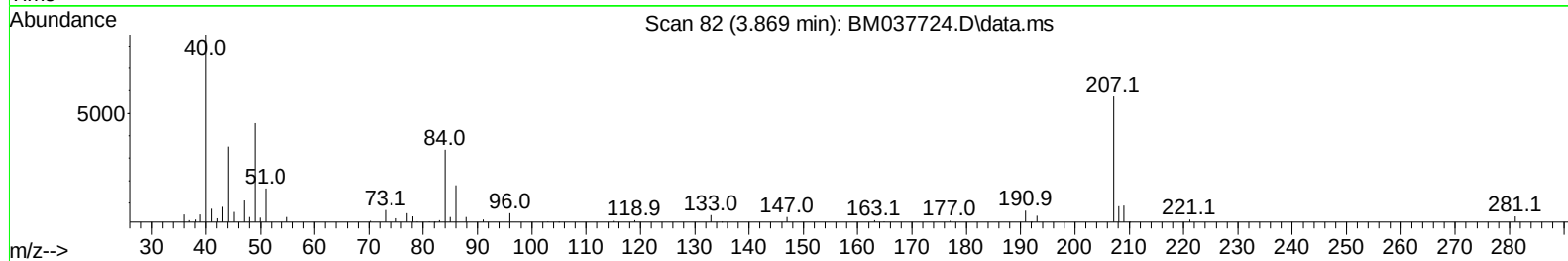
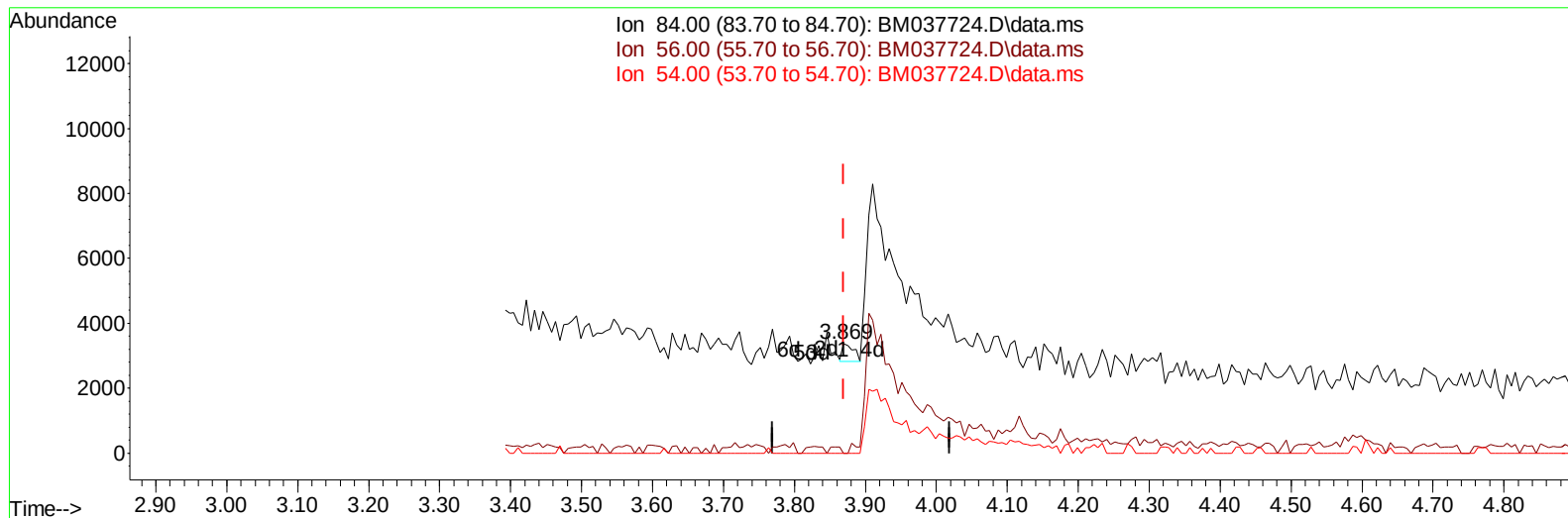
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037724.D
Acq On : 26 Nov 2022 09:22
Operator : CG/JU
Sample : N5591-17
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C9829

Manual IntegrationsAPPROVED

Quant Time: Nov 27 21:26:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 24 01:27:15 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/28/2022
Supervised By :mohammad ahmed 11/28/2022



TIC: BM037724.D\data.ms

(4) Pyridine-d5 (S)

3.869min (0.000) 0.04 ng/ul

response 603

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	74.90	0.00#
54.00	37.60	0.00#
0.00	0.00	0.00

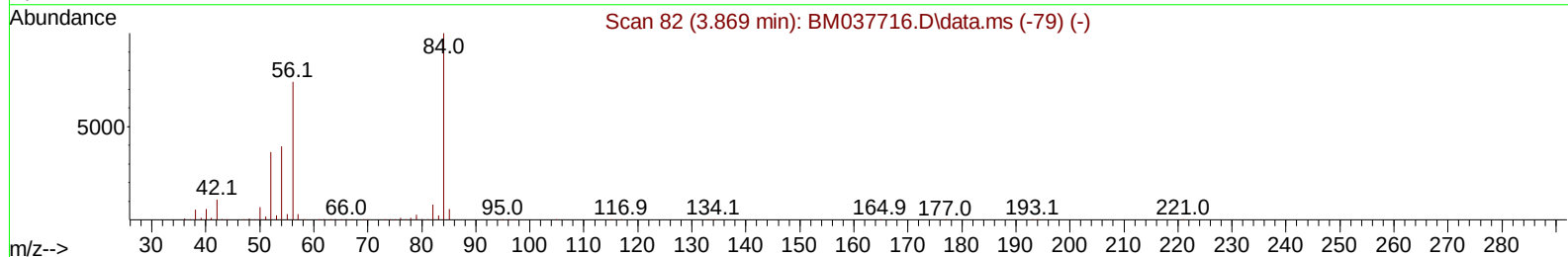
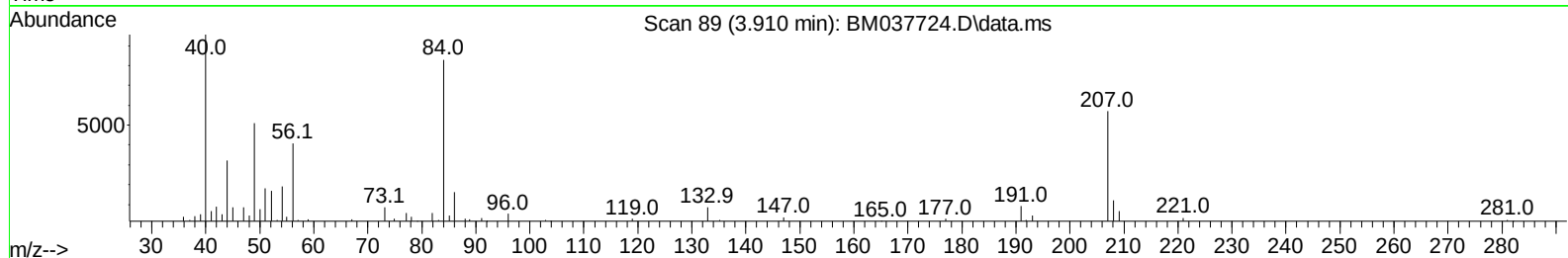
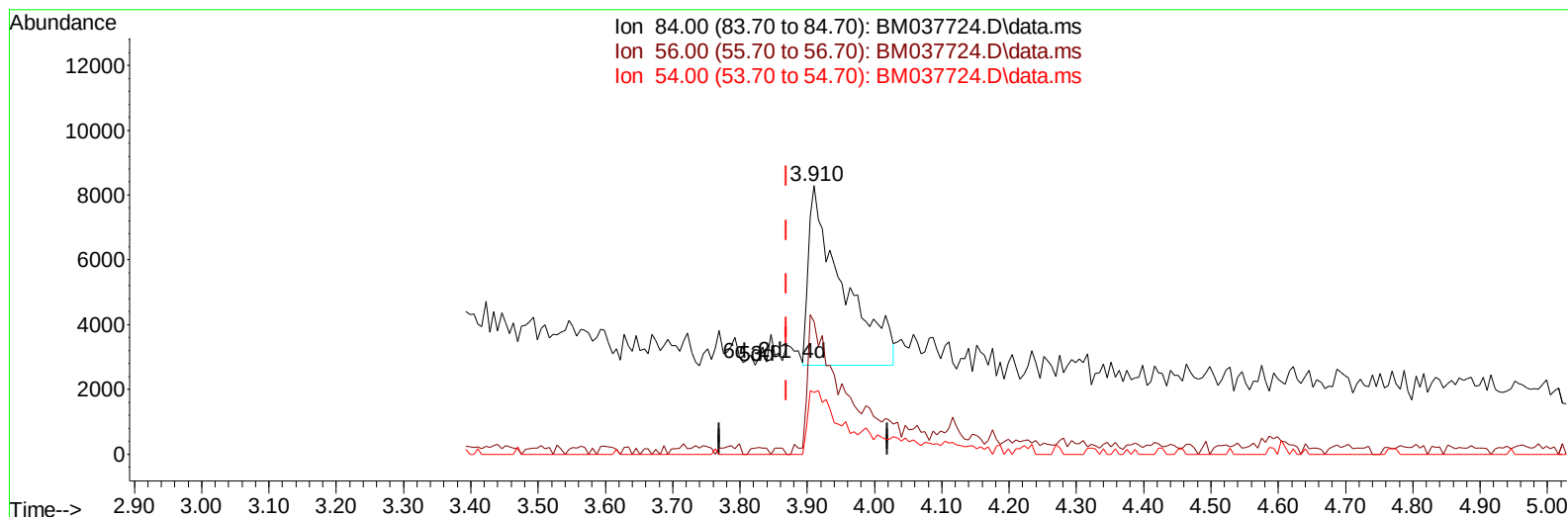
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(4) Pyridine-d5 (S)

3.910min (+ 0.041) 1.21 ng/ul m

response 19756

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	74.90	49.29#
54.00	37.60	22.95#
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	8.110	152	208793	20.000	ng/ul	0.00
20) Naphthalene-d8	10.934	136	991068	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	650375	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.474	188	1359106	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	1257983	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1167887	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	19834	3.940	ng/uL	0.00
4) Pyridine-d5	3.910	84	19756m	1.206	ng/ul	0.04
7) Phenol-d5	7.257	99	400236	19.214	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	282597	20.337	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	298022	20.590	ng/ul	0.00
15) 4-Methylphenol-d8	8.816	113	317092	19.130	ng/ul	0.00
21) Nitrobenzene-d5	9.281	128	164657	21.193	ng/ul	0.00
24) 2-Nitrophenol-d4	10.010	143	160473	20.961	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.545	165	281487	20.092	ng/ul	0.00
31) 4-Chloroaniline-d4	11.069	131	396566	17.058	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	1025805	22.267	ng/ul	0.00
49) Acenaphthylene-d8	14.433	160	1145146	21.060	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	171936	16.944	ng/ul	-0.01
60) Fluorene-d10	15.721	176	890676	22.076	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	106537	14.572	ng/ul	-0.01
73) Anthracene-d10	17.568	188	1399964	22.027	ng/ul	0.00
81) Pyrene-d10	19.851	212	1639286	25.643	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.962	264	1398808	23.269	ng/ul	0.00

Target Compounds Qvalue

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

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