

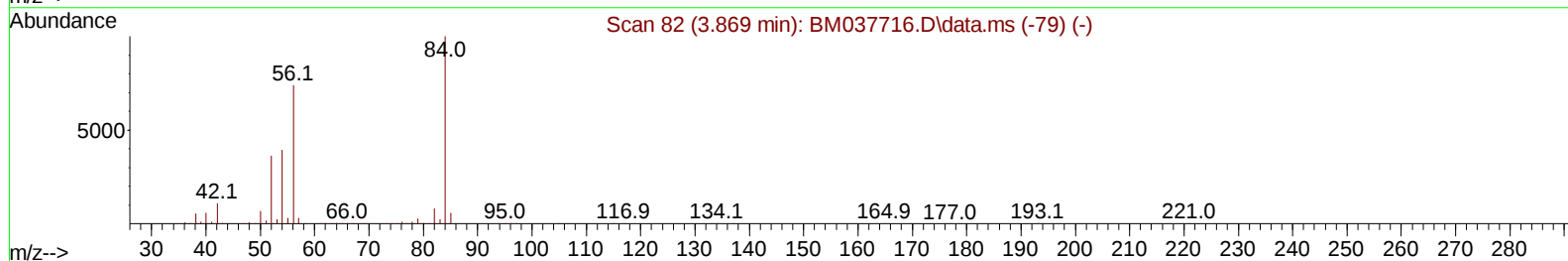
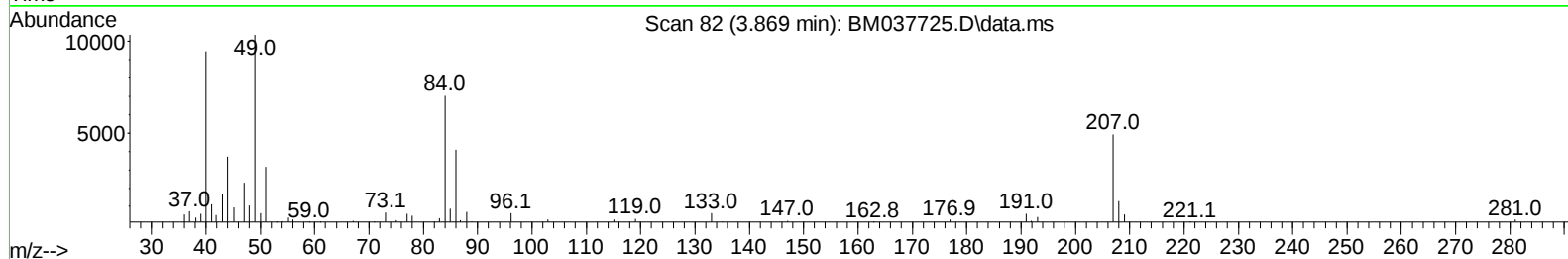
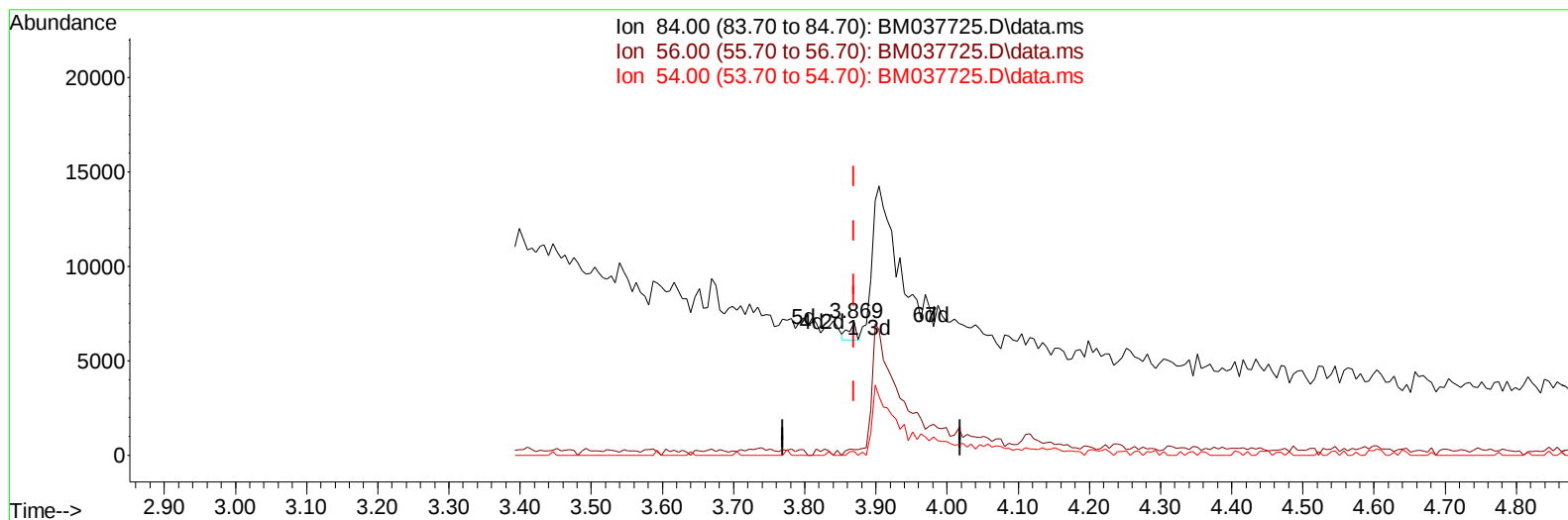
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037725.D
Acq On : 26 Nov 2022 09:59
Operator : CG/JU
Sample : N5591-15
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C9827

Manual IntegrationsAPPROVED

Quant Time: Nov 27 21:26:43 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: BM037725.D\data.ms

(4) Pyridine-d5 (S)

3.869min (0.000) 0.04 ng/u1

response 659

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	74.90	4.35#
54.00	37.60	2.69#
0.00	0.00	0.00

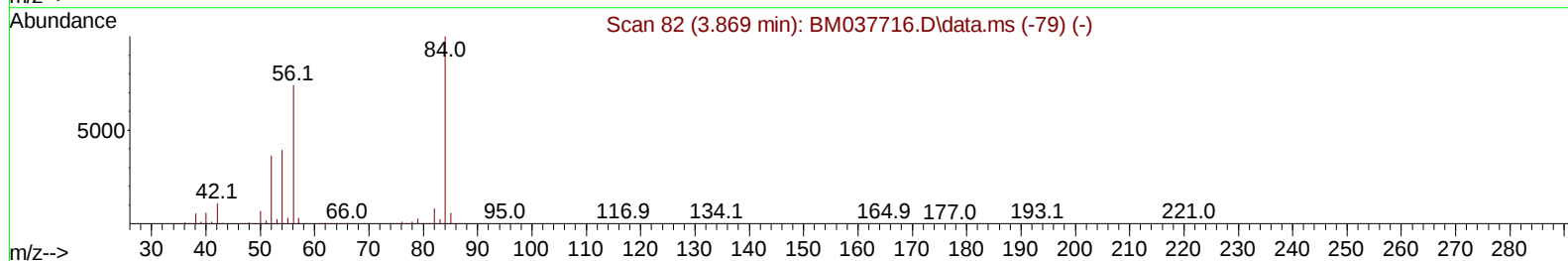
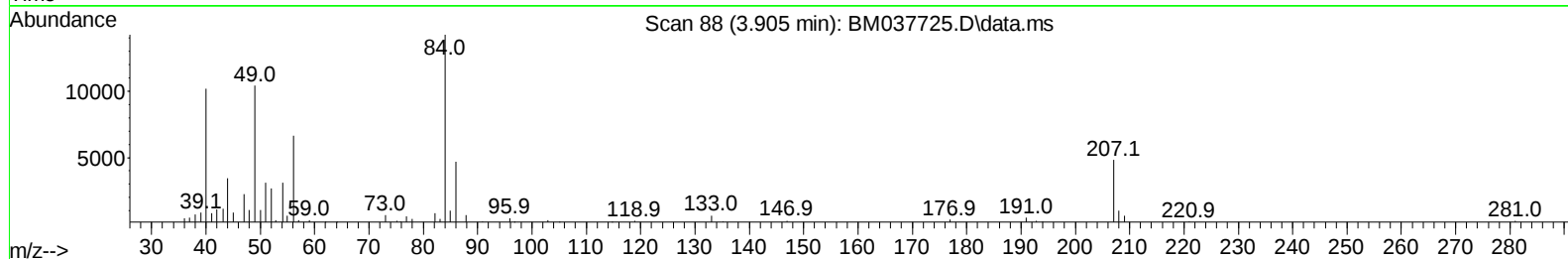
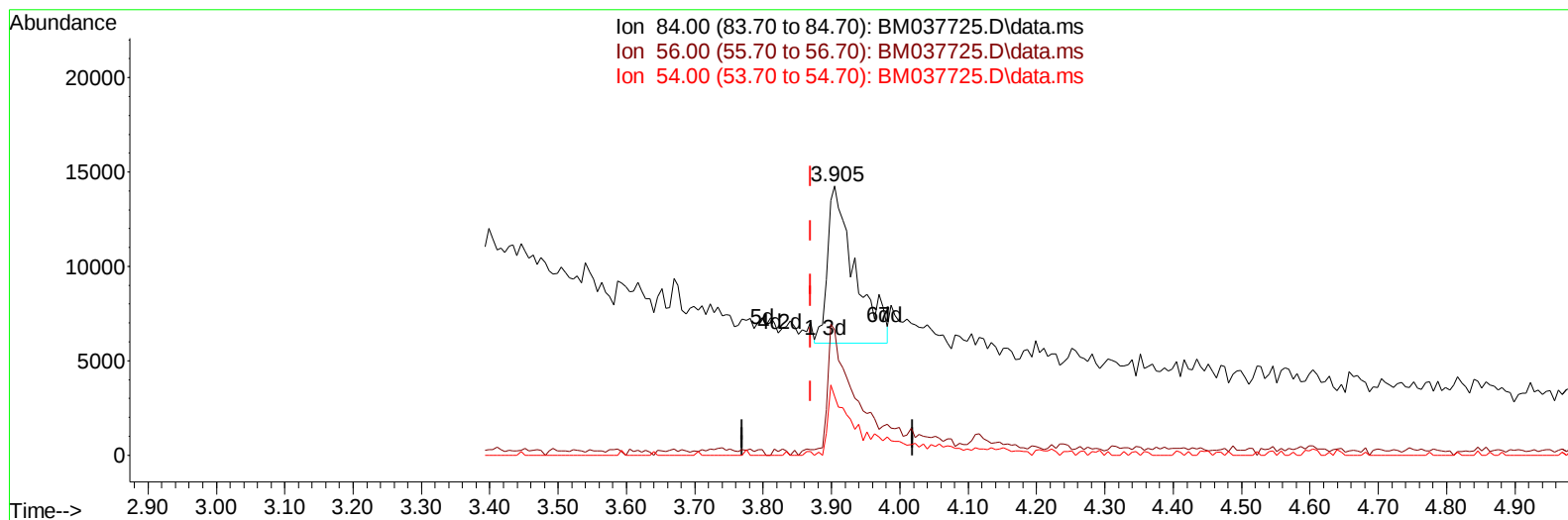
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(4) Pyridine-d5 (S)

3.905min (+ 0.035) 1.36 ng/ul m

response 22976

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	74.90	46.59#
54.00	37.60	21.72#
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	214653	20.000	ng/ul	0.00
20) Naphthalene-d8	10.933	136	1033359	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.739	164	648315	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.474	188	1313595	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	1235107	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1164347	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	24590	4.751	ng/uL	0.00
4) Pyridine-d5	3.905	84	22976m	1.364	ng/ul	0.04
7) Phenol-d5	7.257	99	513644	23.985	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	363022	25.412	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	381302	25.624	ng/ul	0.00
15) 4-Methylphenol-d8	8.816	113	377701	22.164	ng/ul	0.00
21) Nitrobenzene-d5	9.281	128	207879	25.661	ng/ul	0.00
24) 2-Nitrophenol-d4	10.010	143	200640	25.135	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.545	165	345313	23.639	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	488248	20.142	ng/ul	-0.01
46) Dimethylphthalate-d6	14.139	166	1228479	26.752	ng/ul	0.00
49) Acenaphthylene-d8	14.433	160	1384285	25.539	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	226359	22.378	ng/ul	0.00
60) Fluorene-d10	15.721	176	1082144	26.907	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	139100	19.685	ng/ul	-0.01
73) Anthracene-d10	17.568	188	1627479	26.494	ng/ul	0.00
81) Pyrene-d10	19.851	212	1911186	30.450	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.962	264	1653546	27.590	ng/ul	0.00

Target Compounds Qvalue

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

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