

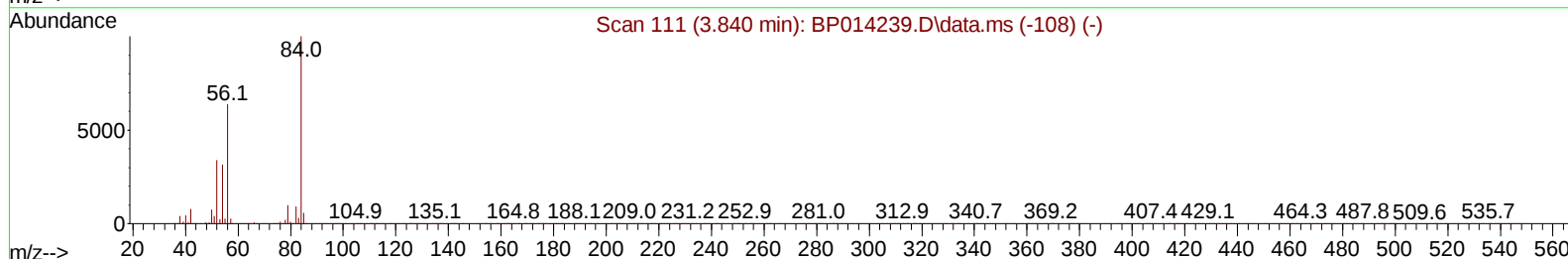
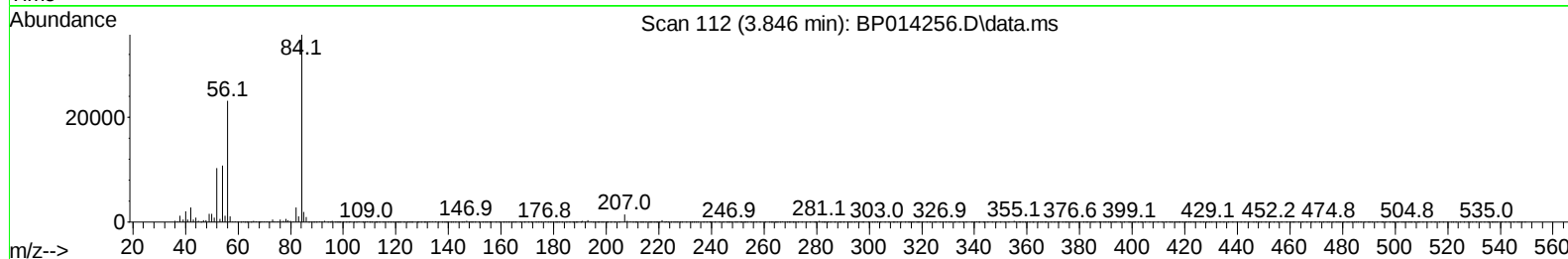
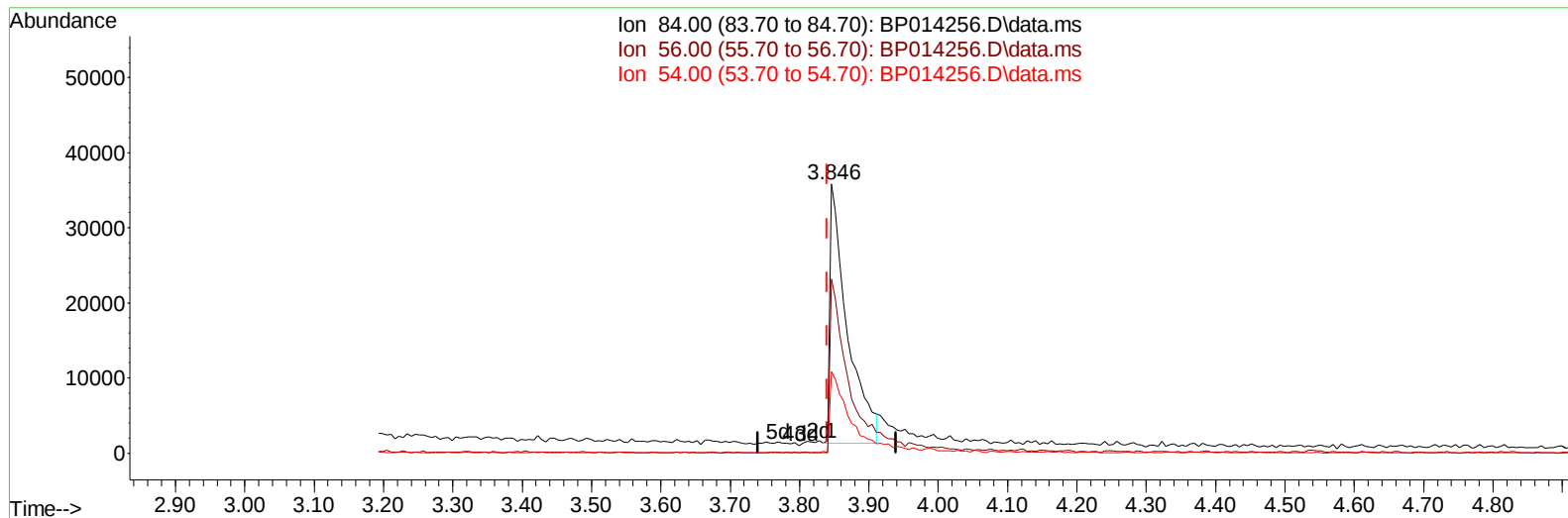
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014256.D
Acq On : 23 Mar 2023 14:08
Operator : CG/JU
Sample : 02010-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB952

Manual IntegrationsAPPROVED

Quant Time: Mar 24 02:02:35 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 23 04:56:06 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/24/2023
Supervised By :Jagrut Upadhyay 03/24/2023



TIC: BP014256.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 5.33 ng/u1

response 60210

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	64.72
54.00	33.20	30.29
0.00	0.00	0.00

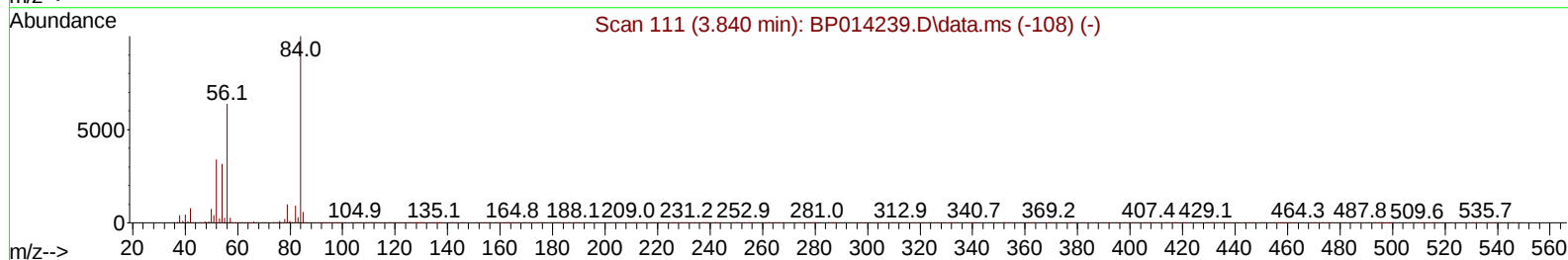
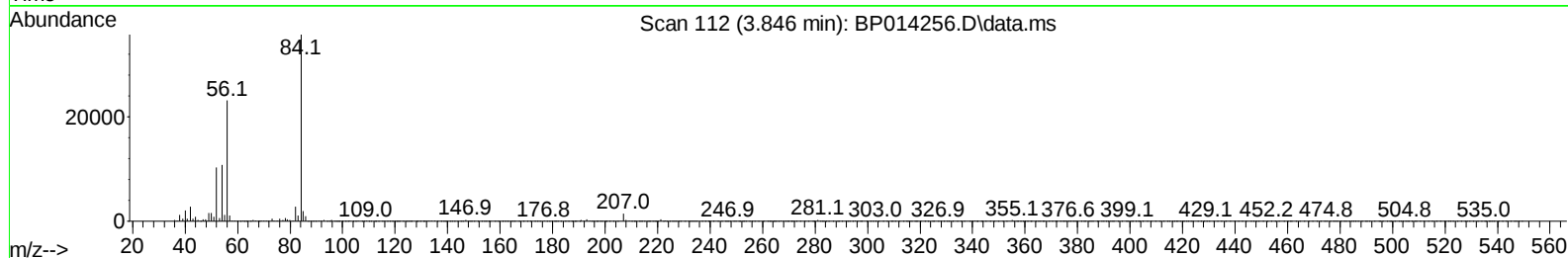
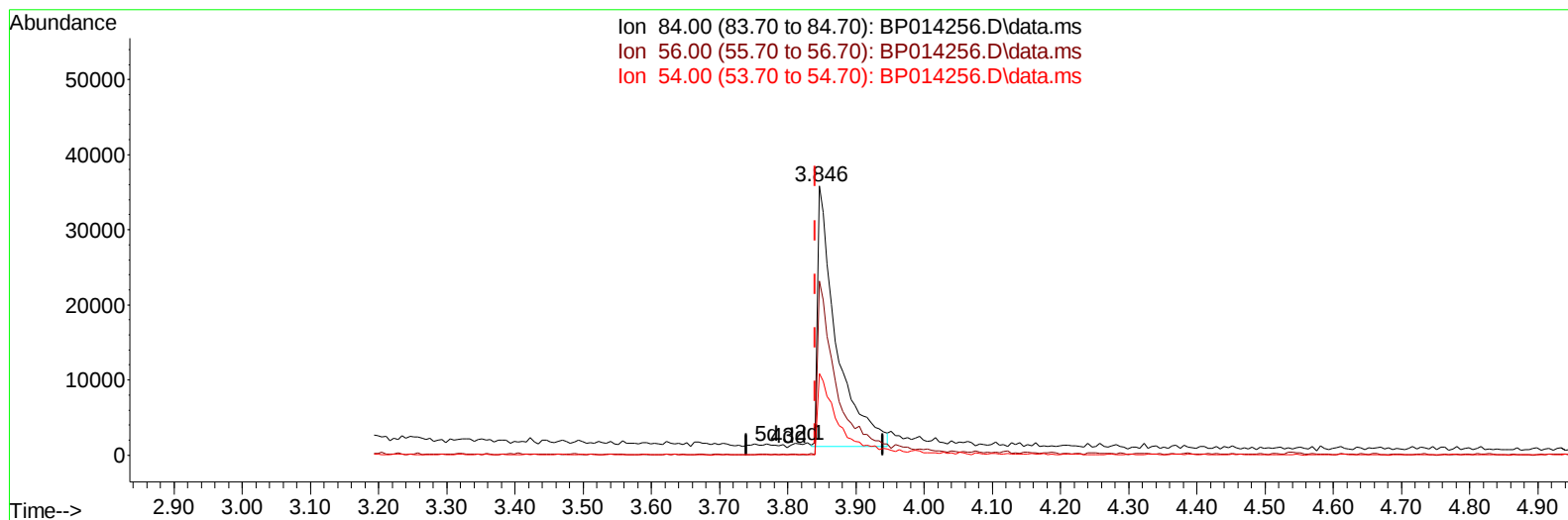
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TIC: BP014256.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 5.89 ng/u1 m

response 66569

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	64.72
54.00	33.20	30.29
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	8.034	152	154156	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	708031	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.681	164	515950	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1226614	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.521	240	1219419	20.000	ng/ul	# 0.00
88) Perylene-d12	24.045	264	1217584	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	15530	3.806	ng/uL	0.00
4) Pyridine-d5	3.846	84	66569m	5.893	ng/ul	0.00
7) Phenol-d5	7.199	99	76832	5.648	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	218012	27.479	ng/ul	0.00
11) 2-Chlorophenol-d4	7.563	132	201522	19.443	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	154380	13.893	ng/ul	0.00
21) Nitrobenzene-d5	9.210	128	133849	23.693	ng/ul	0.00
24) 2-Nitrophenol-d4	9.940	143	154309	22.818	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	273764	21.515	ng/ul	0.00
31) 4-Chloroaniline-d4	11.004	131	285367	16.947	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	1135195	25.925	ng/ul	0.00
49) Acenaphthylene-d8	14.375	160	1179268	25.423	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	27215	3.527	ng/ul	0.02
60) Fluorene-d10	15.675	176	991042	26.706	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	167597	18.078	ng/ul	0.00
73) Anthracene-d10	17.539	188	1639007	27.525	ng/ul	0.00
81) Pyrene-d10	19.763	212	1991098	29.605	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1809675	27.942	ng/ul	0.00

Target Compounds Qvalue

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

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