

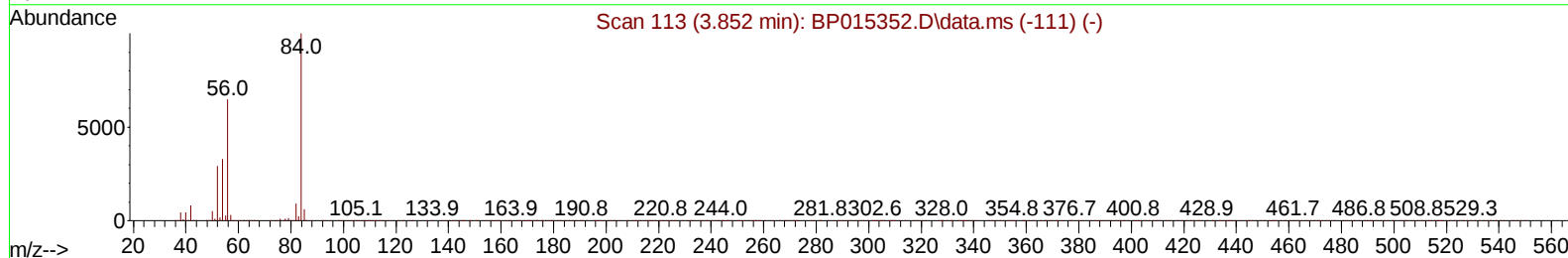
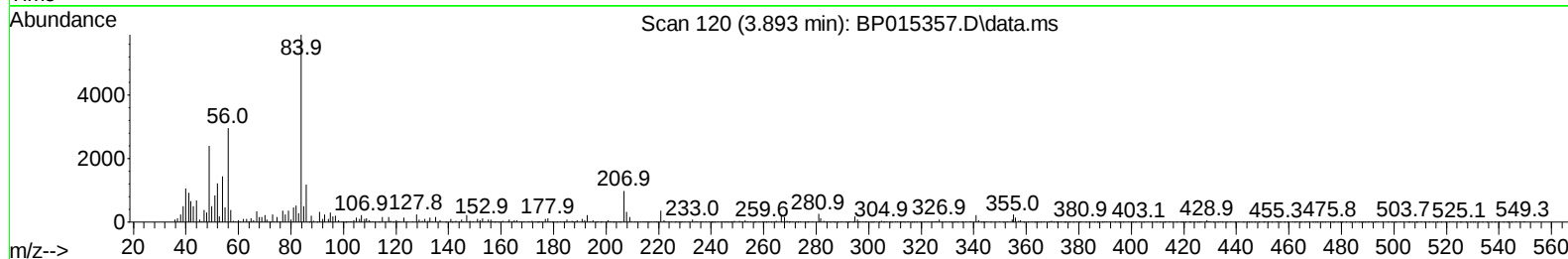
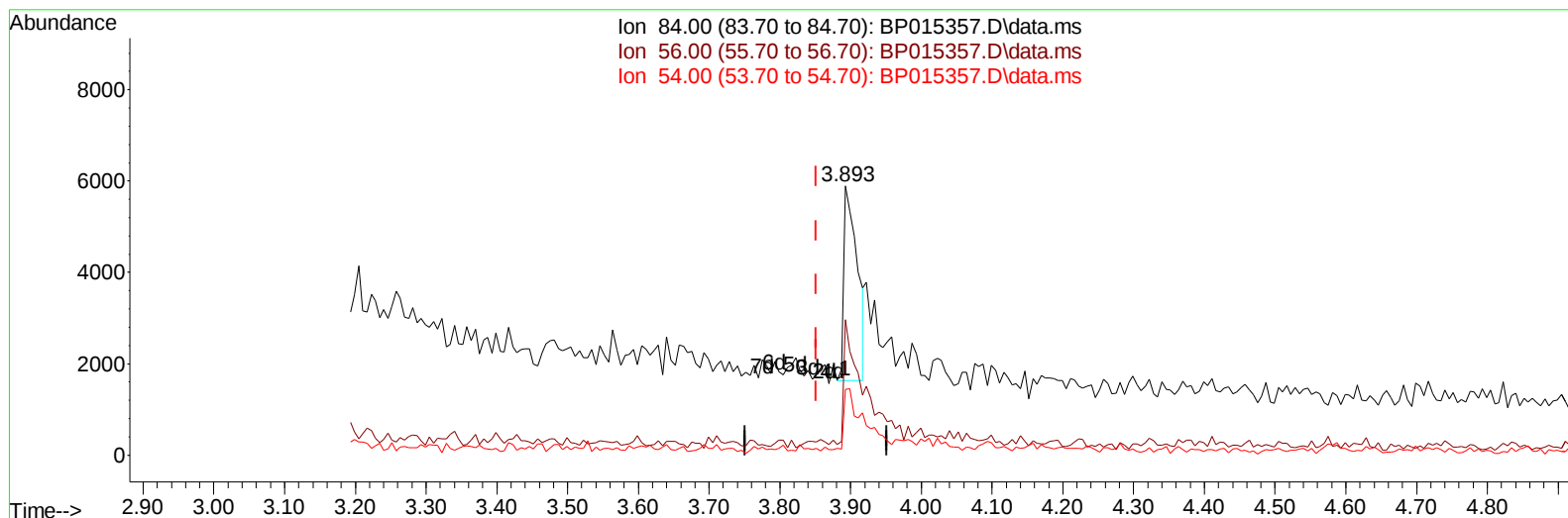
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
 Data File : BP015357.D
 Acq On : 31 May 2023 13:45
 Operator : CG/JU
 Sample : 02851-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0FK4

Manual IntegrationsAPPROVED

Quant Time: May 31 22:42:15 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 31 14:24:45 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/01/2023
 Supervised By :mohammad ahmed 06/01/2023



TIC: BP015357.D\data.ms

(4) Pyridine-d5 (S)

3.893min (+ 0.041) 0.55 ng/u1

response	5565	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	50.31
54.00	29.10	24.54
0.00	0.00	0.00

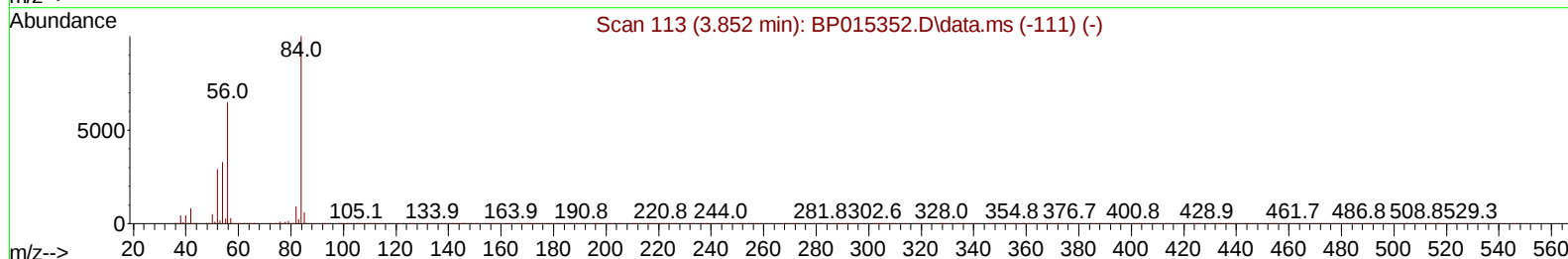
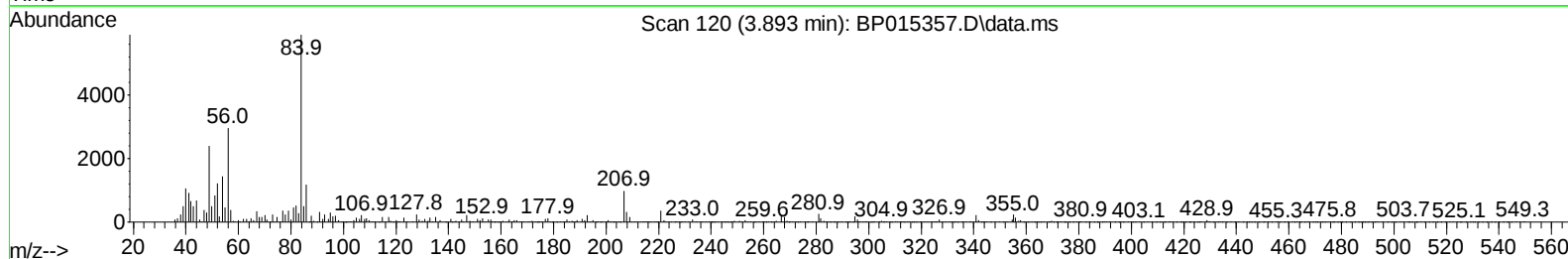
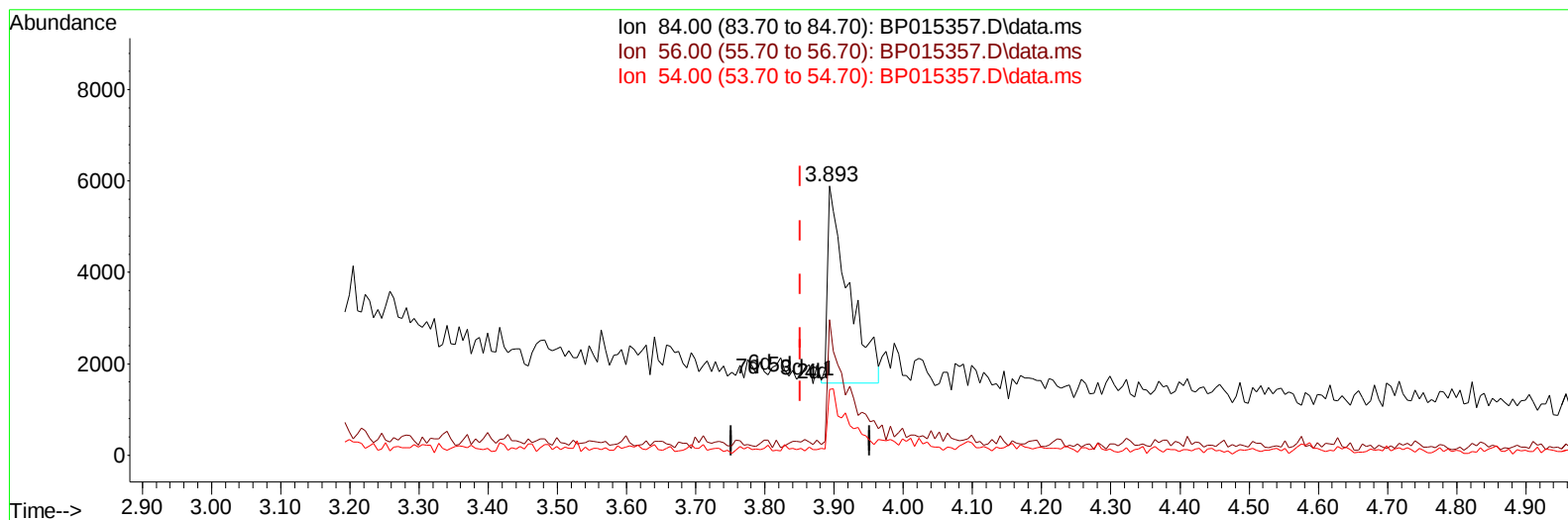
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TIC: BP015357.D\data.ms

(4) Pyridine-d5 (S)

3.893min (+ 0.041) 0.88 ng/u1 m

response 8918

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	50.31
54.00	29.10	24.54
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/01/2023
 Supervised By :mohammad ahmed 06/01/2023

Quant Time: May 31 22:55:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 31 14:24:45 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.105	152	139331	20.000	ng/ul	0.00
20) Naphthalene-d8	10.934	136	610139	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.751	164	449318	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188	1109419	20.000	ng/ul	0.00
79) Chrysene-d12	21.580	240	1209885	20.000	ng/ul	0.00
88) Perylene-d12	24.139	264	1349988	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.417	96	11914	3.322	ng/uL	0.00
4) Pyridine-d5	3.893	84	8918m	0.877	ng/ul	0.04
7) Phenol-d5	7.258	99	228715	17.680	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	141232	18.903	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	189688	19.943	ng/ul	0.00
15) 4-Methylphenol-d8	8.816	113	199830	19.385	ng/ul	0.00
21) Nitrobenzene-d5	9.287	128	90676	18.913	ng/ul	0.00
24) 2-Nitrophenol-d4	10.010	143	114893	21.537	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	198689	20.640	ng/ul	0.00
31) 4-Chloroaniline-d4	11.081	131	82616	5.801	ng/ul	0.00
46) Dimethylphthalate-d6	14.151	166	846071	24.919	ng/ul	0.00
49) Acenaphthylene-d8	14.445	160	872392	22.466	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	107310	16.276	ng/ul	0.00
60) Fluorene-d10	15.739	176	714084	24.924	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	126688	18.120	ng/ul	0.00
73) Anthracene-d10	17.604	188	1244903	24.505	ng/ul	0.00
81) Pyrene-d10	19.822	212	1652668	23.552	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264	1721739	25.692	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.940	105	68691	4.270	ng/ul	95

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

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