

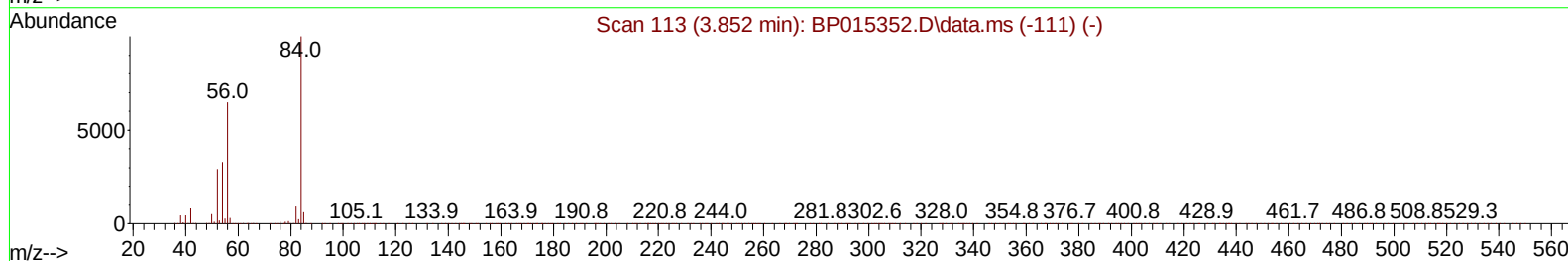
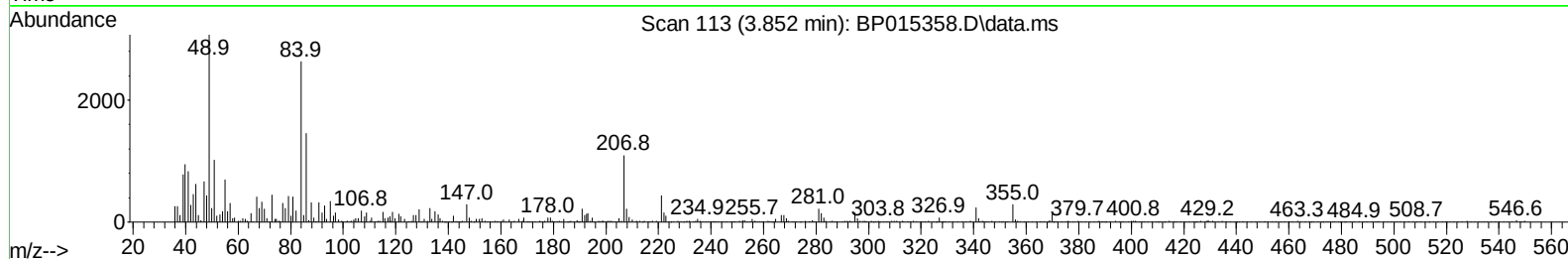
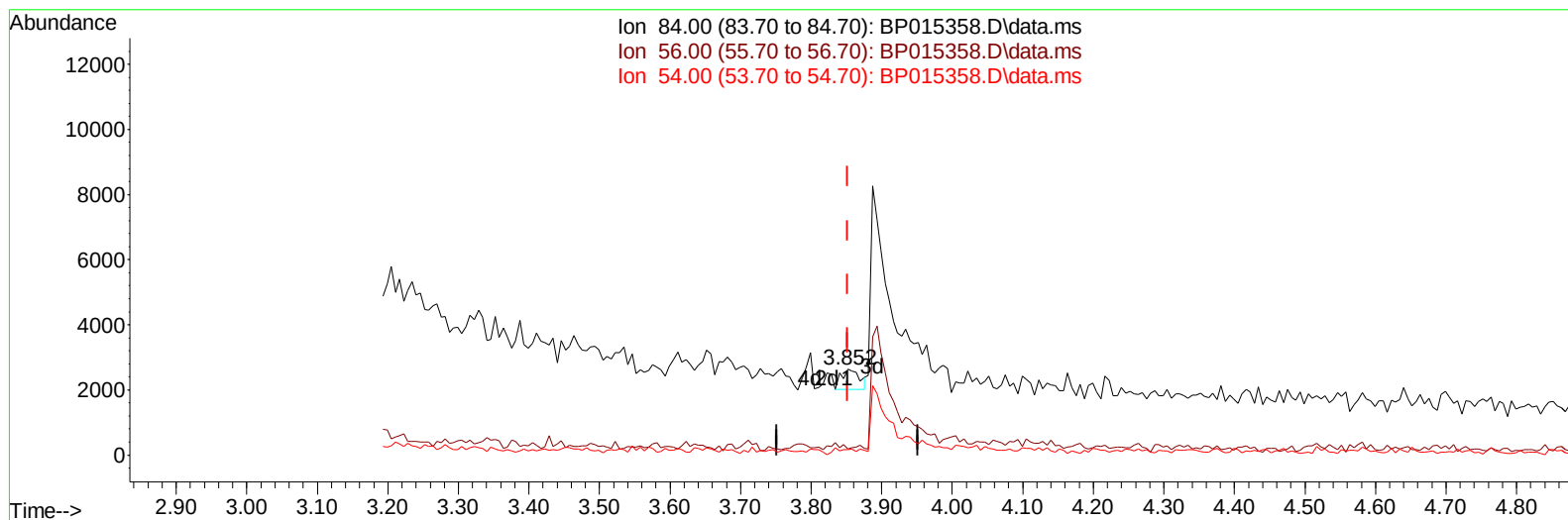
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
 Data File : BP015358.D
 Acq On : 31 May 2023 14:19
 Operator : CG/JU
 Sample : 02851-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0FK5

Manual IntegrationsAPPROVED

Quant Time: May 31 22:42:25 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: BP015358.D\data.ms

(4) Pyridine-d5 (S)

3.852min (+ 0.000) 0.09 ng/u1

response 1119

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	6.91#
54.00	29.10	6.69#
0.00	0.00	0.00

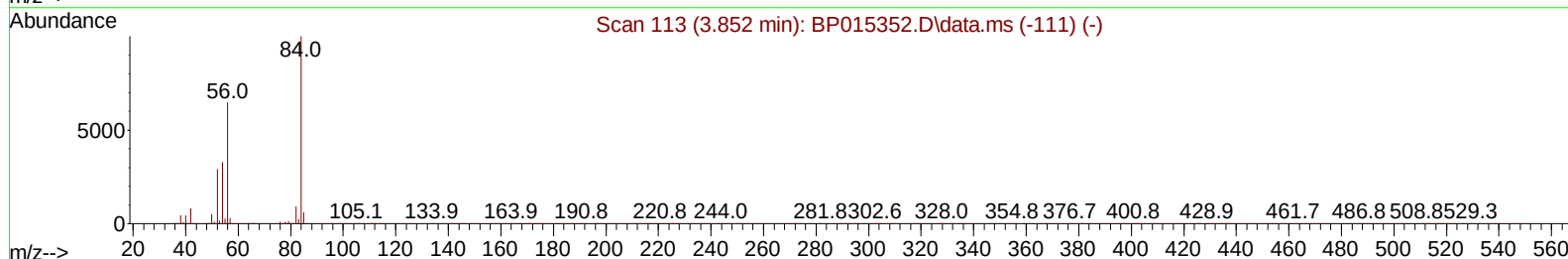
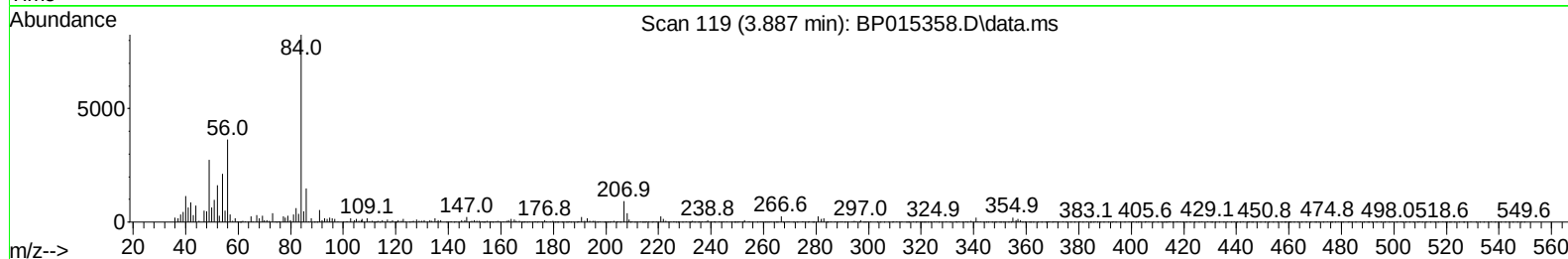
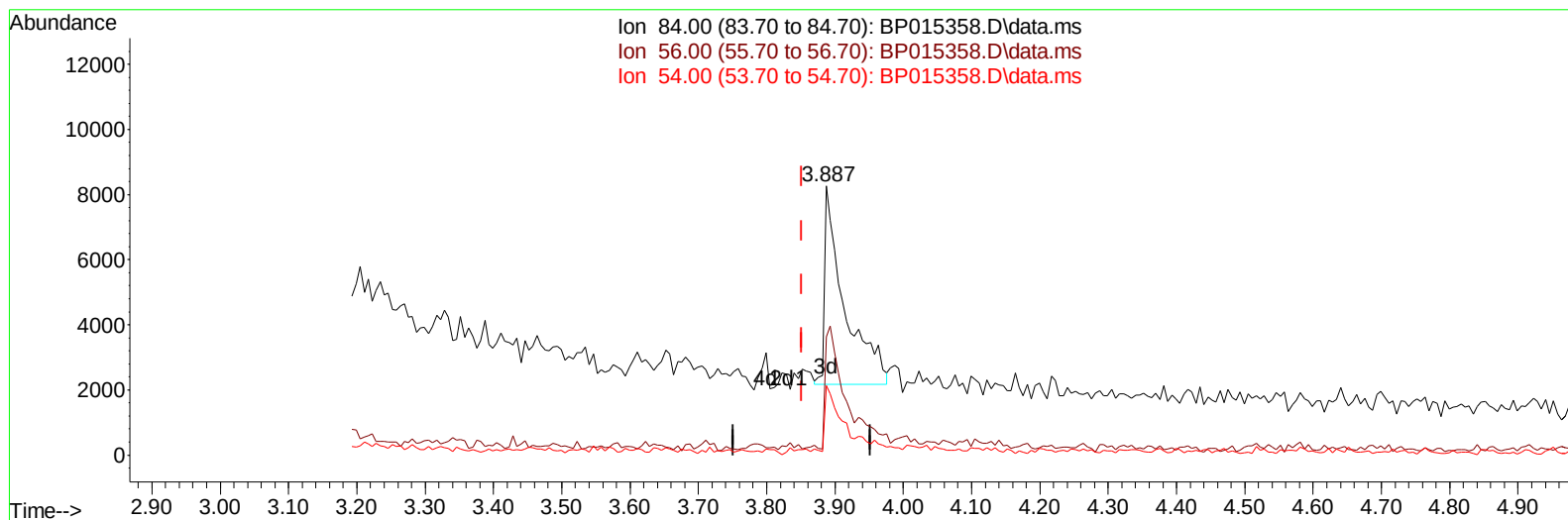
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(4) Pyridine-d5 (S)

3.887min (+ 0.035) 1.00 ng/ul m

response 12382

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	44.02#
54.00	29.10	25.89
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 22:57:37 2023
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.105	152		170273	20.000	ng/ul	0.00
20) Naphthalene-d8	10.934	136		730222	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.751	164		526643	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188		1309273	20.000	ng/ul	0.00
79) Chrysene-d12	21.580	240		1410819	20.000	ng/ul	0.00
88) Perylene-d12	24.139	264		1452500	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.417	96		15898	3.628	ng/uL	0.00
4) Pyridine-d5	3.887	84		12382m	0.997	ng/ul	0.04
7) Phenol-d5	7.258	99		299666	18.955	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67		183756	20.125	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132		247205	21.267	ng/ul	0.00
15) 4-Methylphenol-d8	8.822	113		256995	20.400	ng/ul	0.00
21) Nitrobenzene-d5	9.281	128		131860	22.980	ng/ul	0.00
24) 2-Nitrophenol-d4	10.010	143		158055	24.756	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.552	165		262369	22.773	ng/ul	0.00
31) 4-Chloroaniline-d4	11.075	131		252335	14.803	ng/ul	0.00
46) Dimethylphthalate-d6	14.151	166		1094992	27.515	ng/ul	0.00
49) Acenaphthylene-d8	14.445	160		1134047	24.916	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143		170968	22.124	ng/ul	0.00
60) Fluorene-d10	15.740	176		909409	27.082	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200		200242	24.268	ng/ul	0.00
73) Anthracene-d10	17.604	188		1595361	26.610	ng/ul	0.00
81) Pyrene-d10	19.828	212		2135669	26.100	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264		2081258	28.865	ng/ul	0.00
Target Compounds							
16) Acetophenone	8.940	105		48784	2.481	ng/ul	Qvalue 94

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

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