

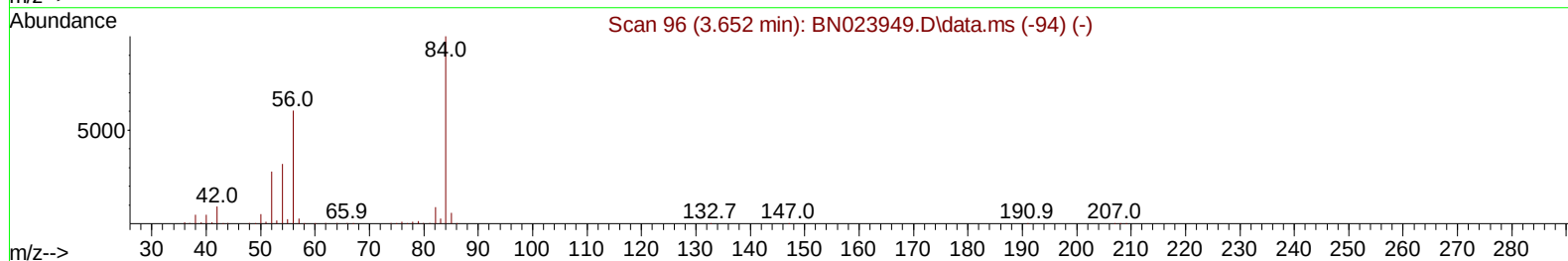
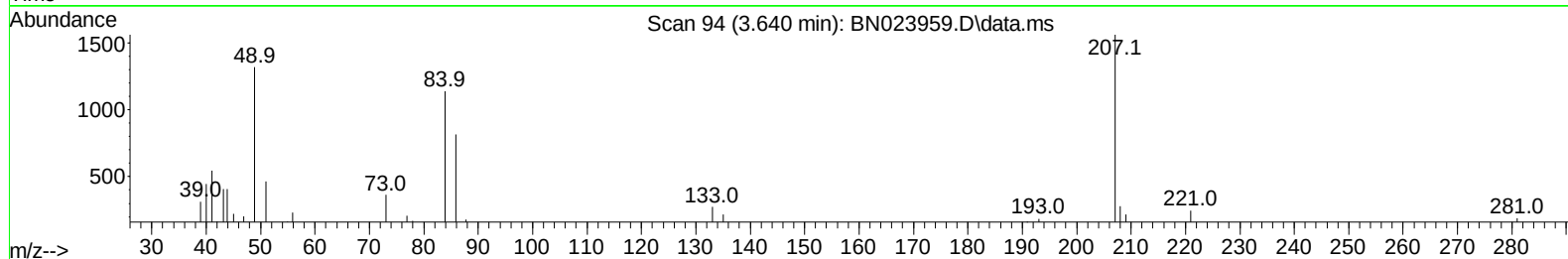
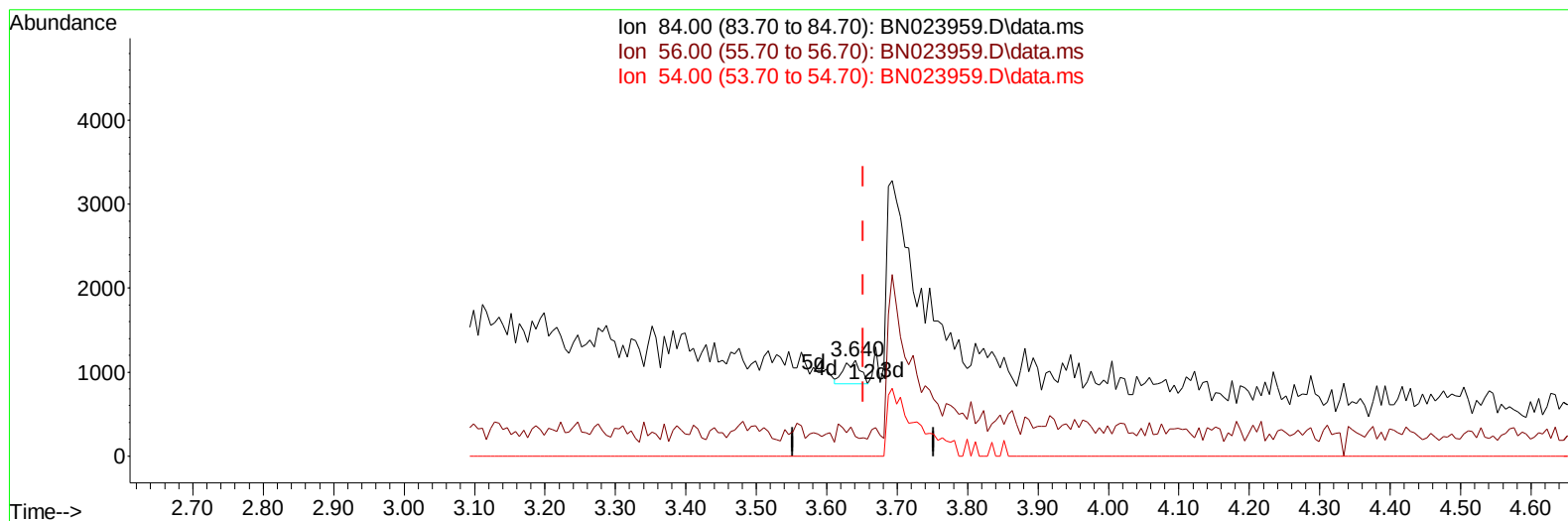
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020623\
Data File : BN023959.D
Acq On : 06 Feb 2023 19:02
Operator : CG/JU
Sample : 01362-05MEDL 20X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4443MEDL

Manual IntegrationsAPPROVED

Quant Time: Feb 07 03:01:26 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020623.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 07 02:28:26 2023
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/07/2023
Supervised By :mohammad ahmed 02/07/2023



TIC: BN023959.D\data.ms

(4) Pyridine-d5 (S)

3.640min (-0.012) 0.05 ng/u1

response 440

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.40	20.46#
54.00	31.70	0.00#
0.00	0.00	0.00

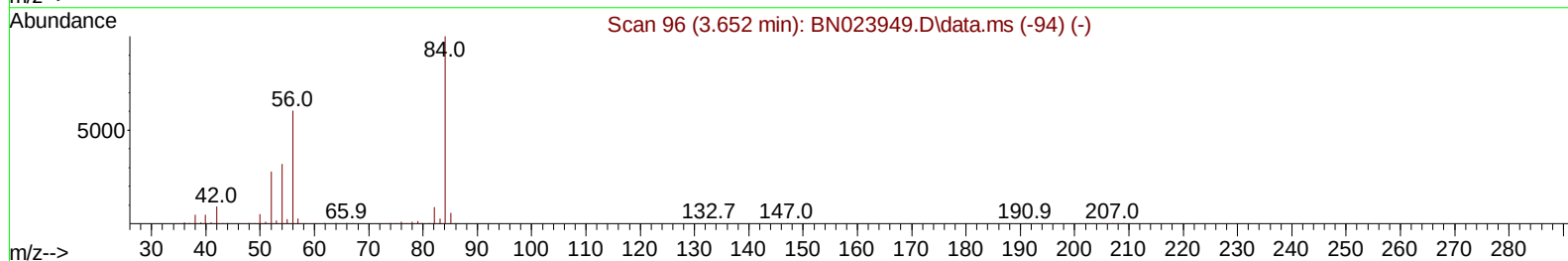
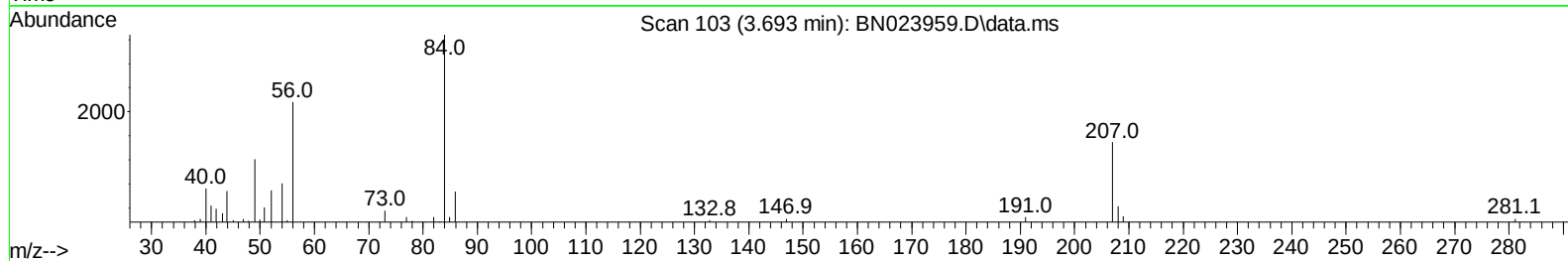
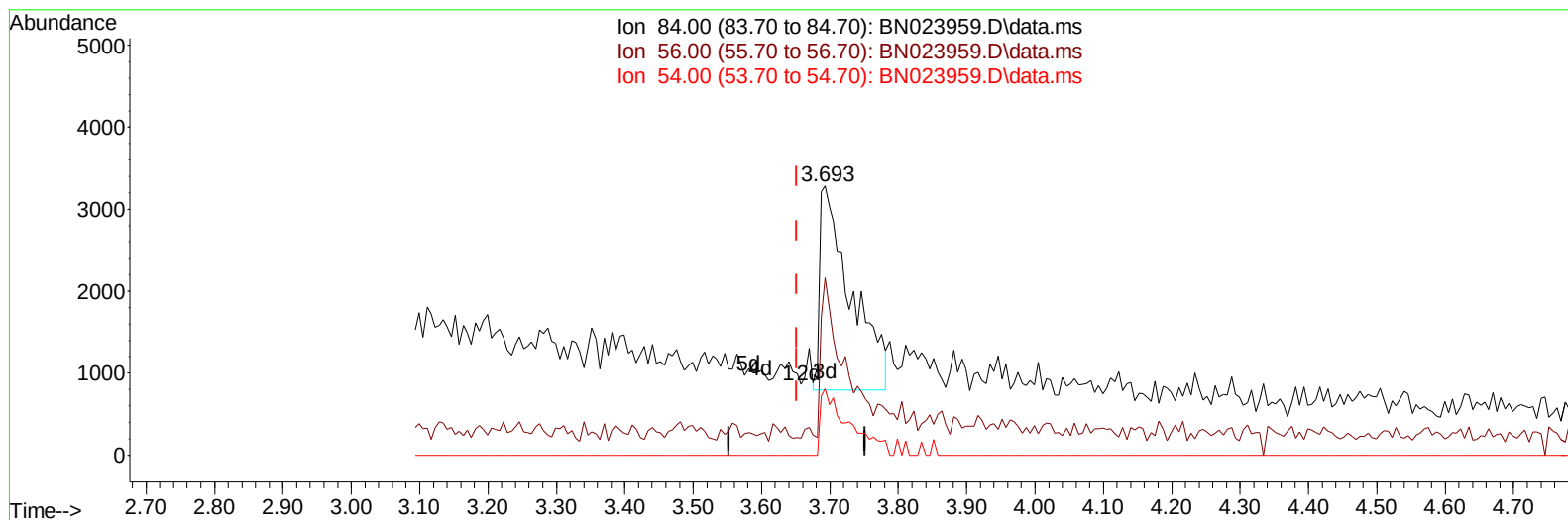
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(4) Pyridine-d5 (S)

3.693min (+ 0.041) 0.87 ng/ul m

response 7875

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.40	65.89
54.00	31.70	24.72#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.840	152		132514	20.000	ng/ul	0.00
20) Naphthalene-d8	10.646	136		547124	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164		349959	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188		765205	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240		649234	20.000	ng/ul	0.00
88) Perylene-d12	23.816	264		662626	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	0.000	96		0d	0.000	ng/uL	
4) Pyridine-d5	3.693	84		7875m	0.866	ng/ul	0.04
7) Phenol-d5	7.011	99		13946	1.231	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67		10694	1.496	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132		11676	1.325	ng/ul	0.00
15) 4-Methylphenol-d8	8.558	113		10323	1.166	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128		5412	1.239	ng/ul	0.00
24) 2-Nitrophenol-d4	9.728	143		4960	1.011	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165		10465	1.145	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131		16502	1.336	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166		41580	1.537	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160		36912	1.232	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143		0d	0.000	ng/ul	
60) Fluorene-d10	15.487	176		36746	1.613	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200		0d	0.000	ng/ul	
73) Anthracene-d10	17.339	188		49012	1.406	ng/ul	0.00
81) Pyrene-d10	19.633	212		58265	1.490	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.669	264		46763	1.379	ng/ul	0.00
Target Compounds							
						Qvalue	
86) Bis(2-ethylhexyl)phtha...	21.345	149		708454	25.728	ng/ul	96

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

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