

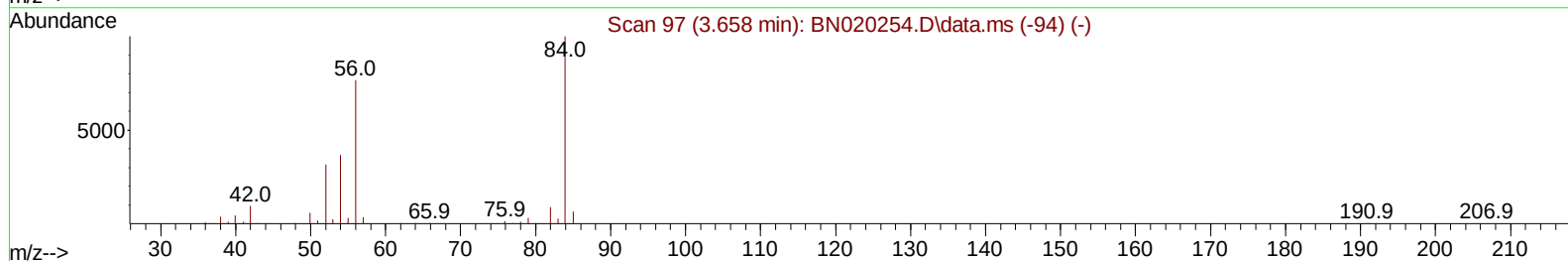
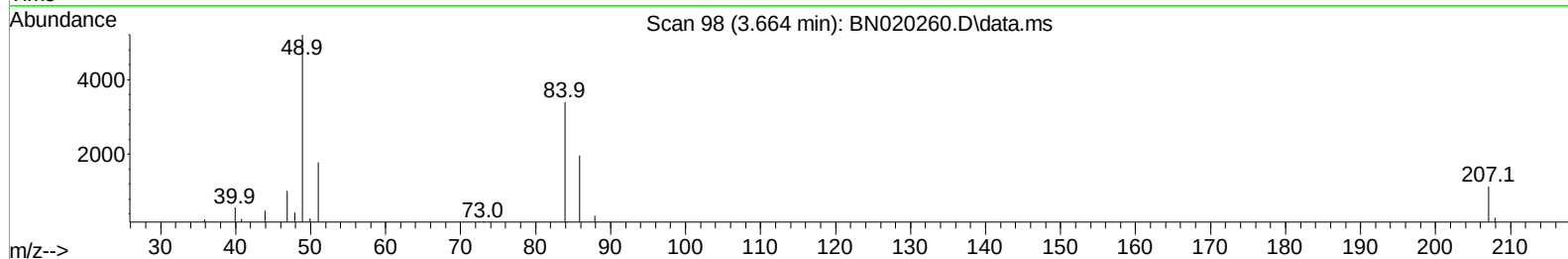
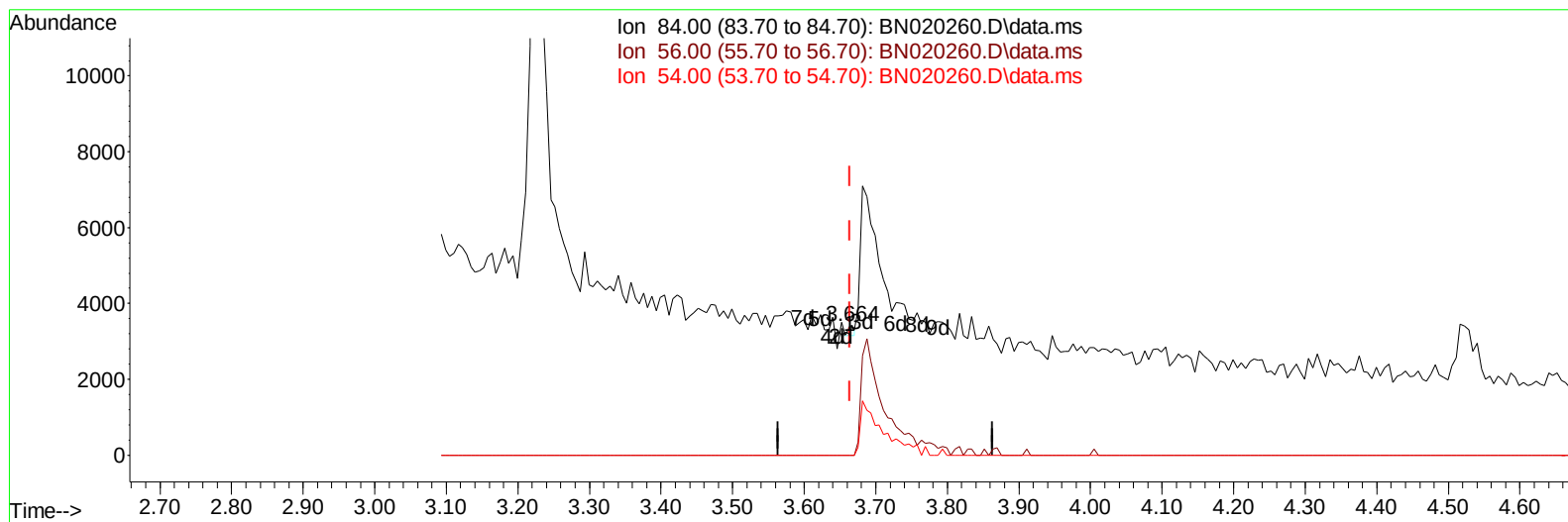
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061722\
 Data File : BN020260.D
 Acq On : 18 Jun 2022 13:20
 Operator : CG/JU
 Sample : N3313-07DL 10X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YB7L5DL

Manual IntegrationsAPPROVED

Quant Time: Jun 20 02:23:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 17 22:53:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/20/2022
 Supervised By :mohammad ahmed 06/22/2022



TIC: BN020260.D\data.ms

(4) Pyridine-d5 (S)

3.664min (-0.000) 0.02 ng/u1

response 143

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	0.00#
54.00	33.00	0.00#
0.00	0.00	0.00

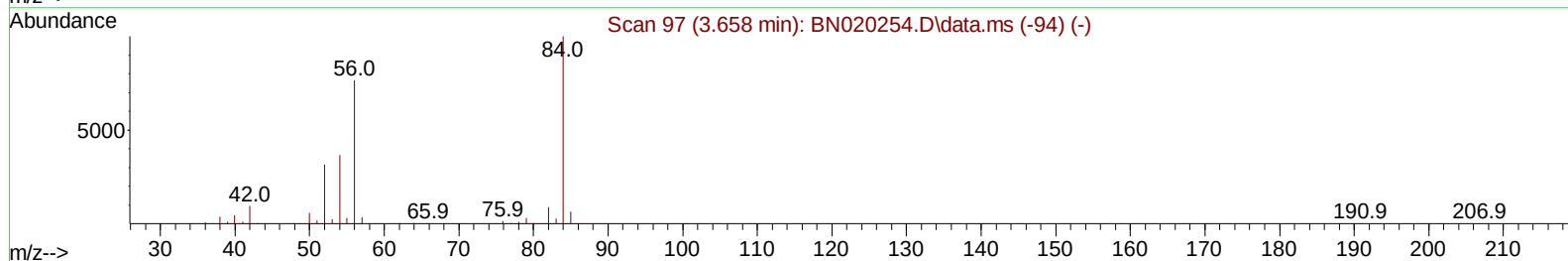
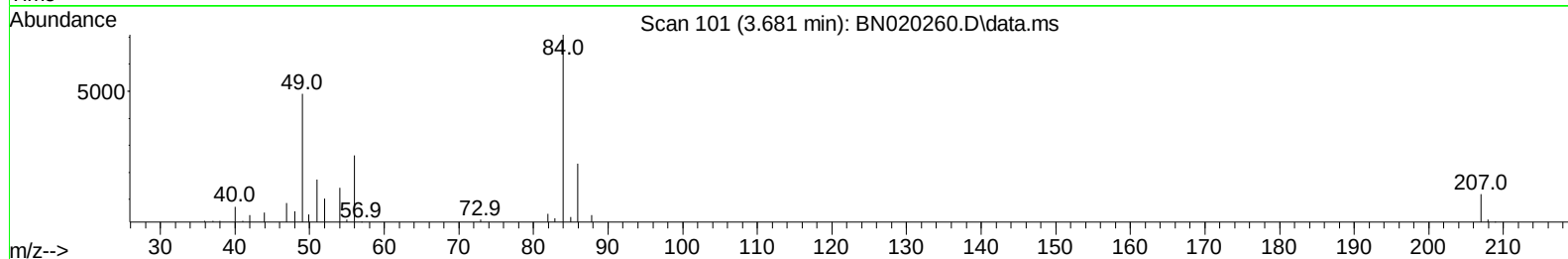
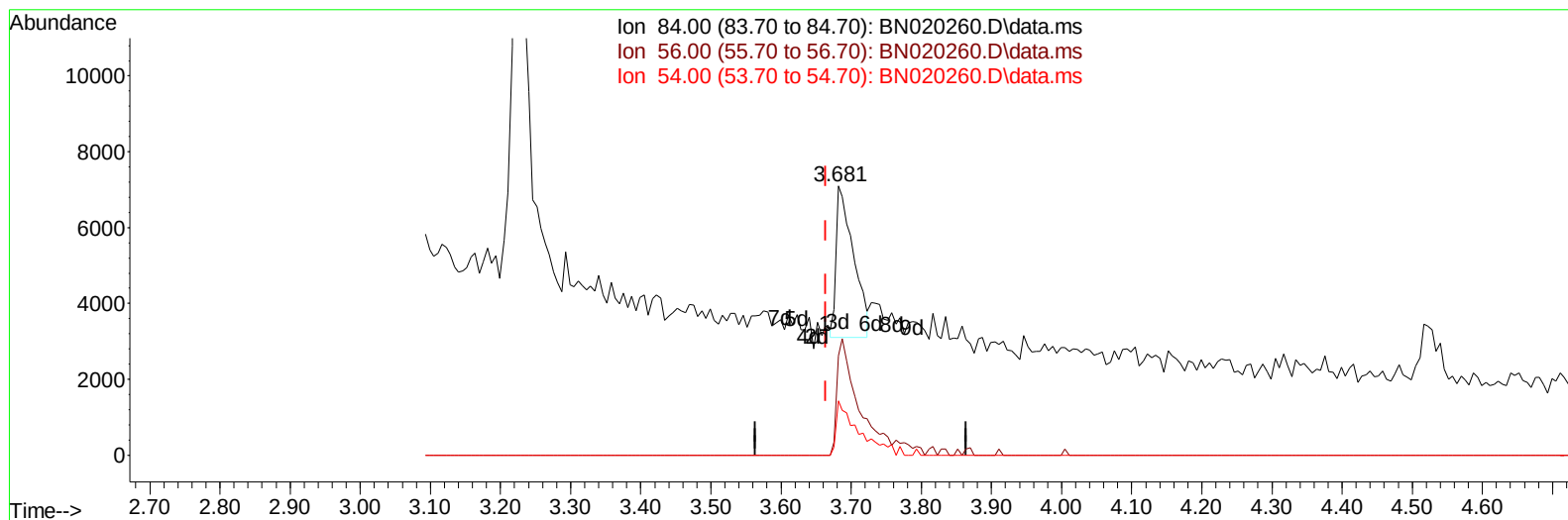
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(4) Pyridine-d5 (S)

3.681min (+ 0.018) 0.93 ng/ul m

response 6849

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	36.92#
54.00	33.00	20.10#
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	7.787	152	117276	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	583121	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	423630	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.157	188	876360	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	702850	20.000	ng/ul	0.00
88) Perylene-d12	23.663	264	571888	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	64592	23.855	ng/uL	-0.02
4) Pyridine-d5	3.681	84	6849m	0.925	ng/ul	0.02
7) Phenol-d5	6.987	99	5624	0.580	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	19542	3.223	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	19009	2.411	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	11843	1.411	ng/ul	0.00
21) Nitrobenzene-d5	8.940	128	10953	2.469	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	11046	2.346	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	22366	2.721	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	30300	2.277	ng/ul	0.01
46) Dimethylphthalate-d6	13.828	166	101610	3.296	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	115116	3.220	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.410	176	97756	3.833	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.257	188	155480	4.065	ng/ul	0.00
81) Pyrene-d10	19.557	212	168250	4.416	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.516	264	101698	3.635	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.287	88	47086	16.473	ng/uL	93

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

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