

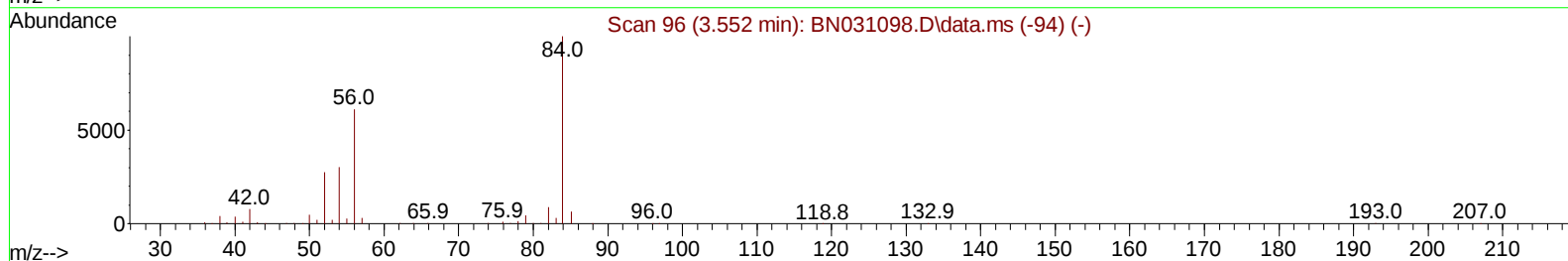
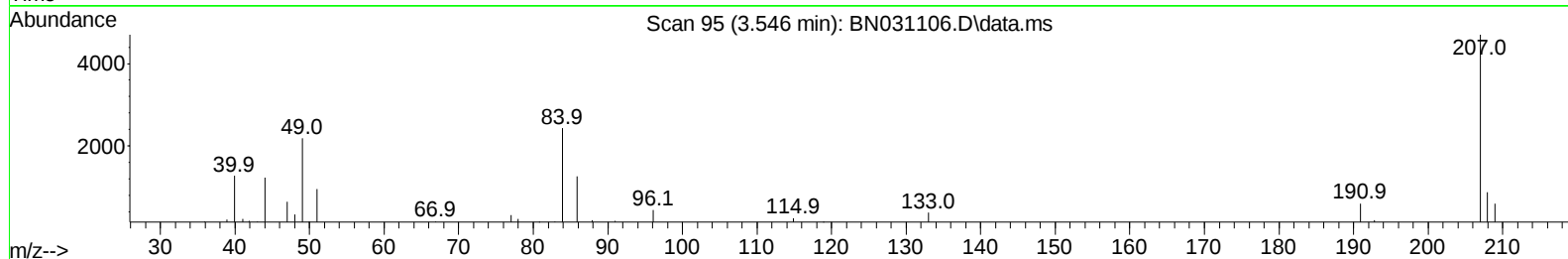
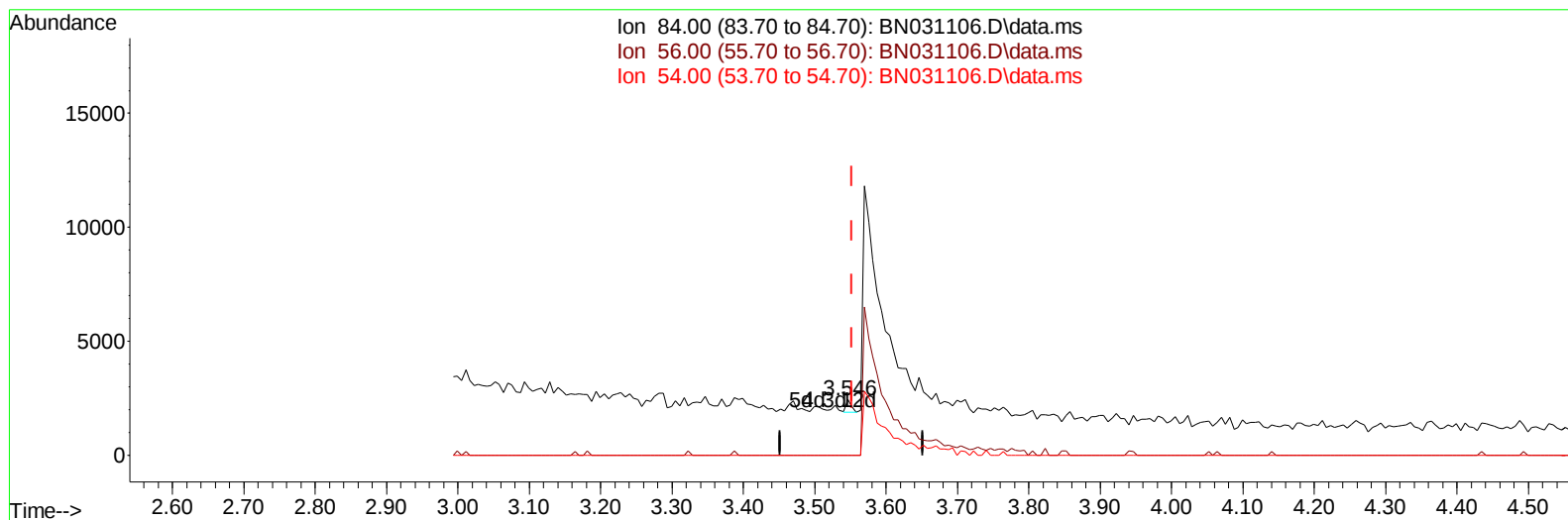
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031106.D
Acq On : 21 Apr 2024 13:09
Operator : MA/JU
Sample : P2161-02
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORD2

Manual IntegrationsAPPROVED

Quant Time: Apr 21 14:44:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 19 22:56:25 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/22/2024
Supervised By :mohammad ahmed 04/22/2024



TIC: BN031106.D\data.ms

(4) Pyridine-d5 (S)

3.546min (-0.006) 0.04 ng/u1

response 277

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.00	0.00#
54.00	30.10	0.00#
0.00	0.00	0.00

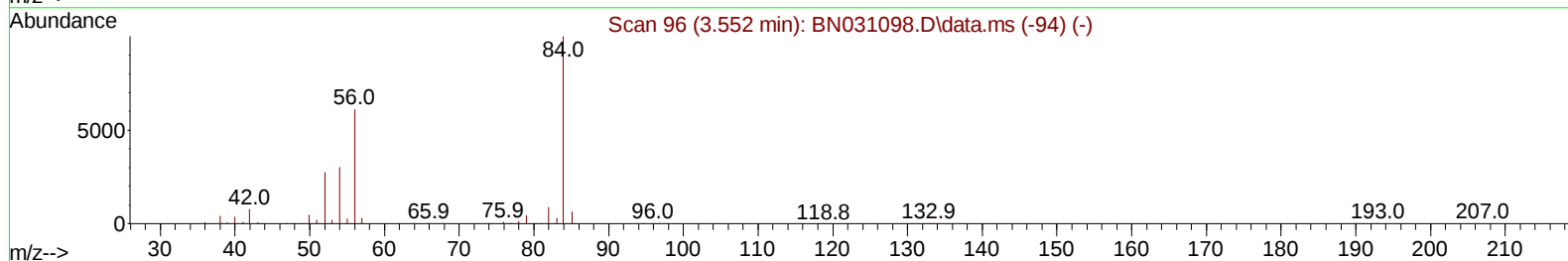
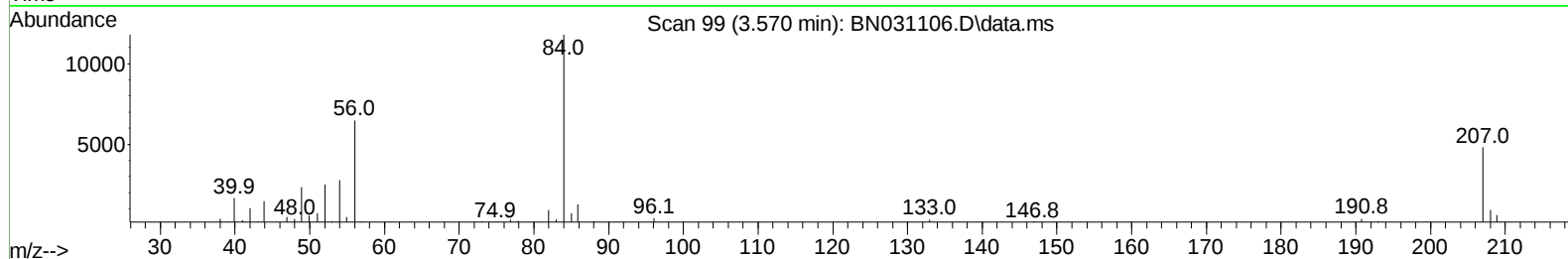
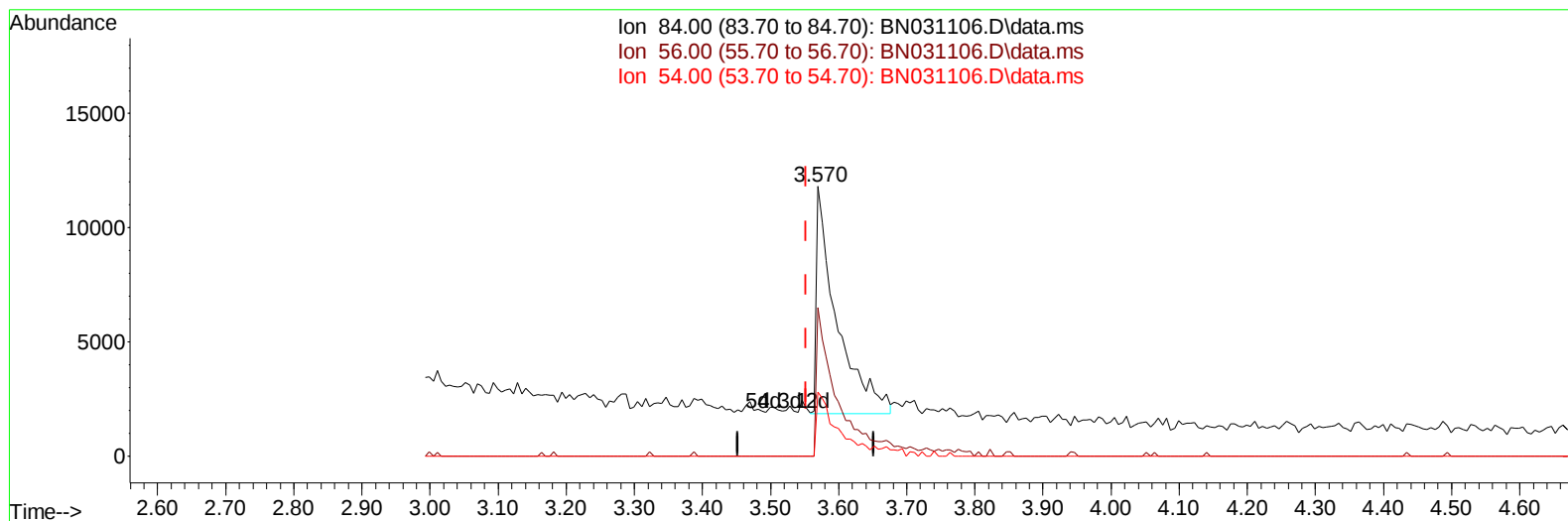
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(4) Pyridine-d5 (S)

3.570min (+ 0.018) 2.80 ng/u1 m

response 20368

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.00	55.03
54.00	30.10	23.67#
0.00	0.00	0.00

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 Sample : P2161-02
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 CORD2

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/22/2024
 Supervised By :mohammad ahmed 04/22/2024

Quant Time: Apr 21 14:45:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 19 22:56:25 2024
 Response via : Initial Calibration

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.634	152		117833	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136		531375	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.281	164		356971	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188		820985	20.000	ng/ul	0.00
79) Chrysene-d12	21.263	240		1012618	20.000	ng/ul	0.00
88) Perylene-d12	24.169	264		1301982	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.146	96		13479	5.344	ng/uL	0.00
4) Pyridine-d5	3.570	84		20368m	2.804	ng/ul	0.02
7) Phenol-d5	6.805	99		104338	10.497	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67		187076	32.348	ng/ul	0.00
11) 2-Chlorophenol-d4	7.164	132		212458	27.451	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113		165135	19.754	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128		123297	32.295	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143		121612	30.644	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.034	165		233527	29.839	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131		47010	3.844	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166		834943	33.612	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160		952882	33.851	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143		46703	9.008	ng/ul	0.00
60) Fluorene-d10	15.275	176		767196	35.478	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200		99192	23.948	ng/ul	0.00
73) Anthracene-d10	17.128	188		1341673	37.704	ng/ul	0.00
81) Pyrene-d10	19.433	212		1735421	34.876	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.969	264		2360382	38.230	ng/ul	0.00
Target Compounds							
						Qvalue	
75) 1,2,3,4-Tetrachloroben...	13.063	216		12187	1.266	ng/uL	97

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

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