

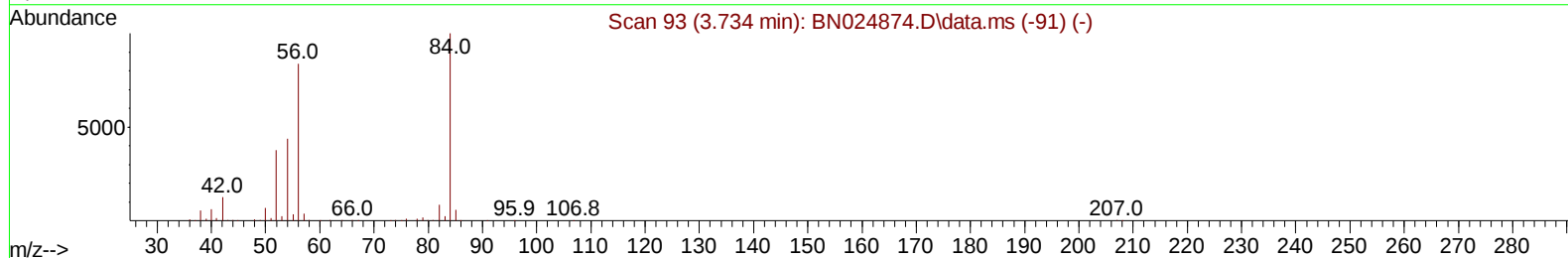
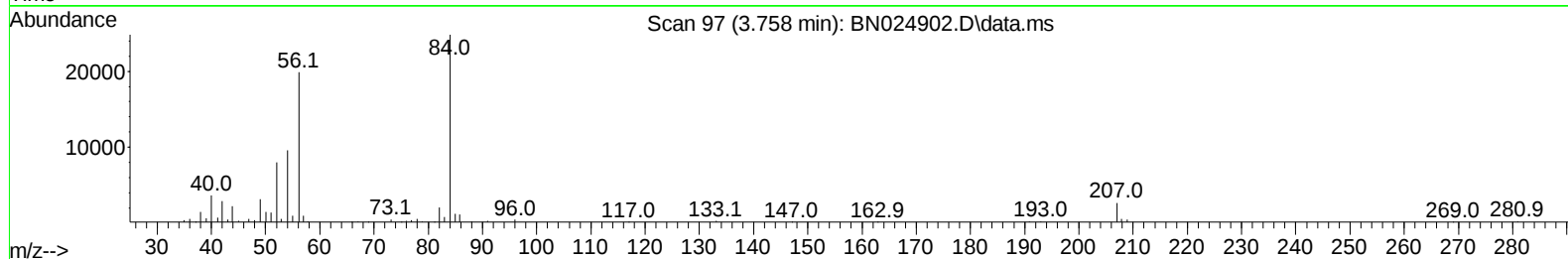
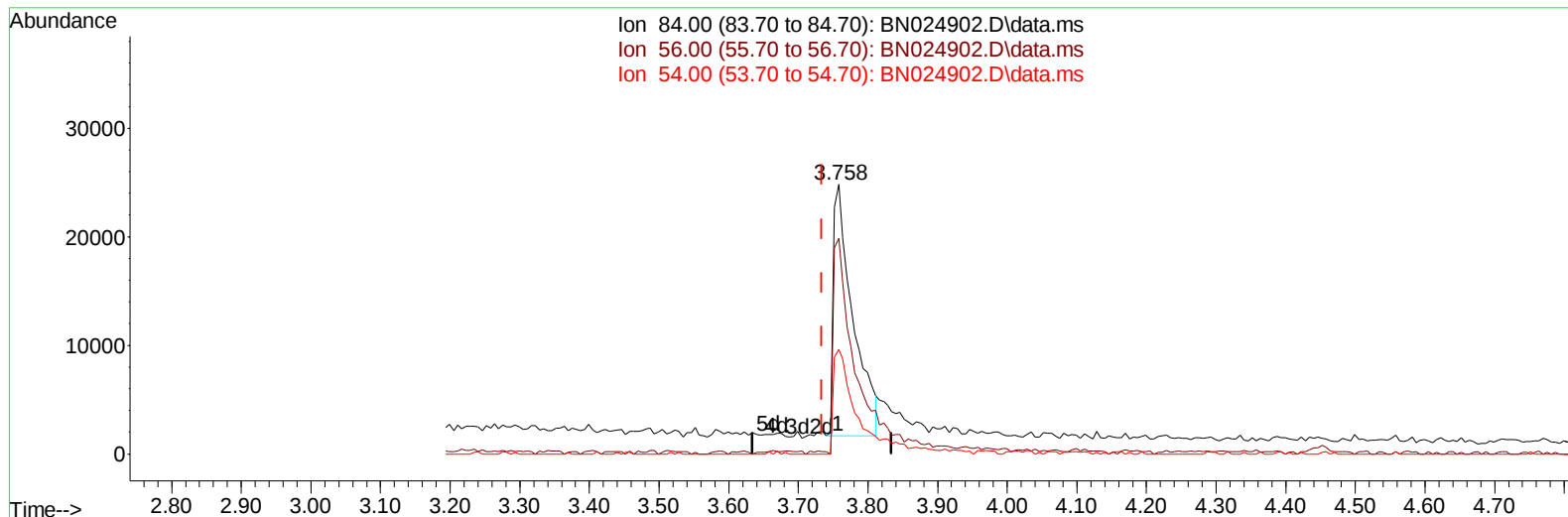
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024902.D
Acq On : 14 Apr 2023 17:46
Operator : CG/JU
Sample : 02252-10
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
COME6

Manual IntegrationsAPPROVED

Quant Time: Apr 14 22:48:21 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 14 22:34:56 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/17/2023
Supervised By :Jagrut Upadhyay 04/17/2023



TIC: BN024902.D\data.ms

(4) Pyridine-d5 (S)

3.758min (+ 0.024) 2.09 ng/u1

response 44740

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	80.11
54.00	40.20	38.73
0.00	0.00	0.00

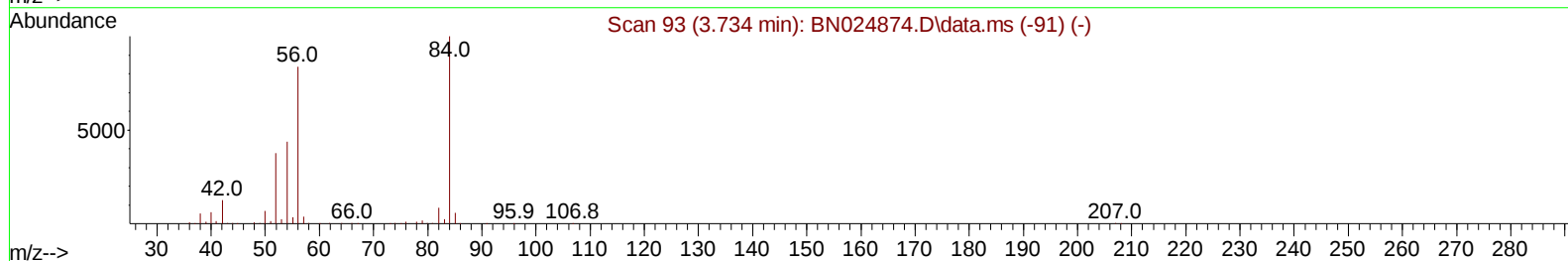
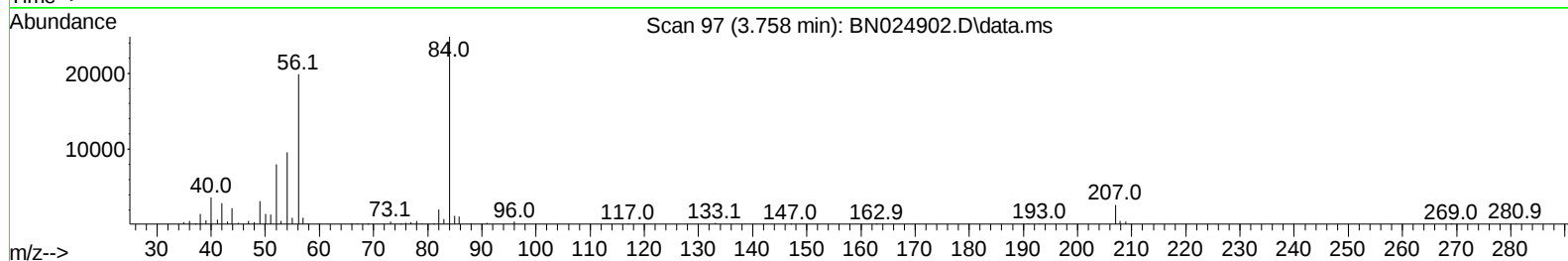
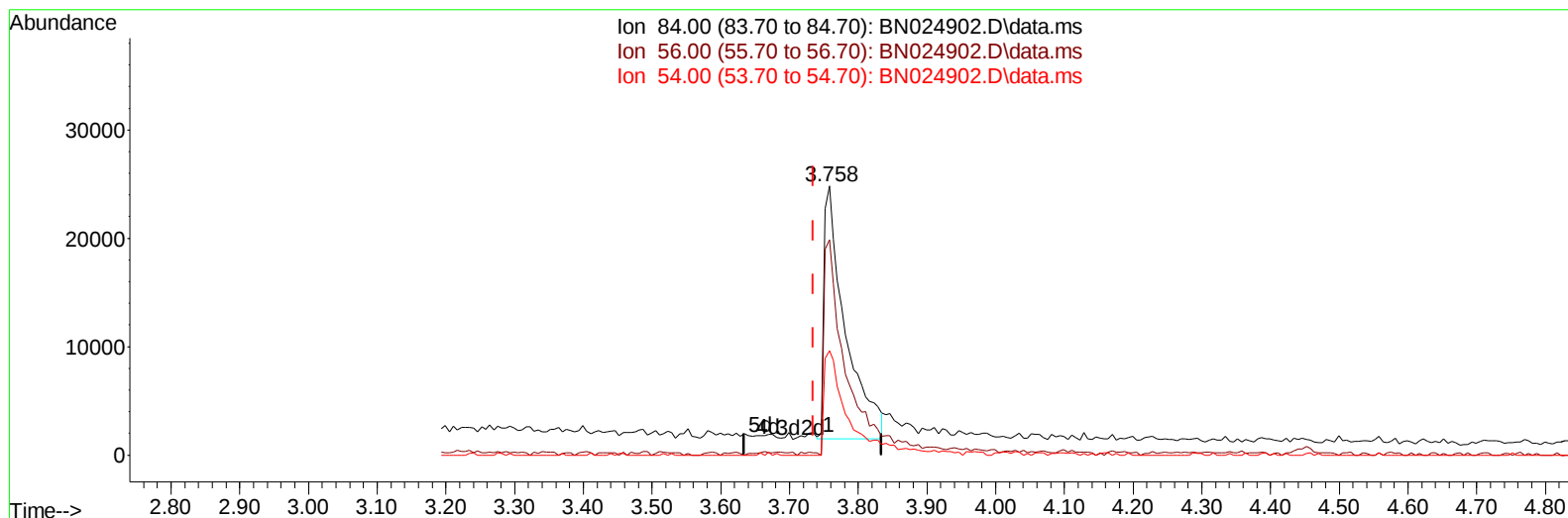
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(4) Pyridine-d5 (S)

3.758min (+ 0.024) 2.34 ng/u1 m

response 49924

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	80.11
54.00	40.20	38.73
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	7.940	152	271628	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	1196541	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	753714	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.334	188	1697744	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	1610188	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264	1739852	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	27271	3.821	ng/uL	0.00
4) Pyridine-d5	3.758	84	49924m	2.336	ng/ul	0.02
7) Phenol-d5	7.099	99	101367	3.828	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	384462	21.705	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	304882	15.666	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	200242	9.525	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	209090	21.582	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	208611	19.879	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	358582	19.118	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	444977	15.633	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1382091	23.654	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1533532	22.904	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	29059	2.426	ng/ul	0.00
60) Fluorene-d10	15.575	176	1231493	24.963	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	163613	14.808	ng/ul	0.00
73) Anthracene-d10	17.434	188	2012936	25.245	ng/ul	0.00
81) Pyrene-d10	19.728	212	2423715	25.922	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	2246248	24.692	ng/ul	-0.01

Target Compounds Qvalue

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

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