

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
 Data File : BN024930.D
 Acq On : 15 Apr 2023 12:13
 Operator : CG/JU
 Sample : 02252-05DL 5X
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

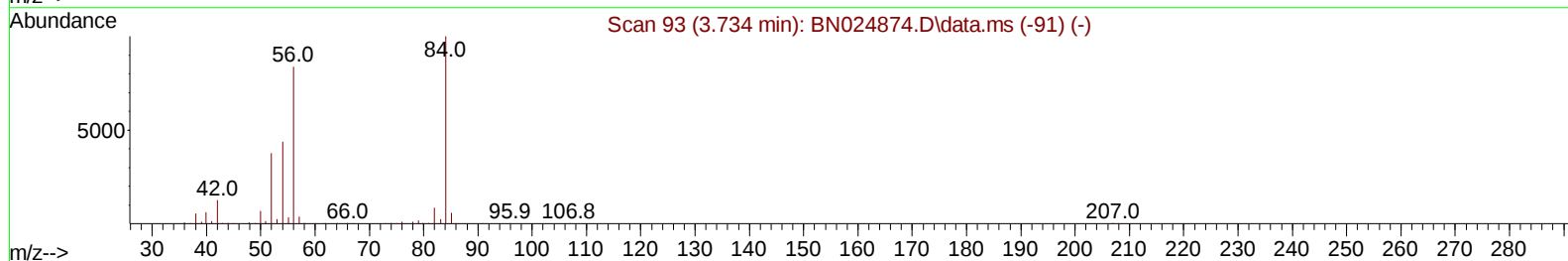
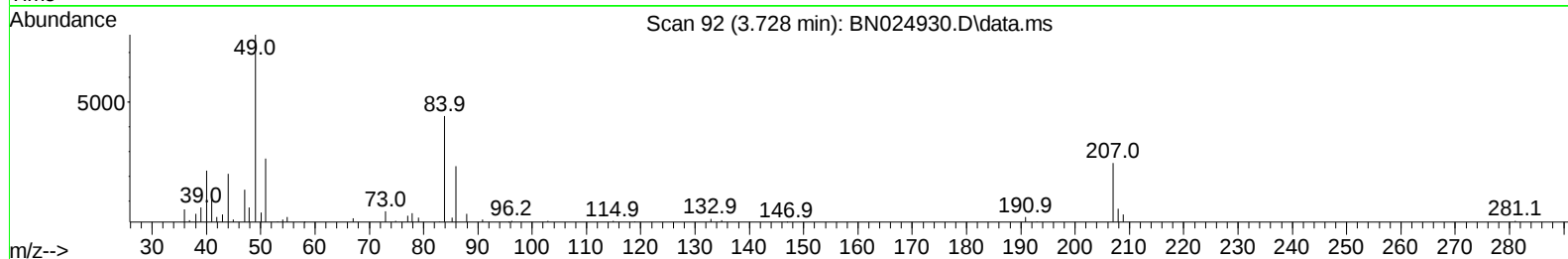
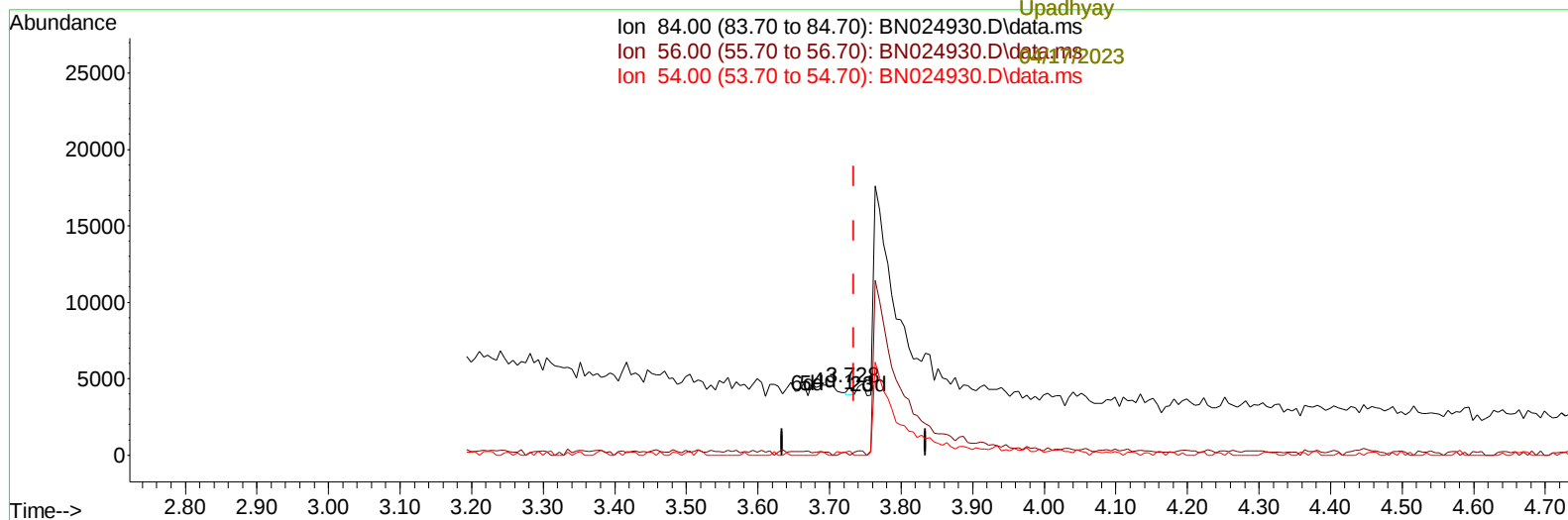
Instrument :
 BNA_N
 ClientSampleId :
 COME1DL

Manual IntegrationsAPPROVED

Quant Time: Apr 17 03:01:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 17 02:58:51 2023
 Response via : Initial Calibration

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

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 Upadhyay



TIC: BN024930.D\data.ms

(4) Pyridine-d5 (S)

3.728min (-0.006) 0.01 ng/u1

response 167

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	0.00#
54.00	40.20	5.91#
0.00	0.00	0.00

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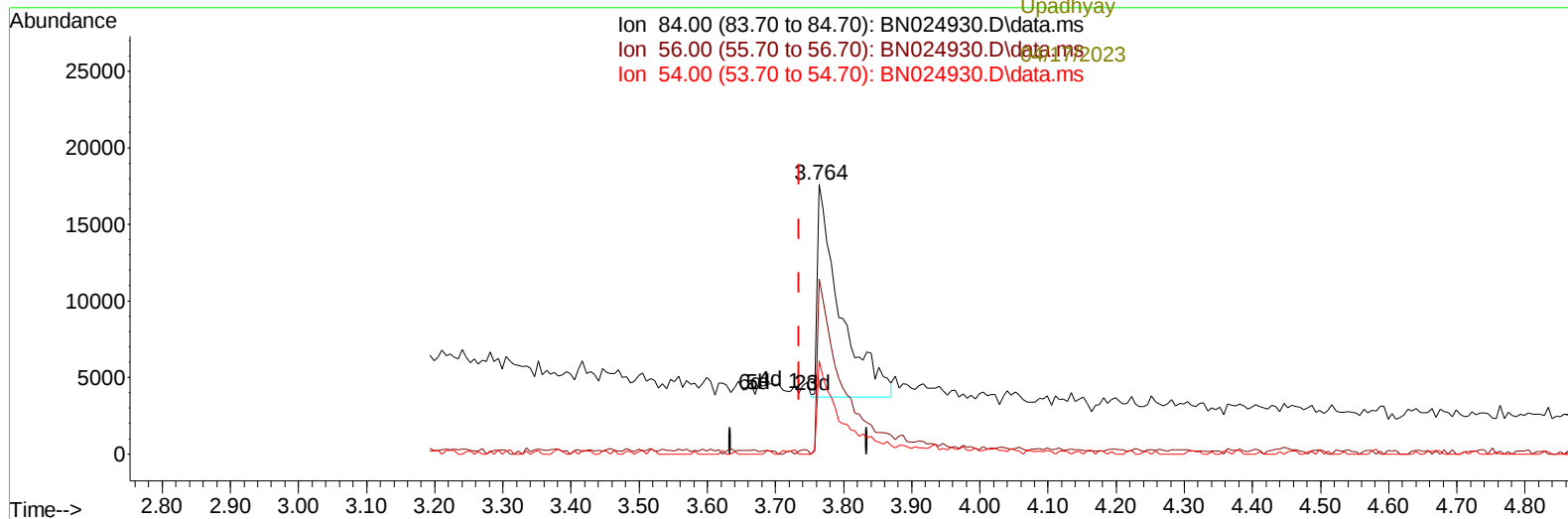
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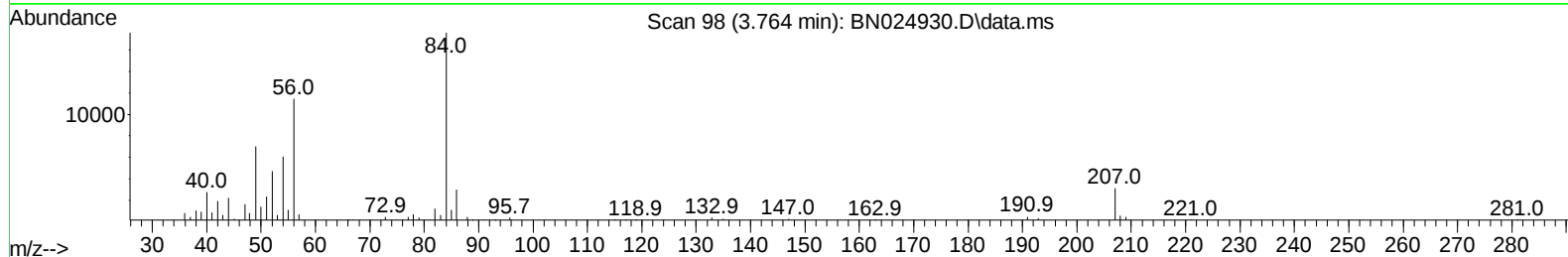
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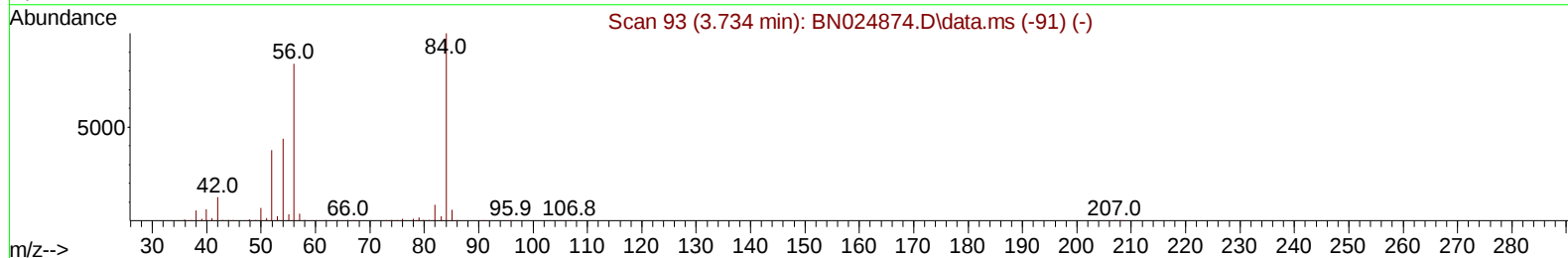
Ion 84.00 (83.70 to 84.70): BN024930.D\data.ms
 Ion 56.00 (55.70 to 56.70): BN024930.D\data.ms
 Ion 54.00 (53.70 to 54.70): BN024930.D\data.ms



Scan 98 (3.764 min): BN024930.D\data.ms



Scan 93 (3.734 min): BN024874.D\data.ms (-91) (-)



TIC: BN024930.D\data.ms

(4) Pyridine-d5 (S)

3.764min (+ 0.029) 1.22 ng/ul m

response 31944

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	65.01#
54.00	40.20	34.44
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----04/17/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	333733	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	1438332	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	886477	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1953376	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	1912905	20.000	ng/ul	0.00
88) Perylene-d12	23.968	264	2079307	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	6529	0.745	ng/uL	0.00
4) Pyridine-d5	3.764	84	31944m	1.217	ng/ul	0.03
7) Phenol-d5	7.099	99	29130	0.895	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	229800	10.559	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	92701	3.877	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	58194	2.253	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	58511	5.024	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	60430	4.791	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	104416	4.631	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	135324	3.955	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	404821	5.891	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	410918	5.218	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.569	176	356760	6.149	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.428	188	569022	6.202	ng/ul	0.00
81) Pyrene-d10	19.716	212	736512	6.631	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.810	264	683023	6.282	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.122	94	42079	1.242	ng/ul	99
22) Nitrobenzene	9.146	77	864266	27.618	ng/ul	98

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

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