

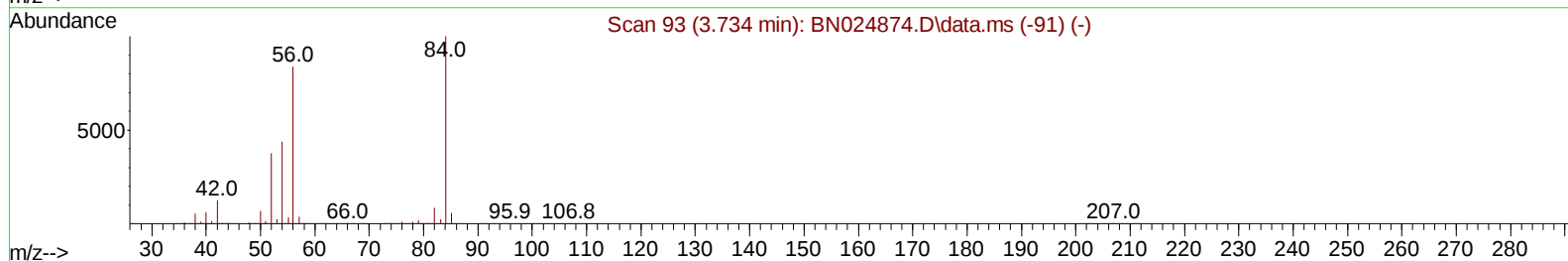
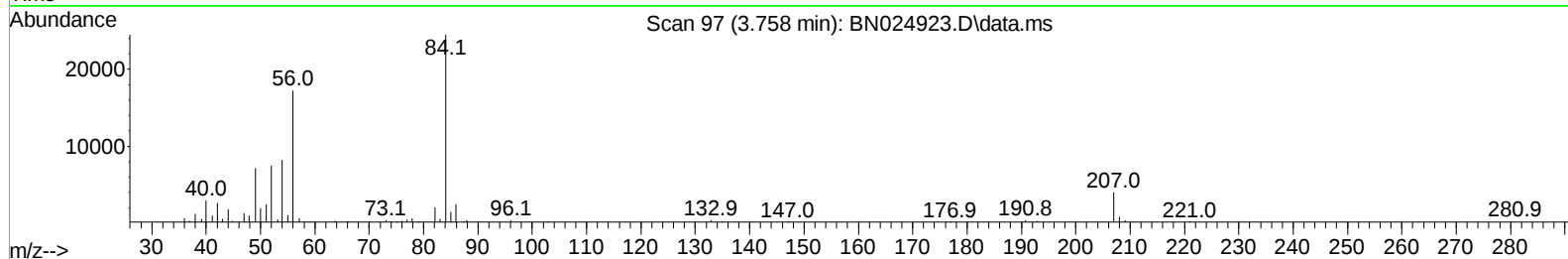
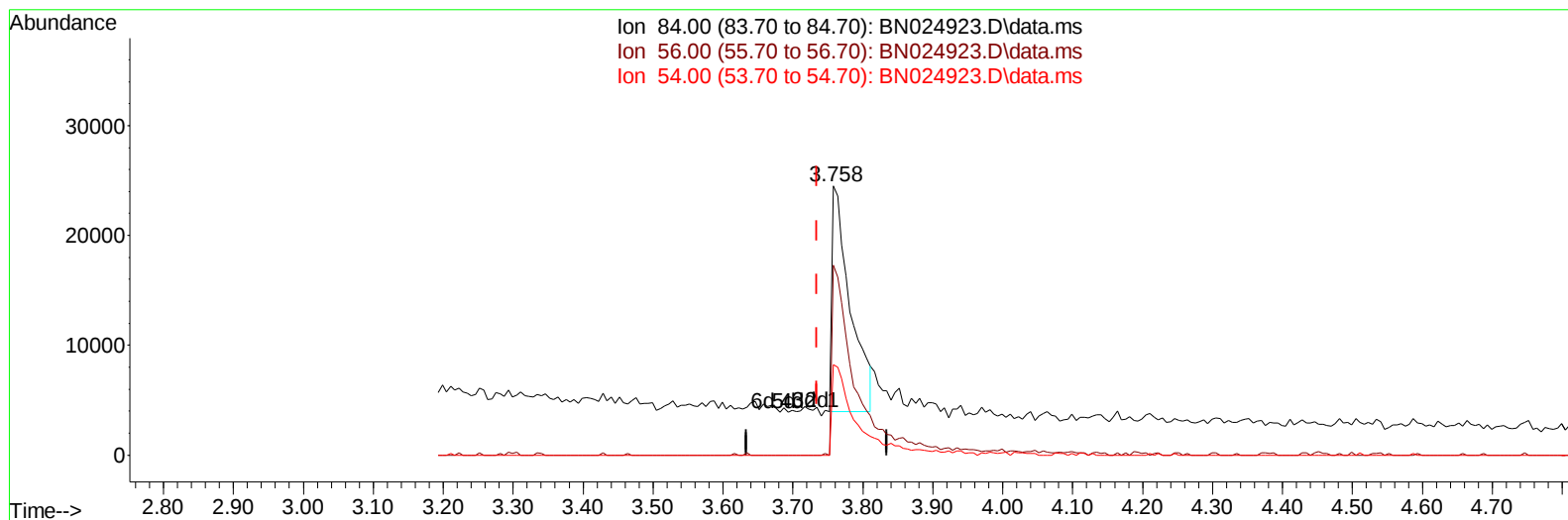
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
 Data File : BN024923.D
 Acq On : 15 Apr 2023 07:21
 Operator : CG/JU
 Sample : PB152131BL
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK131

Manual IntegrationsAPPROVED

Quant Time: Apr 17 02:51:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 15 00:01:14 2023
 Response via : Initial Calibration

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



TIC: BN024923.D\data.ms

(4) Pyridine-d5 (S)

3.758min (+ 0.023) 1.39 ng/u1

response 37438

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	70.53
54.00	40.20	33.71
0.00	0.00	0.00

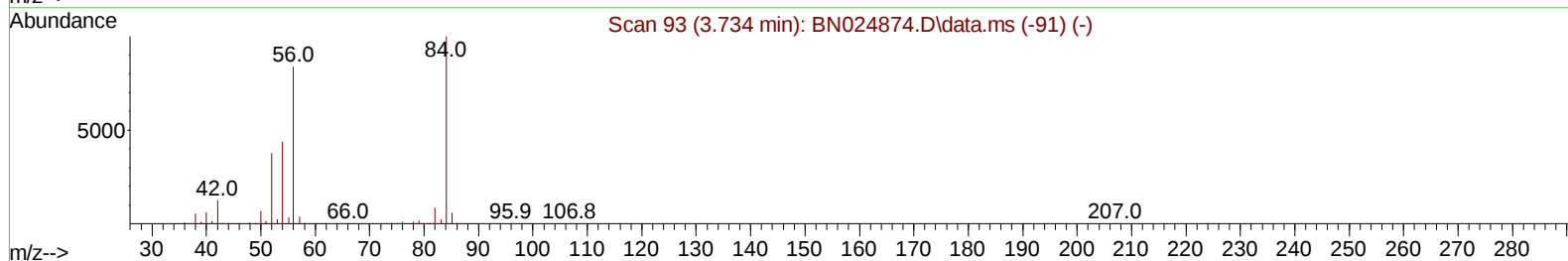
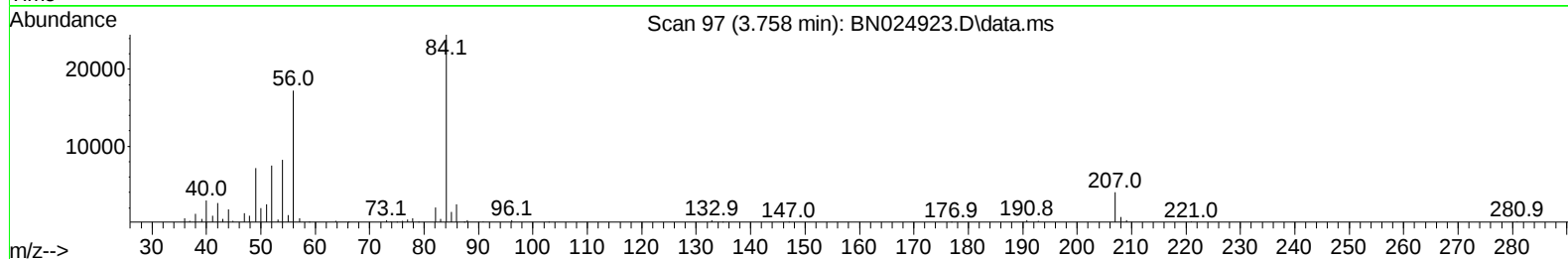
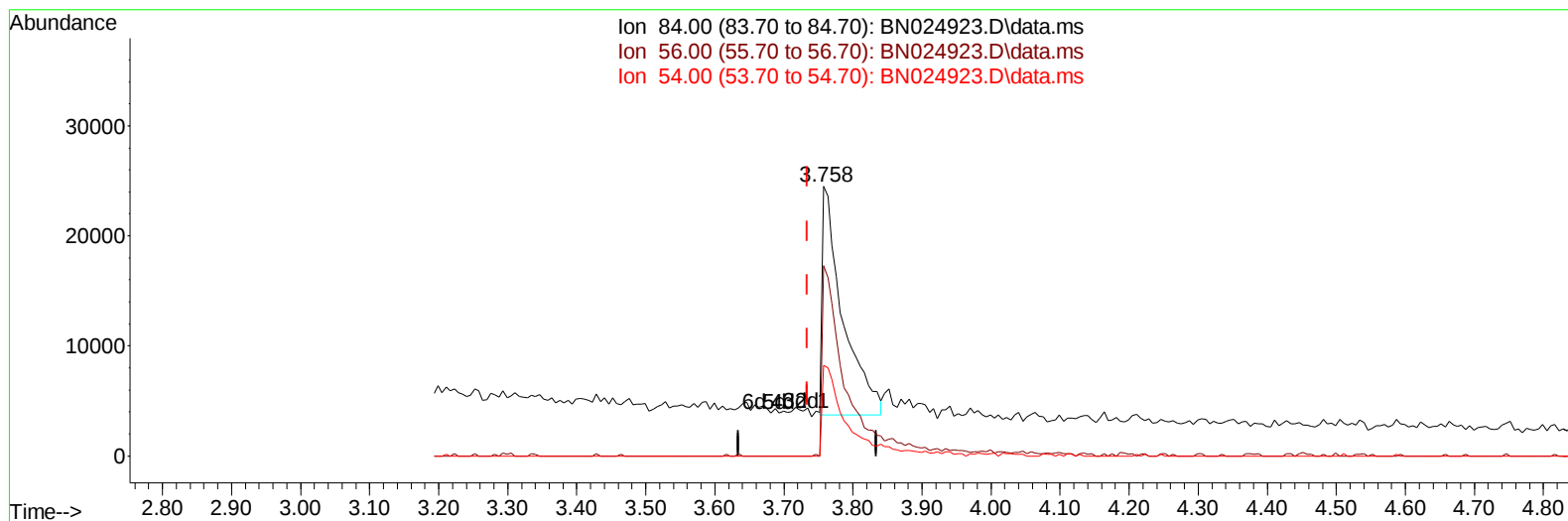
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TIC: BN024923.D\data.ms

(4) Pyridine-d5 (S)

3.758min (+ 0.023) 1.57 ng/u1 m

response 42406

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	70.53
54.00	40.20	33.71
0.00	0.00	0.00

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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	342387	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	1505703	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	955030	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	2125446	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	2000698	20.000	ng/ul	0.00
88) Perylene-d12	23.968	264	2044352	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	44986	5.001	ng/uL	0.00
4) Pyridine-d5	3.758	84	42406m	1.574	ng/ul	0.02
7) Phenol-d5	7.099	99	696634	20.869	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	567596	25.422	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	640169	26.096	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	522341	19.711	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	321643	26.383	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	351331	26.605	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	596878	25.288	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	17546	0.490	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	2007889	27.121	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1905267	22.457	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	358002	23.589	ng/ul	0.00
60) Fluorene-d10	15.575	176	1708720	27.335	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	315800	22.831	ng/ul	0.00
73) Anthracene-d10	17.428	188	2672771	26.775	ng/ul	0.00
81) Pyrene-d10	19.722	212	3120543	26.860	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	2446357	22.886	ng/ul	0.02

Target Compounds Qvalue

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

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