

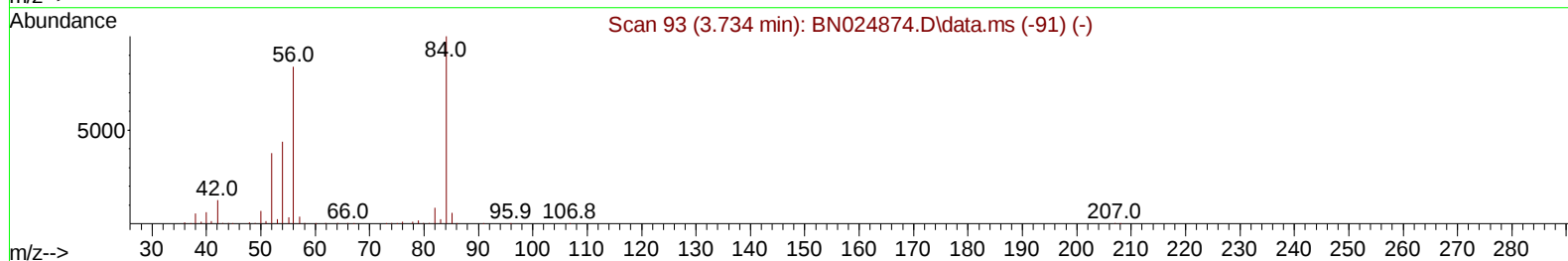
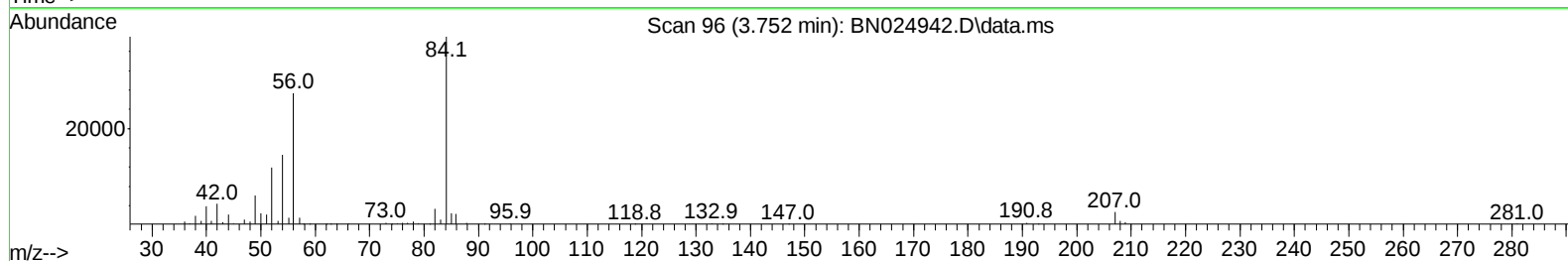
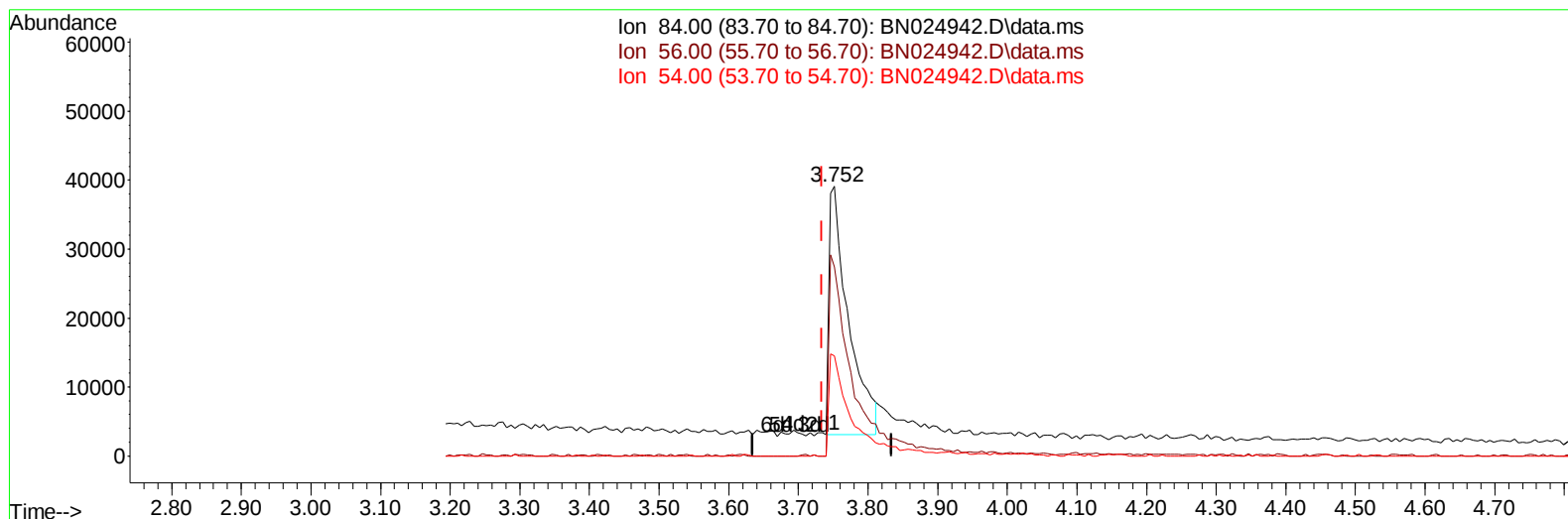
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024942.D
Acq On : 15 Apr 2023 20:07
Operator : CG/JU
Sample : PB152129BL
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK129

Manual IntegrationsAPPROVED

Quant Time: Apr 17 04:04:45 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



TIC: BN024942.D\data.ms

(4) Pyridine-d5 (S)

3.752min (+ 0.018) 2.76 ng/u1

response 69094

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	70.06
54.00	40.20	37.20
0.00	0.00	0.00

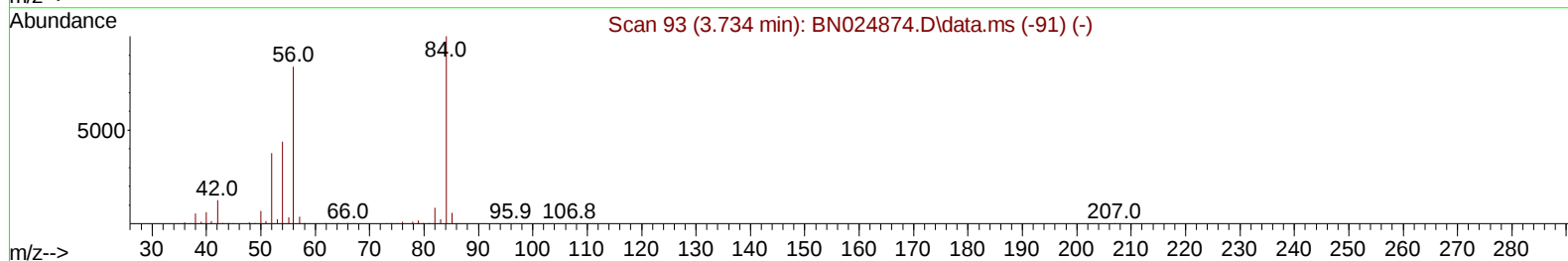
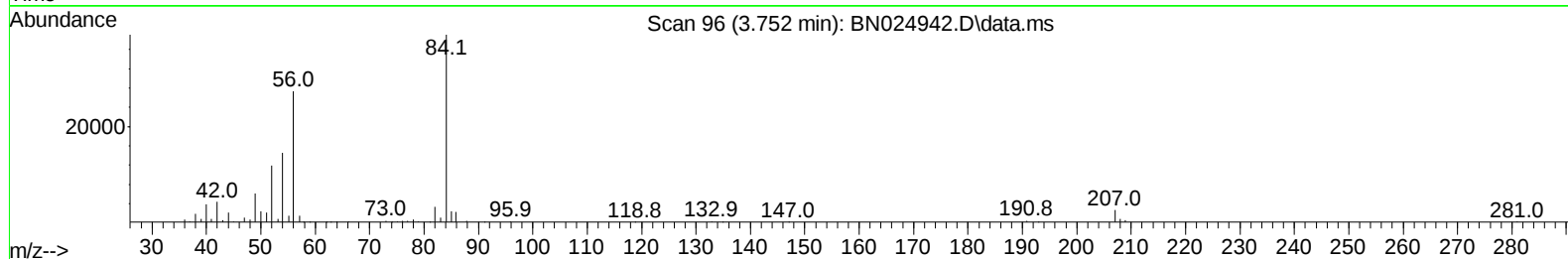
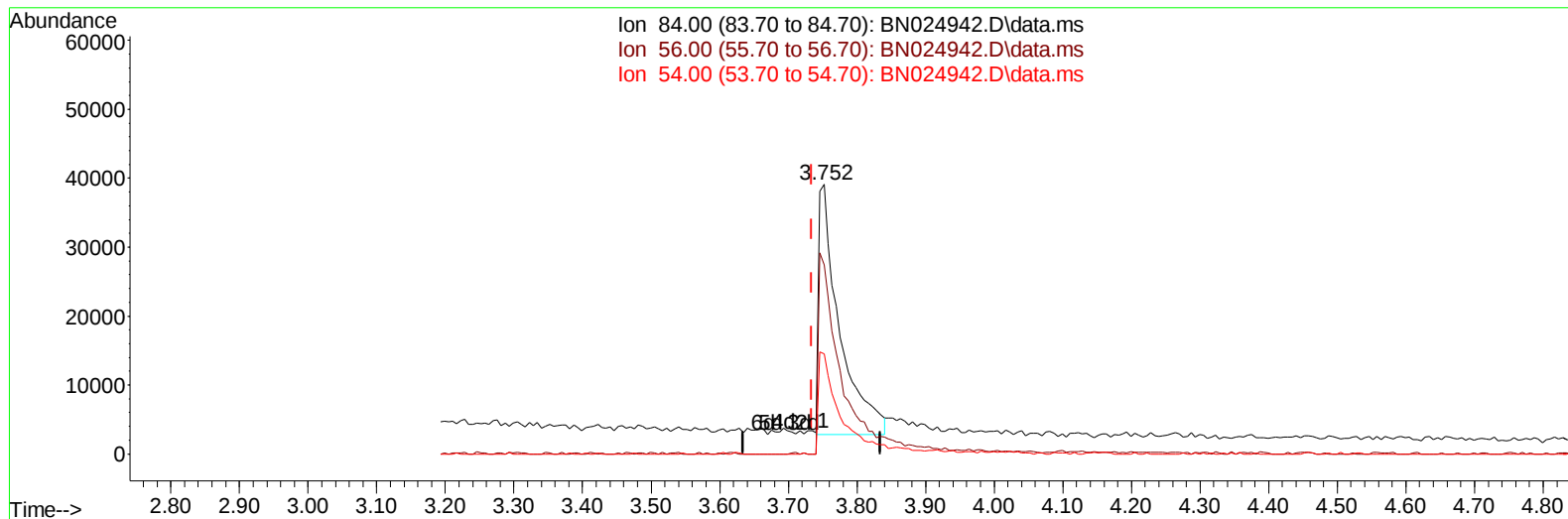
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(4) Pyridine-d5 (S)

3.752min (+ 0.018) 3.06 ng/u1 m

response 76695

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	70.06
54.00	40.20	37.20
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	7.934	152	318086	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	1395295	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	875884	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1929460	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	1865775	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264	1973322	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	46940	5.616	ng/uL	0.00
4) Pyridine-d5	3.752	84	76695m	3.065	ng/ul	0.02
7) Phenol-d5	7.093	99	843852	27.211	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	601390	28.994	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	676620	29.689	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	671938	27.293	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	338139	29.930	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	378523	30.933	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	633013	28.941	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	70341	2.119	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	2129274	31.359	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	2316221	29.768	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	379157	27.241	ng/ul	0.00
60) Fluorene-d10	15.569	176	1773644	30.938	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	336302	26.782	ng/ul	0.00
73) Anthracene-d10	17.428	188	2902736	32.032	ng/ul	0.00
81) Pyrene-d10	19.722	212	3338491	30.814	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	3174173	30.764	ng/ul	0.00

Target Compounds Qvalue

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

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