

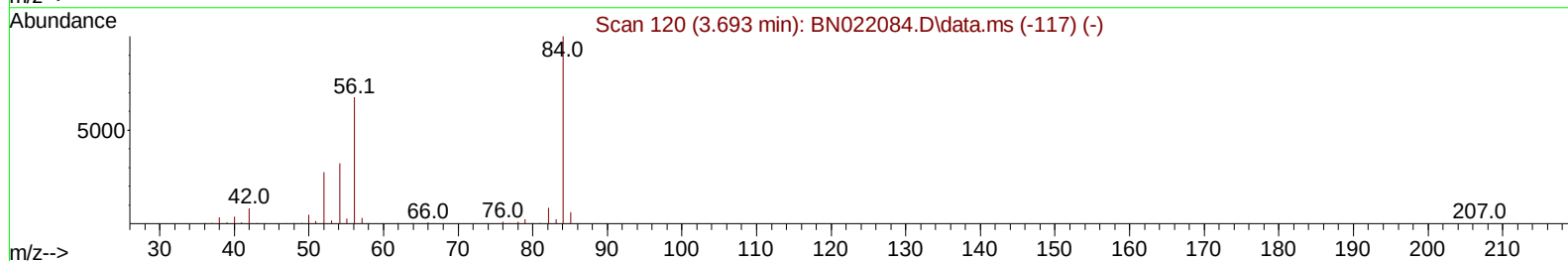
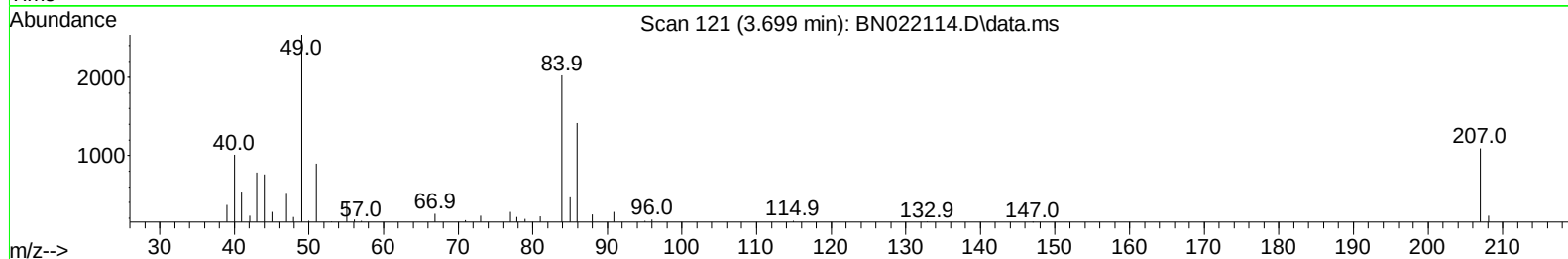
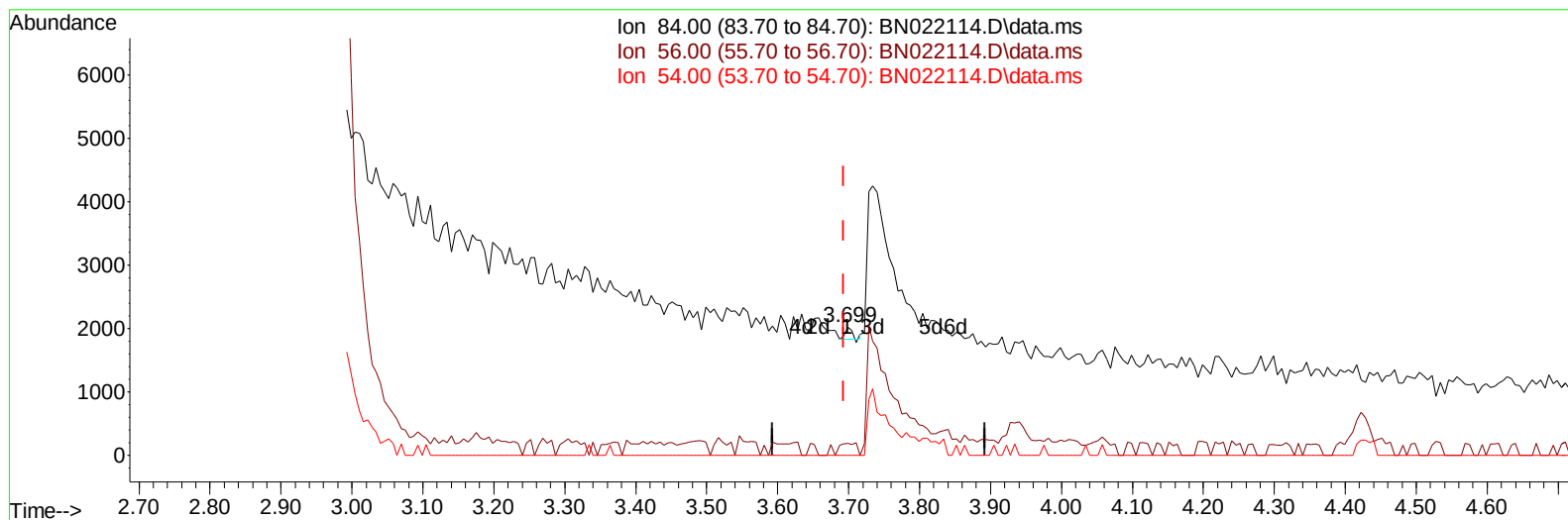
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101422\
 Data File : BN022114.D
 Acq On : 14 Oct 2022 08:00
 Operator : CG/JU
 Sample : N5049-12
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BP8

Manual IntegrationsAPPROVED

Quant Time: Oct 14 08:30:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 13 03:08:08 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/14/2022
 Supervised By :mohammad ahmed 10/16/2022



TIC: BN022114.D\data.ms

(4) Pyridine-d5 (S)

3.699min (+ 0.006) 0.01 ng/ul

response 138

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.60	9.15#
54.00	30.80	0.00#
0.00	0.00	0.00

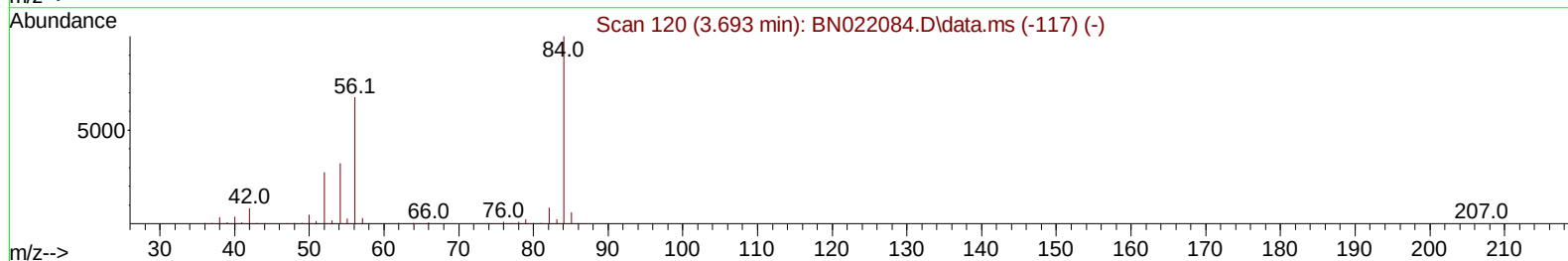
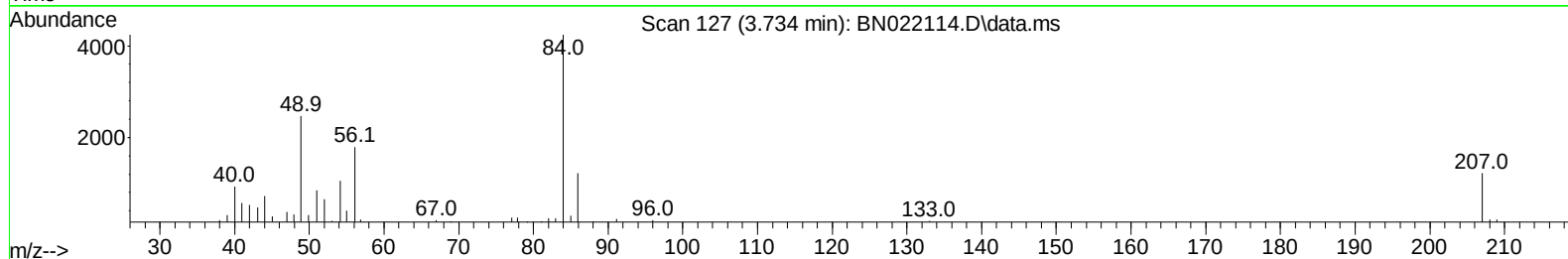
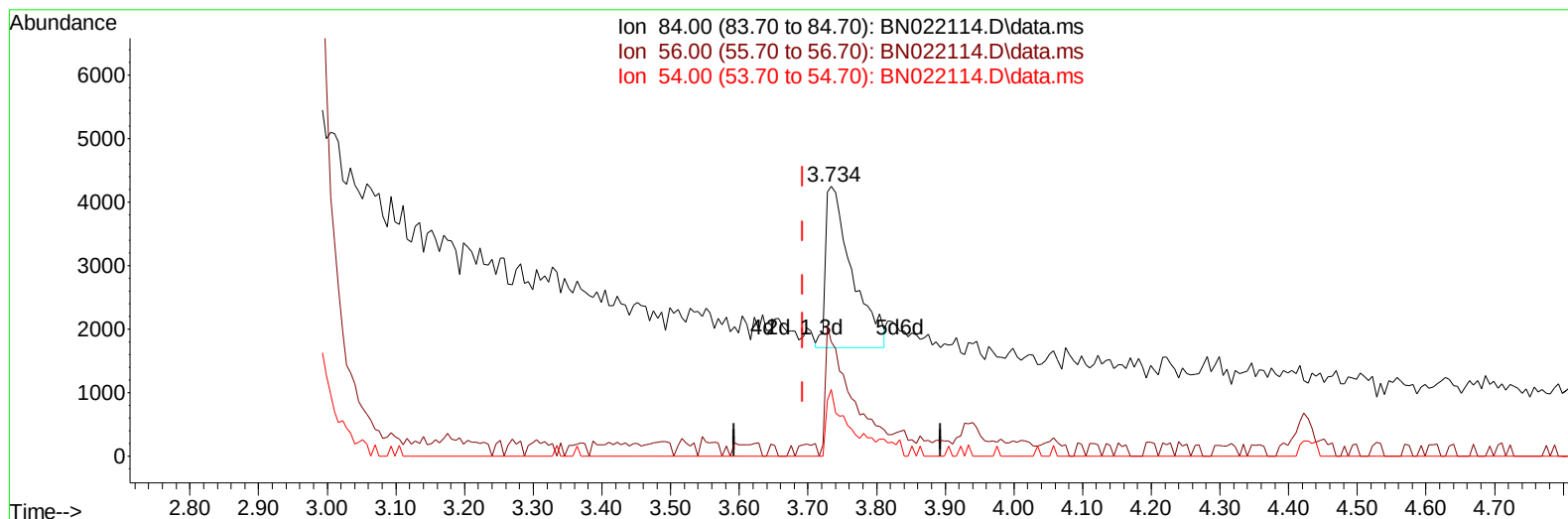
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(4) Pyridine-d5 (S)

3.734min (+ 0.041) 0.42 ng/ul m

response 6762

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.60	42.23#
54.00	30.80	24.75
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	7.951	152	194136	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	925018	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	595912	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.380	188	1251307	20.000	ng/ul	-0.01
79) Chrysene-d12	21.580	240	1174652	20.000	ng/ul	-0.01
88) Perylene-d12	24.051	264	1140502	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	18616	3.634	ng/uL	0.00
4) Pyridine-d5	3.734	84	6762m	0.424	ng/ul	0.04
7) Phenol-d5	7.110	99	391445	20.625	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	243320	20.460	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	290292	22.060	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	302403	20.052	ng/ul	-0.01
21) Nitrobenzene-d5	9.134	128	152615	22.500	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	169438	24.372	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	284353	21.790	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.928	131	358588	16.807	ng/ul	0.00
46) Dimethylphthalate-d6	14.039	166	937601	22.571	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	1049223	22.619	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	177361	20.003	ng/ul	0.00
60) Fluorene-d10	15.622	176	794465	22.675	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	125627	17.598	ng/ul	0.00
73) Anthracene-d10	17.480	188	1209818	22.248	ng/ul	0.00
81) Pyrene-d10	19.780	212	1426475	24.391	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.886	264	1302926	23.295	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.134	94	21063	1.046	ng/ul	95

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

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