

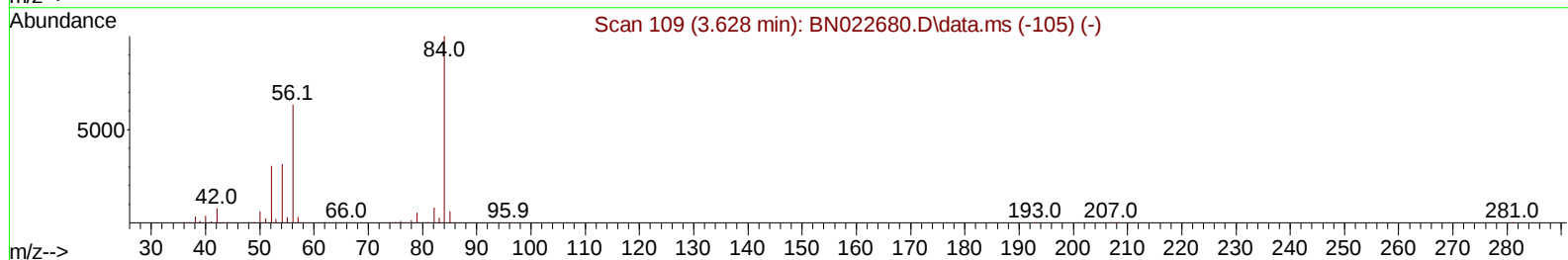
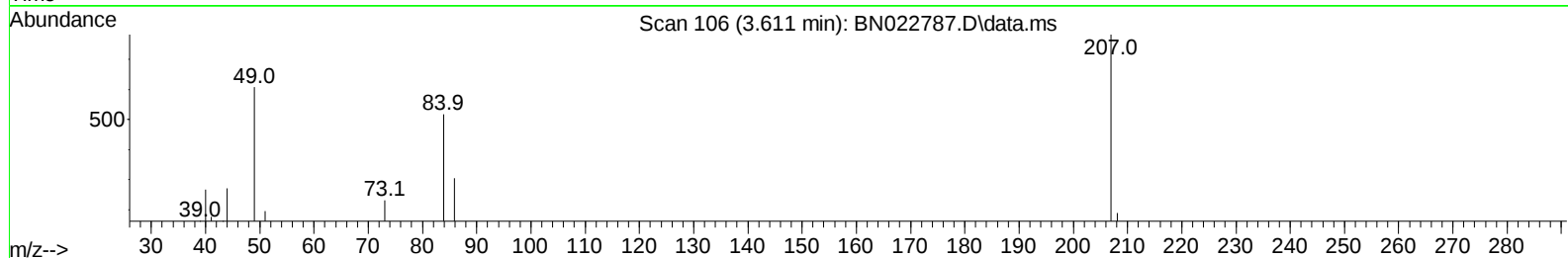
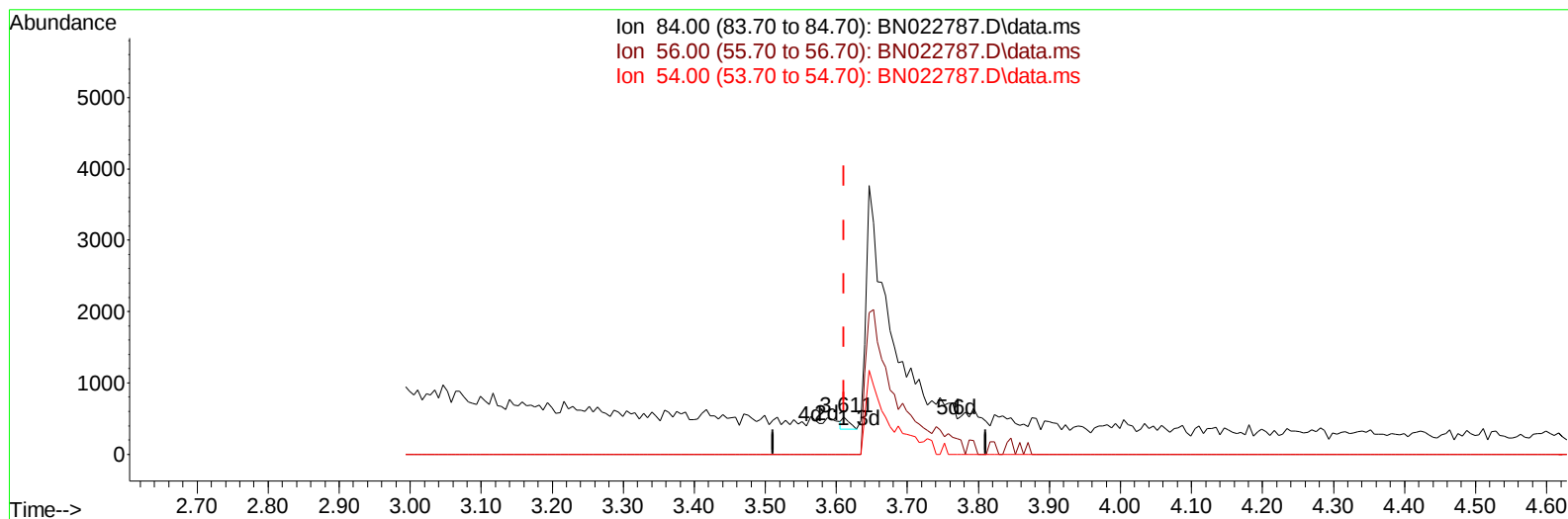
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
 Data File : BN022787.D
 Acq On : 21 Nov 2022 15:07
 Operator : CG/JU
 Sample : N5713-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGC79

Manual IntegrationsAPPROVED

Quant Time: Nov 21 15:43:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN11522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 19 05:51:00 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/22/2022
 Supervised By :mohammad ahmed 11/25/2022



TIC: BN022787.D\data.ms

(4) Pyridine-d5 (S)

3.611min (-0.000) 0.03 ng/u1

response 121

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	0.00#
54.00	31.90	0.00#
0.00	0.00	0.00

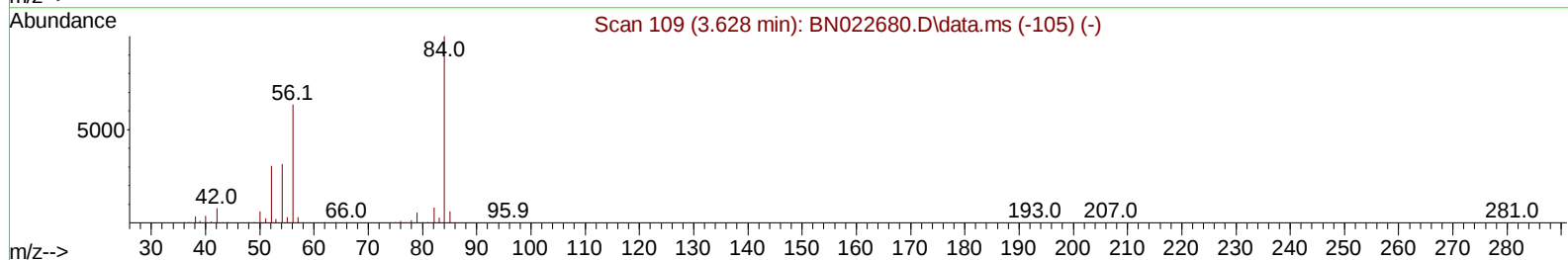
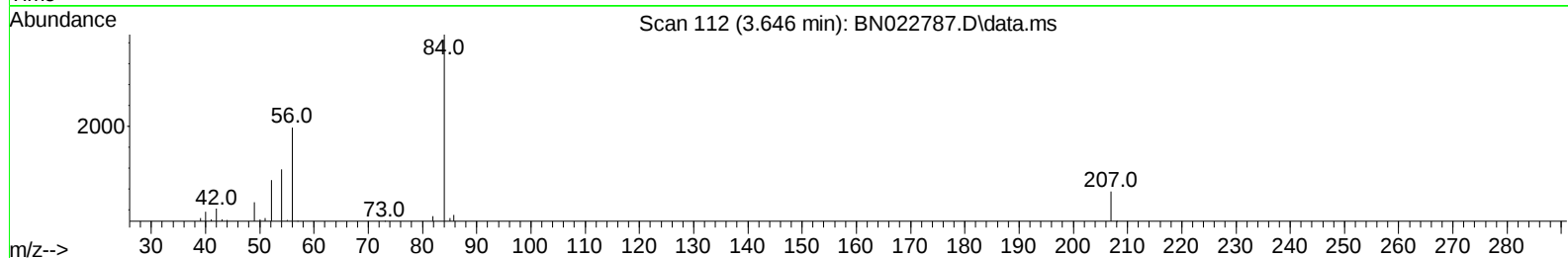
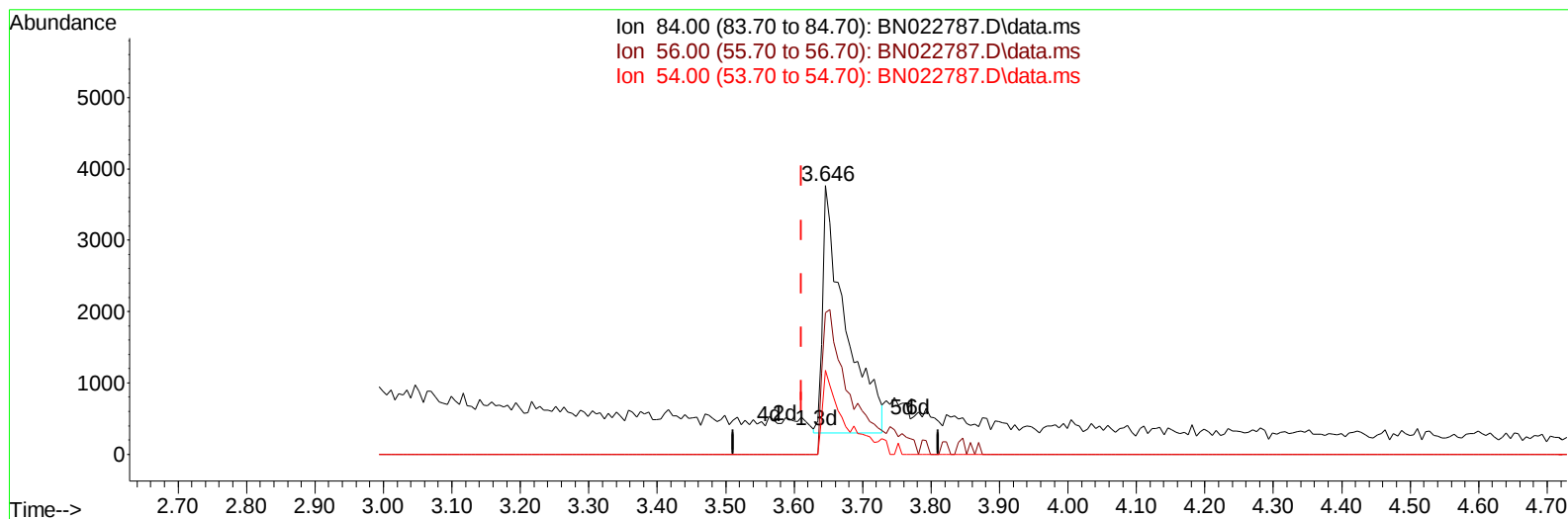
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(4) Pyridine-d5 (S)

3.646min (+ 0.035) 1.72 ng/u1 m

response 8016

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	52.74#
54.00	31.90	31.26
0.00	0.00	0.00

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 Acq On : 21 Nov 2022 15:07
 Operator : CG/JU
 Sample : N5713-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGC79

Manual IntegrationsAPPROVED

Quant Time: Nov 21 15:43:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 19 05:51:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	56342	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	274316	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	178845	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	381280	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	377925	20.000	ng/ul	0.00
88) Perylene-d12	23.869	264	375259	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	2727	1.652	ng/uL	0.00
4) Pyridine-d5	3.646	84	8016m	1.718	ng/ul	0.04
7) Phenol-d5	7.016	99	41963	7.886	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	33224	10.413	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	33130	8.501	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	34382	8.184	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	18386	8.659	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	19731	8.328	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	31814	7.877	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	34222	5.562	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	138742	10.666	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	124154	8.662	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	15973	6.159	ng/ul	0.01
60) Fluorene-d10	15.516	176	111370	10.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	14706	6.068	ng/ul	0.00
73) Anthracene-d10	17.375	188	162534	9.735	ng/ul	0.00
81) Pyrene-d10	19.675	212	179428	9.161	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	164083	8.698	ng/ul	0.00

Target Compounds Qvalue

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

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