

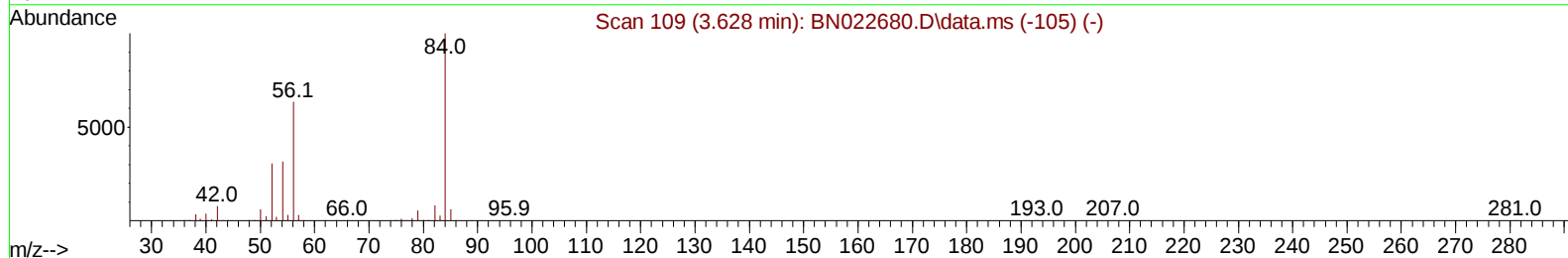
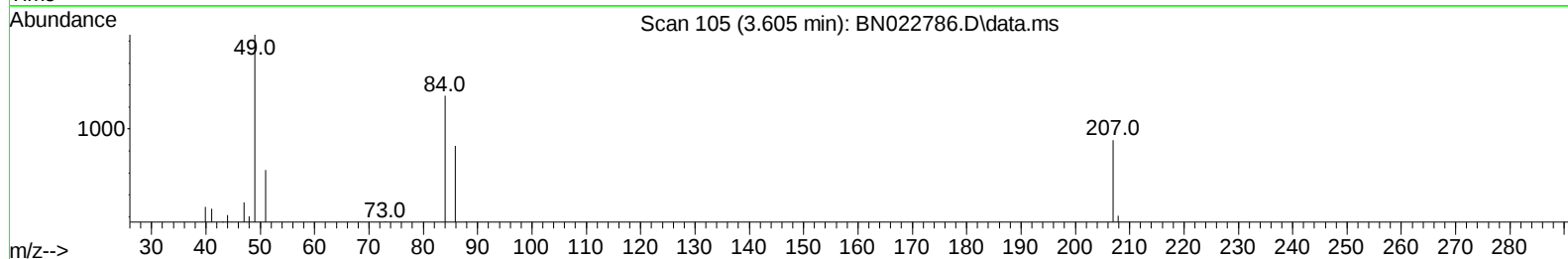
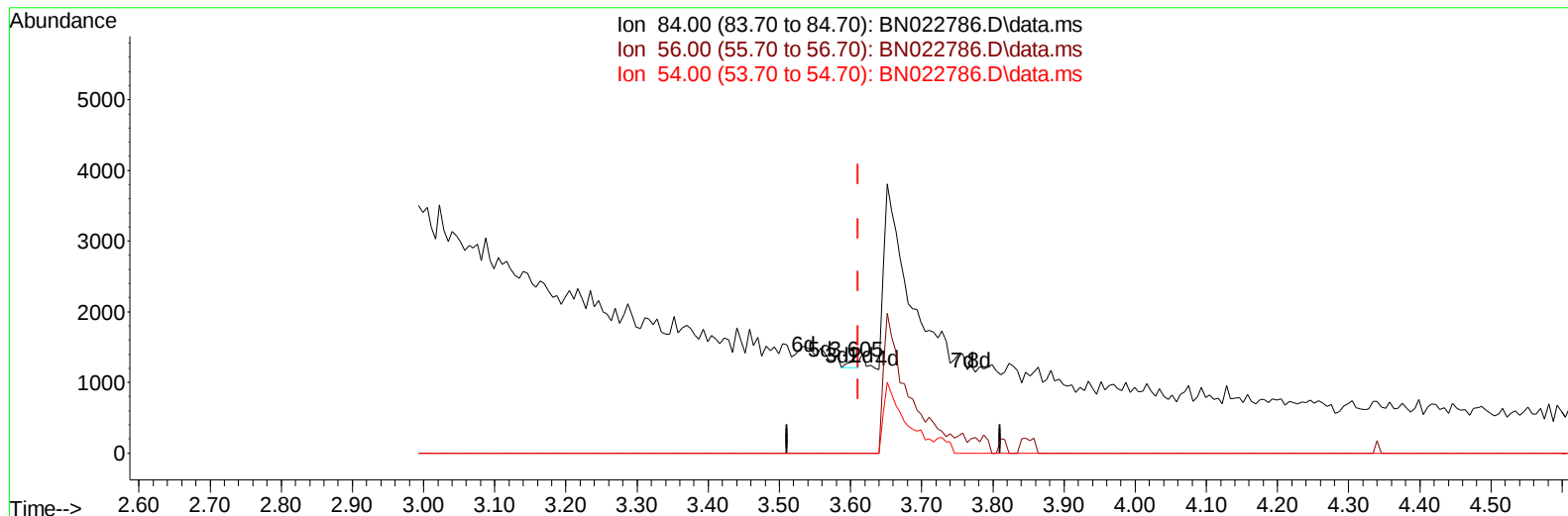
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
Data File : BN022786.D
Acq On : 21 Nov 2022 14:31
Operator : CG/JU
Sample : N5713-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGC80

Manual IntegrationsAPPROVED

Quant Time: Nov 21 15:26:10 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: BN022786.D\data.ms

(4) Pyridine-d5 (S)

3.605min (-0.006) 0.02 ng/u1

response 104

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

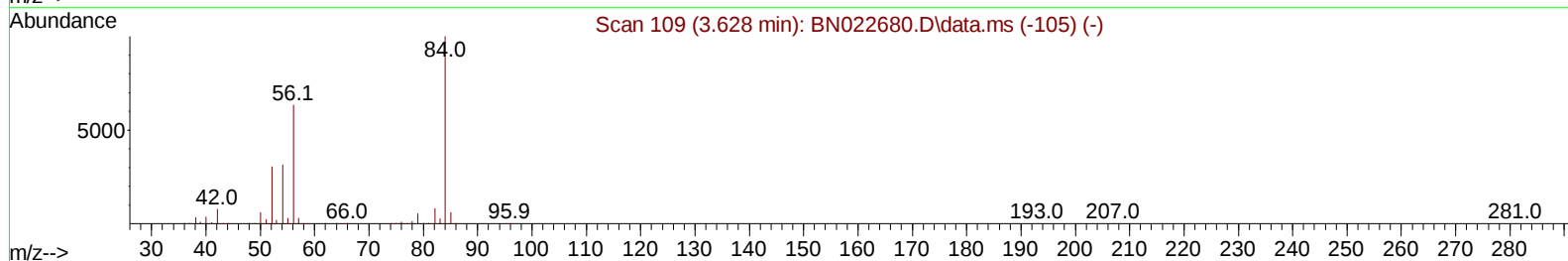
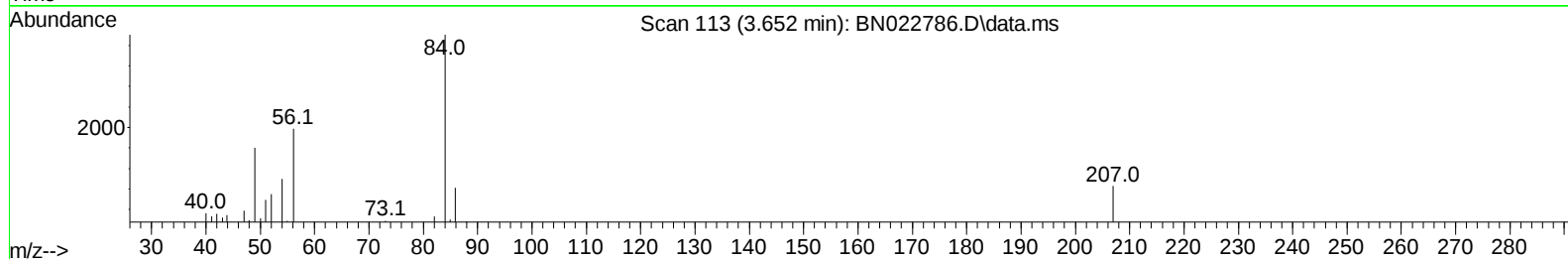
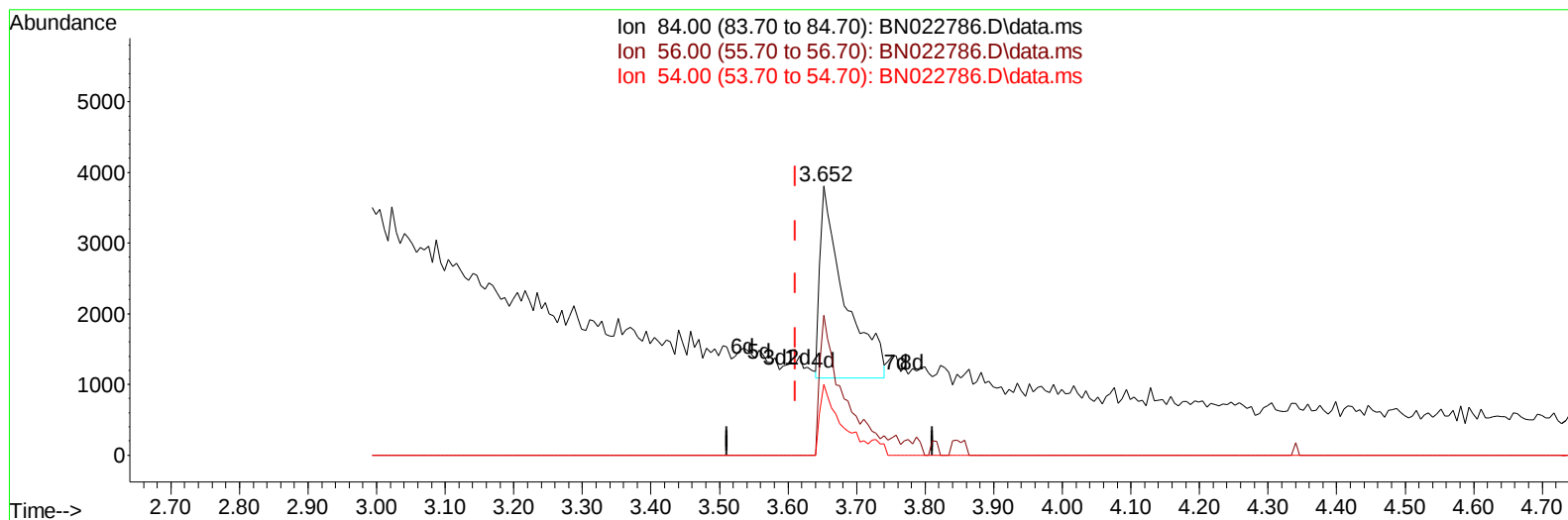
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TIC: BN022786.D\data.ms

(4) Pyridine-d5 (S)

3.652min (+ 0.041) 1.20 ng/ul m

response 6724

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	51.94#
54.00	31.90	26.22
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
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 Acq On : 21 Nov 2022 14:31
 Operator : CG/JU
 Sample : N5713-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGC80

Manual IntegrationsAPPROVED

Quant Time: Nov 21 15:26:10 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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 Supervised By :mohammad ahmed 11/25/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	67395	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	326483	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	215438	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	446068	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	442126	20.000	ng/ul	0.00
88) Perylene-d12	23.874	264	414804	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5497	2.785	ng/uL	0.00
4) Pyridine-d5	3.652	84	6724m	1.204	ng/ul	0.04
7) Phenol-d5	7.016	99	88454	13.896	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	67905	17.793	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	69539	14.917	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	61112	12.161	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	39976	15.819	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	43533	15.439	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	67650	14.073	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	39317	5.370	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	282928	18.055	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	278410	16.125	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	36093	11.554	ng/ul	0.00
60) Fluorene-d10	15.516	176	226432	17.618	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	37077	13.077	ng/ul	0.00
73) Anthracene-d10	17.375	188	334181	17.109	ng/ul	0.00
81) Pyrene-d10	19.675	212	365246	15.941	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	322211	15.452	ng/ul	0.00

Target Compounds Qvalue

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

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