

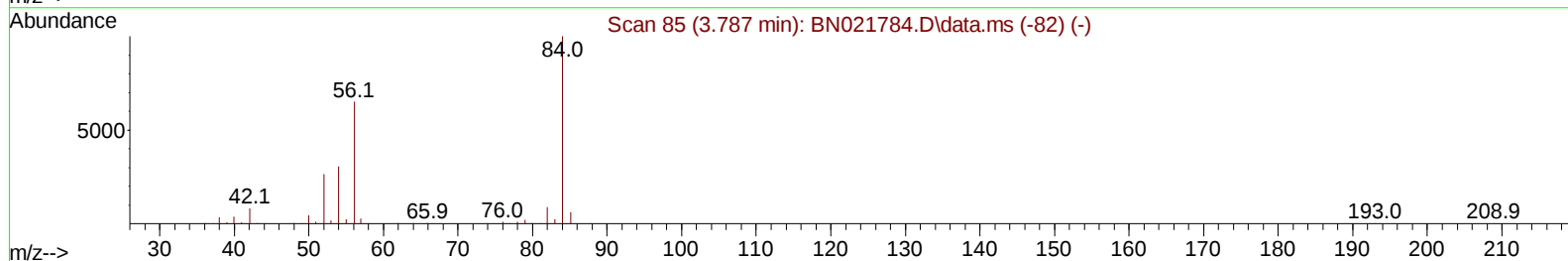
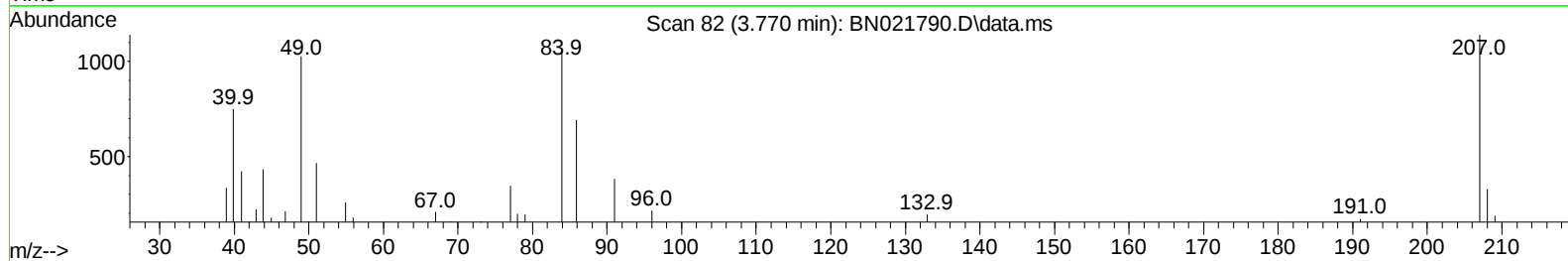
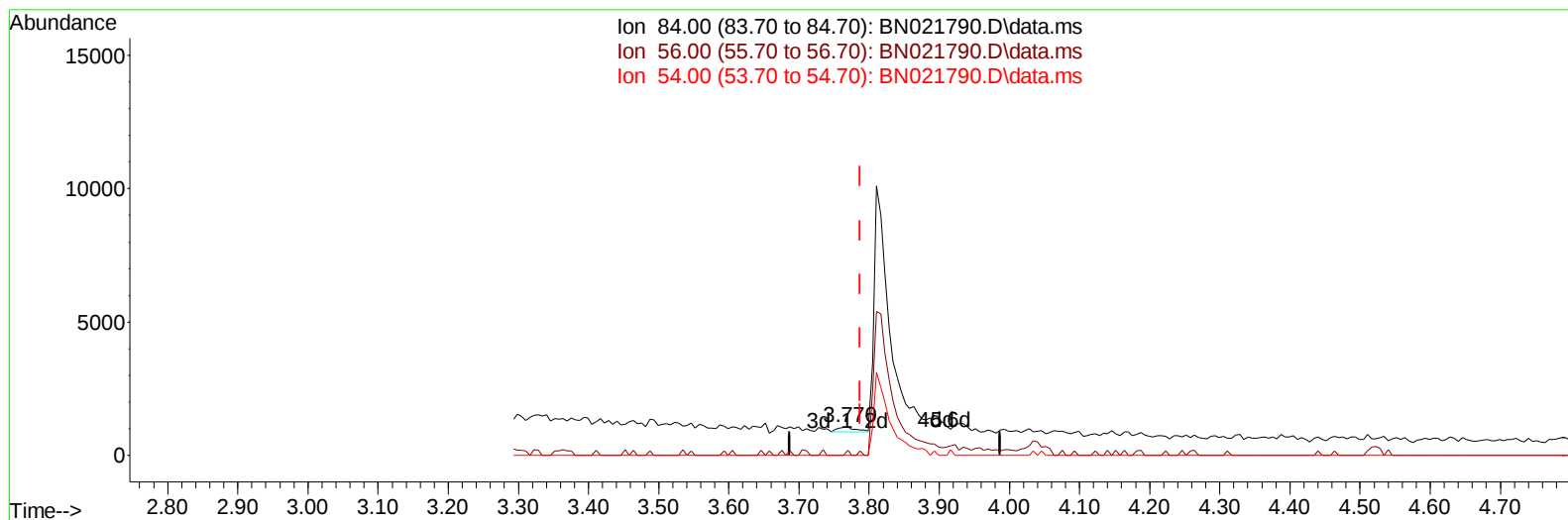
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091622\
 Data File : BN021790.D
 Acq On : 16 Sep 2022 15:54
 Operator : CG/JU
 Sample : N4601-15
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5756

Manual IntegrationsAPPROVED

Quant Time: Sep 17 00:22:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 17 00:17:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/19/2022
 Supervised By :mohammad ahmed 09/19/2022



TIC: BN021790.D\data.ms

(4) Pyridine-d5 (S)

3.770min (-0.018) 0.03 ng/u1

response 369

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	16.46#
54.00	33.00	0.00#
0.00	0.00	0.00

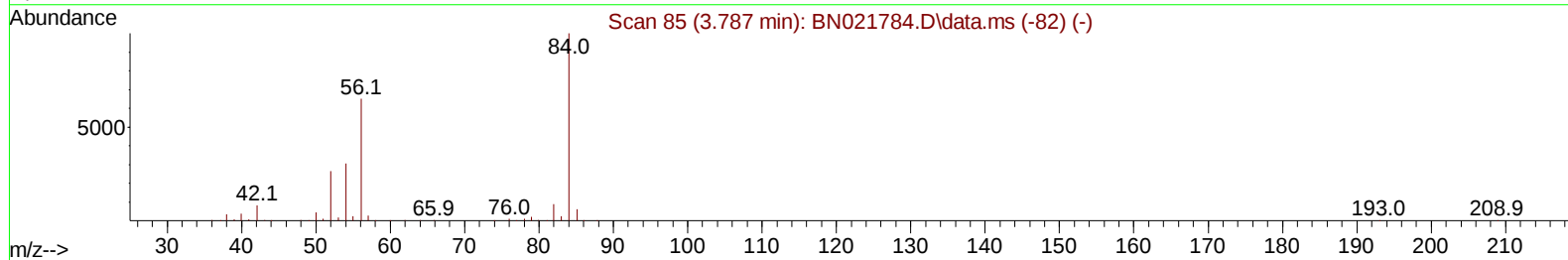
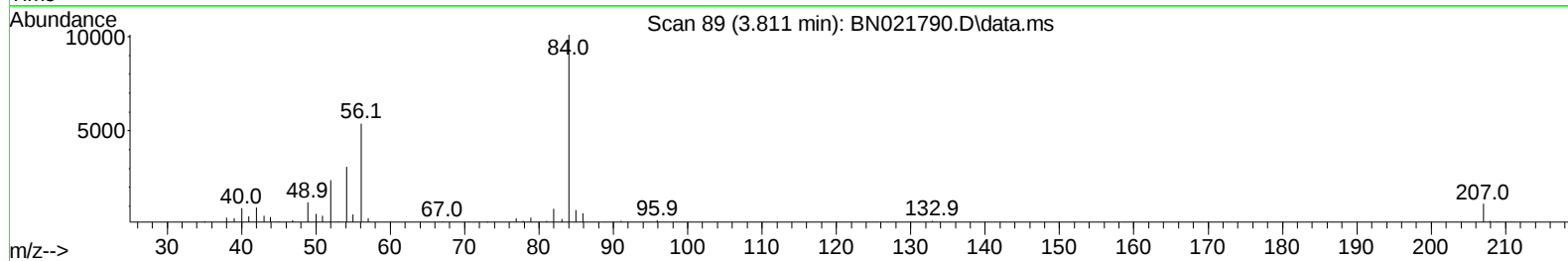
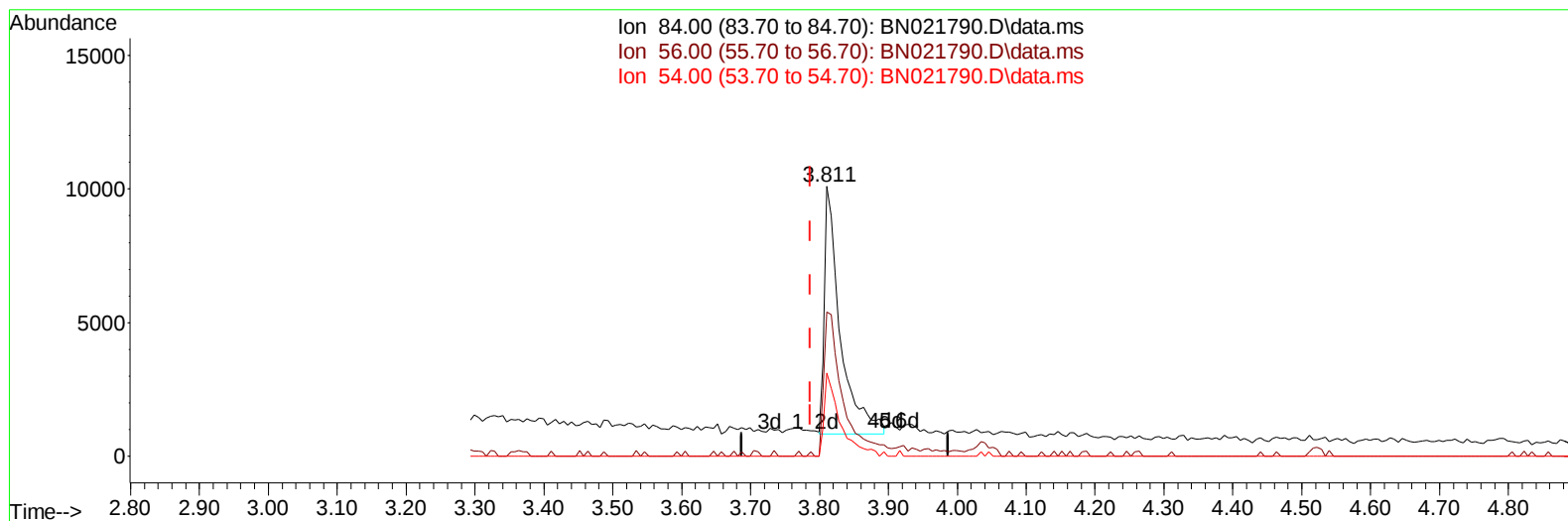
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(4) Pyridine-d5 (S)

3.811min (+ 0.024) 1.19 ng/u1 m

response 14843

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	53.43#
54.00	33.00	30.74
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	171278	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	768941	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	485319	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.457	188	981442	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240	725208	20.000	ng/ul	0.00
88) Perylene-d12	24.157	264	629279	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	18830	4.128	ng/uL	0.00
4) Pyridine-d5	3.811	84	14843m	1.194	ng/ul	0.02
7) Phenol-d5	7.205	99	365577	23.533	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	223266	23.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	293618	25.979	ng/ul	0.00
15) 4-Methylphenol-d8	8.763	113	303010	24.860	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	149523	28.541	ng/ul	0.00
24) 2-Nitrophenol-d4	9.952	143	166150	35.507	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.499	165	273035	26.347	ng/ul	0.00
31) 4-Chloroaniline-d4	11.022	131	391779	22.260	ng/ul	0.00
46) Dimethylphthalate-d6	14.116	166	869817	26.222	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	1031006	26.661	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	122505	18.559	ng/ul	0.00
60) Fluorene-d10	15.698	176	716968	24.840	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	94196	23.995	ng/ul	0.00
73) Anthracene-d10	17.557	188	1107669	25.611	ng/ul	0.00
81) Pyrene-d10	19.851	212	1181289	34.252	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.992	264	809644	27.145	ng/ul	0.00

Target Compounds Qvalue

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

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