

Fluorescent detection of silver ions in water with organic nano-aggregates

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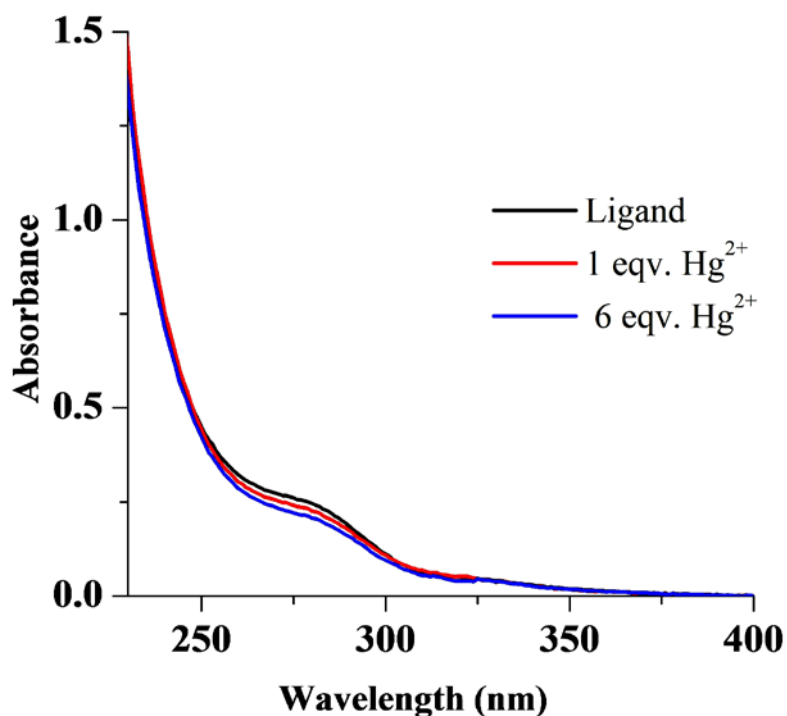


Figure (S1). UV-vis spectra : **1** with Hg^{2+} in $\text{H}_2\text{O}/\text{DMSO}$ (99:1).

Job's Plot (fluorescence method): Stock solutions of **1** (10 μM) and silver nitrate (1 mM) were prepared separately in respective solvents. The fluorescence spectra was recorded for each of the 11 solutions containing the two solutions (**1** and AgNO_3) at a total concentration of 80 μM in $\text{CH}_3\text{OH}/\text{DMSO}$ (99:1) and 8 μM for $\text{H}_2\text{O}/\text{DMSO}$ (99:1) in the following volume ratios. 2.0:0, 1.8:0.02, 1.6:0.04, 1.4:0.06, 1.2:0.08, 1.0:0.10, 0.8:0.12, 0.6:0.14, 0.4:0.16, 0.2:0.18, and 0.0:0.20.

S-3

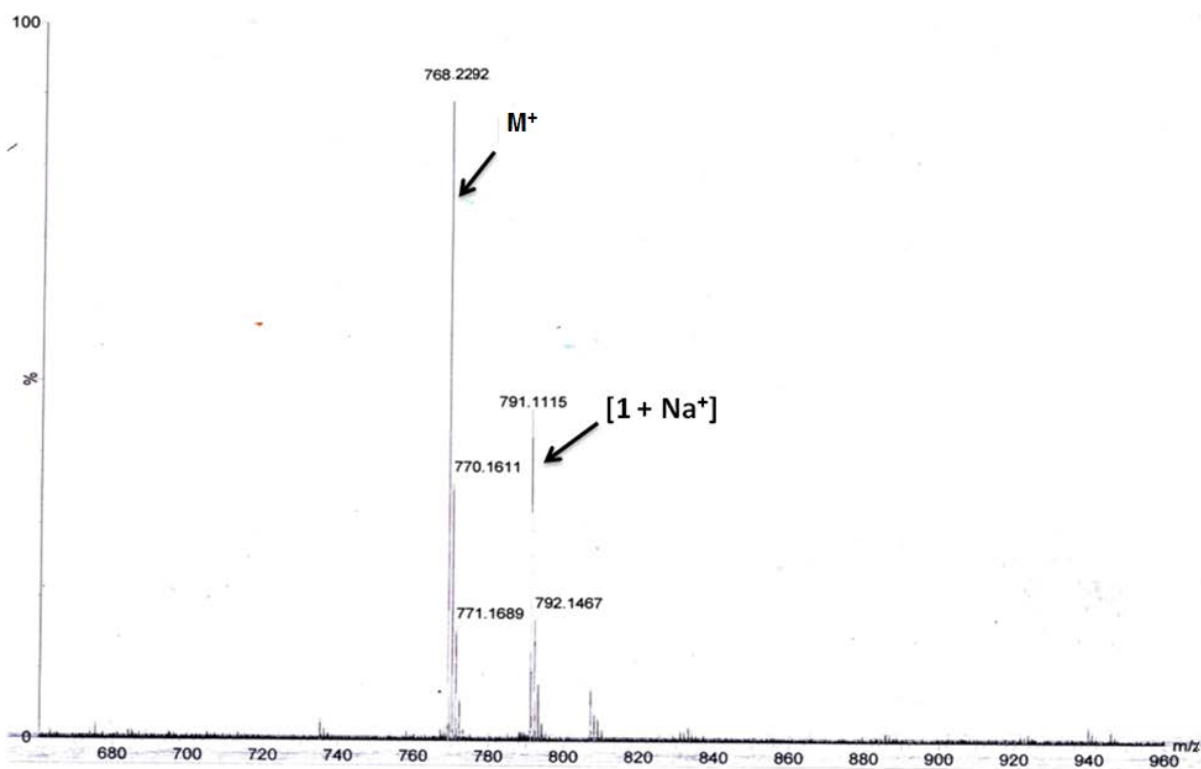


Figure (S2). ESI-MS of the chemosensor **1**.

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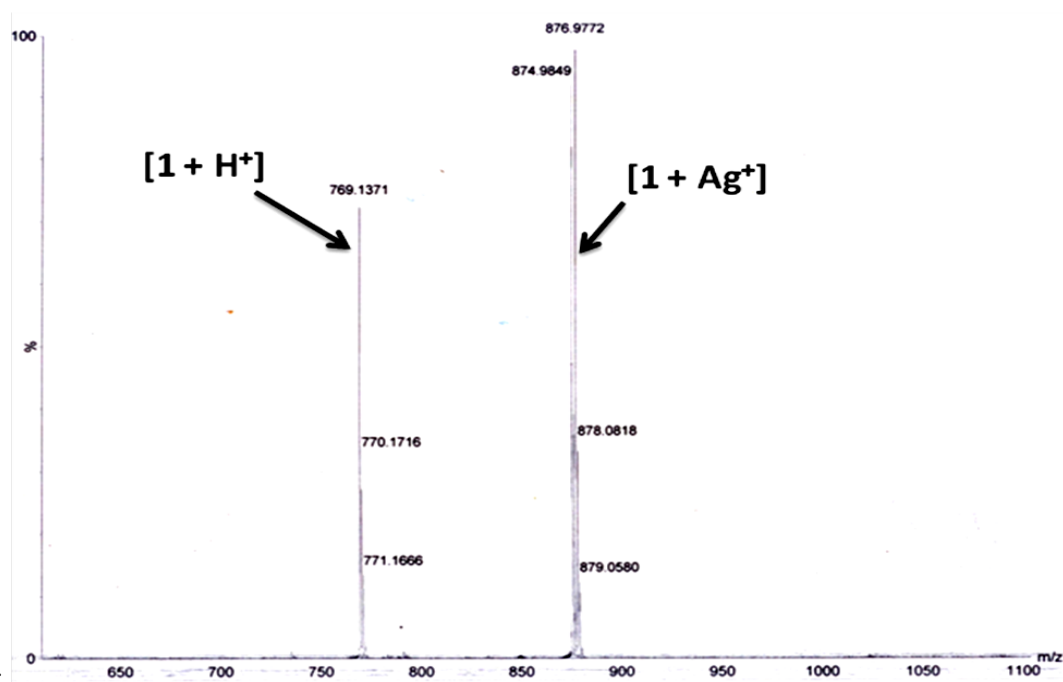
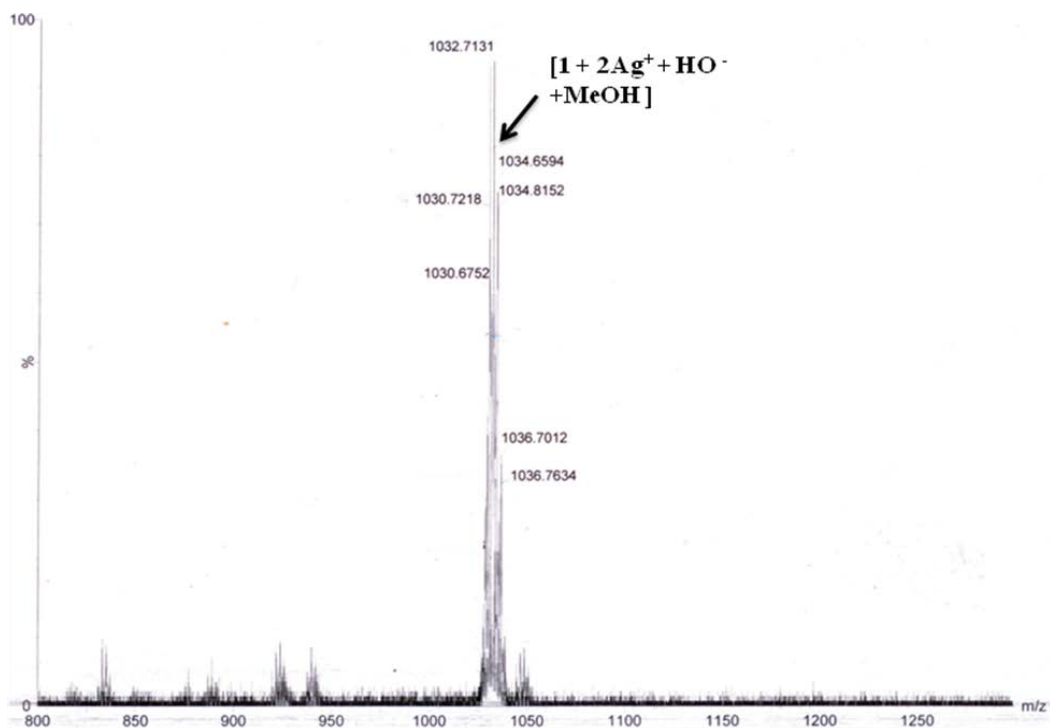


Figure (S3). ESI-MS spectrum of complex [1 + Ag⁺] from an aqueous sample.



Figure(S4). ESI-MS spectrum of complex [1 + 2Ag⁺ + MeOH + HO⁻].

S-5

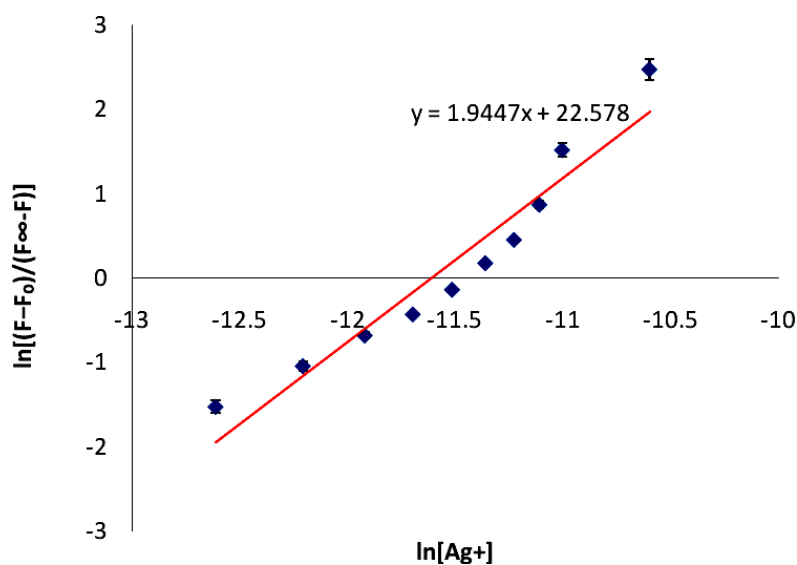


Figure (S5). Binding constant in MeOH:DMSO (99:1, v/v) was determined from the plot of $\ln[(F-F_0)/(F_\infty-F)]$ against $\ln[Ag^+]$; the stoichiometry of 1- Ag^+ association, obtained directly from the slope, is $1.94 \cong 2$. Following equation 1, the intercept gave an association constant of $8.0 \times 10^5 M^{-2}$. The error calculated from experiments performed in triplicate was 6%.

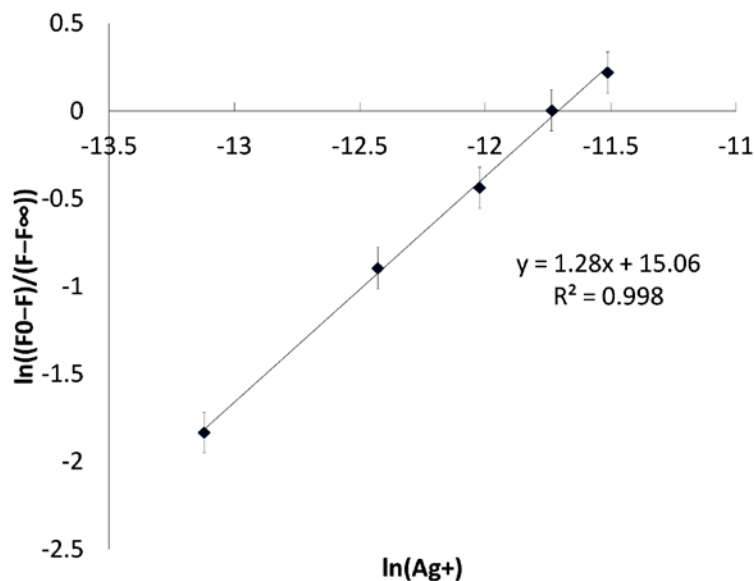
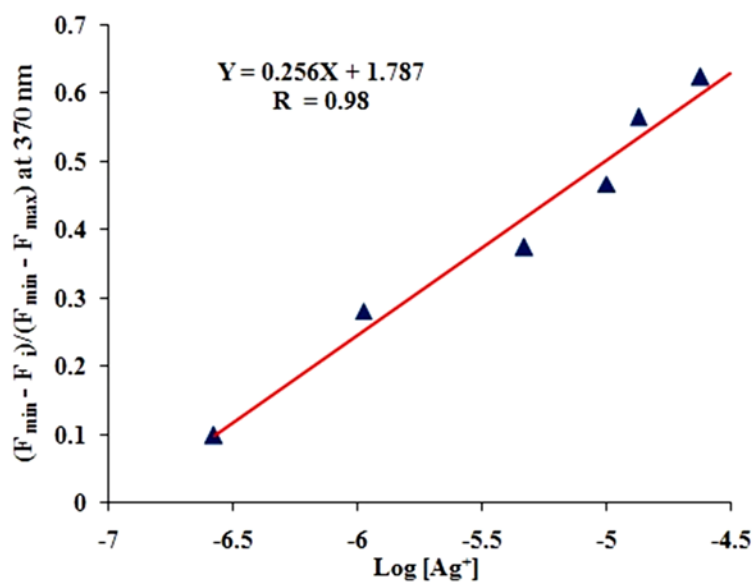


Figure (S6). Binding constant in H₂O:DMSO (99:1, v/v) was determined from the plot of $\ln[(F_0-F)/(F-F_\infty)]$ against $\ln[Ag^+]$; the stoichiometry of 1- Ag^+ association, obtained from the slope, is 1.28. Following equation 1, the intercept gave an association constant of $1.37 \times 10^6 M^{-1}$. The error calculated from experiments performed in triplicate was 4%.

S-6



Figure(S7). Determination of detection limit from fluorescence intensity of chemosensor **1** in H₂O/DMSO (99:1, v/v).

S-7

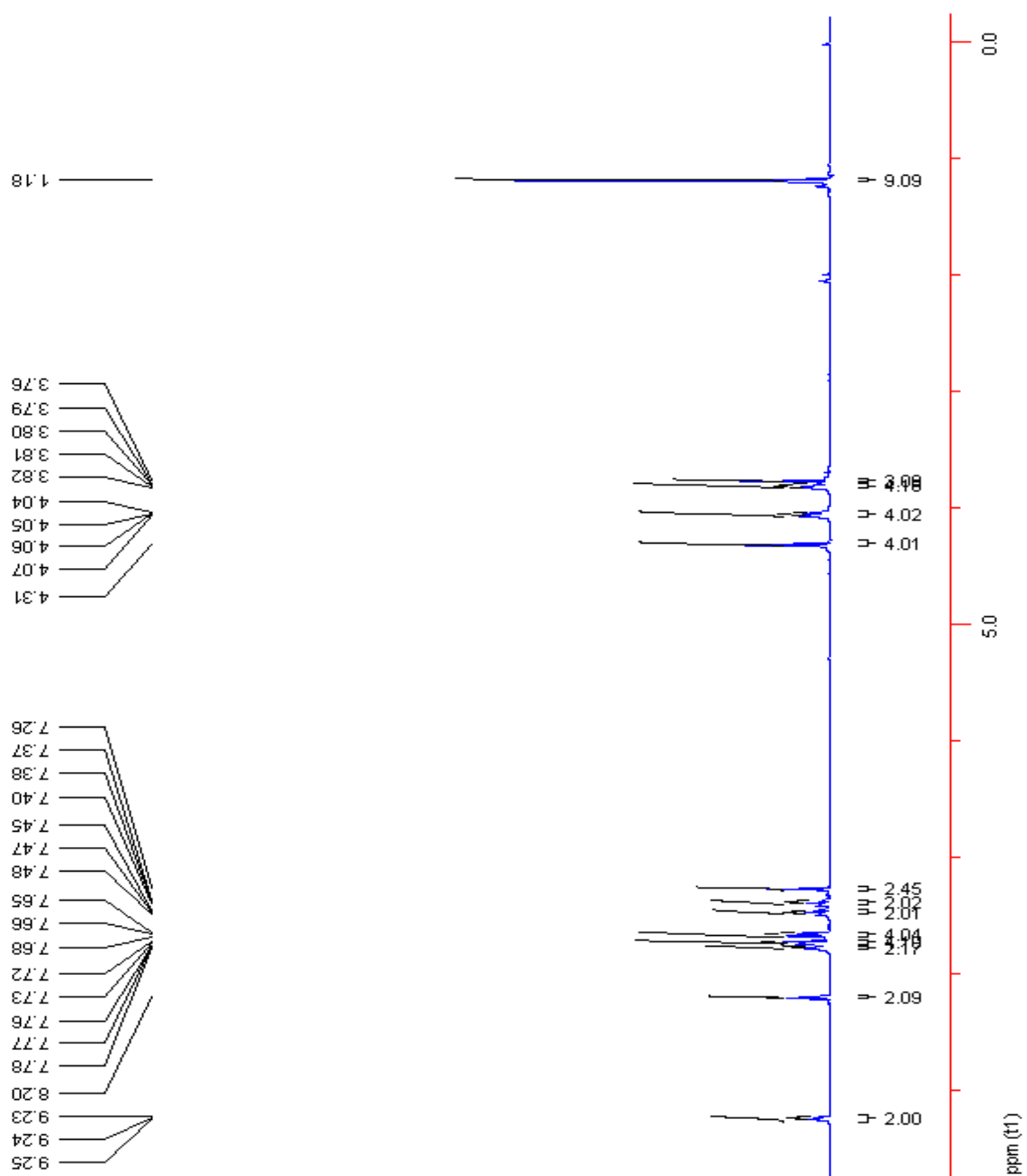


Figure (S8). $^1\text{H-NMR}$ spectrum of **1**.

S-8

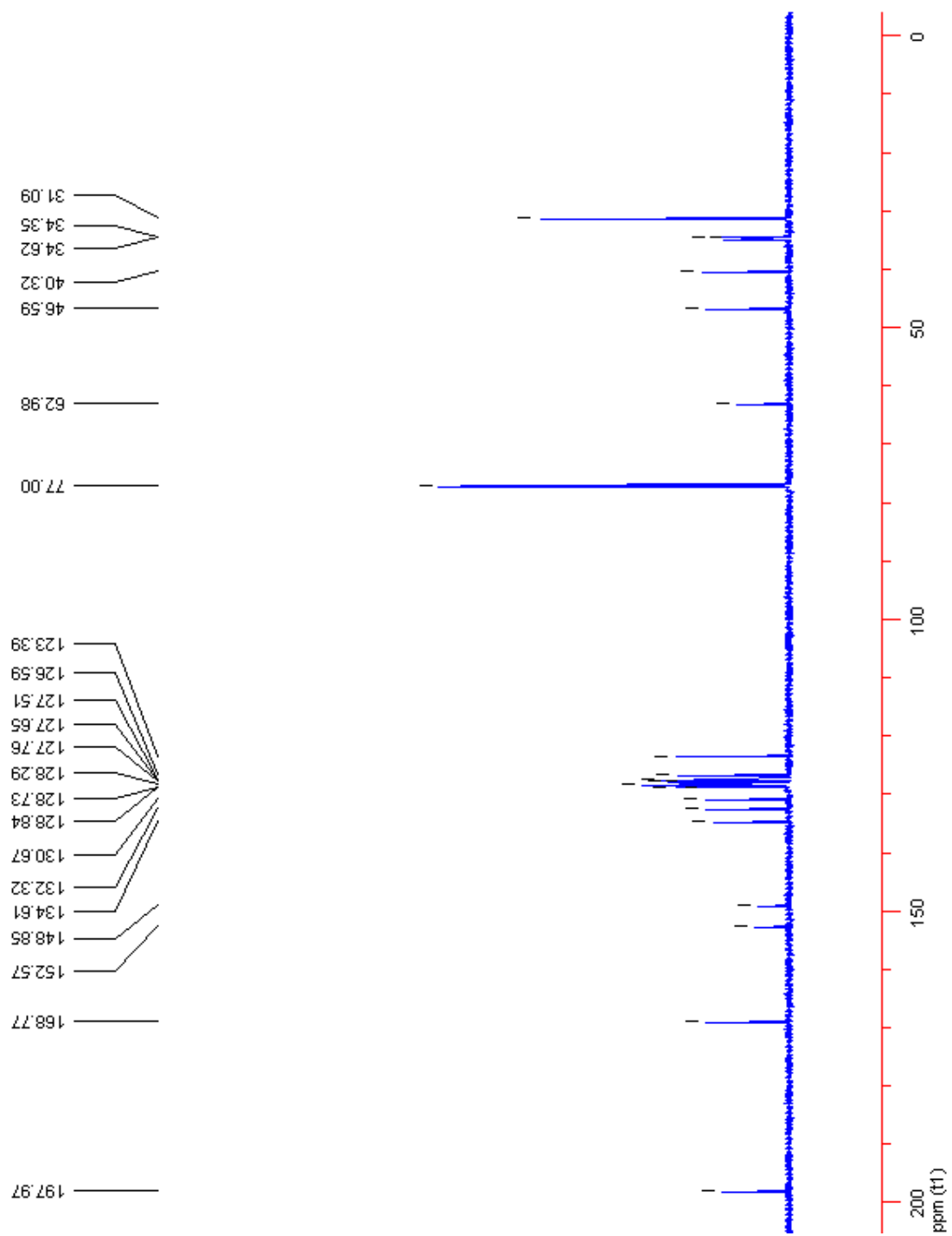


Figure (S9). ^{13}C -NMR spectrum of **1**.