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Supporting Information

Synthesis, "turn-on" fluorescent signals towards Zn²⁺ and Hg²⁺ and monoamine oxidase A inhibitory activity using molecular docking approach of morpholine analogue Schiff base linked organosilanes.

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(Fig.S2) ¹H NMR spectrum of compound 4b























(Fig.S10)¹³C NMR spectrum of compound 5a



(Fig.S12)¹³C NMR spectrum of compound 5c



(Fig.S13) Mass spectrum of compound 4c



(Fig.S14) Mass spectrum of compound 4b















(Fig.S18) Mass spectrum of compound 5c



(Fig.S19) (a) Absorption spectra of receptor 5a in the presence of various metal ions in MeOH– H_2O (9:1 v/v) solvent. UV-Visible absorption spectra after addition of increasing amount of (b) Zn^{2+} (c) Hg^{2+} in MeOH- H_2O (9:1 v/v) solution.



(Fig.S20) Linear regression plot (S20a-S20b) and Benesi–Hildebrand plot (S20c-S20d) of receptor 5a with Zn^{2+} and Hg^{2+} , using UV-Visible titration data, respectively.



(Fig.S21) Job's Plot for receptor 5a (a) Zn^{2+} and (b) Hg^{2+} .



(Fig.S23) ¹H NMR spectrum of compound 5a and Hg²⁺ complex