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Synthesis, Characterization, EPR Spectroelectrochemistry Studies and Theoretical Calculations of Manganese(II) Complexes with the Ligands H_3bpeten and H_3bnbpeten

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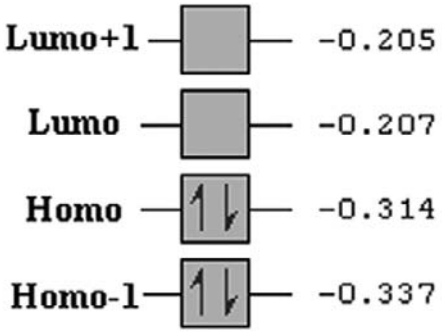

Figure S1. Molecular orbital energy levels for the frontier orbitals of the ligand H_3bnbpeten.

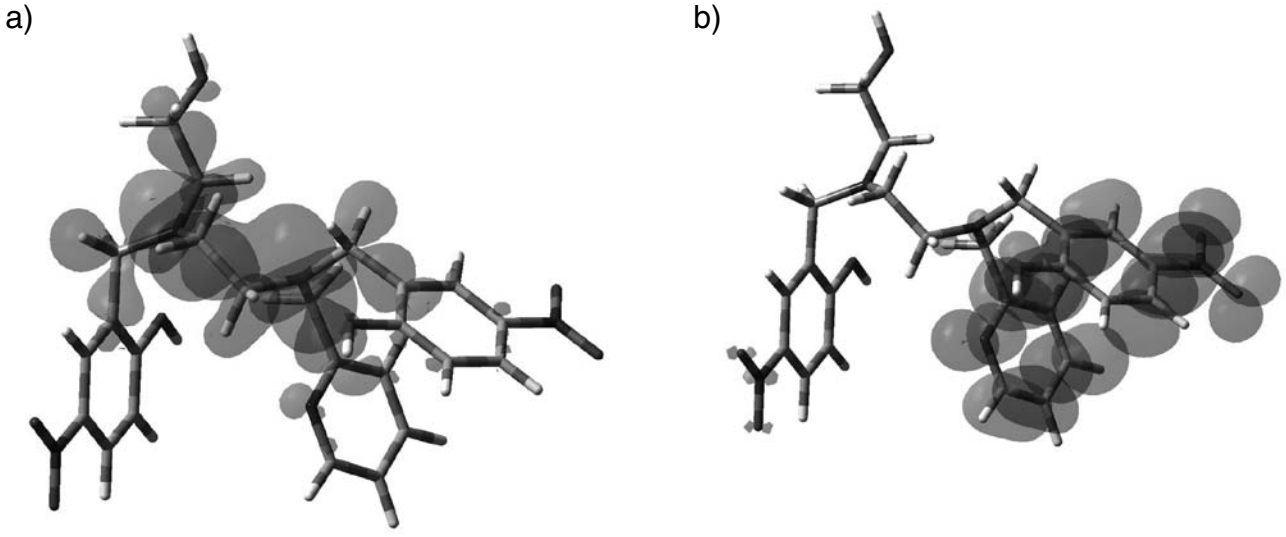

Figure S2. Graphical representation of the HOMO (a) and LUMO (b) of the H_3bnbpeten using Hartree-Fock.

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