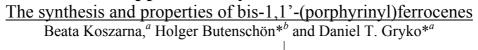
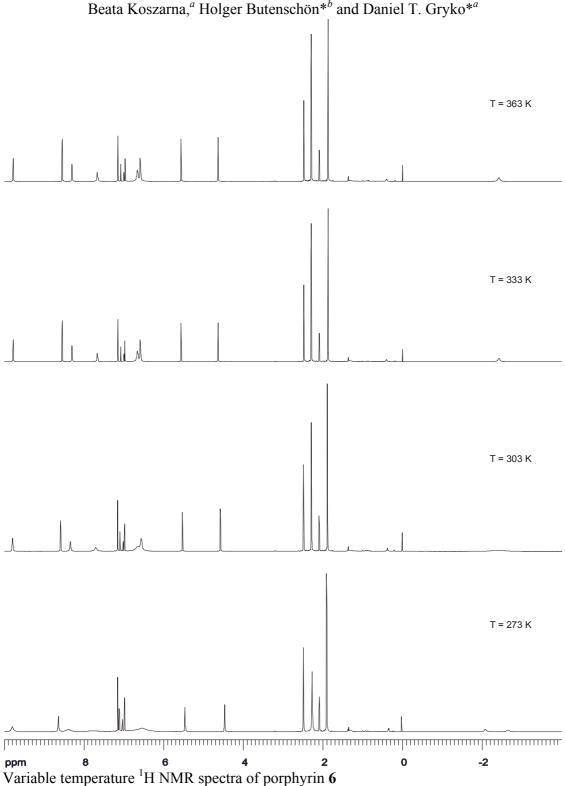
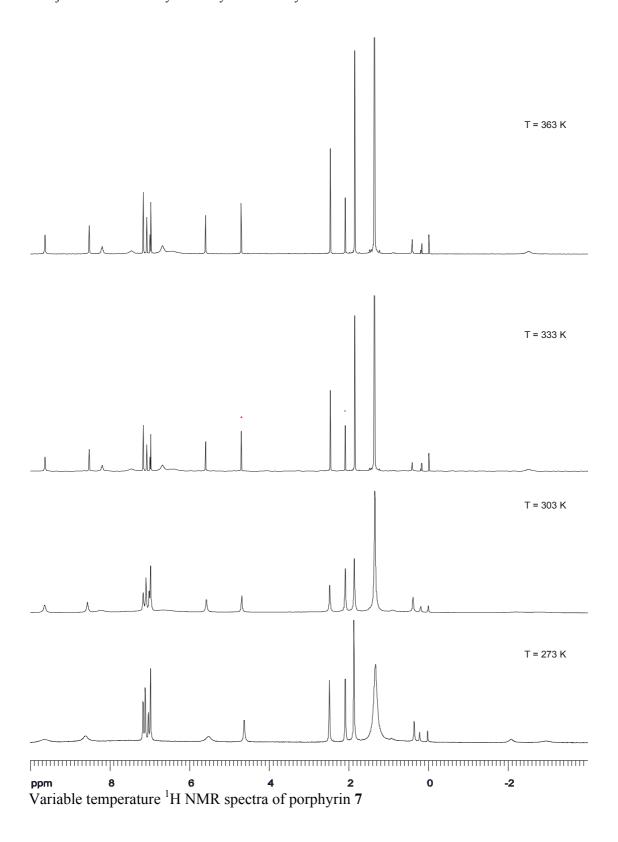
Supplementary information:



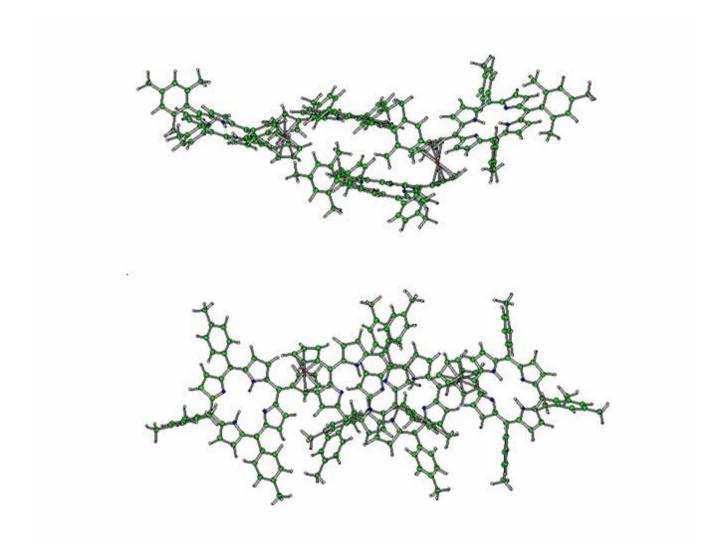




Molecular modeling of porphyrin 7 obtained from calculations performed within the HyperChem® Package by extensive energy minimizations (Polak-Ribiere conjugate gradient) alternating between molecular mechanics (MM+ force field) and semiempirical calculations (PM3).

Linear conformation

Kinked conformation



Dimer formation from linear conformation.