Supporting information

Binuclear Zinc Complexes with Radical-Anionic Diimine Ligands

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The geometry of the model system was optimized with the restricted (singlet) and unrestricted (triplet) formalism at the B3LYP/6-31G* level of theory with use of the GAUSSIAN 03 package.[1] After optimization NBO analysis was performed on the basis of the NBO 5.0 code.[2]

References

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