## Crystal Structure and Magnetic Properties of a New Hetero-Dinuclear Cu<sup>II</sup>Mn<sup>II</sup> Schiff Base Complex

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Mn[Cu(L)(O<sub>2</sub>CMe)<sub>2</sub>]·H<sub>2</sub>O (L = N,N'-bis(2-hydroxy-3-methoxybenzylidene)-1,3-diaminopropane) was synthesized and the crystal structure determined. (C<sub>23</sub>H<sub>26</sub>CuMnN<sub>2</sub>O<sub>8</sub>).H<sub>2</sub>O, monoclinic, space group  $P2_1/c$ , a = 12.017(3), b = 8.217(3), c = 24.786(4) Å,  $\beta = 92.10(2)^\circ$ , V = 2446(1) Å<sup>3</sup>, Z = 4. The crystal structure consists of ordered dinuclear units with Cu<sup>II</sup> and Mn<sup>II</sup> ions bridged by two oxygen atoms of the Schiff base ligand. The Cu<sup>II</sup> coordination sphere is a slightly distorted square-plane formed by the N<sub>2</sub>O<sub>2</sub> donor set of the Schiff base ligands. The average Cu—O and Cu—N distances are 1.920(1) and 1.957(4) Å, respectively. The coordination around the Mn<sup>II</sup> ion is a distorted tetrahedron with the donor oxygen atoms of the Schiff base ligands and oxygen atoms of the acetate anions. The Cu<sup>...</sup>Mn separation is 3.327(4) Å. There is also one non-coordinating water molecule in the crystal structure. The  $\chi$  and  $\chi$ T versus T plots,  $\chi$  being the molar magnetic susceptibility per Cu<sup>II</sup>Mn<sup>II</sup> unit and T the temperature, has been measured in the 4.9–301 K temperature range. The values of the interaction parameters are J = -28.3 cm<sup>-1</sup>, g<sub>Mn</sub> = 2.01, g<sub>Cu</sub> = 2.07. This indicates an intramolecular antiferromagnetic interaction between Cu<sup>II</sup> and Mn<sup>II</sup> ions.

*Key words:* Hetero-Dinuclear Cu<sup>II</sup>Mn<sup>II</sup> Complex, Super-Exchange Interactions, Antiferromagnetic Interaction