

# The Crystal Structures of [N,N'-Bis(3-methoxysalicylidene)-1,3-diaminopropane]nickel(II) and -copper(II)

Ayhan Elmali<sup>a</sup>, Celal T. Zeyrek<sup>b</sup>, and Yalcin Elerman<sup>a</sup>

<sup>a</sup> Department of Engineering Physics, Faculty of Engineering, University of Ankara, 06100 Besevler-Ankara, Turkey

<sup>b</sup> Ankara Nuclear Research and Training Center, Turkish Atomic Energy Authority, 06100 Besevler-Ankara, Turkey

Reprint requests to A. Elmali. E-mail: elmali@eng.ankara.edu.tr

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[N,N'-Bis(3-methoxysalicylidene)-1,3-diaminopropane]nickel(II) dihydrate [Ni(C<sub>19</sub>H<sub>20</sub>N<sub>2</sub>O<sub>4</sub>)·2(H<sub>2</sub>O)] **1** and [N,N'-bis(3-methoxysalicylidene)-1,4-diaminobutane]copper(II) [Cu(C<sub>20</sub>H<sub>22</sub>N<sub>2</sub>O<sub>4</sub>)] **2** have been synthesized and their crystal structures determined. Crystals of compound **1** are orthorhombic, space group *Pnma*,  $a = 7.509(3)$ ,  $b = 22.070(7)$ ,  $c = 11.532(4)$  Å,  $V = 1611.1(12)$  Å<sup>3</sup>,  $Z = 4$  and  $D_c = 1.498$  g·cm<sup>-3</sup>. The molecule **1** has mirror symmetry, but the ligand is not planar. The two parts of the Schiff base moieties are folded so as to form an angle of 21.6(1)°. The Ni atom is in a distorted octahedral geometry and coordinated by the donor atoms of the ligand in the horizontal plane and of two water molecules. Crystals of compound **2** are monoclinic, space group *P2<sub>1</sub>/c*,  $a = 9.488(1)$ ,  $b = 21.918(3)$ ,  $c = 8.413(1)$  Å,  $\beta = 91.45(1)^\circ$ ,  $V = 1749.0(4)$  Å<sup>3</sup>,  $Z = 4$  and  $D_c = 1.587$  g·cm<sup>-3</sup>. The Cu atom is coordinated by an N<sub>2</sub>O<sub>2</sub> donor set from the imine-phenol ligand in a distorted planar geometry, with the two phenolate O atoms deprotonated. The Cu–O bond lengths are 1.854(3) and 1.868(3) Å. The Cu–N bond lengths are 1.931(3) and 1.950(3) Å, the dihedral angle between the two 3-methoxysalicylidene groups is 43.4(1)°.

**Key words:** Schiff Base Complexes, Nickel(II) Complex, Copper(II) Complex, Square-Planar Coordination, Octahedral Coordination