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SYNTHESIS AND MOLECULAR STRUCTURE OF NEW COMPLEX OF COPPER

Javadzade T.A.^{1,2}, Musayev F.N.³, Aliyeva F.S.¹, Sujayev A.R.⁴, Chiragov F.M.¹

⁽¹⁾ *Baku State University*

AZ 1148, Azerbaijan, Baku, Z. Xalilov 23

⁽²⁾ *Khazar University*

AZ1096, Baku, Azerbaijan, Mabsati 41

⁽³⁾ *Virginia Commonwealth University*

VA 23284, USA, Richmond, 907 Floyd Avenue

⁽⁴⁾ *Institute of Chemistry of Additives*

AZ1029, Baku, Azerbaijan, H. Javid Avenue 102

The synthesis of the copper complex of 4-((2,4-dioxopentan-3-yl) diazenyl) benzoic acid was carried out according to the following methodology: The ligand (4-((2,4-dioxopentan-3-yl) diazenyl) benzoic acid) and the metal salt ((CH₃COO)₂Cu) were taken from the same concentration solutions in a 1:2 ratio and mixed with a magnetic stirrer for 30 minutes, then a solution of 2 mol of ethylene diamine in alcohol was added to it and stirred for 1 hour in a heater equipped with a reflux condenser. After the synthesis, the colored solution obtained was poured into a beaker and stored. After 48 hours, needle-shaped crystals were formed at the bottom of the container.

The structure of the complex was studied on an XtalLAB Synergy, Dual ex, HyPix, "BRUKER D8 ADVANCE" X-ray diffractometer ($\lambda = 1.54184 \text{ \AA}$, $T = 100 \text{ K}$). Analysis of crystallographic parameters shows that C₂₆H₃₄CuN₆O₁₀ belongs to the P-1 space group and crystallizes in the triclinic syngony. The degree of correspondence of the analyzed crystal to the true structure is $R = 0.0603$, which indicates that the structure is accurately determined. The lattice parameters and molecular structure of complexes is shown below:

$$a = 7.3645(3) \text{ \AA}$$

$$b = 14.2102(8) \text{ \AA}$$

$$c = 15.2334(9) \text{ \AA}$$

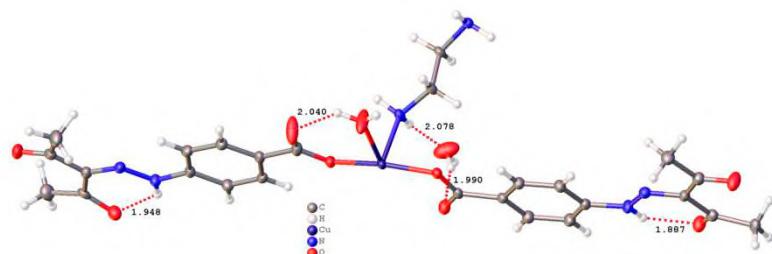
$$\alpha = 110.626(6)^\circ$$

$$\beta = 96.088(4)^\circ$$

$$\gamma = 99.886(4)^\circ$$

$$\text{Volume (V)} =$$

$$1445.67(15) \text{ \AA}^3$$



Molecular structure of C₂₆H₃₄CuN₆O₁₀

During the crystallization of the complex compound, in addition to two molecules of C₁₂H₁₃TN₂O₄, one molecule of ethylenediamine and two molecules of water are coordinated around the complexing agent.

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