MS30-P8 Three dimensional Zinc(II) and Cadmium(II)-Metal Organic Frameworks (MOFs) based on

3,3',5,5'-Azobenzenetetracarboxylate and bis(imidazole) ligands

Okan Zafer YEŞİLEL $^{\rm I},$ Mürsel ARICI $^{\rm I},$ Murat TAŞ $^{\rm 2},$ Hakan ERER $^{\rm I}$

- 1. Department of Chemistry, Eskişehir Osmangazi University, 26480 Eskişehir, Turkey
- 2. Department of Science Education, Ondokuz Mayıs University, 55139, Samsun, Turkey

email: yesilel@ogu.edu.tr

The design and synthesis of coordination polymers have attracted attention due to their potential applications in the fields such as gas adsorption/separation, catalysis, conductivity, luminescence, magnetism and fascinating topologies [1]. Organic linkers and metal clusters containing secondary building units (SBUs) are very important to obtain three dimensional (3D) coordination polymers with the desired structures [2]. In this study, novel coordination polymers, namely { $[Zn_2(abtc)(1,4-betib)]\cdot DMF$ } (1) { $[Cd_2(abtc)(1,4-betix)]\cdot DMF$ }_n (H₄abtc=3,3',5,5'-azobenzenetetracarboxylic and **(2)** acid. 1,4-betib = 1,4-bis(2-ethylimidazole-1yl)butane and 1,4-betix = 1,4-bis(2-ethylimidazole-1ylmethyl)benzene) were synthesized and structurally characterized by elemental analysis, IR spectra, single crystal X-ray diffraction. Complexes 1 and 2 were synthesized by the solvothermal methods at 100 °C in the presence of HNO₂. Single crystal X-ray diffraction of analyses reveal that 1 and 2 crystallize in the monoclinic and orthorhombic systems with the space groups C2/c, P2₁2₁2₁, respectively. The crystal structures of 1 and 2 with the atom labeling are shown in Fig. 1. As seen in Fig 1, the Zn(II) and Cd(II) ions are five coordinated with a distorted square pyramidal geometry by four oxygen atoms from four different abtc ligands and one nitrogen atom from bis(imidazole) ligands. Each abtc ligand acts as an octadentate ligand, 3,3'-carboxyl and 5,5'-carboxyl groups display bis(monodentate) bridging modes to connect to eight metal centers. Topological analysis revealed that both complexes are 3,6 connected sqc5381 net. Thermal and photoluminescence properties of the complexes were investigated in detail.

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References

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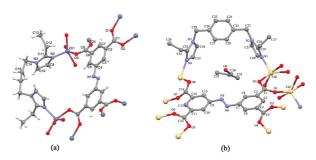


Figure 1. The crystal structures of 1 (a) and 2 (b) with the atom labeling

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