

**CRYSTAL STRUCTURES OF $[\text{Hg}(\text{N-ETHYLTHIOUREA})_2(\text{CN})_2]$
AND $[\text{Hg}(\text{N-PROPYLTHIOUREA})_2(\text{CN})_2]$**

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Two mercury(II) cyanide complexes of *N*-ethylthiourea (Ettu) and *N*-propylthiourea (Prtu) ligands, $[\text{Hg}(\text{Ettu})_2(\text{CN})_2]$ (**1**) and $[\text{Hg}(\text{Prtu})_2(\text{CN})_2]$ (**2**), were prepared and their crystal structures were determined by X-ray crystallography. In both structures, the mercury atom is coordinated to two sulfur atoms of thioureas and two cyanide carbon atoms in a pseudo-tetrahedral mode with the bond angles in the range of 90.52(11)-162.2(3)°. The structures are stabilized by N-H---S, N-H---N, and C-H---N intramolecular and intermolecular hydrogen bonds.

Keywords: mercury(II) cyanide, *N*-ethylthiourea, *N*-propylthiourea, X-ray structures.