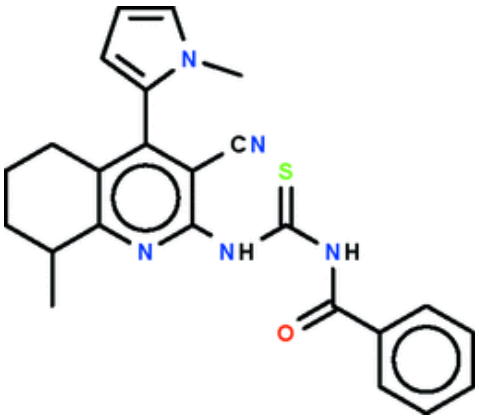

Acta Cryst. (2011). E**67**, o2430 [doi:10.1107/S1600536811033046]

1-Benzoyl-3-[3-cyano-8-methyl-4-(1-methyl-1*H*-pyrrol-2-yl)-5,6,7,8-tetrahydroquinolin-2-yl]thiourea

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Abstract: In the *N*-substituted benzoylthiourea, $C_{24}H_{23}N_5OS$, the benzoylthiourea unit is non-planar (r.m.s. deviation = 0.126 Å). The aliphatic part of the tetrahydroquinoline fused-ring system is disordered over two positions in a 0.592 (5):0.408 (5) ratio. The pyridine and pyrrole rings are twisted by 55.2 (1)° in order to avoid crowding of their respective substituents. Pairs of molecules are linked by N-H...N hydrogen bonds, forming centrosymmetric dimers. Furthermore, an intramolecular N-H...O hydrogen bond stabilizes the molecular conformation.