

# Crystallography of 3,3,6,6-Tetramethyl-9-(3,4-dimethylphenyl)-3,4,6,7,9,10-hexahydroacridine-1,8-dione

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**Abstract**—The title compound 3,3,6,6-Tetramethyl-9-(3,4-dimethylphenyl)-3,4,6,7,9,10-hexahydroacridine-1,8-dione ( $C_{25}H_{31}NO_2$ ) was synthesized and crystallized by the Chemistry Department of Shivaji University, Kohlapur. The compound crystallizes in the monoclinic space group  $P 2_1/c$  with unit cell parameters:  $a = 18.616(3)$ ,  $b = 9.9747(13)$ ,  $c = 11.8093(15)$  Å,  $\beta = 91.89(13)^\circ$ ,  $Z = 4$ . The crystal structure was solved by direct methods and refined by full-matrix least-squares procedures to a final  $R$ -value of 0.075 for 3088 observed reflections. The central ring of the acridinedione system adopts a boat conformation. The four essentially planar atoms (C10/C11/C13/C14) of this ring [maximum deviations =  $0.2047(3)$  Å] forms dihedral angles of  $87.32(13)$ , with the benzene ring. The two outer rings of the acridinedione system adopt sofa conformations. The crystal packing is stabilized by intermolecular  $N-H...O$  interactions.

**Keywords:** Crystal structure. Direct method. Boat and Sofa Conformations. Intermolecular interactions