

STRUCTURAL STUDIES OF 3-CHLORO-*N*-(8'-QUINOLYL)× BENZO[*b*]THIOPHENE-2-CARBOXAMIDE

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3-Chloro-*N*-(8'-quinolyl)benzo[*b*]thiophene-2-carboxamide was synthesized from 3-chlorobenzo[*b*]thiophene -2-carboxyl chloride and 8-aminoquinoline in the presence of triethylamine. The single crystal X-ray structure determination confirmed the earlier proposed structure and also characterized by ¹HNMR, and Mass spectroscopy. Crystallographic study reveals that the structure crystallizes in monoclinic system, $a = 14.878(4) \text{ \AA}$, $b = 8.4292(15) \text{ \AA}$, $c = 25.461(7) \text{ \AA}$, $\beta = 112.022(18)^\circ$, $Z = 8$, $V = 2960.20(12) \text{ \AA}^3$ with space group $C2/c$ (No. 15). In the structure packing, three kinds of interactions are responsible for the stability of the structure. Infinite two-dimensional stair-like layered chains are formed by relatively strong intermolecular hydrogen bonds [C14–H14...O1]. These parallel chains are connected by several π – π and CH– π interactions, alternatively. There are two such parallel chains with 70.53° , which are in contact by van der Waals interactions.

Keywords: benzothiophene, carboxamide, crystal structure, hydrogen bonding, CH– π interaction, π – π interaction.