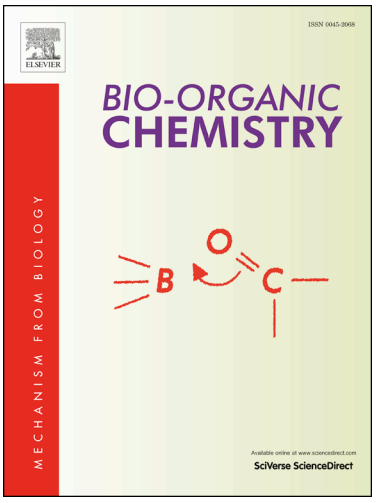

Synthesis, Computational Molecular Docking Analysis and Effectiveness on Tyrosinase Inhibition of Kojic Acid Derivatives

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