

Accepted Manuscript

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

Gülşah Karakaya, Aslı Türe, Ayşe Ercan, Selin Öncül, Mutlu Dilsiz Aytemir

PII: S0045-2068(19)30078-1
DOI: <https://doi.org/10.1016/j.bioorg.2019.102950>
Article Number: 102950
Reference: YBIOO 102950

To appear in: *Bioorganic Chemistry*

Received Date: 17 January 2019
Revised Date: 16 April 2019
Accepted Date: 23 April 2019

Please cite this article as: G. Karakaya, A. Türe, A. Ercan, S. Öncül, M. Dilsiz Aytemir, Synthesis, Computational Molecular Docking Analysis and Effectiveness on Tyrosinase Inhibition of Kojic Acid Derivatives, *Bioorganic Chemistry* (2019), doi: <https://doi.org/10.1016/j.bioorg.2019.102950>

This is a PDF file of an unedited manuscript that has been accepted for publication. As a service to our customers we are providing this early version of the manuscript. The manuscript will undergo copyediting, typesetting, and review of the resulting proof before it is published in its final form. Please note that during the production process errors may be discovered which could affect the content, and all legal disclaimers that apply to the journal pertain.

