

## Review of: "Expansion of the Experimental Antifungal Activities Through in Silico Docking Study of Compounds From Albizia Lebbeck"

Mohamed A.A. Radwan<sup>1</sup>

1 National Research Center, Egypt

Potential competing interests: No potential competing interests to declare.

This publication presents the results of an in silico molecular docking investigation aimed at determining if chemicals isolated from Albizia lebbeck. The work is presented in an effective manner overall, and the computational technique is pertinent to the study's goals. A few small changes are recommended to increase clarity and fix some constraints:

- -More information about the docking protocol parameters, such as the scoring functions and docking algorithms that are employed, is required.
- -For clarity, Table 1 may include compound names, and the Figure quality could be improved.

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