Review of: "Design and Molecular Screening of Various Compounds by Molecular Docking as BACE-1 Inhibitors"

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Potential competing interests: No potential competing interests to declare.

1. The manuscript is well written by selecting a suitable disease and applying the computational approach.
2. The manuscript has to address the following queries:
3. How the compounds were designed for the study. What is the rationale?
4. Provide the molecular docking interactions for the lead compounds.
5. Concise the conclusion part.
6. Why the computational approach is selected for the study.