## **IJARSCT**



International Journal of Advanced Research in Science, Communication and Technology (IJARSCT)

International Open-Access, Double-Blind, Peer-Reviewed, Refereed, Multidisciplinary Online Journal

Volume 4, Issue 1, December 2024

## Review of Computer-Aided Drug Design and its Implications in Drug Discovery and Development

Machhindra M Sarode<sup>1</sup>, Bhagyashree B Randhwan<sup>2</sup>, Snehal Fand<sup>3</sup>, Abhishek A Chougule<sup>4</sup>, Gaurav P Dhanwade<sup>5</sup>, Aditya D Dhiwar<sup>6</sup>, Sohel C Shaikh<sup>7</sup>, Naman K Gandhi<sup>8</sup>
Students, Arihant College of Pharmacy, Kedgaon, Ahilyanagar, Maharashtra, India<sup>1,4,5,6,7,8</sup>
Professor, Arihant College of Pharmacy, Kedgaon, Ahilyanagar, Maharashtra, India<sup>3,4</sup>

**Abstract:** Computer-aided drug design (CADD) is a rapidly growing field of research that uses computers to help design new drugs. It's a fascinating area that combines different aspects of both basic and applied science. The core principles of CADD rely on quantum mechanics and molecular modeling techniques. These techniques include designing drugs based on a known structure, designing drugs based on existing molecules (ligands), searching databases for potential drug candidates, and predicting how well a drug might bind to a specific biological target. This review will explore how CADD tools are used to support the drug discovery process.

DOI: 10.48175/568

Keywords: Molecular modeling, Target molecules, Drug discovery process

