# Chapter 26 Role of Molecular Docking in Computer-Aided Drug Design and Development

Rahul Agarwal Shiv Nadar University, India

Ashutosh Singh Shiv Nadar University, India

Subhabrata Sen Shiv Nadar University, India

## ABSTRACT

Molecular Docking is widely used in CADD (Computer-Aided Drug Designing), SBDD (Structure-Based Drug Designing) and LBDD (Ligand-Based Drug Designing). It is a method used to predict the binding orientation of one molecule with the other and used for any kind of molecule based on the interaction like, small drug molecule with its protein target, protein – protein binding or a DNA – protein binding. Docking is very much popular technique due to its reliable prediction properties. This book chapter will provide an overview of diverse docking methodologies present that are used in drug design and development. There will be discussion on several case studies, pertaining to each method, followed by advantages and disadvantages of the discussed methodology. It will typically aim professionals in the field of cheminformatics and bioinformatics, both in academia and in industry and aspiring scientists and students who want to take up this as a profession in the near future. We will conclude with our opinion on the effectiveness of this technology in the future of pharmaceutical industry.

### 1. INTRODUCTION

Computer aided drug designing or CADD, is a strategy that harnesses state of the art technology to expedite the drug development process. Traditionally the drug development /discovery (that includes random screening, serendipitous discovery and process optimization) takes nearly a decade to complete with an

DOI: 10.4018/978-1-5225-1762-7.ch026

average expense of ~300 million dollar. CADD tends to curtail this expenditure and timeline by providing a holistic view of a drug discovery project and enables a SWOT (strength-weakness-opportunity-threat) analysis to evaluate the viability of the program.

Molecular docking is one of the fundamental pillars of CADD. It analyzes the binding interaction between a target protein and small molecules (Lengauer & Rarey, 1996). These small molecules also referred as ligands are the potential drug candidates that are being developed against the phenotypes and the therapeutic model for which the protein in the CADD is the target (Sliwoski, Kothiwale, Meiler, & Lowe, 2014).

## 2. HISTORY OF CADD

The evolution of CADD began in 1900s when Emiel Fischer (in 1894) and later Paul Ehlrich (in 1909) propagated the concept of receptors and lock and key mechanism (Fischer, 1894). Lock and key concept explains the interaction of a drug and its receptor. Just as there are innumerable pairs of lock and key, where each key inserts into a specific lock to open it, drugs are designed to be tailor made for a particular receptor. Depending on the perfection of the design, the drugs its binding capacity of the drugs with the receptor excels. But just as duplicate keys can be made for a single lock, which makes the lock vulnerable, there is a chance that multiple drugs can work on single receptor thereby giving rise to undue side effects. It was Paul Ehlrich, who proposed the concept of "magic bullet" the elusive drug that only bind to the chosen receptor thereby exhibiting no side effects (Elrich P, 1909,1957).

Figure 1. Flowsheet of the operations in CADD



26 more pages are available in the full version of this document, which may be purchased using the "Add to Cart" button on the publisher's webpage: <u>www.igi-global.com/chapter/role-of-molecular-docking-in-computer-aided-</u> drug-design-and-development/174146

## **Related Content**

#### Industrial Enzyme Technology: Potential Applications

Michael Bamitale Osho (2018). *Research Advancements in Pharmaceutical, Nutritional, and Industrial Enzymology (pp. 375-394).* www.irma-international.org/chapter/industrial-enzyme-technology/203824

#### Liposomes: Properties and Therapeutic Applications

Ljiljana Djekic (2017). *Novel Approaches for Drug Delivery (pp. 27-51).* www.irma-international.org/chapter/liposomes/159653

#### Molecular Docking at a Glance

Maryam Hamzeh-Mivehroud, Babak Sokoutiand Siavoush Dastmalchi (2016). *Methods and Algorithms for Molecular Docking-Based Drug Design and Discovery (pp. 1-38).* www.irma-international.org/chapter/molecular-docking-at-a-glance/151881

#### Effect of Ocimum sanctum on Cancer Progression and Its Effective Therapeutic Activity

Chittipolu Ajaykumar, K. Swathiand Pavithra Dasari (2023). *Natural Products as Cancer Therapeutics (pp. 1-23).* 

www.irma-international.org/chapter/effect-of-ocimum-sanctum-on-cancer-progression-and-its-effective-therapeuticactivity/329152

#### Secondary Metabolites From Plants for Cardiovascular Disease

Jose Prakash (2023). *Pharmacological Benefits of Natural Agents (pp. 155-171).* www.irma-international.org/chapter/secondary-metabolites-from-plants-for-cardiovascular-disease/327308