

# **Drug Design: Structure- And Ligand-Based Approaches**

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The first computational structure-based drug design methods came into existence in the early 1980s and are, to an extent, still in their infancy. There have been

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## **Ligand- and structure- based drug design**

Owing to the biological role of PPARs, these receptors have become important therapeutic targets for antidiabetic and antimetabolic syndrome therapies . However

## **Figure 2: structure and ligand based drug design**

Figure 2: Basic principles and types of drug design. The term SBDD explicates various approaches wherein the structural knowledge of the drug target is exploited and

## **Ligand docking and structure- based drug design |**

The Center for Computational Biology The University of Kansas 2030 Becker Drive Lawrence, KS 66047

## **Structure and ligand based drug design strategies**

Structure Based Drug Design. SBDD is the approach where the structural information of the drug target is exploited for the development of its inhibitor.

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The first is referred to as ligand-based drug design and the second, structure-based drug design. Ligand bound to a potent ligand. This approach to drug discovery

## **Computational and structural approaches to drug**

Computational methods impact all aspects of modern drug discovery and most notably these methods move rapidly from academic exercises to becoming drugs in clinical

## **Structure- based drug design: nmr- based approach**

Structure-based drug design: NMR-based approaches will undoubtedly be helpful in high throughput screening and in providing structural Ligand-based experiments.

## **Combining ligand- based and structure- based drug**

In recent years there has been considerable interest in developing new and improved approaches for the identification of novel therapeutic agents.

## **Structure- based drug design (sbdd) : positive**

Structure-based drug design is one The role of structure-based ligand design and 3 comments on Structure-Based Drug Design (SBDD) : Positive Approach

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Structure-based approaches to a new approach to computer-aided ligand- and receptor-based drug design which seek compounds with structure

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1. Future Med Chem. 2011 Apr;3(6):735-50. doi: 10.4155/fmc.11.18. Integrating structure-based and ligand-based approaches for computational drug design.

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