Ligands and Receptors: A Love Story

Binding Affinities Between Opioids and the μ -opioid Receptor

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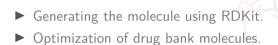
https://github.com/JoshuaKSt

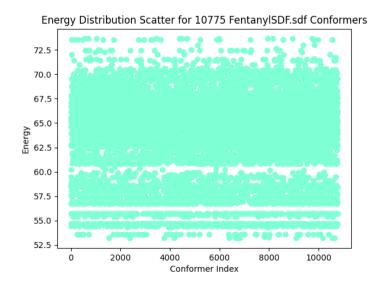
August 1, 2024

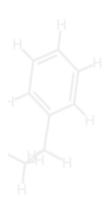
Research Goal

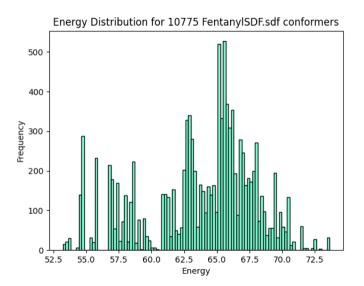
- ▶ Direction.
- ► Understand opioid receptors.
- ► Create a molecule.
- Simulate its interaction with the μ -opioid receptor and compare.

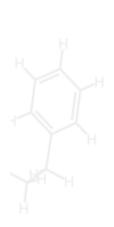
Molecule Creation and Optimization Method



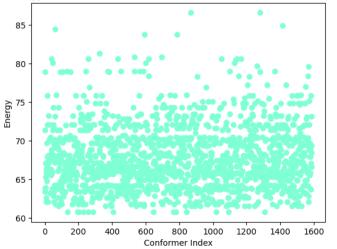


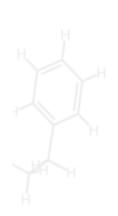


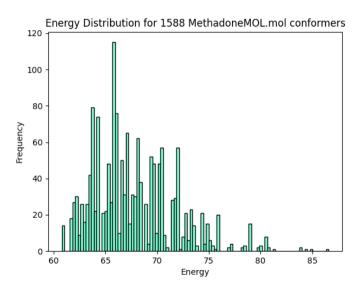


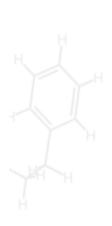


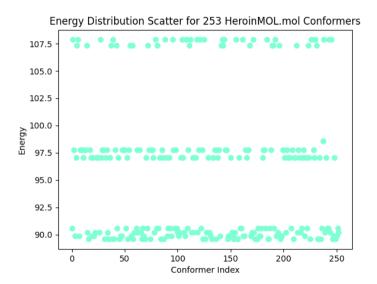


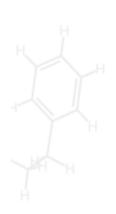


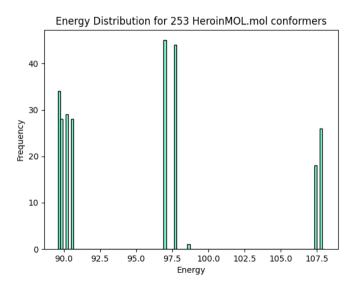


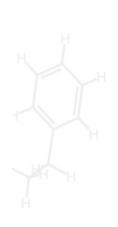




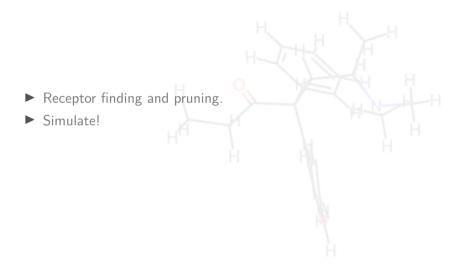








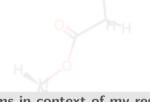
Experiment Setup



What are: "Binding Affinities"?

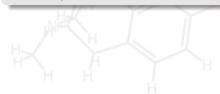
They are:

- ► How tightly bound an interaction is (i.e. a ligand and protein)
- ightharpoonup Often referred to as the free energy of binding (ΔG)
- Used in drug optimization, computations of biological systems etc.



Simple terms in context of my research:

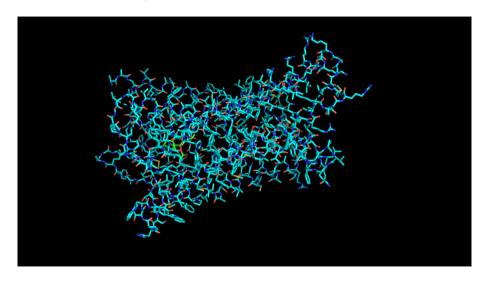
The total energy decrease when a ligand binds to a receptor



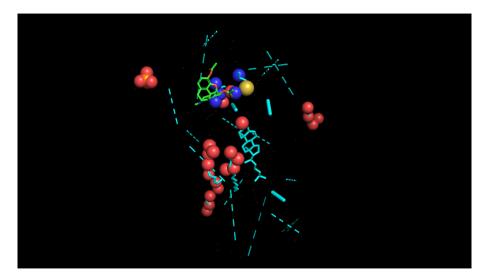
Briefing

- ightharpoonup $\mu ext{-opioid Receptors}$
 - ► 6DDF (Human, Mouse, Removed DAMGO)
 - ► 8E0G (Mouse, Llama, Removed BU72)
 - ► 8K9K (Human, Removed DAMGO)
- ► Ligands
 - ► Heroin
 - ► Morphine
 - ► Fentanyl
 - ▶ Methadone
 - ► Carfentanil
 - ► BU72
- ► .SDF and .MOL
 - ► Structure Data Files

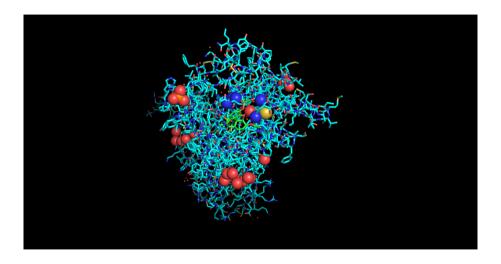
Heroin 6DDF Docking Output



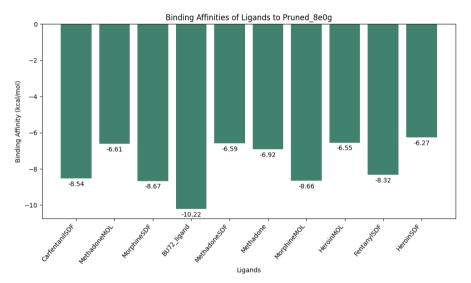
Heroin 8E0G Docking Output



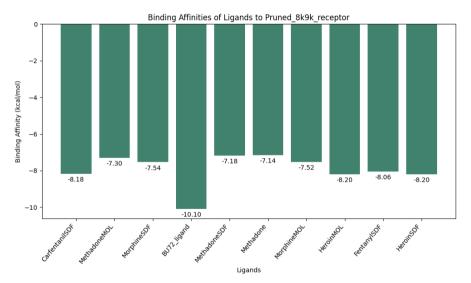
Heroin 8E0G Docking Output



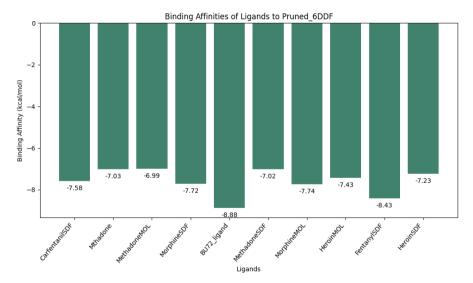
Affinity Graphs



Affinity Graphs



Affinity Graphs



Analysis

- ▶ BU72 Consistently had the strongest binding.
- ► Fentanyl and Carfentanil consistently showed strong binding.
- Morphine was stronger bound to a higher quality model.
- ► The gap in affinity lessened when the structure was lower quality.
- Heroin and Methadone varied greatly across structure and file format.

Thank You!

