



## Al-assisted Drug Discovery

**Enabling and Accelerating Candidate Nomination** 

- Aurigene.Al: An Al-assisted drug discovery platform
- Aurimine: A virtual database comprising of 12+ billion compounds

## **Key Capabilities**

- Aurigene. Al consists of a meticulously curated compound database, which is ever-expanding using both in-house and public domain information.
- This database currently consists of 180 million compounds, and 1.6 million validated bioassay data points and serves as a training data for the Aurigene. Al platform.
- This evolving platform consists of several top-tier Al algorithms 4 predictive, 11 generative and 8 CADD models.
- This modular platform enables our scientific experts to cherry-pick the right set of Al models for a given application. The model output can help prioritize the chemical designs with appropriate properties, thereby leading to increased probability of success.
- We are well equipped to quickly synthesize and run a variety of experimental assessments both in-vitro and in-vivo. The data output can be used for training the AI models to further improve the chemical design.
- The accuracy and reliability of the data output from Aurigene. Al results from an iterative DMTA cycle followed by training the Al models with generated data in real-time.

## Top 5 reasons to partner with Aurigene.Al

- Interpretable AI platform
- Enables "fail fast and iterate faster" model
- Scientific experts can cherry-pick the right set of AI models for a given application
- Customizable and ever-expanding data base and algorithms
- An end-to-end solution combining AI with Aurigene's core expertise in synthesizing and experimentally testing the molecules