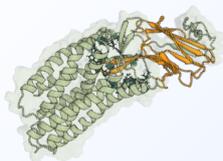
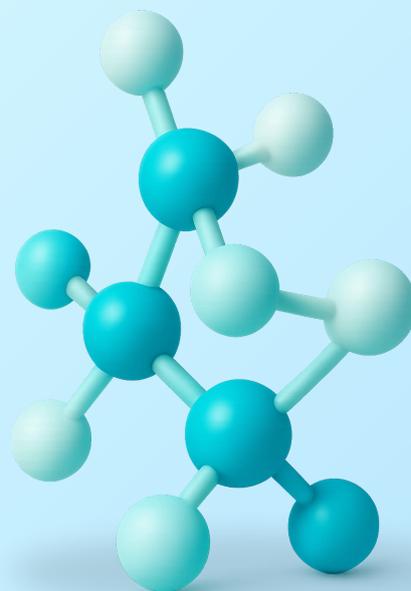


MedGraph Topaaz™

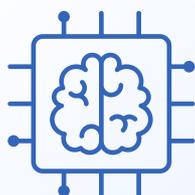
AI-Powered De Novo Design Platform for Next-Gen Molecule Discovery

Precision Discovery | Scientifically
Grounded | Experimentally Relevant



Generative Design

AI-driven compound generation using curated fragment libraries to build novel, target-specific scaffolds



AI-Driven Optimization

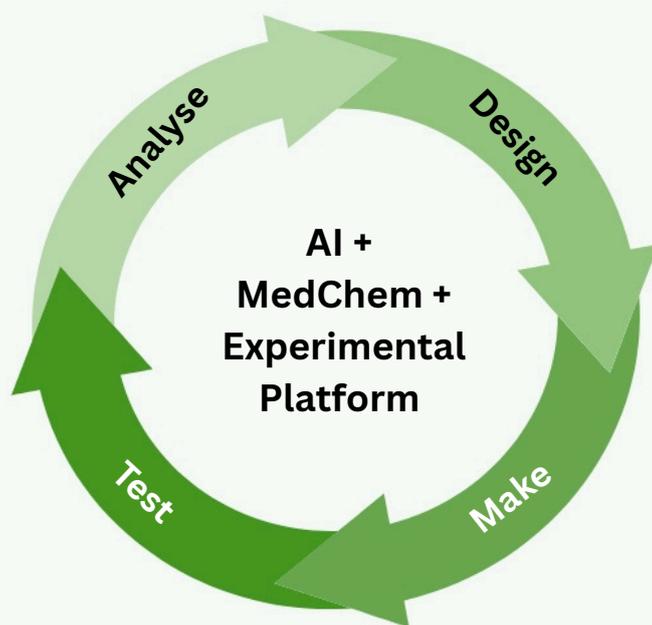
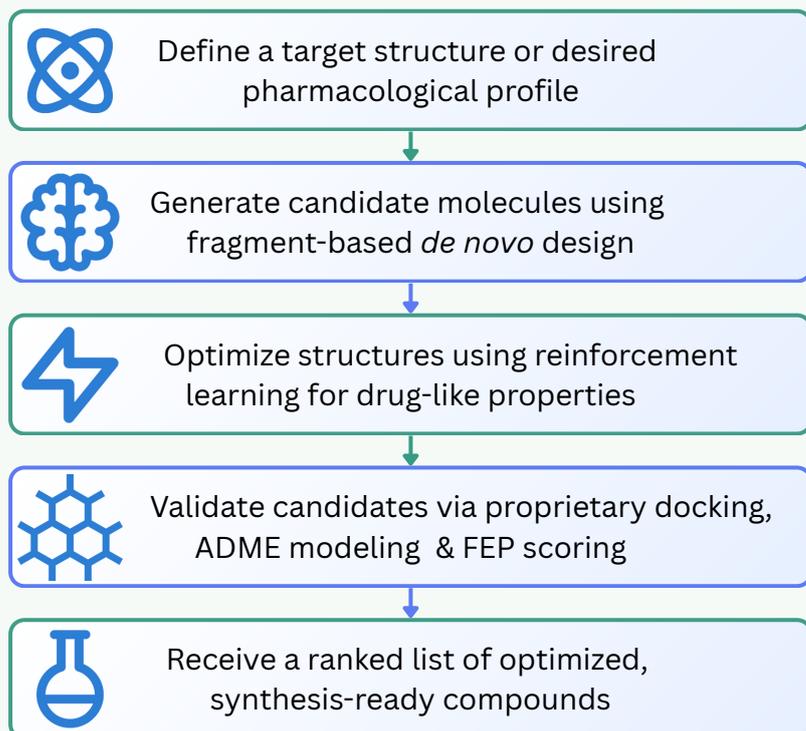
Refines molecules for ADME, efficacy, safety and synthetic feasibility using reinforcement learning



AI+Physics Screening

Ranks drug-like candidates through docking, AI-driven scoring and FEP-based binding affinity scoring

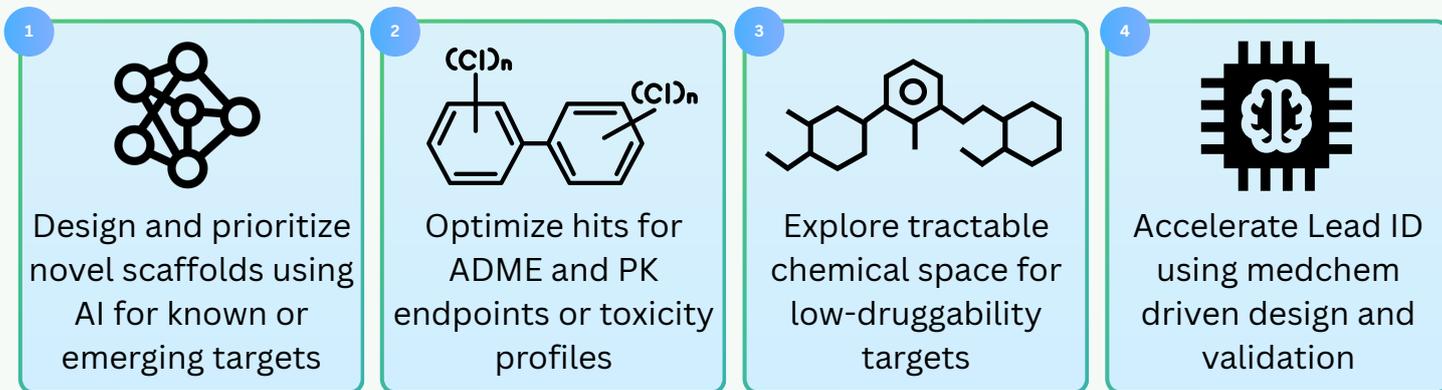
Workflow Snapshot



**AI meets chemistry to unlock
breakthrough candidates**



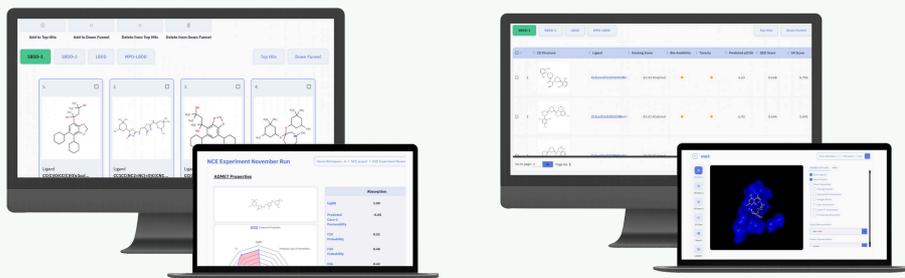
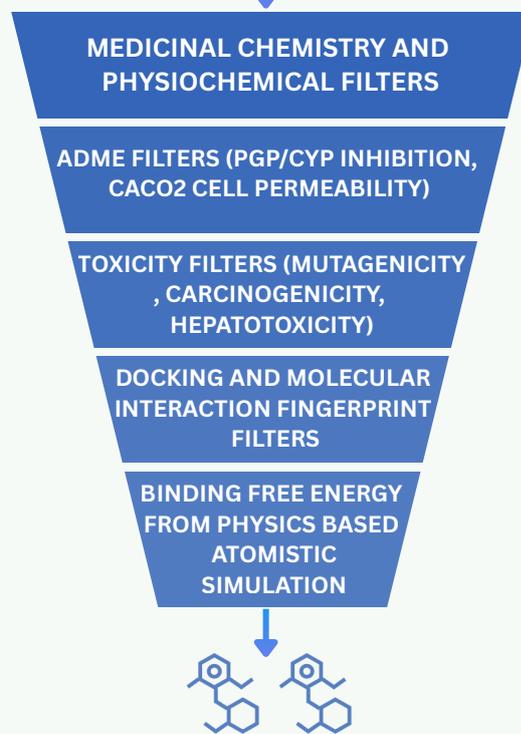
Applications of MedGraph Topaaz™



Why Choose MedGraph - Topaaz™

Revolutionizing Drug Discovery with Generative AI

- End-to-end AI-powered molecule design: From ideation to preclinical candidate in a single pipeline
- Hybrid approach: Generative AI + FEP: Combines creativity and precision for confident decision-making
- High scalability and secure deployment: Available as cloud, on-premise or hybrid deployment
- Customization for every program: Define ADME, novelty, selectivity and synthetic feasibility criteria
- Rapid turnaround and modular access: Reduced TAT and flexible engagement



Engagement Model

Platform Module Licensing

Lead Candidate-as-a-Service

Outcome Driven Productized Solutions

Co-development

Contact Us

Let's explore how we can support your next discovery efforts



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Topaaz is built to accelerate your discovery goals without compromising scientific integrity



Contact Us



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