





REAL-WORLD INNOVATION

Al-first discovery platform generating multi-parameter molecules

November 27, 2025

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THERE IS A GAP

BETWEEN REAL-WORLD CHALLENGES AND INNOVATIVE SOLUTIONS





UTILIZING

AI AND ADVANCED PROPRIETARY ALGORITHMS

TO DELIVER INNOVATIVE SMALL MOLECULES, TAILORED TO REAL-WORLD COMMERCIAL PRODUCT NEEDS



WE CALL IT: REAL-WORLD INNOVATION



EVOGENE PIONEERS **REAL-WORLD INNOVATION**

Using a proprietary generative AI engine, we design highly potent and novel small molecules, optimized across multiple-parameters, for the pharmaceutical and ag-chemical industries







OUR STORY IN 3 SENTENCES

GEN AI ENGINE

Design of small molecules with high probability of success



REVOLUTIONIZING AG

Proven results with strategic collaborations



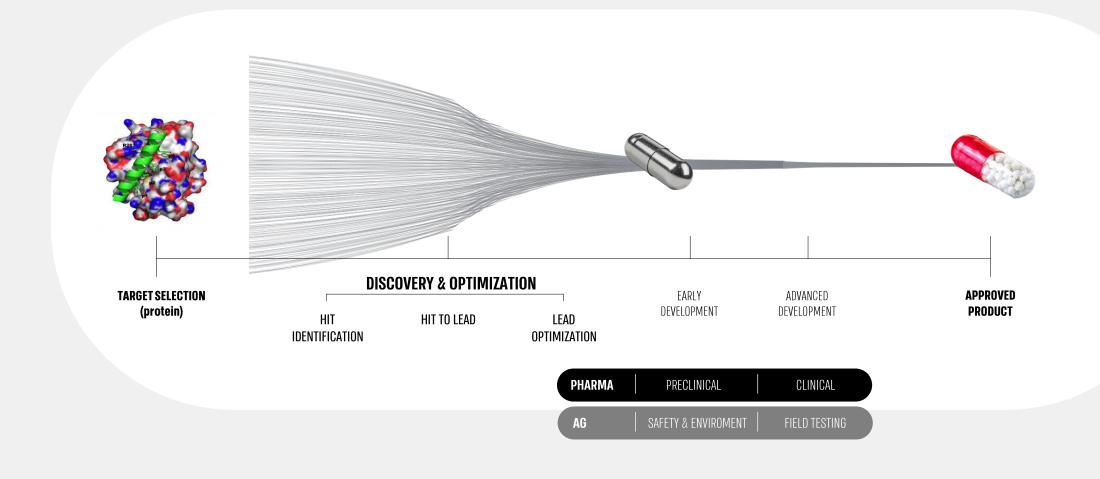
PIONEERING IN PHARMA

Differentiated tech offering that fits market demand



THE DISCOVERY & OPTIMIZATION CHALLENGE

DESIGN THE RIGHT CANDIDATES TO REACH THE MARKET



LESS THAN
10% OF
MOLECULES
TESTED IN
CLINICAL TRIALS
MAKE IT TO
MARKET





Discovery & optimization are the critical stages for increasing market success probability

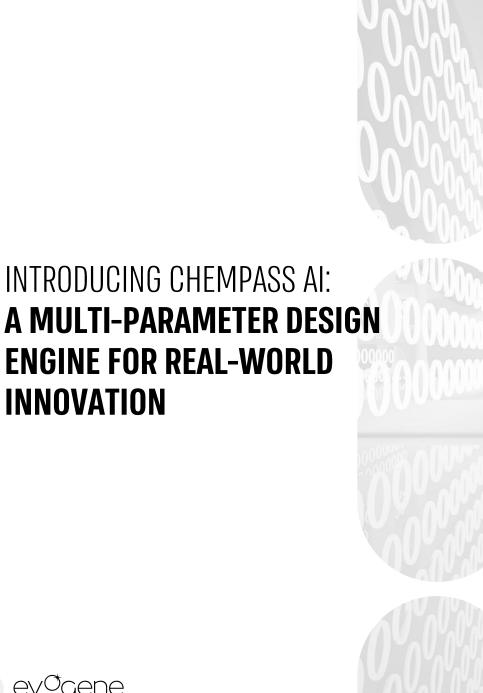
THE MULTI-PARAMETER OPTIMIZATION **CHALLENGE**

Creating a molecule that will succeed through all development phases requires balancing in advance multiple, often competing, parameters

THE CHALLENGE FOR TRADITIONAL METHODS:

- Optimization typically limited to a small number of parameters
- Cannot address multipleparameters simultaneously
- Requires lengthy, resourceintensive iterative cycles







MULTI-PARAMETER OPTIMIZATION

Optimizes multiple-parameters simultaneously, adapted to project specific chemical, biological, and physical constraints, increasing probability of success

NOVEL MOLECULAR STRUCTURES

Based on 38B molecules, the algorithm generates completely novel, synthesizable and effective compounds, extending beyond explored chemical space.

HIGHLY POTENT

Al-first designed molecules optimized through targeted experimental validation

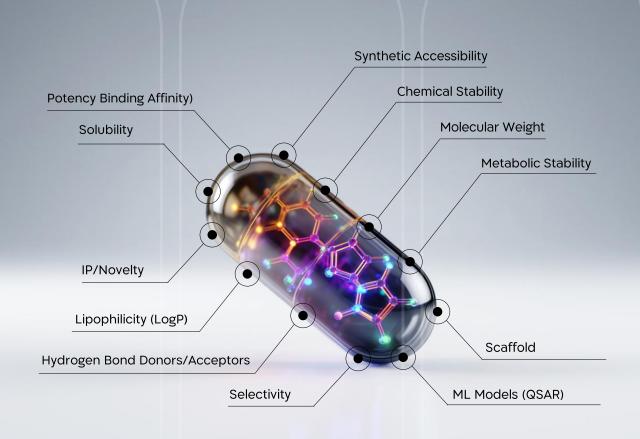


INNOVATION

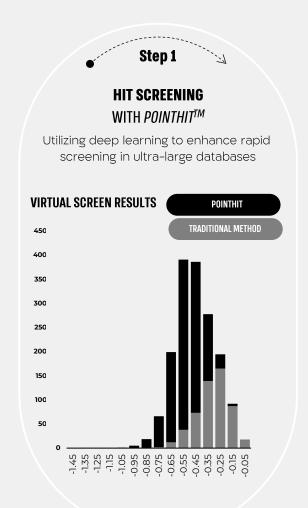
CHEMPASS AI: A MULTI-PARAMETER GENERATIVE DESIGN ENGINE FOR REAL-WORLD INNOVATION

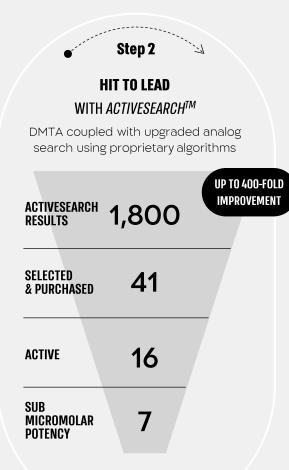
CHEMPASS AI TRANSFORMS TRADITIONAL METHODS:

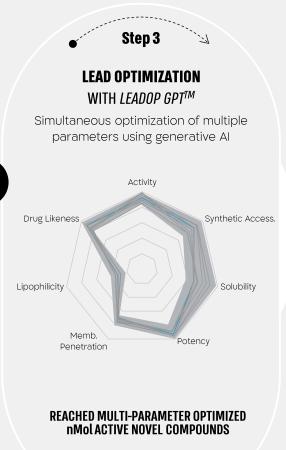
- Optimization of diverse product specific parameters
- Addresses multiple constraints simultaneously
- Minimizes lengthy, resource-intensive iterative cycles



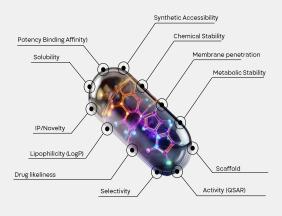
CHEMPASS AI FOUNDED BY PROPRIETARY TECHNOLOGY BOOSTING PROBABILITY OF SUCCESS





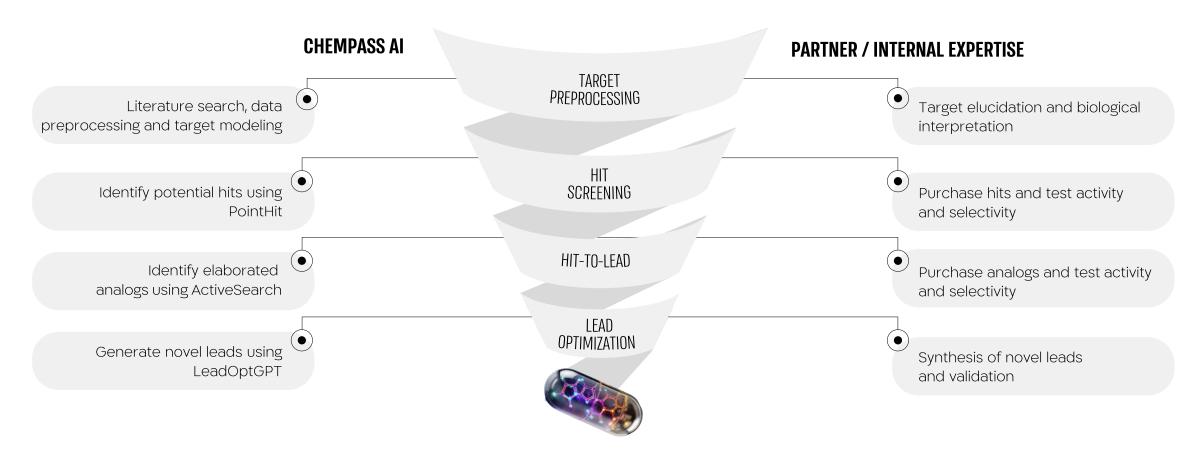


Result PRODUCT CANDIDATE



THE KEY FOR SUCCESS -STRATEGIC PARTNERSHIPS / INTERNAL EXPERTISE

TAILORED TO EACH CHALLENGE & OPTIMIZED FOR SUCCESS





Making breakthroughs in life sciences smarter and faster

GOOGLE CLOUD X EVOGENE

BREAKING NEW GROUND IN AI-FIRST MOLECULE DESIGN

TRANSFORMATIVE PARTNERSHIP

Combining Evogene's proprietary AI foundation model, with Google Cloud's world-class infrastructure to scale revolutionary molecular discovery

ACCELERATED INNOVATION

Google's infrastructure enabled efficient use of a massive dataset starting with millions and ultimately reaching a training set of approximately 38 billion molecular structures

3x IMPROVED ACCURACY

Optimization across multiple-parameters reached 90% precision (vs. 29% in traditional GPT Al-model)

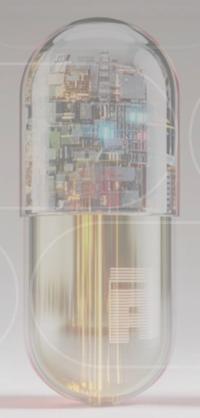
"We're pleased to collaborate with Evogene's innovation in Alpowered molecule design. Their progress with ChemPass AI highlights the strength of pairing advanced AI infrastructure with deep scientific insight. We look forward to seeing the impact of this new model in drug discovery and agriculture."

BOAZ MAOZ, MANAGING DIRECTOR, GOOGLE CLOUD ISRAEL

OUR STORY IN 3 SENTENCES

GEN AI ENGINE

Design of small molecules with high probability of success



REVOLUTIONIZING AG

Proven results with strategic collaborations



PIONEERING IN PHARMA

Differentiated tech offering that fits market demand





A GROWING MARKET WITH PLENTY OF ROOM FOR INNOVATION

- Increase of pest resistance & regulatory requirements
- Urgent need for new Modes of Action (MoAs)
- Decreased rate in discovery of new pesticides due to lack of innovation



\$43.3B BROAD SPECTRUM HERBICIDES¹

\$22.2B FUNGICIDES FOR MAJOR DISEASES²

\$22.3B BROAD SPECTRUM INSECTICIDES³







STRATEGIC COLLABORATIONS



"Bringing together AgPlenus' expertise with Bayer's CropKey approach to crop protection innovation will help accelerate the delivery of essential, sustainable, and affordable solutions to farmers and set a new benchmark in the industry."

Rachel Rama, Head of Small Molecules, Crop Science division



Nontreated



Treated

Development of new sustainable Mode-of-Action broad spectrum, herbicide

Demonstrated high weed control efficacy and good tolerance in corn



"The collaboration with AgPlenus has accelerated the identification of a class of herbicide chemistry that targets a new modeof-action for weed control, something the industry has been lacking for decades."

Vid Hegde, Former VP of Crop Protection Discovery and Development



Treated nontreated

Development of novel herbicides

Demonstrated weed growth reduction in post emergence application



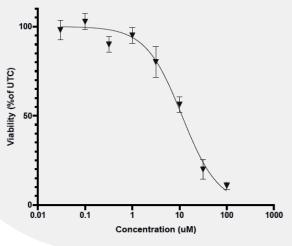


INTERNAL PIPELINE WHEAT BLOTCH

DISRUPTING A \$1.2B PROBLEM AT RECORD SPEED

- 70% of EU fungicide usage in wheat is for Wheat Blotch¹
- EU market alone >\$1.2B annually1
- Widespread resistance to current top products ('Strobilurins') with 2024 sales of \$4.59B²

APTF-4 PRESENTS STRONG POTENTIAL FOR REAL-WORLD INNOVATION

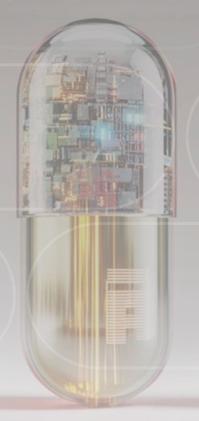


- Shows clear concentration– dependent antifungal efficacy
- Ongoing optimization to enhance potency
- 18 months from target to optimized hit

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GROWING OPPORTUNITY IN AI-DRUG DISCOVERY







SMALL MOLECULES

Account for **58%** Of total pharmaceutical market (\$1,344B)

AI-based candidates in pipeline

>60

150% CAGR in last 3 years

Discovery market expected to reach

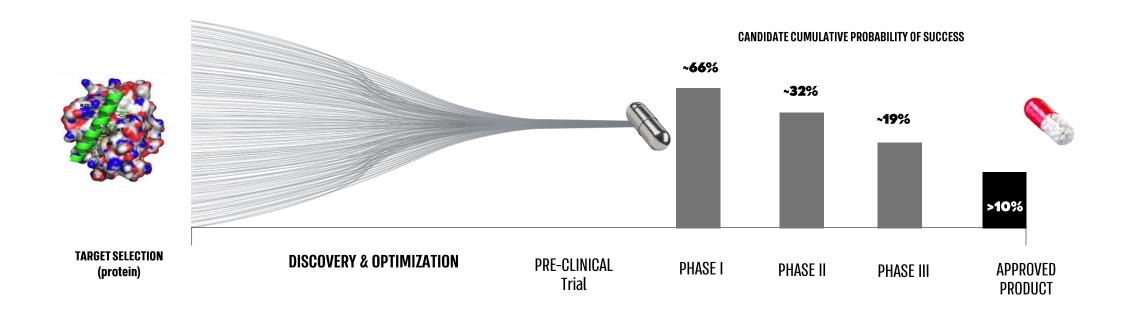
\$190.68B

by 2034

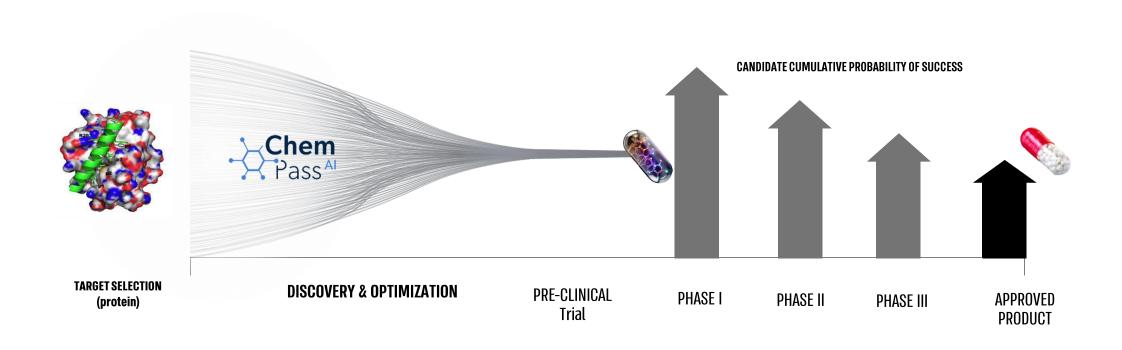




THE DISCOVERY AND OPTIMIZATION CHALLENGE DEVELOPING CANDIDATES THAT WILL REACH THE MARKET



CHEMPASS AI: DRIVING HIGHER CUMULATIVE PROBABILITY OF SUCCESS



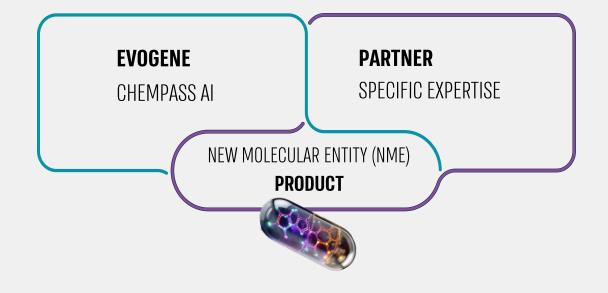
BUSINESS STRATEGY

CAPTURE THE VALUE OF CHEMPASS AI

through diverse collaborative partnerships to accelerate small molecule-based drug development.

PARTNERING WITH PHARMA AND BIOTECH COMPANIES

that complement our technology, enabling groundbreaking innovations that benefit both parties.



STRATEGIC COLLABORATIONS

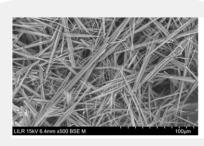


Evogene and Professor Ehud Gazit of Tel Aviv University Collaborate to develop new therapeutics for metabolite Disease

The collaboration aims to design novel small molecules that can effectively inhibit pathological self-assembly processes, leading to transformative new therapeutics with the potential to improve the lives of millions of patients suffering from diseases including Tyrosinemia, Gout and Maple Syrup Urine Disease (MSUD).

The collaboration brings together Evogene's ChemPass Al's state-of-the-art computational capabilities for generative molecular design with Professor Gazit's world-renowned expertise in the characterization and manipulation of molecular self-assembly.

Scanning electron microscopy of crystals formed by the self-assembly of a metabolite



* Costa-Bauza, A., & Grases, F. (2023). Biomolecules, 13(12), 1769.



DELIVERING REAL-WORLD INNOVATION

THROUGH TWO STRATEGIC MARKET OPERATIONS DIVISIONS



PHARMA DIVISION



EXTERNAL COLLABORATIONS



INTERNAL PIPELINE

Undisclosed

STRATEGIC COLLABORATION





INTERNAL PIPELINE

Wheat Blotch

evøgene



Creating

REAL-WORLD INNOVATION

for a better future