

app

advanced features

analytics

drug explorer

fda dashboard


interaction network


molecule visualizer

trade dashboard


Navigation

☒ Analysis

☐  Drug Explorer

☐  Advanced AI

Features

☐  Drug Interactions

☐ Analytics

☐ Email Reports


Analysis Progress

Waiting to start...




Welcome to PharmaGenie AI

Your AI-Powered Drug Repurposing Assistant

▼  How It Works

1. Enter a drug name you're interested in
2. Select a therapeutic area (optional)
3. Click *Analyze* to generate insights
4. View detailed reports and export results


>  Example Queries

Drug Name

Enter drug name

Therapeutic Area (optional)

e.g., Oncology

 Start Analysis



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Select Feature

 Drug Repurposi... 



Advanced AI Features

Patent-Worthy Innovations powering next-generation pharmaceutical intelligence. These features combine cutting-edge AI, real-time data, and novel algorithms.



AI-Powered Drug Repurposing Engine

Patent-Worthy Innovation: Multi-modal drug discovery with confidence scoring

- Molecular similarity analysis using graph neural networks
- Disease pathway overlap detection
- Clinical evidence aggregation with temporal weighting


Enter Drug Name



Target Disease (Optional)



Metformin

All Diseases 



Analyze Repurposing Opportunities

Found 5 repurposing opportunities!

Total Opportunities

5

High Confidence

1

Avg Confidence

63.7%

Repurposing Candidates

> #1: Cardiovascular (Confidence: 77%)

> #2: Autoimmune (Confidence: 61%)

> #3: Alzheimer'S (Confidence: 61%)

> #4: Diabetes (Confidence: 60%)

> #5: Depression (Confidence: 59%)



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Analytics Dashboard



Drug Comparison

Enter first drug name

Metformin

Enter second drug name

Aspirin

Compare Drugs



Statistics Overview

Total Analyses

150

Success Rate

46.7%

Unique Drugs

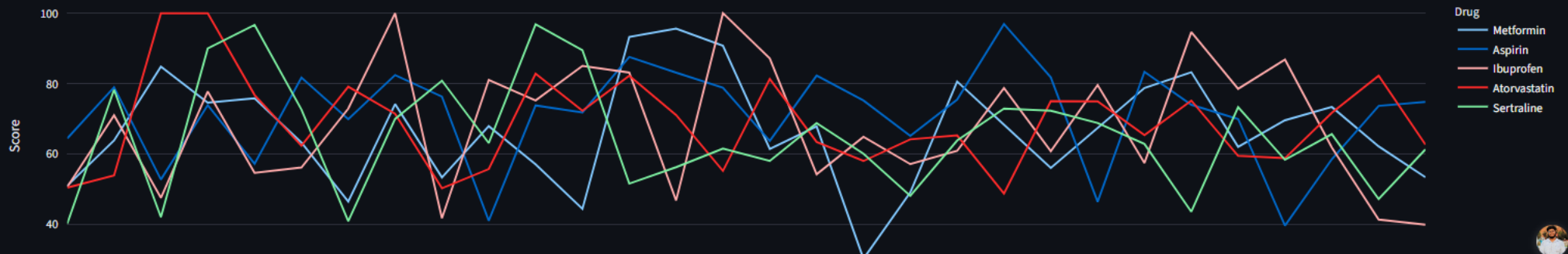
5

Date Range

2025/11/17 – 2025/12/16


Score Trends Over Time

Scores Over Time



Comprehensive Drug Explorer

Explore detailed drug information including classification, mechanism of action, molecular structure, safety information, and interactions - all powered by free APIs.

 Enter Drug Name: ?

e.g., Aspirin, Ibuprofen, Metformin

 Search Drug

Quick Examples:

Aspirin


Ibuprofen

Paracetamol

Metformin

Omeprazole

Amoxicillin






 Drug information retrieved successfully!



Aspirin

RxCUI

1191

 Overview  Molecular Info  Safety  Interactions  Structure



Molecular Structure Visualization

Molecular Weight

180.04 g/mol

H-Bond Donors

1

Rotatable Bonds

2

LogP

1.31

H-Bond Acceptors

3

Heavy Atoms

13

☐ Generate 3D structure (slower) ?

2D Structure 3D Structure

2D Structure



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

interaction network

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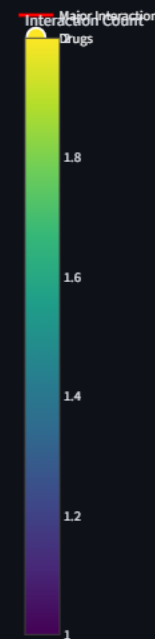
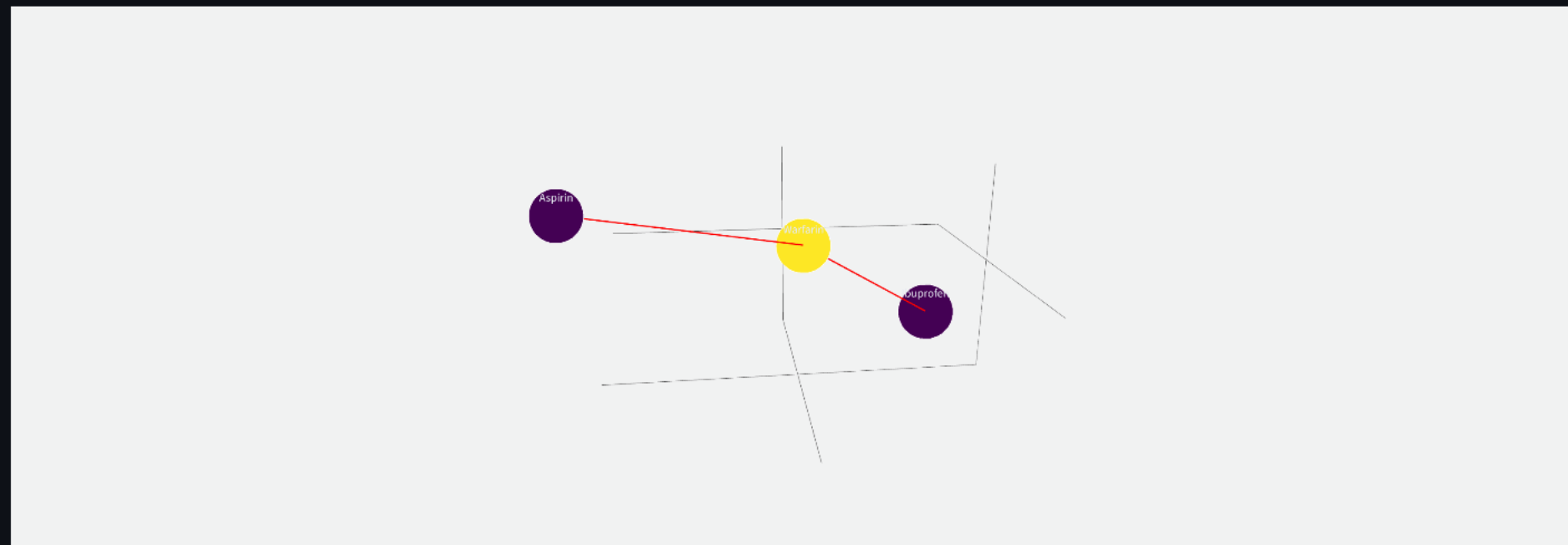
AI Drug Interaction Network

Patent-Worthy Feature: Dynamic 3D network visualization with real-time pharmacokinetic simulation and temporal interaction analysis.

 Interaction Network  Pharmacokinetic Profiles  Timing Analysis

3D Interaction Network

3D Drug Interaction Network



Select Drugs

Choose drugs to analyze

Warfarin ×

Aspirin ×



Ibuprofen ×

Detected Interactions

> ● warfarin ↔ aspirin - MAJOR

> ● warfarin ↔ ibuprofen - MAJOR



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

molecule visualizer

trade dashboard



AI Drug Interaction Network

Patent-Worthy Feature: Dynamic 3D network visualization with real-time pharmacokinetic simulation and temporal interaction analysis.

 Interaction Network  Pharmacokinetic Profiles  Timing Analysis

Temporal Interaction Analysis

First Drug

Warfarin

Second Drug

Aspirin

Warfarin Dose (mg)

100.00

Aspirin Dose (mg)

100.00

Administration Time

18:16

Administration Time


18:16

Analyze Timing

 **Innovation:** This tool uses advanced network theory and pharmacokinetic modeling to predict drug-drug interactions in real-time. Patent-pending algorithm.

Select Drugs

Choose drugs to analyze

Warfarin Aspirin Ibuprofen 



Molecule Visualizer

Visualize molecular structures in 2D and 3D. Enter a drug name or SMILES string to get started.

Enter a drug name (e.g., Aspirin, Ibuprofen) or SMILES string:

Aspirin

Try these examples:

Aspirin

Ibuprofen

Paracetamol

Metformin

Caffeine

Visualize Molecule

Molecular Weight

180.04 g/mol

H-Bond Donors

1

Rotatable Bonds

2

LogP


1.31

H-Bond Acceptors

3

Heavy Atoms

13

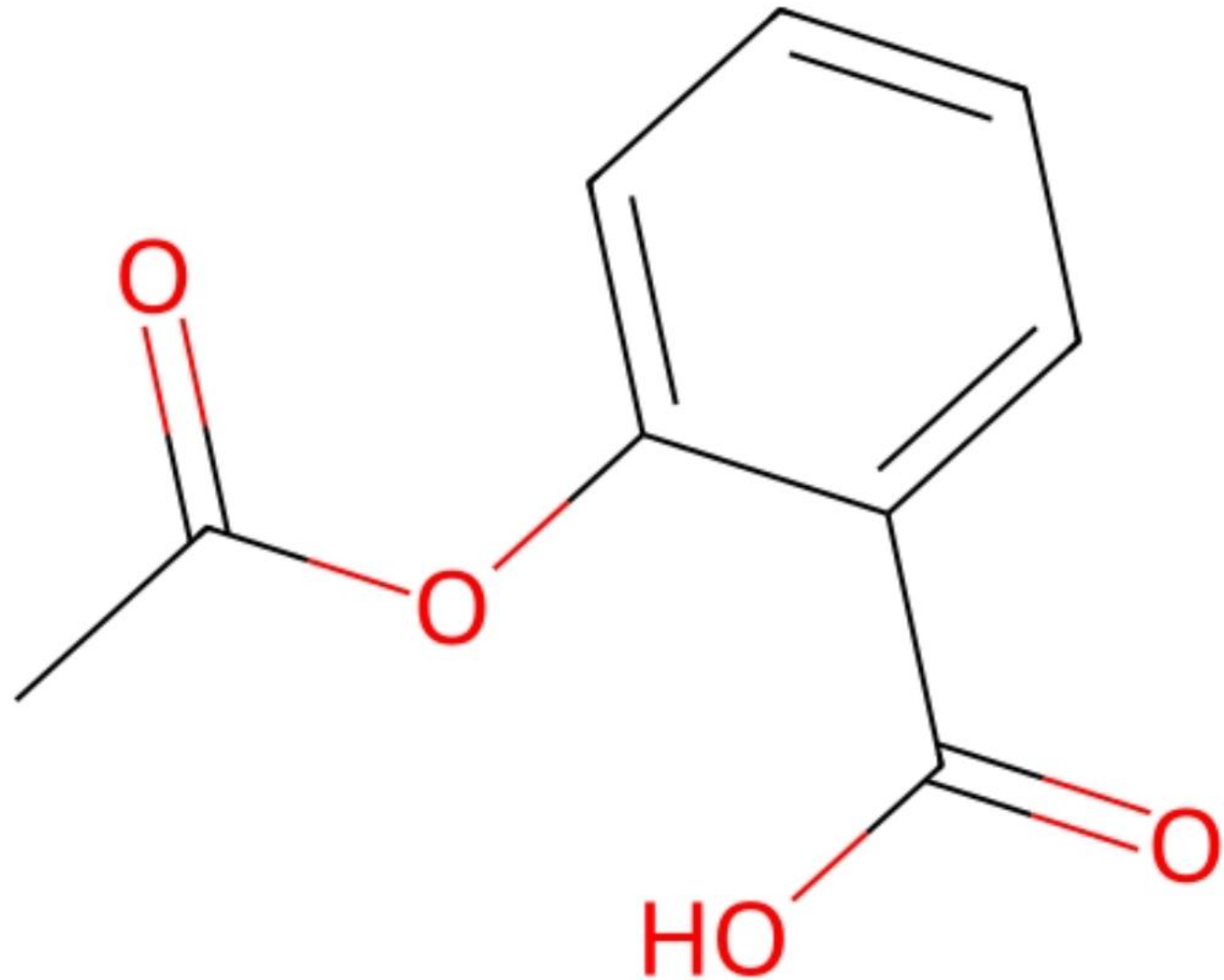
☐ Generate 3D structure (slower) 

2D Structure 3D Structure



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2D Structure



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Filters

Select Country

India ▼

Select Trade Indicator

Exports of goods and ... ▼

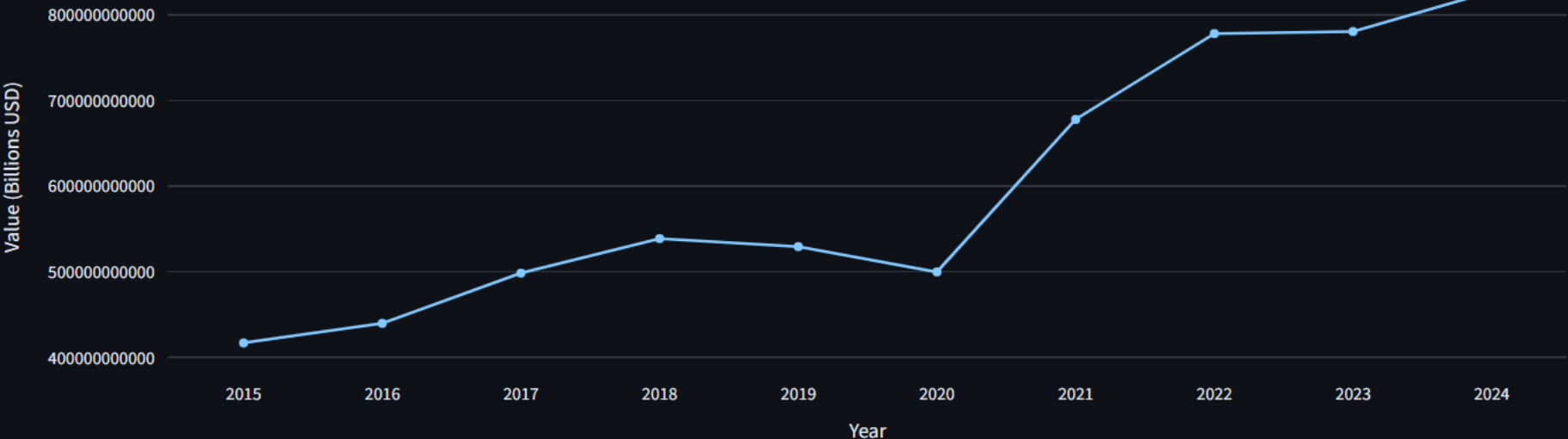
Select Year Range

2015 2024

↑ +6.0% from previous year

Exports of goods and services (current US\$) - India

Exports of goods and services (current US\$) - India (2015 - 2024)



Trade Balance

Trade Balance	Exports	Imports
\$-90.93B	\$827.41B	\$918.33B
↑ Deficit		

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Filters

Select Country

India

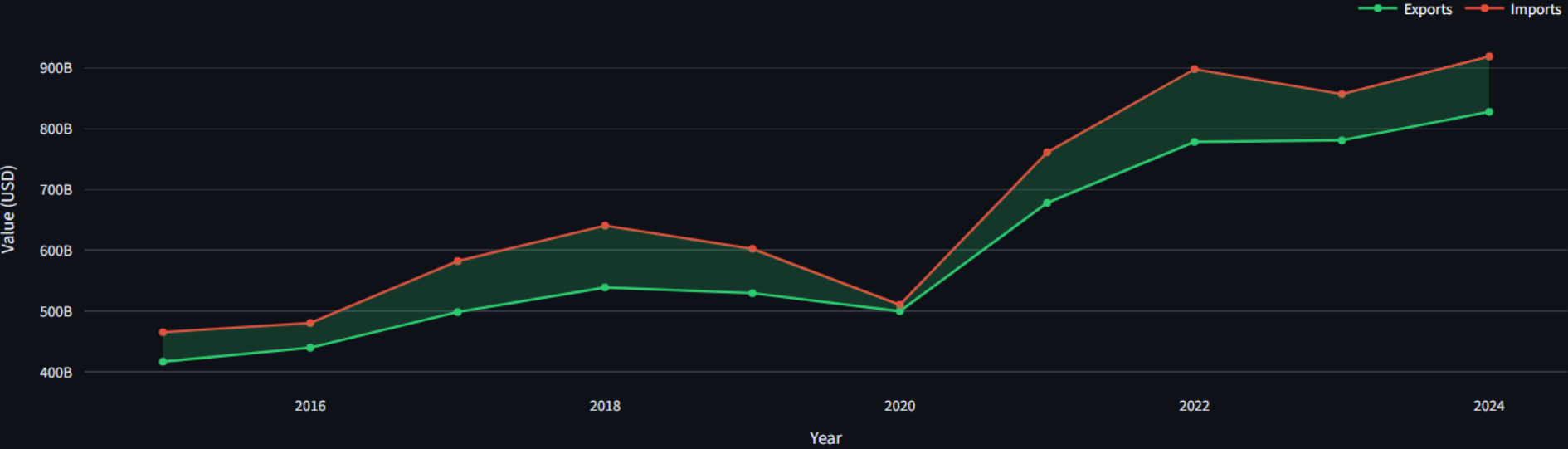
Select Trade Indicator

Exports of goods an...

Select Year Range

2015 2024

Exports vs Imports Over Time



View Raw Data

	Year	Exports of goods and services
9	2015	416787940847.812
8	2016	439642578076.358
7	2017	498258808231.521
6	2018	538635186135.438
5	2019	529245063167.215
4	2020	499728529072.629
3	2021	677769255428.552
2	2022	778021514754.531
1	2023	780408530088.596

