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drug discovery is the process through which potential new medicines are identified it involves a wide range of scientific disciplines including biology chemistry and pharmacology

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mit researchers have identified a new algorithmic framework that automatically identifies the best molecules to test for more streamlined drug discovery credits credit istock the use of ai to streamline drug discovery is exploding researchers are deploying machine learning models to help them identify molecules among billions of options

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a new ai powered program will allow researchers to level up their drug discovery efforts the program called topoformer was developed by an interdisciplinary team led by guowei wei a michigan

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phenotypic drug discovery has re emerged over the past decade as an approach to systematically pursue drug discovery based on therapeutic effects in realistic disease models

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the new model known as conplex can match target proteins with potential drug molecules without having to perform the computationally intensive step of calculating the molecules structures using this method the researchers can screen more than 100 million compounds in a single day much more than any existing model

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