



Your Quantum Prescriptions

How quantum algorithms can drive innovation in life sciences



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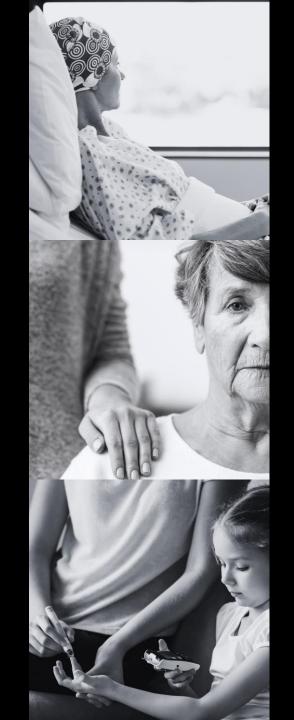
c.90% of drugs in the market

only work for 30-50% of the population

7% of all hospital admissions

Are due to adverse drug reactions

The current efficacy rate of standard drug treatments commands improvements



Cancer: 25%

Alzheimer's disease: 30%

Diabetes:

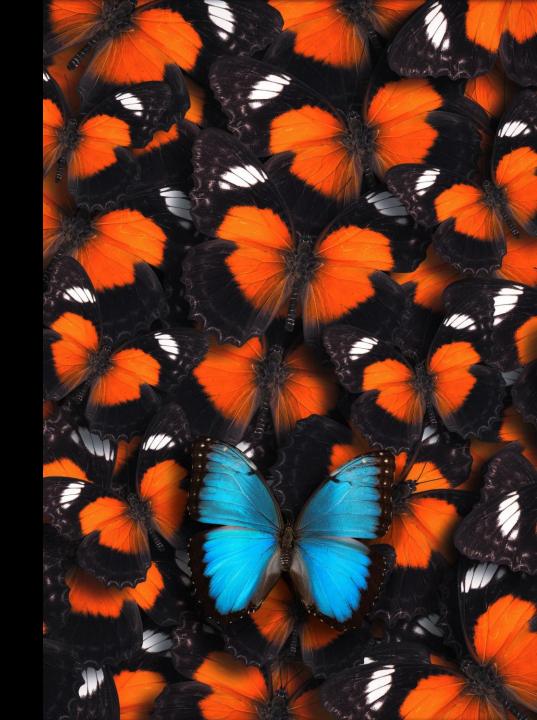
57%

Source: FDA

We need to focus our efforts on powering the paradigm shift in medical treatments

From "one size fits all"

To personalised medicine.



To truly understand each individual, we need to understand them at the deepest level

This journey begins with simulations at the molecular level, followed by interconnected systems, then organs, leading to a truly unique digital representation of the human body.



Quantum computing is driven by two key properties which enable exponential scale...



Quantum Superposition

Quantum Entanglement

... breaking the glass ceiling of traditional information processing and tapping into the power of nature at the smallest scales.

Quantum Bit
0 and 1



Hybrid Quantum Computing: Delivering Impact Today and Exponential Improvement in the Future









GPUs Transformed AI.

Now They're Here For Quantum.

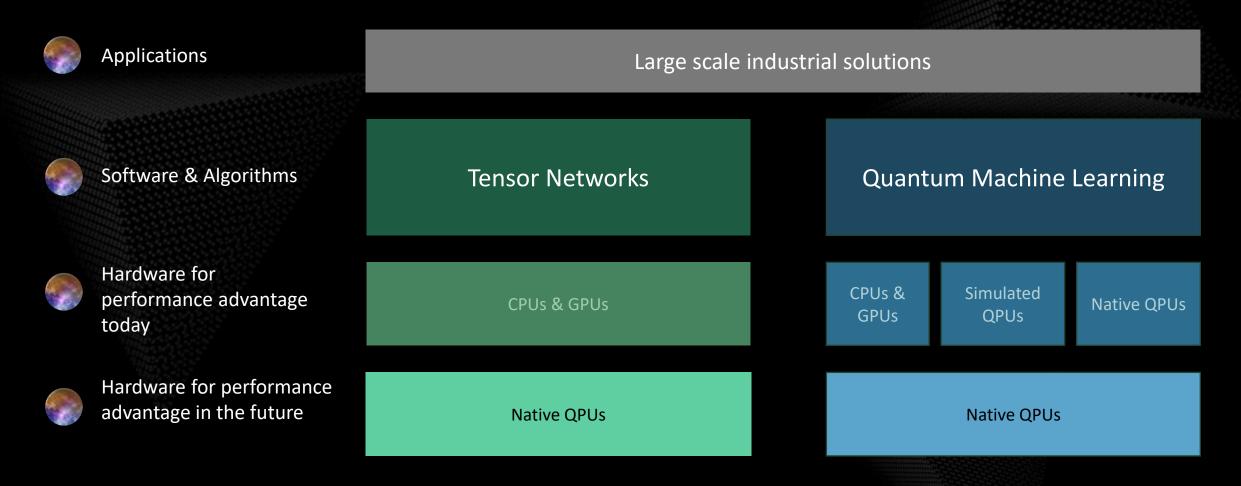


BEST COMPUTE PLATFORMS OF TODAY

Performance advantage on CPUs, GPUs, Simulated Qubits, etc NATIVE QUANTUM COMPUTERS

Improved performance on large scale QPUs of tomorrow

Leveraging our capabilities in Tensor Networks & Quantum Machine Learning



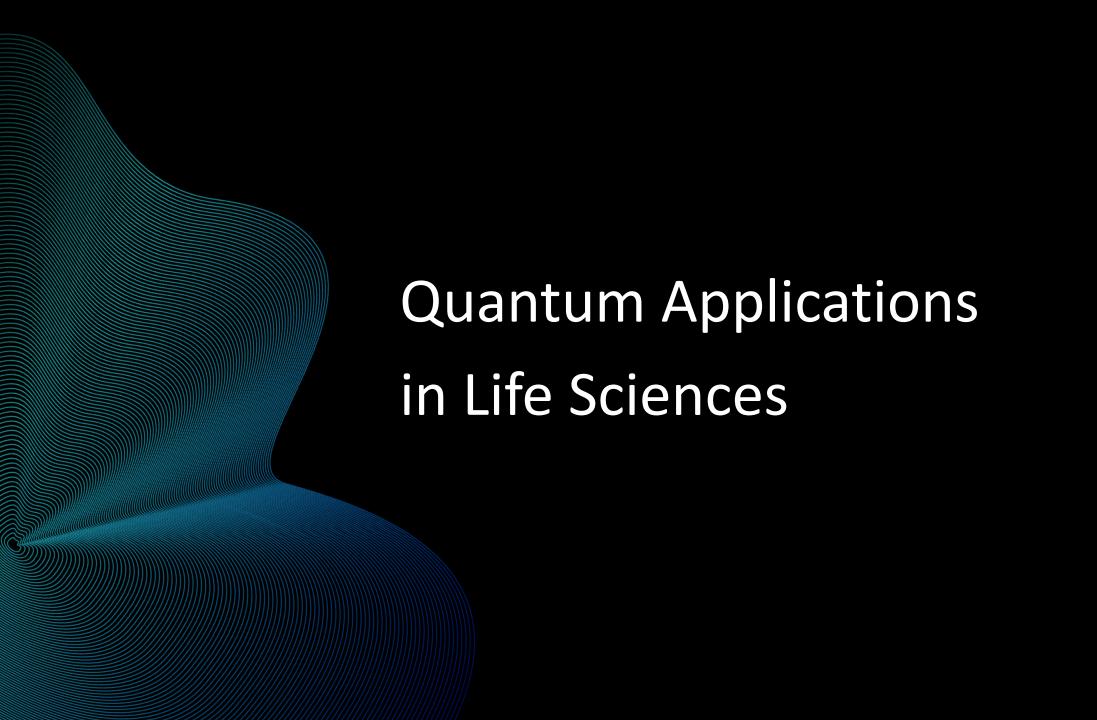
Quantum Software can deliver impactful solutions today

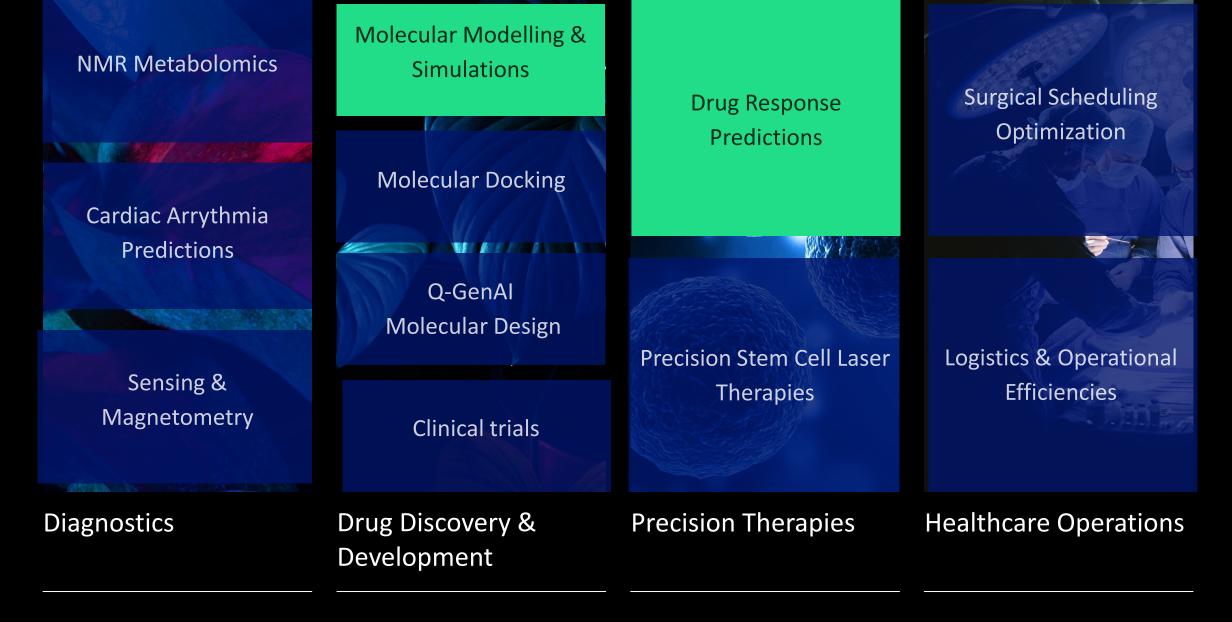








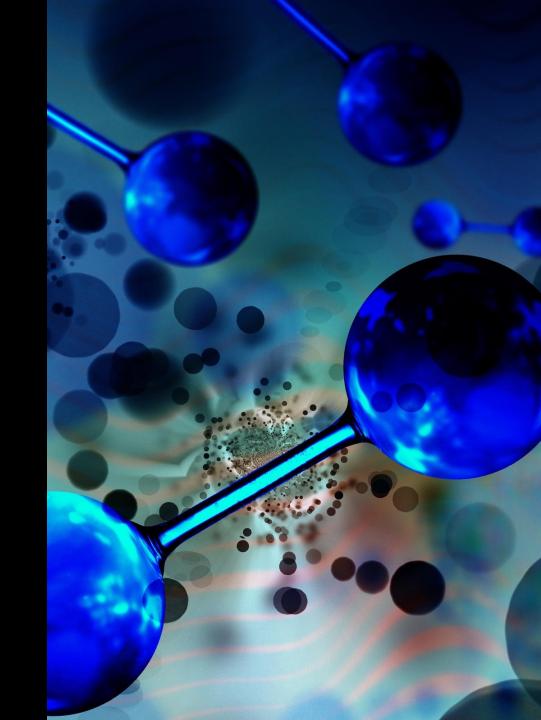




Our work spans various life sciences & healthcare applications

Shortening the Path
Towards Hypertargeted Treatments

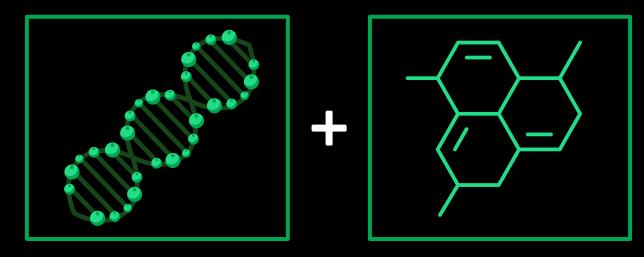
More accurate predictions with less data



Even when considering a particular type of cancer, there are many subtypes of cancer that exist, expressed in a large range of cell lines.

By understanding the structure of the cancer cell line and of the related drug, we are able to predict their interactions and hence anticipate the drug's effectiveness.

We ran a study of such interactions

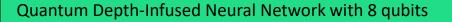


Encoded into a Neural Network of 256 nodes



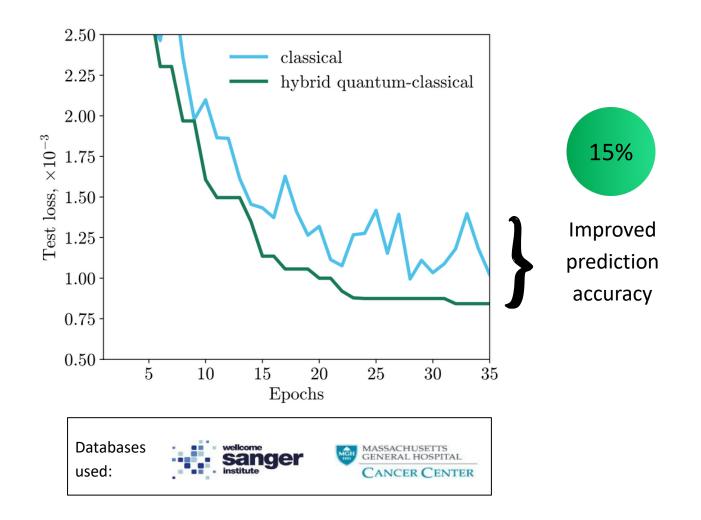




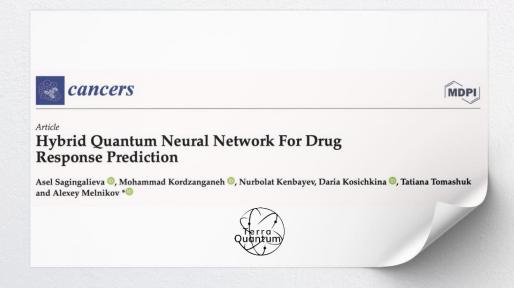




The Terra Quantum quantum-enhanced approach delivered 15% accuracy improvement in predicting the cancer drug response

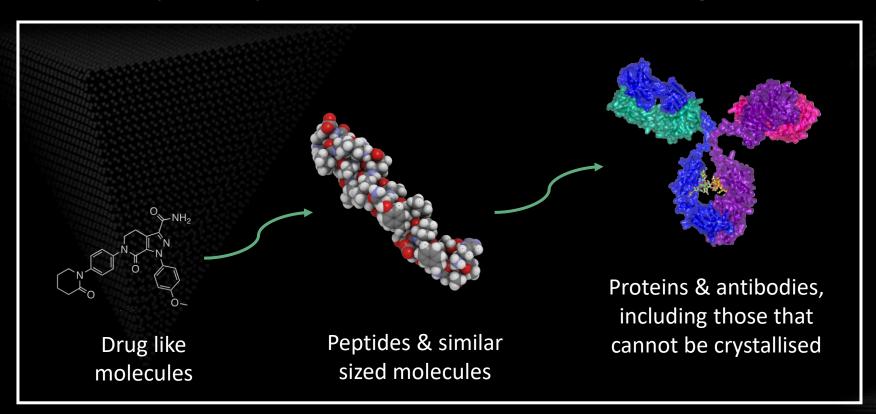


The results and detailed methodology are published in the Cancers Journal



Our vision is to create the world's most advanced molecular structure modelling tool that can accelerate drug development

Our pathway for molecular structure modelling



Additional future functionality

Optimisation of antibody sequences

Protein – protein interaction modelling

Broader optimisation of antibodies

Our methods can work with existing molecular structure prediction tools (such as Alphafold) but are not limited to the molecules considered here and can also produce greater levels of accuracy.



Key reasons why our solutions deliver benefits versus alternatives



Better quality solutions: Our methods are highly suited to navigating complex landscapes for optimisation problems and achieve better quality solutions with diversified answers.



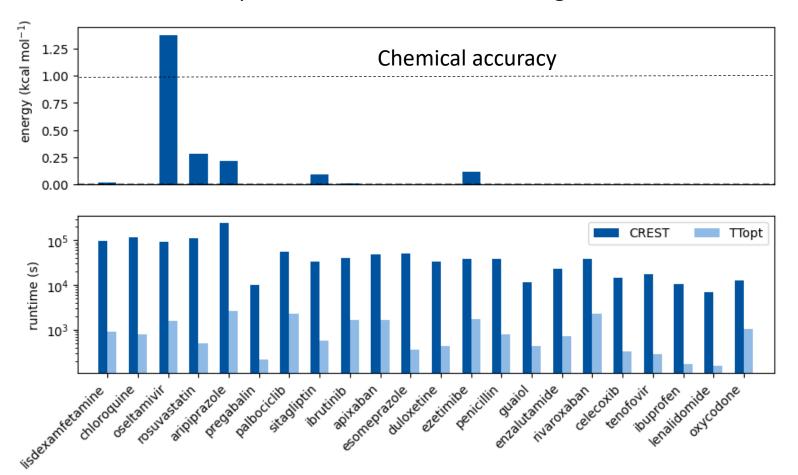
Significant speed up: The improved speed delivered by our solutions allows deployment of methods that were previously computationally too expensive or considered unfeasible.



Greater generalisability: Our solutions do not require training data, unlike many other machine learning / Al based methods. This allows our solutions to be generalised to cases where there is limited training data or of poor quality.

TTConf: Accelerating Conformer Sampling for Druglike Molecules

TTConf vs CREST: comparison of runtimes and energies



TTConf achieves chemical accuracy 10-100x faster than CREST

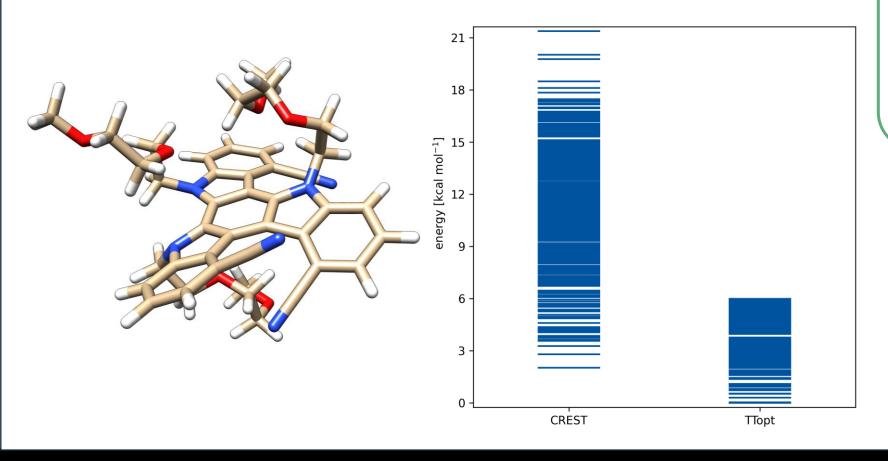
system specifications
Intel® Core™ i7-11700 @ 2.50 GHz
24 GB | 8 Threads

TTConf settings
2 sweeps | rank = 4
grid in degree = [30, 60, ..., 330]



TTConf: Optimizing Flexible Molecules and Beyond

Case Study: Optimizing a nonlinear Catalyst



TTConf achieves chemical accuracy in 2:38 h using 8 threads

CREST takes 12:30 h using 24 threads

Potential applications:

- Screening various databases
- Extension to tree tensor networks

Online-Demo of TTConf available



At Terra Quantum, we recognize the importance of advancing this field and have therefore made the healthcare domain a key part of our focus





Thank you!

Let's connect on LinkedIn







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