

# Scalable Quantum Simulation of Molecular Energies

James Mills<sup>1</sup>, Shuhao Yang<sup>1</sup>, and Shanice St John<sup>1</sup>

<sup>1</sup>MSc Quantum Technologies, UCL



## Introduction

Quantum computing is a rapidly advancing field predicted to revolutionise many areas of science and technology. It is predicted to have important real-world applications in encryption and communication systems and in the development of new medicines and materials to name but a few. An important point is that these quantum computers are not considered a replacement to classical computers, they will only be brought to bear on certain types of problems which are too difficult for classical computers to solve. Classical computers is a term used to describe all everyday computational computing devices which rely on classical (non-quantum) phenomena to compute.

Simulating systems in quantum chemistry is a problem too difficult for classical computers. Efficient simulation of quantum chemistry experiments would enable a dramatic leap forward in our understanding of fundamental chemistry. For example it would significantly reduce the need for cumbersome and expensive trial-and-error techniques in the development of new medicines and materials.

## Key terms

**Qubit** For representation, we describe information as qubit in quantum mechanics. It is an analogous concept of the classical bit in which classical information is encoded in 0s and 1s. For the information stored in a qubit can be represented by a state.

**Superposition** A qubit has extraordinary property, which can be in both states of '0' and '1' at the same time, but a classical bit is either '0' or '1'.

**Quantum simulations in chemistry** Quantum theory is our best description of phenomena happening at the smallest scales imaginable. Simulating quantum systems with classical computers is extremely difficult, and becomes increasingly difficult as the size of the system increases. Quantum simulations with quantum computing hardware are much easier, as we are using quantum phenomena (the same phenomena that make the system difficult to describe with a classical computer) to boost our computing power.

**Algorithm** A set of instructions used to solve a problem, especially by a computer. The instructions are created so that it can be understood by the computer. This is then sent to the quantum computer. The computer then follows the instructions using some program or software. The end of the process provides an answer or even many possible answers.

## Techniques in Paper

The work by the researchers on the paper 'Scalable Quantum Simulation of Molecular Energies' uses supercooled quantum qubits to run algorithms to find out information about the fundamental properties of molecules. The variational quantum eigensolver (VQE) and the phase estimation algorithm (PEA) are two types of algorithm which are performed on the qubit hardware to find molecular energies. A special type of quantum hardware known as Superconducting Quantum Interference Devices (SQUIDs) is used to run the algorithms. This type of qubit has a superposition of "charged" states and can be used in quantum circuits.

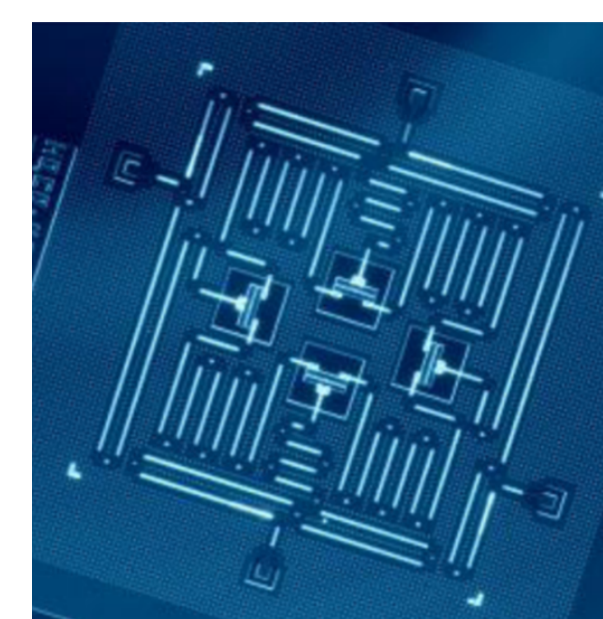


Figure 1: IBM qubit

The qubits used to run the algorithms were a type of charge qubit (the qubit is in a superposition of two different charges - see Superposition box) that is kept at very low temperatures. It is similar to the IMB qubit shown in Figure 1.

## VQE

The VQE algorithm uses the averages of the total energy so that the energy of a molecule can be calculated.

- Firstly, we start by preparing a state using a "special" method. This method makes it possible for the algorithm run in the way that we want.
- Secondly, we measure the total energy for each distance between the molecules. We must select the lowest value as we are focusing on the ground energy.
- After that, a tool is then used to find more lower values of the total energy.
- All steps are repeated until we reach the smallest value possible.

## PEA

PEA is the algorithm estimate the value of index over exponential coefficient in order to work out energy of a molecule.

- Similarly to the VQE algorithm, we must prepare a state using another "special" method since these are two different algorithms.
- Next, some "magical" equipment must be set up in a superposition of states - this will make it possible to measure the state.
- Then, we use some operations to set up the state and the apparatus so that it is ready to be measured.
- Finally, we measure the state to find the lowest value of the total energy. We repeat this process until find an accurate value of the total energy.

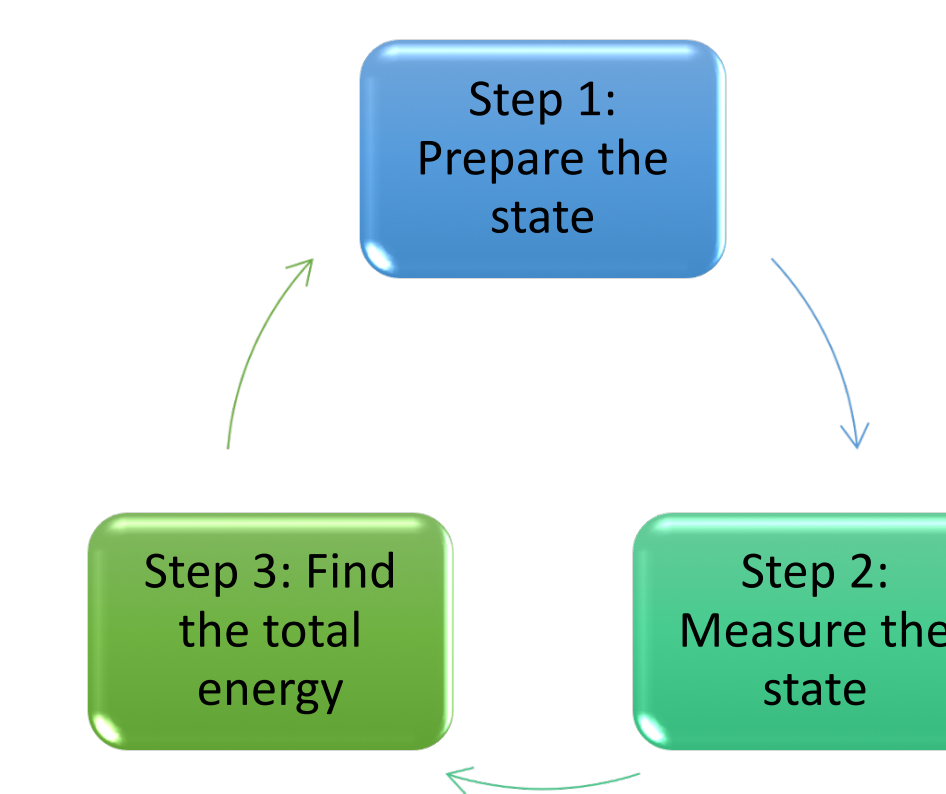


Figure 2: VQE

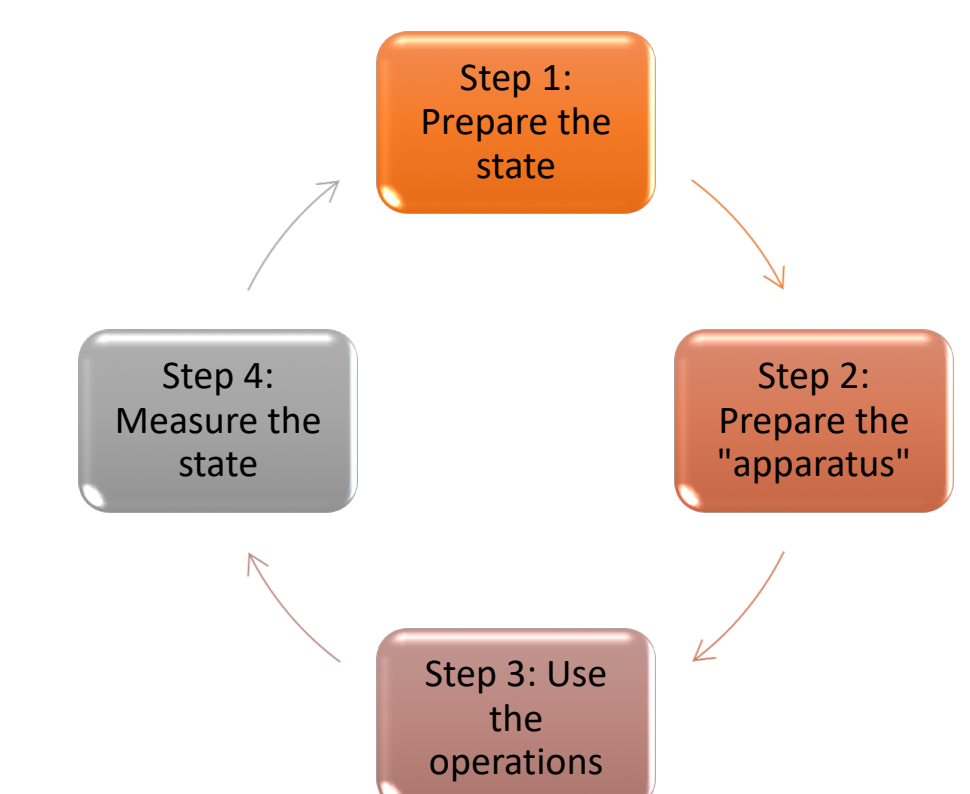


Figure 3: PEA

## Conclusion and Outlook

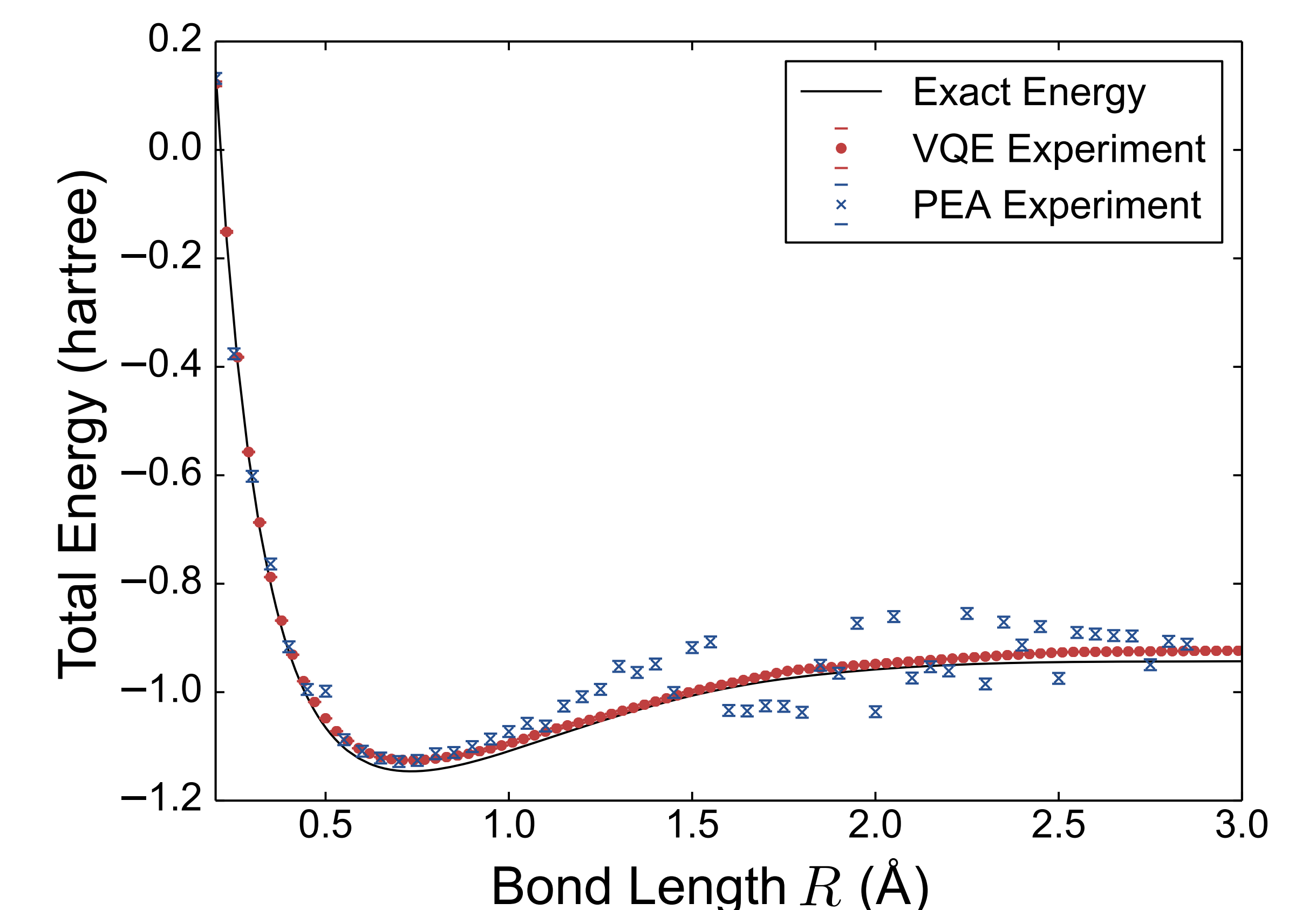


Figure 4: Computed  $H_2$  energy curve, energy surface of molecular hydrogen as determined by both VQE and PEA



VQE was proven to work better than PEA because it gives more accurate results. Therefore, it is used more than PEA. Creating more algorithms such as VQE can

lead to making new medicines, materials and other cool applications!