

# Using the variational quantum eigensolver to study a lattice fermion model

Ronak Desai<sup>1</sup> and Michael McGuigan<sup>2</sup>

<sup>1</sup>Department of Physics and Astronomy, Rowan University, Glassboro, New Jersey 08028, USA <sup>2</sup>Computational Science Initiative, Brookhaven National Laboratory, Upton, New York 11973, USA

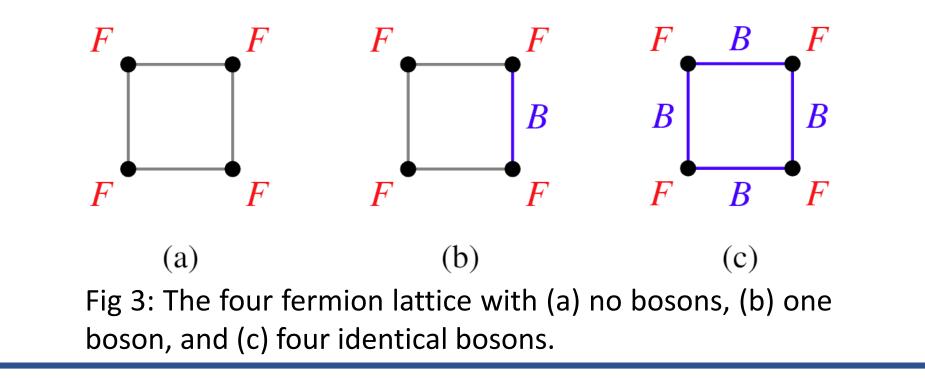
### ABSTRACT

Although computing power has been steadily rising over the years, there are some problems that classical computers simply cannot solve. Many lattice gauge theories used in nuclear physics contain such computationally complex problems that cannot be solved in a reasonable time. However, advances in low-temperature physics allow these models to be simulated by ultracold atoms. These cold atom systems can be studied by quantum computers which have recently been made available to the public via the cloud from IBM. Since quantum computers have the capability to solve certain problems exponentially faster than classical computers, we decided to study a lattice fermion model using the IBM's Qiskit. Specifically, we used the variational quantum eigensolver algorithm to find the ground state energy of several similar systems. We report our findings using both a local simulator (Qiskit) and an actual quantum computer (IBMQ).

#### MODEL

Our model describes four fermions in a square lattice where nearest neighbors can interact through a boson link. The hamiltonian is given by  $\hat{H} = (\bigotimes^{j-1} 1_3) \otimes \frac{\hat{P}^2}{2} \otimes (\bigotimes^{n_B - j} 1_3) \otimes 1_{16} - \sum_{i=i}^4 c_{Fi}^{\dagger} A_{i,j} c_{Fj}$ 

Here,  $\otimes$  is the tensor product,  $\hat{P}$  is the position basis representation of momentum,  $A_{i,j}$  is the adjacency matrix accounting for fermion-boson interactions, and  $c_{Fi}$  is the fermion operator.





## QUANTUM COMPUTING

Quantum computers are machines that rely on the physics of quantum mechanical two-level systems to perform computations. They operate in a fundamentally different way than classical computers by using qubits instead of bits.

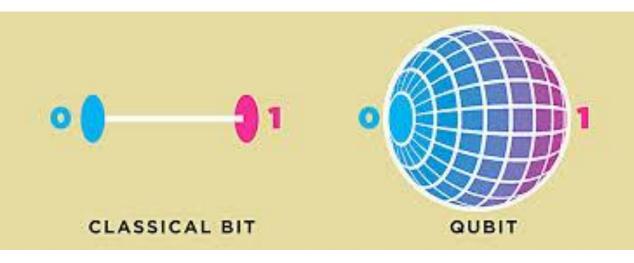


Fig 1: Qubits exhibit superposition and entanglement which quantum computers leverage to perform computations that have the capability to outclass modern supercomputers.

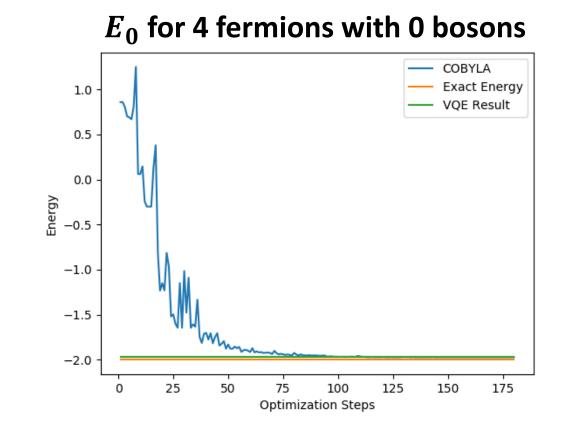
Image Credit: https://www.austinchronicle.com/screens/2019-04-19/quantum-computing-101-abeginners-guide-to-the-mind-bending-new-technology/

## VARIATIONAL QUANTUM EIGENSOLVER

We use IBM's quantum computing platform Qiskit [1] to run the variational quantum eigensolver (VQE) algorithm [2] on a lattice model in nuclear physics. The VQE finds the lowest eigenvalue (ground state energy) of a Hermitian matrix (hamiltonian). For each run, we specify the number of shots (repeated quantum calculation to generate adequate statistics) and trials (optimization steps).

## RESULTS

Using the local simulator, we report the VQE converge plots which show the estimate of the ground state energy  $(E_0)$  as the optimization subroutine progresses.



#### $E_0$ for 4 fermions with 1 boson

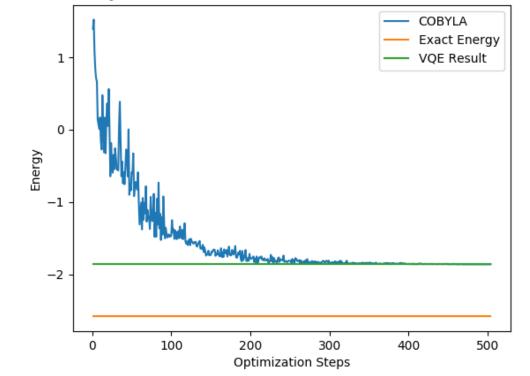
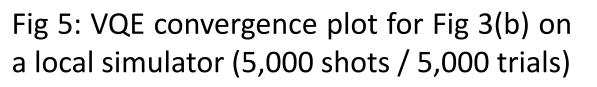


Fig 4: VQE convergence plot for Fig 3(a) on a local simulator (500 shots / 500 trials)

Fig 6 shows that IBMQ cannot run through many optimization steps and is not able to converge



#### $E_0$ for 4 fermions with 0 bosons (IBMQ)



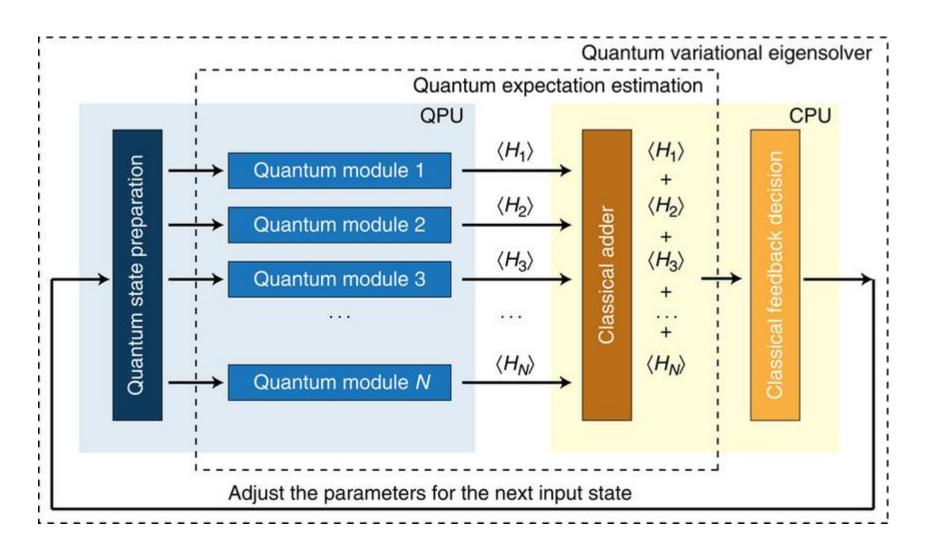


Fig 2: The flow chart of the VQE algorithm (from Peruzzo et al.) which is based on the quantum mechanical variational method. The VQE takes advantage of quantum hardware that can perform parallel computation and is guaranteed to find an upper bound for the ground state energy.

to correct value. This is most likely due to limitations on what the public can access. The quantum computer ran at a rate around 60x slower than the local simulator for obtaining ground state energy of Fig 3(a).

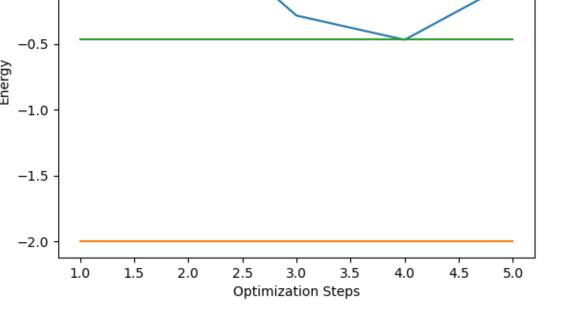


Fig 6: VQE convergence plot for Fig 3(a) on IBM quantum computer (500 shots / 5 trials)

#### FUTURE WORK

We hope that we can find better results for more complicated systems in the future by finding a good trial wave function for the VQE and finding a way to simulate many qubits without creating large matrices.

#### REFERENCES

[1] A. Peruzzo et al., "A variational eigenvalue solver on a photonic quantum processor," Nature Communications 5, 4213 (2014).

[2] G. Aleksandrowicz et al., "Qiskit: An Open-source Framework for Quantum Computing," (2019).



ACKNOWLEDGMENTS

This project was supported in part by the U.S. Department of Energy, Office of Science,

**U.S. DEPARTMENT OF** 

Office of Workforce Development for Teachers and Scientists (WDTS) under the Science

Undergraduate Laboratory Internships Program (SULI). We acknowledge use of the IBM

Q for this work. The views expressed are those of the authors and do not reflect the

official policy or position of IBM or the IBM Q team.

