## Jack Ceroni

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<b>Education</b> September 2020— Present	<ul> <li>Degree: Bachelor of Science in Pure Mathematics</li> <li>Where: University of Toronto, Pure Math Specialist Program</li> <li>Cumulative GPA: 4.0/4.0</li> </ul>
Employment May 2023— Present	<ul><li><b>Position:</b> Visiting Research Student</li><li><b>Where:</b> Massachusetts Institute of Technology</li></ul>
	• Working in Isaac Chuang's research group in the Research Lab- oratory of Electronics.
	• Collaborating with Zane M. Rossi on research in fault-tolerant quantum algorithms.
September 2022— Present	<b>Position:</b> Research Scientist <b>Where:</b> Stealth Startup
	• Was the first (non-founder) employee at a quantum comput- ing/machine learning startup, started by former team leads at Google, IBM, and BMW Quantum. Took a gap year from school to help build the initial versions of the company's core technol- ogy.
	• Built the core code-base utilized for quantum simulation and probabilistic machine learning applications.
	• Led the development of novel theoretical work and error bounds related to quantum algorithms for simulating large thermal states.
April 2021— September 2021	<b>Position:</b> Quantum Algorithms Research Intern <b>Where:</b> Xanadu Quantum Technologies
	• Conducted research on quantum machine learning for generat- ing approximate molecular ground states. Developed sample bounds for gradient calculations, Cramer-Rao lower bounds on data requirements for the model, and performed numerical sim- ulations.
	• Explored other research directions in the regime of fault-tolerant quantum algorithms for quantum chemistry, namely QSVT-based algorithms for simulation of molecular vibronic structure.

April 2021— September 2021	<ul><li>Position: Quantum Research Resident</li><li>Where: Xanadu Quantum Technologies</li></ul>		
	<ul> <li>Conducted research on performing end-to-end quantum chemistry simulations on quantum computers, and new techniques for computing accurate energy derivatives on quantum devices.</li> <li>Built software on top of the PennyLane library for simulating</li> </ul>		
	quantum computational chemistry calculations.		
	• Developed the initial version of an automatically differentiable Hartree-Fock solver, for computing derivatives of molecular Hamil- tonians and atomic basis set parameters.		
June 2020— October 2020	<ul><li>Position: Quantum Software Development Intern</li><li>Where: Xanadu Quantum Technologies</li></ul>		
	• Focused on development of Xanadu's quantum machine learning software library, PennyLane.		
	• Published tutorials on cutting-edge variational quantum algo- rithms implemented using PennyLane, such as the Variational Quantum Thermalizer, the Quantum Graph Neural Network, and the Quantum Approximate Optimization Algorithm (QAOA).		
	• Led the development effort to incorporate native QAOA func- tionality into PennyLane, in the pennylane.qaoa module.		
<b>Other Experience</b> September 2022— Present	<b>Position:</b> Research Student <b>Where:</b> University of Toronto		
	• Working with Prof. Nathan Wiebe on research related to ex- tensions of Quantum Signal Processing and Quantum Singular Value Transforms, and novel techniques for Hamiltonian simu- lation.		
July 2019— July 2019	<ul><li>Position: Summer Student</li><li>Where: The Perimeter Institute for Theoretical Physics</li></ul>		
	• Participated in the International Summer School for Young Physicists at the Perimeter Institute.		
	• Took classes on basic concepts in modern physics (quantum mechanics, special relativity, etc.).		
	• Participated in a small seminar course on introductory Standard Model particle physics.		
	• Worked on a quantum information project with one of the IS-SYP mentors, Prof. Jamie Sikora, for several months after the program ended.		

October 2019—	<b>Position:</b> Qiskit Advocate	
Present	Where:	IBM Quantum

• Credited with writing two sections for the Qiskit textbook on introductory linear algebra for quantum computing, and the Variational Quantum Linear Solver algorithm.

## Selected Preprints and Publications

- Jack Ceroni, Torin F Stetina, Maria Kieferova, Carlos Ortiz Marrero, Juan Miguel Arrazola, Nathan Wiebe "Generating Approximate Ground States of Molecules Using Quantum Machine Learning." arXiv preprint arXiv:2210.05489 (2022).
- Jack Ceroni, Alain Delgado, Soran Jahangiri, Juan Miguel Arrazola "Tailgating quantum circuits for high-order energy derivatives." arXiv preprint arXiv:2207.11274 (2022).
- Juan Miguel Arrazola, Soran Jahangiri, Alain Delgado, **Jack Ceroni**, Josh Izaac, Antal Száva, Utkarsh Azad et al. "Differentiable quantum computational chemistry with PennyLane." arXiv preprint arXiv:2111.09967 (2021).

## Invited Talks

- North Carolina State University Quantum Workshop (January 2023). Talk title: Generating Approximate Ground States of Molecules Using Quantum Machine Learning.
- Quantum Technology and Application Consortium Journal Club (December 2022). Talk title: Generating Approximate Ground States of Molecules Using Quantum Machine Learning.

## Honors and awards

- Weston Youth Innovation Award (2018)
- University of Toronto Scholars Award (2020)
- MITACS Accelerate Grant (2021, 2022)
- Haylon Chan Memorial Award (2022)
- University of Toronto Dean's List (2021, 2022, 2023)