

## EDUCATION

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### California Institute of Technology (Caltech)

*Ph. D. in Theoretical Chemistry*

*Advisor:* Garnet Kin-Lic Chan

*Thesis:* Finite temperature simulations of strongly correlated systems.

**Pasadena, CA**

*Jul. 2016 - Dec. 2020*

### Peking University

*B. S. in Chemistry with honor*

*Thesis advisor:* Wenjian Liu

*Thesis:* Density matrix embedding theory in terms of localized molecular orbitals.

**Beijing, China**

*Sep. 2011 - Jul. 2015*

### Princeton University

*Centennial Fellowship graduate Student*

**Princeton, NJ**

*Sep. 2015 - Jun. 2016*

### Massachusetts Institute of Technology (MIT)

*Exchange undergraduate Student*

**Cambridge, MA**

*Sep. 2013 - Dec. 2013*

## SKILLS

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**Research Topics:** quantum chemistry, numerical simulation, quantum computing, machine learning.

**Coding & Software:** Python, C/C++, PySCF, OpenFermion, PyTorch, JAX.

## RESEARCH & WORK EXPERIENCE

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### Research scientist, Zapata Computing Inc.

*May. 2022 - Jan. 2023*

*Supervisor:* Dr. Peter Johnson ([URL for website](#))

- Developed quantum algorithms and software for ground-state energy estimation and quantum resource estimation.

### Postdoctoral fellow, University of Toronto

*Feb. 2021 - Oct. 2022*

*Supervisor:* Prof. Alán Aspuru-Guzik ([URL for website](#))

- Extended normalizing flows to be able to preserve the anti-symmetric feature of fermionic wavefunctions, which provided an alternative view to encoding antisymmetry with the Slater determinant.
- Co-developed quantum iterative power algorithm (QIPA) to accelerate solving large number factorization and ground-state energy estimation.

### Graduate research assistant, Caltech

*Jul. 2016 - Dec. 2020*

*Supervisor:* Prof. Garnet Kin-Lic Chan ([URL for website](#))

- Developed the finite-temperature density matrix embedding theory (FT-DMET) algorithm for strongly-correlated systems.
- Implemented the finite temperature extension of existing numerical algorithms, including exact diagonalization, Lanczos algorithm, density matrix renormalization group (DMRG), and spin-symmetry restored coupled cluster (CC).
- Developed quantum imaginary time evolution (QITE) algorithm and quantum minimally entangled typical thermal states (QMETTS) algorithm.

### Undergraduate research assistant, Peking University

*Jan. 2015 - Jun. 2015*

*Supervisor:* Prof. Wenjian Liu ([URL for website](#))

- Implemented localized molecular orbitals with the Foster-Boys localization method in the density matrix embedding theory (DMET) method to study molecular systems.

### Undergraduate research assistant, Peking University

*Jul. 2014 - Nov. 2014*

Supervisor: Prof. Hong Jiang ([URL for website](#))

- Studied the low-spin to high-spin (LS-HS) phase transition of  $[\text{Fe}(\text{pz})\text{Pt}(\text{CN})_4]\cdot 2\text{H}_2\text{O}$  with Monte Carlo and density functional theory (DFT).

## AWARDS & HONORS

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| • Barbara J. Burger Fellowship                                   | <i>Caltech, 2019</i>              |
| • CCE Teaching (TA) Award  | <i>Caltech, 2018</i>              |
| • Centennial Fellowship in the Natural Science and Engineering   | <i>Princeton University, 2015</i> |
| • Chun-Tsung Scholarship   | <i>Peking University, 2014</i>    |
| • Merit Student  | <i>Peking University, 2014</i>    |
| • Pan Gu Fellowship  | <i>Peking University, 2013</i>    |
| • First Prize in the 29th National Undergraduate Physics Contest | <i>Beijing, 2012</i>              |

## INVITED TALKS

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|---|--------------------------------|
| • Density matrix embedding theory and quantum imaginary time evolution. | <i>ByteDance, 2021</i>         |
| • Quantum imaginary time evolution.                                     | <i>QIP, 2020</i>               |
| • Quantum computing for quantum chemistry.                              | <i>Peking University, 2020</i> |

## TEACHING EXPERIENCE

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### University of Toronto

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| • CHM427H1S Statistical Mechanics - course instructor. | <i>Spring 2021-2022</i> |
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### Californian Institute of Technology

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|---|-------------------------|
| • CH125a Quantum Mechanics (A) - recitation instructor.             | <i>Fall 2016-2017</i>   |
| • CH21b Physical Chemistry (B) - recitation instructor.             | <i>Winter 2016-2017</i> |
| • CH125b The Elements of Quantum Chemistry - recitation instructor. | <i>Winter 2017-2018</i> |

## SELECTED PUBLICATIONS

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- L. Thiede, **C. Sun**, A. Aspuru-Guzik. Waveflow: Enforcing boundary conditions in smooth normalizing flows with application to fermionic wave functions. Pre-print: [arXiv:2211.14839](#) Submitted to ICML 2023.
- K. Gratsea, **C. Sun**, P. Johnson. When to Reject a Ground-State Preparation Algorithm. Pre-print: [arXiv:2212.09492](#).
- **C. Sun**, U. Ray, Z.-H. Cui, M. Stoudenmire, M. Ferrero and G. K. Chan. Finite temperature density matrix embedding theory. *Phys. Rev. B*, **101**, 075131 (2020)
- M. Motta, **C. Sun**, A. T. K. Tan, M. J. O' Rourke, E. Ye, A. J. Minnich, F. G. S. L. Brandao, G. K. Chan. Determining eigenstates and thermal states on a quantum computer using quantum imaginary time evolution. *Nature Physics*, **16**, 205 (2020)
- Z.-H. Cui, **C. Sun**, U. Ray, B.-X. Zheng, Q. Sun, G. K. Chan, Ground-state phase diagram of the three-band Hubbard model in various parametrizations from density matrix embedding theory. *Phys. Rev. Research*, **2**, 043259 (2020)
- **C. Sun**, Finite Temperature Simulations of Strongly Correlated Systems. *Dissertation (Ph.D.)*, *California Institute of Technology*. doi:10.7907/dchn-p020 (2021)

- H.-Z. Ye, C. **Sun** and H. Jiang. Monte-Carlo Simulation of Spin-Crossover Phenomena Based on a Vibronic Ising-like Model with Realistic Parameters. *Phys. Chem. Chem. Phys.*, **17**, 6801 (2015)