

# George Lai Gie Ho

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eigenfoo.github.io  
georgeho1618@gmail.com

Education	<b>The Cooper Union, New York, NY.</b> Bachelor of Science in Engineering, General Engineering. Projected May 2019. <i>Relevant Coursework:</i> Artificial Intelligence, Machine Learning, Data Structures and Algorithms, Numerical Analysis, Linear Algebra	
Experience	<b>Quantitative Research and Investments Intern</b>	Summer 2017
	Quantopian Inc., Boston <ul style="list-style-type: none"><li>Developed open-source and in-house Python libraries for portfolio risk analysis and performance attribution</li><li>Maintained, developed and managed 3 open-source Python libraries: collaborated with worldwide team to coordinate significant feature additions</li><li>Operationalized risk analysis and performance attribution of institutional fund portfolio, leading to significant restructuring of fund-level portfolio</li><li>Contributed to crowd-sourced algorithm selection process using machine learning techniques</li></ul>	
	<b>Undergraduate Research Fellow</b>	Summer-Fall 2017
	Complex Fluid Physics and Engineering (CoFPhE) Laboratory <ul style="list-style-type: none"><li>Performing direct numerical simulation of complex ABC fluid flow in Python and Fortran to determine roles of density, Stokes number and Coriolis forces on inertial transport</li><li>Work will be presented at 2017 annual meeting of the American Physical Society Division of Fluid Dynamics at Denver, Colorado</li></ul>	
Projects	<b>Independent Projects in Algorithmic Trading</b>	Fall 2016
	Quantopian, Cornell University <ul style="list-style-type: none"><li>Analyzed minutely market pricing data using machine learning algorithms to research and develop a pairs-trading strategy in Python</li><li>Won 3rd place in Cornell University's Sparkstone Algorithmic Trading Challenge (out of hundreds of competitors) using a momentum long-short strategy</li></ul>	
	<b>Molecular Modeling of Ammonium Halide Nanoparticles</b>	Spring 2016
	The Cooper Union <ul style="list-style-type: none"><li>Wrote MATLAB scripts and performed quantum calculations to develop a molecular model for ammonium fluoride nanoclusters</li><li>Presented findings at the 2016 Meeting of the American Chemical Society</li></ul>	
Skills	<ul style="list-style-type: none"><li><i>Programming Languages:</i> Fluent in Python. Functional in C++ and MATLAB.</li><li><i>Software Packages and Libraries:</i> NumPy, pandas, matplotlib.</li><li><i>Languages:</i> Functional in Mandarin, Cantonese and French</li></ul>	
Awards	<ul style="list-style-type: none"><li>Cooper Union Half-Tuition Scholarship</li><li>Cooper Union Innovator Merit Scholarship</li><li>Dean's List (School Honors)</li></ul>	<ul style="list-style-type: none"><li>2015 - 2019</li><li>2015 - 2019</li><li>All semesters</li></ul>
References	Available upon request.	