


CHRISTOPHER R. COLLINS

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 <http://github.com/crcollins>

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EDUCATION

- **Carnegie Mellon University** **Pittsburgh, PA**
Ph.D. in Chemistry 2014 – 2019 Expected
- **University of North Georgia** **Dahlonega, GA**
B.S. in Chemistry with Chemical Physics Concentration 2010 – 2014

RESEARCH

- **Carnegie Mellon University** **Pittsburgh, PA**
Graduate Student; Advisor: David Yaron 2014 – Present
 - Developed parallel and distributed computing code for optimizing INDO parameters
 - Predicted optoelectronic properties of various polymeric systems using Deep Learning
 - Developed new descriptor for machine learning models in chemistry
 - Advised undergraduate students in machine learning and Python project
- **University of North Georgia** **Dahlonega, GA**
Undergraduate Researcher; Advisor: Aimée Tomlinson 2011 – 2014
 - Researched Benzobisazole Cruciform structures using computational chemistry for organic solar cells
 - Designed and wrote set of tools to automate research process (Chemtools/Chemtools-webapp)
 - Instructed other group members about Unix usage methods
 - Led group members in structure studies (multicore systems, donor-acceptor systems, increased rings systems)
 - Developed machine learning model for predicting Benzobisazole properties
- **Carnegie Mellon University** **Pittsburgh, PA**
Undergraduate Researcher; Advisor: David Yaron Summer 2013
 - Researched decomposition of Thiaphospholes using computational methods
 - Wrote MATLAB code to automate running Gaussian/AMPAC calculations and parsing results
 - Wrote code to calculate molecular overlaps and generate molecular orbital diagrams
 - Created routines to draw structures with molecular orbital overlaps indicated on each atom

WORK EXPERIENCE

- **Tripwire, Inc.** **Alpharetta, GA**
Engineering Intern 2012 – 2014
 - Rewrote entire test suite for Benchmark product using Selenium and PyTest (Python)
 - Aided in migration of critical IP360 product and wrote automated test suite for API
 - Maintained and expanded installation automation for CCM product
 - Wrote automated script to test for CSRF and XSS vulnerabilities in IP360 product

TEACHING

- Teaching Assistant
 - 09-221 Laboratory I: Introduction to Chemical Analysis
 - 09-214 Physical Chemistry
 - 09-101 Introduction to Experimental Chemistry (2 times)
 - 09-103 Atoms, Molecules, and Chemical Change
- Weekly Lecture Series in Theory Suite (Organizer)
 - Intro to Python
 - Intro to Bash Scripting (2 lectures)
 - Intro to Object Oriented Programming
 - Intro to NumPy (2 lectures)

SIDE PROJECTS

- **Chemtools/Chemtools-Webapp** (<http://gauss.crcollins.com>)
 - Implemented a set of tools for use in Benzobisazole research
 - * Gaussian log output parser to collect molecular properties
 - * Machine learning model to predict polymer properties from monomers
 - * Benzobisazole structure generator and machine learning predictor for optoelectronic properties
 - * Supercomputer job submitter and curator
 - Full automated testing suite (95% code coverage)
 - Python/Django Application with a Bootstrap frontend
 - Vagrant VM build system
- **pyOS**
 - Implementation of a *nix like operating system using Python
 - Created shell similar to Bash Shell
 - Reimplemented standard unix programs (cp, mv, rm, ls, and etc)
 - Included permissions, pipes, multiple users, and a virtual file system
- **Other**
 - 8 bit CPU (Logisim)
 - Mandelbrot/Buddhabrot Generator (Python, Javascript, C/C++ , CUDA C, TI-BASIC)
 - Various Quadtree Visualizations (Python, C/C++ , Javascript)
 - Cellular Automata Based Electronic Simulation (Javascript)
 - Object Relation Management Library (Python/SQL)
 - University Course Scraper (Python)
 - Cryptography Library (Python, Logisim)
 - Automated Peer Review System Web Application (Python/Django)
 - Virtual Cluster and Server Architecture Vagrant Builds

PUBLICATIONS

- **Collins, C. R.** ; Gordon, G. J.; Yaron, D. J. Constant Size Molecular Descriptors For Use With Machine Learning. *Submitted.*
- **Collins, C. R.** ; Tomlinson, A. L. Application of Machine Learning to Predict the Optoelectronic Properties of Benzobisazoles. *In Progress.*
- **Collins, C. R.** ; Yaron, D. J. Developing Coarse-Grained Site Models for Excited Electronic States of Conjugated Polymers. *SPIE 2015.*

- Qiu, Y.; Worch, J. C.; Chirdon, D. N.; Kaur, A.; Maurer, A. B.; Amsterdam, S.; **Collins, C. R.** ; Pintauer, T.; Yaron, D.; Bernhard, S.; Noonan, K. J. T. Tuning Thiophene with Phosphorus: Synthesis and Electronic Properties of Benzobisthiaphospholes. *Chemistry A European Journal* **2014**.
- Tlach, B. C.; Tomlinson, A. L.; Morgan, K. D.; **Collins, C. R.** ; Jeffries-EL, M. Evaluation of the Impact of Extended Conjugation on the Optoelectronic Properties Benzo[1,2-d:4,5-d']bisoxazole Polymers. *Aust. J. Chem.* **2013**.

PRESENTATIONS AND POSTERS

- | | |
|--|-----------------------------------|
| • Predicting Chemical Properties Using Machine Learning Methods
<i>Department Retreat Poster Session</i> | Farmington, PA
2015 |
| • The Influence of Chemical Representations on the Efficiency of Molecular Screening
<i>Graduate Seminar</i> | Pittsburgh, PA

2015 |
| • A Mission to Multivariate Adaptive Regression Splines (MARS) and Its Other-worldly Neighbors
<i>10-705 Statistical Machine Learning Term Project</i> | Pittsburgh, PA

2015 |
| • Predicting Chemical Properties Using Machine Learning Methods
<i>10-701 Intro to Machine Learning Term Project</i> | Pittsburgh, PA
2014 |
| • Application of Machine Learning to Predict the Optoelectronic Properties of Benzobisazoles
<i>Senior Seminar</i> | Dahlonega, GA

2014 |
| • Machine Learning in Chemical Compound Space
<i>Junior Seminar</i> | Dahlonega, GA
2013 |
| • Donor-Acceptor Behavior of Benzobisazole Cruciform Polymers
<i>American Chemical Society Southeastern Regional Meeting</i> | Atlanta, GA
2013 |
| • The Impact of Conjugation Length on Benzobisazole Cruciforms
<i>American Chemical Society Southeastern Regional Meeting</i> | Raleigh, NC
2012 |
| – Runner-up Undergraduate Poster in Organic Chemistry | |

TECHNICAL SKILLS

Languages: Python, Javascript, C, C++ , MATLAB, Octave, Bash, Maple, CUDA C, Mathematica, TI-BASIC, SQL, Go, HTML, CSS, L^AT_EX, Java, Haskell

Software: *nix (Linux, Centos, FreeBSD, AIX, Solaris), Gaussian/AMPAC, LabVIEW, Git, SVN, SSH, rsync, coreutils, Make, Logisim, ab, Vagrant, PostgreSQL, Varnish, nginx, Apache, Torque

Libraries: Django, NumPy, SciPy, Matplotlib, Scikit-Learn, PyBrain, Selenium, Paramiko, PyTest, Twitter Bootstrap, jQuery, jQueryUI, SDL, Caffe, OpenMP, MPI