# Dakota Folmsbee

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## Education

**University of Pittsburgh** 

Ph.D. STUDENT IN PHYISCAL CHEMISTRY

Aug. 2016 - PRESENT

Pittsburgh, PA

**Clarkson University** 

Aug. 2012 - May 2016

B.S. IN CHEMISTRY Potsdam NY

### Skills

Chemistry NMR, HPLC, Mass Spectrometry, FTIR, UV/Vis

Programming Python, Scikit-Learn, Tensorflow, Keras, PyTorch, Bash, ŁT-X, C++, Julia

General GNU/Linux, VIM, Microsoft Office Suite

## **Professional Experience**

#### **Computational/Physical Chemistry Graduate Student**

Jan. 2017 - PRESENT

University of Pittsburgh

- · Instituted fast property prediction models to aid a genetic algorithm in rapid material screening
- Developed a molecular machine learning representation for chemical property predictions

#### **General Chemistry Teaching Assistant/Fellow**

Aug. 2016 - Dec. 2017

University of Pittsburgh

- · Recitation & Lab Honors General Chemistry
- Recitation & Lab General Chemistry
- · Lab General Chemistry for Engineers

#### **Undergraduate Researcher**

Aug. 2013 - May 2016

CLARKSON UNIVERSITY

- Synthesized carriers for cancer detecting molecules and chemotherapy drugs
- Analyzed compounds using techniques such as NMR, TOF-MS, and HPLC
- Researched procedures and applications for Gold nanoparticles and nanorods
- Synthesized Gold nanorods and analyzed with thermogravimetric analysis

### **General Chemistry Teaching Assistant/Mentor**

Aug. 2013 - May. 2016

**CLARKSON UNIVERSITY** 

- Recitation & Lab General Chemistry for Engineers
- Recitation & Lab General Chemistry for Chemistry and Chemical Engineering

### **Publications**

D. Folmsbee, S. Upadhyay, A. Dumi, D. Hiener, & D. Mulvey. (2019, July 12). chemreps/chemreps: Molecular Machine Learning Representations (Version 0.1.1). Zenodo. http://doi.org/10.5281/zenodo.3333856

### Presentation

#### Frederick Kaufman Memorial Lecture Series, University of Pittsburgh

Pittsburgh, PA

Oct. 2019

Oct 2019

POSTER PRESENTATION

Assessing Conformer Energies: Machine Learning vs Conventional Quantum Chemistry

### Science 2019, University of Pittsburgh

Pittsburgh, PA

POSTER PRESENTATION

Assessing Conformer Energies: Machine Learning vs Conventional Quantum Chemistry

DAKOTA FOLMSBEE · CURRICULUM VITAE

Advancing Research through Computing 2019, University of Pittsburgh	Pittsburgh, PA
POSTER PRESENTATION	Mar. 2019
Rapid Predictive Methods to Aid in Screening of Organic Dielectric Materials	
Science 2018, University of Pittsburgh	Pittsburgh, PA
POSTER PRESENTATION  Rapid Predictive Methods to Aid in Screening of Organic Dielectric Materials	Oct. 2018
Frederick Kaufman Memorial Lecture Series, University of Pittsburgh	Pittsburgh, PA
Poster Presentation	Oct. 2018
Rapid Predictive Methods to Aid in Screening of Organic Dielectric Materials	
Covestro Lecture Series, University of Pittsburgh	Pittsburgh, PA
POSTER PRESENTATION	Sept. 2018
Rapid Predictive Methods to Aid in Screening of Organic Dielectric Materials	
Simulators Meeting 2018, Carnegie-Mellon University	Pittsburgh, PA
ORAL PRESENTATION  Machine Learning to Aid in Screening for Organic Dielectric Materials	May 2018
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Covestro Lecture Series, University of Pittsburgh  Poster Presentation	Pittsburgh, PA Oct. 2017
Genetic Algorithms & Machine Learning for Rapid Materials Screening	Oct. 2011
Frederick Kaufman Memorial Lecture Series, University of Pittsburgh	Pittsburgh, PA
POSTER PRESENTATION	Oct. 2017
Genetic Algorithms & Machine Learning for Rapid Materials Screening	
Programming Projects	
chemreps	Aug. 2018 - PRESENT
Developer	
<ul><li>https://github.com/chemreps/chemreps</li><li>Developed a molecular representation library for machine learning in chemistry.</li></ul>	
QM/MM Study Group	July 2018 - Dec. 2018
Instructor & Organizer	
<ul> <li>https://github.com/shivupa/QMMM_study_group</li> <li>Organized and taught various lessons surrounding computational chemistry.</li> </ul>	
Honors & Awards	

**Safford Teaching Award**, University of Pittsburgh

**First Year Graduate Teaching Assistant Mentor**, University of Pittsburgh

**Walsh Fellow,** Clarkson University