

# Chase Clark

Postdoctoral Research Associate

## Curriculum Vitae

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📍 VCRGE/BIOTECH/GENOMICS, School of Pharmacy, University of Wisconsin-Madison  
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## Skills

- Bioinformatics
- Metabolomics
- Genomics/Metagenomics
- Natural product isolation and drug discovery

## Current Bioinformatic Tools

- Advanced: R, Rmarkdown, Shiny, Neo4j, Nextflow
- Intermediate: Python, Flask, HTML
- Familiar: JavaScript

## Education

2020	<b>PhD Pharmacognosy</b> Chicago, IL	University of Illinois at Chicago
2012	<b>BS Biochemistry</b> Mount Berry, GA	Berry College
2010	<b>Principia Consortium Study Abroad</b> Glasgow, Glasgow	University of Glasgow

## Experience

2020-Current	<b>Postdoctoral Research Associate</b> School of Pharmacy, Pharmaceutical Sciences Division	University of Wisconsin-Madison
2021-Current	<b>Computation and Informatics in Biology and Medicine (CIBM) Trainee</b>	National Library of Medicine
2019-2020	<b>F31 Fellow</b> National Center for Complementary and Integrative Health	U.S. National Institutes of Health (NIH)
2015-2019	<b>Ph.D. Student</b> Medicinal Chemistry and Pharmacognosy	University of Illinois at Chicago
2013-2015	<b>Research and Development Technician</b> R&D, Method Development	Deerland Enzymes
2008-2012	<b>Bonner Scholar</b>	Bonner Foundation

## Recent Solo Projects

2021	<b>socialgene</b> Repository-scale genome search platform	Python, Flask app, Nextflow, High-throughput computing
2019	<b>electricShine</b> Shiny packaging with Electron	R, Javascript, Node/Electron
2019	<b>mzEasy</b> Shiny app for converting and visualizing mass spectrometry data	R, Shiny
2019	<b>mzPlotter</b> Automated summary of LC-MS/MS data with interactive Rmarkdown reports	R, Rmarkdown
2019	<b>mzFromImage</b> Predict a mass spectrum's values from a static image file	R, Shiny
2018	<b>IDBac</b> MALDI protein and small molecule bioinformatics platform	R, Shiny, Electron

## Recent Team Projects

2021	<b>Autometa</b> Automated binning pipeline for single metagenomes	Python, Nextflow
2021	<b>metaBenchmarks</b> Benchmark metagenomic profiling/binning software	Nextflow

## Awards

2017	<b>American Society of Pharmacognosy Student Travel Award</b> \$600
2017	<b>University of Illinois at Chicago Provost Deiss Award</b> \$2,500
2019-2020	<b>NIH Predoctoral Individual National Research Service Award (F31)</b> \$90,000.00
2021-2023	<b>Computation and Informatics in Biology and Medicine (CIBM) training</b> \$100,000.00

## Talks

2020-06-01	<b>IDBac: Bioinformatics Software for Microbial Drug Discovery Prioritization and Culturomics Characterization</b> St. Jude National Graduate Student Symposium (cancelled due to Covid19)
2020-05-08	<b>Protein MS isn't the only MS... Programming in R for Metabolomics Mass Spectrometry</b> May Institute 2020: Future developers meeting
2020-02-25	<b>IDBac: Bioinformatics Software for Microbial Drug Discovery Prioritization and Culturomics Characterization</b> Gordon Research Conference on Marine Natural Products
2020-02-23	<b>IDBac: Bioinformatics Software for Microbial Drug Discovery Prioritization and Culturomics Characterization</b> Marine Natural Products Gordon Research Seminar (GRS)
2019-08-23	<b>Your Missing Step in Reproducible R Programming: Continuous Deployment</b> R/Pharma, Harvard University
2019-04-27	<b>Bioinformatics in R</b> satRdays Chicago
2019-01-23	<b>Single Function Lightning Talks: "lengths()", not length()"</b> Chicago R User Group
2018-08-15	<b>IDBac: A New Paradigm in Developing Microbial Libraries for Drug Discovery</b> R/Pharma, Harvard University
2018-01-27	<b>IDBac: A Shiny App to Analyze Bacterial Fingerprints and Aid in the Discovery of Potential New Antibiotics</b> Chicago R User Group
2017-03-09	<b>IDBac: A proteomic &amp; chemometric pipeline for rapid bacterial characterization</b> Center for Biomolecular Sciences, UIC
2016-11-03	<b>Rational design of bacterial strain libraries for drug discovery</b> UIC Specialized Metabolite Community
2016-07-22	<b>Use of MALDI-MS to Create 'Smart' Libraries for Drug-Lead Discovery</b> Chicago Mass Spec Day
2015-04-15	<b>HPLC-DAD Method for Trace Detection of Benzoic and Sorbic Acids in High Protein Matrices</b> SSAOAC Annual Meeting

## Poster Presentations

11/11/2016	<b>Innovating Microbial Libraries for Drug Discovery Using MALDI-TOF-MS and the Cultivable Freshwater Sponge Microbiome.</b> UIC College of Pharmacy Research Day	Chicago, IL
2/10/2017	<b>MALDI-TOF MS to Innovate Microbial Library Generation in Drug Discovery</b> UIC College of Pharmacy Research Day	Chicago, IL
1/1/2018	<b>IDBac as a Tool to Evaluate Environmental Bacterial Collections in a Single Week</b> American Society of Pharmacognosy	Lexington, KY

7/30/2017	<b>Use of MALDI-TOF MS to Generate Low Redundancy Taxonomic and Specialized Metabolite Libraries from Freshwater Sponge Microbiomes</b> American Society of Pharmacognosy	Portland, OR
7/14/2017	<b>MALDI-TOF MS to Innovate Microbial Library Generation in Drug Discovery</b> Chicago Mass Spec Day	Chicago, IL
5/24/2017	<b>MALDI-TOF MS to Innovate Microbial Library Generation in Drug Discovery</b> MBRB Research Day, UIC	Chicago, IL
07/30/2019	<b>Using IDBac to Investigate the Microbial and Natural Product Potential of Freshwater Sponges</b> WI American Society of Pharmacognosy	Madison,

## Publications

1. Leao, T. F., Clark, C. M., Bauermeister, A., Elijah, E. O., Gentry, E. C., Husband, M., Oliveira, M. F., Bandeira, N., Wang, M., & Dorrestein, P. C. (2021). Quick-start infrastructure for untargeted metabolomics analysis in GNPS. *Citations: 0*. 10.1038/s42255-021-00429-0
2. Clark, C. M., Murphy, B. T., & Sanchez, L. M. (2020). A call to action: The need for standardization in developing open-source mass spectrometry-based methods for microbial subspecies discrimination. *Citations: 2*. 10.1128/msystems.00813-19
3. Clark, C. M., Costa, M. S., Conley, E., Li, E., Sanchez, L. M., & Murphy, B. T. (2019). Using the open-source MALDI TOF-MS IDBac pipeline for analysis of microbial protein and specialized metabolite data. *Citations: 7*. 10.3791/59219
4. Braesel, J., Clark, C. M., Kunstman, K. J., Green, S. J., Maienschein-Cline, M., Murphy, B. T., & Eustáquio, A. S. (2019). Genome sequence of marine-derived streptomyces sp. Strain F001, a producer of akashin a and diazaquinomycins. *Citations: 0*. 10.1128/mra.00165-19
5. Costa, M. S., Clark, C. M., Ómarsdóttir, S., Sanchez, L. M., & Murphy, B. T. (2019). Minimizing taxonomic and natural product redundancy in microbial libraries using MALDI-TOF MS and the bioinformatics pipeline IDBac. *Citations: 7*. 10.1021/acs.jnatprod.9b00168
6. Clark, C. M., Costa, M. S., Sanchez, L. M., & Murphy, B. T. (2018). Coupling MALDI-TOF mass spectrometry protein and specialized metabolite analyses to rapidly discriminate bacterial function. *Citations: 34*. 10.1073/pnas.1801247115