Chase Clark

♀ University of Wisconsin-Madison, School of Pharmacy

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Education

2020 PhD Pharmacognosy University of Illinois Chicago

Chicago, IL

2012 BS Biochemistry Berry College

Mount Berry, GA

2010 **Principia Consortium Study Abroad** University of Glasgow

Glasgow, Glasgow

Employment

Aug 2015 - Aug 2020

Jun 2021 - present Computation and Informatics in Biology and Medicine (CIBM) Postdoctoral Fellow

Remote, Chicago, IL

Appointment to UW-Madison/National Library of Medicine T15 training grant

• Developing genomic and metagenomic bioinformatics software

• Currently developing a repository-scale analysis platform for comparative genomics

Sep 2020 - present
Postdoctoral Research Associate
University of Wisconsin-Madison

Remote, Chicago, IL

May 2019 - Aug 2020 NIH F31 Fellow University of Illinois at Chicago

Chicago, IL, US

Ph.D. Student

Chicago, IL, US

Natural product drug dicovery, specifically bacterial-derived antibiotics

 Conceived, learned to code, and built a bioinformatics program (R Shiny app) for rapid bacterial metabolomics analyses

 Isolated and cultured 1000's of bacterial isolates and performed natural product isolation and structure elucidation

• First or co-first author on four studies/manuscripts

• Mentored undergradaute and graduate students leading to publishable studies

Mar 2013 - Aug 2015 Research and Development Technician

Deerland Probiotics & Enzymes

University of Illinois at Chicago

Kennesaw, GA, US

· Assisted in the design and results analysis of three clinical studies

Presented monthly projects to CEO, VP Sales and VP Science & Technology

 Head of method development and identification of raw materials to meet CFR 21, FDA requirements.

 Developed identifications for over 140 materials, from enzymes and botanicals to minerals and additives.

 Overhauled the Method Development Department establishing inter-departmental responsibilities, improving and creating new SOPs, streamlining identification requests by establishing a formal request process and interactive database

 Strengthened and created new relationships with third-party vendors, labs, and universities.

 Controlled budget of method development for purchase of standards, consumables, equipment and third-party testing.

Mentored interns from Kennesaw Mountain High School magnet program

Aug 2008 - May 2012 Bonner Scholar

Rome, GA, US

r **Scholar** Bonner Foundation

Four-year community service scholarship requiring weekly and summer commitments.

Select Individual Software Projects

2022	socialgene Repository-scale genome pattern search platform	Python, Django, Nextflow, Rust, High-throughput computing
2019	electricShine Shiny packaging with Electron	R, Javascript, Node/Electron
2019	mzEasy Shiny app for converting and visualizing mass spectro	R, Shiny metry data
2019	mzPlotter Automated summary of of LC-MS/MS data with inter-	R, Rmarkdown active Rmarkdown reports
2019	mzFromImage Predict a mass spectrum's values from a static image	R, Shiny
2019	mgfparse Low-dependency R package for quickly and efficiently	R parsing mgf files into R
2018	IDBac MALDI protein and small molecule bioinformatics pla	R, Shiny, Electron

Select Team Software Projects

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

Talks

2022-06-23	National Library of Medicine (NLM) T15 Training Conference
	 Large Scale Analysis of Protein Homology for Microbial Drug Discovery
2022-03-09	 Gordon Research Conference on Marine Natural Products Towards near-instant, repository-scale searching for homologous BGCs with Social-gene (Selected alternate speaker, didn't present)
2021-11-02	 Computation & Informatics in Biology & Medicine Seminars BMI 915 New Computational Tools for Top-Down and Bottom-Up Natural Product Drug Discovery
2020-06-01	 St. Jude National Graduate Student Symposium (cancelled due to Covid19) IDBac: Bioinformatics Software for Microbial Drug Discovery Prioritization and Culturomics Characterization
2020-05-08	 May Institute 2020: Future developers meeting Protein MS isn't the only MS Programming in R for Metabolomics Mass Spectrometry
2020-02-25	 Gordon Research Conference on Marine Natural Products IDBac: Bioinformatics Software for Microbial Drug Discovery Prioritization and Culturomics Characterization
2020-02-23	 Gordon Research Seminar (GRS) on Marine Natural Products IDBac: Bioinformatics Software for Microbial Drug Discovery Prioritization and Culturomics Characterization
2019-08-23	R/Pharma, Harvard University
	 Your Missing Step in Reproducible R Programming: Continuous Deployment
2019-04-27	satRdays Chicago
	Bioinformatics in R
2019-01-23	Chicago R User Group
	 Single Function Lightning Talks: "lengths(), not length()"
2018-08-15	R/Pharma, Harvard University
	 IDBac: A New Paradigm in Developing Microbial Libraries for Drug Discovery

2018-01-27	 Chicago R User Group IDBac: A Shiny App to Analyze Bacterial Fingerprints and Aid in the Discovery of Potential New Antibiotics
2017-03-09	Center for Biomolecular Sciences, UIC
	• IDBac: A proteomic & chemometric pipeline for rapid bacterial characterization
2016-11-03	UIC Specialized Metabolite Community
	Rational design of bacterial strain libraries for drug discovery
2016-07-22	Chicago Mass Spec Day
	• Use of MALDI-MS to Create 'Smart' Libraries for Drug-Lead Discovery
2015-04-15	 SSAOAC Annual Meeting HPLC-DAD Method for Trace Detection of Benzoic and Sorbic Acids in High Protein Matrices

In-Progress

IDBacDB Manuscript	Web resource for sharing and searching FAIR MALDI-MS microbial spectra (a resource reque within many papers for many years now)
SocialGene Manuscript	Repository-scale analyses of protein similarity
clusterrule Manuscript	Initiated a multi-national collaboration to create a standardized collection of HMM models
Worshop Manuscript	Artificial Intelligence Approaches to Natural Product Drug Discovery
Autometa 2.0 Manuscript	Code standardization and enhancements including addition of Nextflow pipleine
Metagenomics Review Manuscript	Co-writing a review on genome resolved metagnomics and natural product discovery
Transcriptomics	Mapping transcriptional regulation in Micromonospora
Metagenomics	Independent metagnomics software development project

Funding/Awards

2021-present	Computation and Informatics in Biology and Medicine (CIBM) Training Program (T32) \$100,000.00
2020	St. Jude National Graduate Student Symposium (cancelled due to Covid19) One of 40 participants selected from 1,000 invite-only applications
2019-2020	NIH Predoctoral Individual National Research Service Award (F31) \$90,000.00
2018	R/Pharma, Travel Award \$500
2017	W.E. van Doren Scholar \$1,000
2017	American Society of Pharmacognosy Student Travel Award \$600
2017	University of Illinois at Chicago Provost/Deiss Award \$2,500
2008-2012	Bonner Scholar Four-year community service scholarship for students with need of financial assistance and a commitment to service
2008-2012	Academic Scholarship
2011	McCaleb, Hubert Scholarship
2011	Frank Plummer Scholarship
2010	Griggs Academic Scholarship
2009	John R. Bertrand Scholarship
2008	Sierra Club College Scholarship

2008 Cherokee Area Eagle Scout of the Year Scholarship

Workshops (host)

2022-07-25 ASP Annual Meeting (Co-host)

• Hands-On training of IDBac, an informatics tool for strain prioritization

2019-05-12 **Department of Pharmacognosy**

IDBac tutorial

Workshops (participant)

Sep 2021	Artificial Intelligence for Natural Product Drug Discovery (Invitation-Only; Lorentz Center)
Aug 2019	Integration of Metabolomics and Genomics (University of Chicago)
Dec 2016	PATRIC/RAST Workshop (Argonne National Laboratory)

Publications

- 1. Clark, C. M., Hernandez, A., Mullowney, M. W., Fitz-Henley, J., Li, E., Romanowski, S. B., Pronzato, R., Manconi, R., Sanchez, L. M., & Murphy, B. T. (2022). Relationship between bacterial phylotype and specialized metabolite production in the culturable microbiome of two freshwater sponges. *ISME Communications*. *Citations*: 0. 10.1038/s43705-022-00105-8
- 2. Clark, C. M., Nguyen, L., Pham, V. C., Sanchez, L. M., & Murphy, B. T. (2022). Automated microbial library generation using the bioinformatics platform IDBac. *Molecules. Citations*: 0. 10.3390/molecules27072038
- 3. 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. *Nature Metabolism. Citations*: 2. 10.1038/s42255-021-00429-0
- 4. 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. *mSystems*. *Citations*: 2. 10.1128/msystems.00813-19
- 5. 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. *Journal of Visualized Experiments. Citations: 9.* 10.3791/59219
- 6. 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. *Microbiology Resource Announcements*. *Citations*: 1. 10.1128/mra.00165-19
- 7. 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. *Journal of Natural Products. Citations*: 13. 10.1021/acs.jnatprod.9b00168
- 8. 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. *Proceedings of the National Academy of Sciences*. *Citations*: 52. 10.1073/pnas.1801247115

Submitted Publications

- "Artificial Intelligence Approaches to Natural Product Drug Discovery" (Submitted to Nature Reviews Drug Discovery)
 - Michael W. Mullowney, Katherine R. Duncan, Somayah S. Elsayed, Neha Garg, Justin J.J. van der Hooft, Nathaniel I. Martin, David Meijer, Barbara Terlouw, Friederike Biermann, Kai Blin, Janani Durairaj, Marina Gorostiola González, Eric J.N. Helfrich, Florian Huber, Stefan Leopold-Messer, Kohulan Rajan, Tristan de Rond, Jeffrey A. van Santen, Maria Sorokina, Marcy J. Balunas, Mehdi A. Beniddir, Doris van Bergeijk, Laura M. Carroll, Chase M. Clark, Chris Dejong, Chao Du, Scarlet Ferrinho, Francesca Grisoni, Albert Hofstetter, Willem Jespers, Olga V. Kalinina, Satria A. Kautsar, Tiago F. Leao, Joleen Masschelein, Evan R. Rees, Raphael Reher, Daniel Reker, Philippe Schwaller, Marwin Segler, Michael A. Skinnider, Allison S. Walker, Egon Willighagen, Barbara Zdrazil, Nadine Ziemert, Rebecca J.M. Goss, Pierre Guyomard, Andrea Volkamer, William H. Gerwick, Hyun Uk Kim, Rolf Müller, Gilles P. van Wezel, Gerard van Westen, Anna K.H. Hirsch, Roger Linington, Serina L. Robinson, Marnix H. Medema

Teaching

Invited Guest Lectures

PSCI 522 Advanced Pharmacognosy (two lectures) University of Illinois at Chicago Spring 2022