

Chase Clark

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Education

2015	PhD Pharmacognosy Chicago, IL	University of Illinois Chicago
2008	BS Biochemistry Mount Berry, GA	Berry College
2010	Principia Consortium Study Abroad Glasgow, Glasgow	University of Glasgow

Employment (Only relevant positions listed)

Jun 2021 - present	Computation and Informatics in Biology and Medicine (CIBM) Postdoctoral Fellow Remote, Chicago, IL	UW-Madison
Sep 2020 - present	Postdoctoral Research Associate Remote, Chicago, IL	University of Wisconsin-Madison
May 2019 - Aug 2020	NIH F31 Fellow Chicago, IL, US	University of Illinois at Chicago
Aug 2015 - Aug 2020	Ph.D. Student Chicago, IL, US <ul style="list-style-type: none">Natural product drug discovery, specifically bacterial-derived antibioticsConceived, learned to code, and built a bioinformatics program (R Shiny app) for rapid bacterial metabolomics analysesIsolated and cultured 1000's of bacterial isolates and performed natural product isolation and structure elucidation	University of Illinois at Chicago
Mar 2013 - Aug 2015	Research and Development Technician Kennesaw, GA, US <ul style="list-style-type: none">Head of method development and identification of raw materials to meet CFR 21, FDA requirements.Assisted in the design and results analysis of three clinical studiesPresented monthly projects to CEO, VP Sales and VP Science & TechnologyDeveloped identifications for over 140 materials, from enzymes and botanicals to minerals, additives, and probiotics.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 databaseStrengthened 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	Deerland Probiotics & Enzymes
Aug 2008 - May 2012	Bonner Scholar Rome, GA, US <ul style="list-style-type: none">Four-year community service scholarship requiring weekly and summer commitments.	Bonner Foundation

Data Science (select, not comprehensive)

Bioinformatics: metabolomics () • genomics () • metagenomics ()

Cloud Computing: remote server computing () • high-performance and high-throughput computing (Open science grid, HTCondor, AWS, etc)

Communication: presentations (public speaking, technical writing, peer-reviewed publications, SOPs) • writing (customer-facing documents, peer-reviewed publications, SOPs) • reports (Rmarkdown, Jupyter) • visualization (Base R, ggplot2, plotly, leaflet, cytoscape, networkx, gephi, etc)

Development: full-stack design (Flask, Django, Shiny, Dash)

Programming: general (R, Python) • database (Neo4j, SQL) • automation (Nextflow, Make) • containers (Docker) • containers (Singularity) • pipeline (Nextflow, Make)

Software Development: source control (Git) • CI/CD/DevOps (Github, Travis, Azure, etc.) • automated testing ()

Statistics: machine learning (scikit-learn, keras, tensorflow, Neo4j, etc) • experimental design (DOE, etc)

Lab Science (select, not comprehensive)

- microbial genomics and transcriptomics
- targeted and untargeted small molecule mass spectrometry and NMR
- biochemistry and biosynthetic gene cluster discovery and characterization
- method development

Select Individual Software Projects

2021-2023	SocialGene Repository-scale genomic graph database	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 spectrometry data	R, Shiny
2019	mzPlotter Automated summary of LC-MS/MS data with interactive Rmarkdown reports	R, Rmarkdown
2019	mzFromImage Predict a mass spectrum's values from a static image file	R, Shiny
2019	mgfparse Low-dependency R package for quickly and efficiently parsing mgf files into R	R
2016-2023	IDBac MALDI protein and small molecule bioinformatics platform	R, Shiny, Node/Electron

Select Team Software Projects

2021	metaBenchmarks Benchmark metagenomic profiling/binning software	Nextflow
2020-2023	Autometa Automated binning pipeline for single metagenomes	Python, Nextflow

Talks

2024-03-14	Gordon Research Conference on Marine Natural Products <ul style="list-style-type: none"> • A Multi-Repository Scale Genomic and Chemical Search Engine to Enable the Discovery, Production, and Function of Natural Products (selected GRS talk)
2024-03-10	Gordon Research Seminar (GRS) on Marine Natural Products <ul style="list-style-type: none"> • A Multi-Repository Scale Genomic and Chemical Search Engine to Enable the Discovery, Production, and Function of Natural Products
2023-09-21	Chicago R User Group <ul style="list-style-type: none"> • How to Write Reproducible Code for Data Science

- 2023-08-10 **Chicago Python User Group**
- Reproducibly Building and Analyzing Knowledge Graphs for Drug Discovery with Nextflow, Neo4j and Python
- 2023-03-07 **Computation and Informatics in Biology and Medicine Seminars**
- SocialGene: Large Scale Knowledge Graphs for Microbial Based Drug Discovery
- 2023-01-30 **Northeastern University Department of Chemistry and Chemical Biology**
- Information-Rich Platforms for Natural Product Antibiotic Drug Discovery and Microbial Characterization
- 2023-01-30 **2023 ASP Younger Members Symposium**
- SocialGene: A Large Scale Search Engine to Find Metagenomic BGCs in Free-Living Organisms
- 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 Socialgene (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 (selected GRS talk)
- 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

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**
Scholarship for students with financial need "in exchange for weekly commitment to intensive and meaningful service with a local community organization over the four years"
- 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**

Publications

1. Rees, E. R., Uppal, S., Clark, C. M., Lail, A. J., Waterworth, S. C., Roesemann, S. D., Wolf, K. A., & Kwan, J. C. (2023). Autometa 2: A versatile tool for recovering genomes from highly-complex metagenomic communities. *Citations: 0*. 10.1101/2023.09.01.555939
2. 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: 4*. 10.1038/s43705-022-00105-8
3. 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: 2*. 10.3390/molecules27072038
4. 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: 11*. 10.1038/s42255-021-00429-0
5. Elfeki, M., Mantri, S., Clark, C. M., Green, S. J., Ziemert, N., & Murphy, B. T. (2021). Evaluating the distribution of bacterial natural product biosynthetic genes across lake huron sediment. *ACS Chemical Biology. Citations: 4*. 10.1021/acscchembio.1c00653
6. 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

7. 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
8. 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: 2. 10.1128/mra.00165-19
9. 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: 16. 10.1021/acs.jnatprod.9b00168
10. 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: 66. 10.1073/pnas.1801247115