## Jeremiah Gaiser

2209 Gilbert Ave, Missoula, MT 59802

jeremiah.gaiser@umontana.edu (406) 407-5409

#### **PROFILE**

I am a doctoral student in the University of Montana's Biochemistry program. My research interests are in development of machine learning approaches to drug discovery and toxicology prediction. I graduated in May of 2020 from the University of Montana with a Bachelors of Science in Computational Biochemistry and minors in Math and Computer Science. As an undergraduate researcher, I developed a foundation in genetic sequencing and annotation methods in genomics research. Before my undergraduate education, I worked as a professional web developer.

#### **EDUCATION**

### THE UNIVERSITY OF MONTANA, Missoula, MT

May 2020

- Bachelors of Science in Computational Biochemistry, GPA 3.63. Minors in Computer Science and Mathematics
- Selected as Outstanding Graduating Computational Biochemistry Senior.

#### **WORK AND RESEARCH EXPERIENCE**

UNIVERSITY OF MONTANA, Missoula, MT
Undergraduate researcher (Wheeler Lab / Grimes Lab)

June 2017-present

- Deep learning models for drug discovery
  - Advisor: Travis Wheeler (University of Montana)
  - Context: Joint European Disruptive Initiative Covid-19 drug discovery contest,
     collaborative with Amit Roy (RML) and Vishwesh Venkatraman (Norwegian University of Science and Technology)
  - Focus: latent-space representation of ligands and proteins as foundation for binding affinity prediction
  - Technology: PyTorch, cluster compute.
- Developed model for annotating protein-coding DNA sequence
  - o Advisor: Travis Wheeler
  - probabilistic model for sequence alignment of amino acid to codon, based on codon neighborhood
  - o Implemented in C programming language and integrated into fork of HMMER codebase

- Investigation of annotation of conserved, non-coding elements across eukaryotic genomes
  - Advisors: Martin Frith (University of Tokyo), Travis Wheeler (University of Montana)
  - Used large volumes of genome-scale metadata (Uniprot, Dfam, fRNAdb) and sequence alignment tools (HMMER, LAST) to estimate conservation in non-coding sequences in eukaryotic genomes.
  - Project initiated during summer internship at the University of Tokyo
- Cytoscape Network Browser (http://cytoscape.cs.umt.edu/)
  - Advisors: Mark Grimes (University of Montana), Travis Wheeler (University of Montana)
  - o Developed graphical user interface for visually navigating protein interaction networks
  - Built and implemented application architecture using several different technologies, including cytoscape and its API, R, python, and front-end web technologies and frameworks such as HTML, CSS, javascript, cytoscape.js, jquery, and react.js.
  - Exercised communication and teamwork with co-developer
  - Manuscript: http://stke.sciencemag.org/content/11/531/eaaq1087.abstract

# PRIME INCORPORATED, Bozeman, MT

#### **Web Developer**

May 2015 - August 2016

- Development of websites and applications for small to medium sized businesses
  - Used frontend and backend technologies including HTML, javascript, css, jquery, PHP, and content-management systems including Expression Engine and CraftCMS.
  - Worked with team of developers, designers, and managers
  - Communicated with clients via email, phone, and in person to discuss projects

#### **RELEVANT COURSEWORK**

Genomics
Biochemistry (including lab)
Nucleic Acids

Genetics and Evolution
Machine Learning
Data Analysis

Organic Chemistry

Design/Analysis of Algorithms

Cellular and Molecular Biology