# Postdoctoral Researcher, Polytechnique

# Research Experience

2022 **Postdoctoral researcher**, LIX in Polytechnique / Sanofi, Paris

Supervised by Maks Ovsjanikov

Geometric learning methods for proteins: Compare protein structure representations and how they can be combined.

Finding antibodies in cryo-EM densities: Partnership with Sanofi to use object detection to automatically identify antibodies in cryo-EM maps.

2022 Visiting Student, La Sapienza Gladia Group, Rome

Supervised by Emanuele Rodolà

*Surface methods for proteins* : Adapt latest surface methods to learning on protein surfaces.

2019-2022 **PhD**, Structural Bioinformatics, Institut Pasteur / CBIO, Mines Paris, Paris Supervised by Michael Nilges and Jean-Philippe Vert.

Reinforcement learning and cryo-EM: Train an agent to trace protein backbones in cryo-EM densities.

Equivariant networks for genomic data: Create a theoretical framework that encompasses existing methods. Implement it and get enhanced results.

Assist Protein-Protein Interaction drug design: Develop the first binding pocket predictor for this data.

Learning on RNA structure as 2.5D graphs: Develop a library for this representation, relevant deep learning layers and motifs mining algorithms.

2018-2019 **Master Thesis**, *McGill Computational Biology Group*, Montréal Supervised by Jérome Waldispühl.

Active ligand generation project: Condition the generation of small molecules on the structure of a protein pocket.

RNA structural drug design: Represent binding pockets as 2.5D graphs and use machine learning to predict active compounds from it.

Spring 2018 **Research Internship**, *McGill Computational Biology Group*, Montréal Developed binding site comparison and embedding for proteins based on graphs vs spatial point sets.

#### Education

- 2018 2019 **M.Sc. Computer Science**, *McGill University*, Machine Learning, Deep Learning, Bioinformatics
- 2015 2018 **M.Eng. Bioinformatics**, *Ecole Polytechnique*, Parallel computing, Big Data, Machine Learning, Bioinformatics
- 2013 2015 Preparatory Program, Louis le Grand, PCSI/PC Track

### Teaching

- 2022-23 **Teacher**, *CRI* (*Centre de Recherche Interdisciplinaire*), 30 hours of classes and practical sessions on Foundations of Al. Joined the board of teachers.
  - 2022 Lecturer, McGill CS Master, Geometric learning for drug design
- 2020-21-22 **Lecturer**, *Université de Paris, Master ISDD*, Generative models for drug design

### **Publications**

#### **Journals**

- 2022 ▷ InDeep: 3D fully convolutional neural networks to assist in silico drug design on protein-protein interactions. Vincent Mallet, Luis Checa Ruano, Alexandra Moine Franel, Michael Nilges, Karen Druart, Guillaume Bouvier, and Olivier Sperandio. *Bioinformatics*, 2022
- 2022 > RNAglib: A python package for RNA 2.5D graphs. Vincent Mallet, Carlos Oliver, Jonathan Broadbent, William L Hamilton, and Jérôme Waldispühl. *Bioinformatics*, 2022
- 2022 **VeRNAI: A Tool for Mining Fuzzy Network Motifs in RNA**. Carlos Oliver\*, Vincent Mallet\*, Pericles Philippopoulos, William L Hamilton, and Jérôme Waldispühl. *Bioinformatics*, 2022
- 2021 puicksom: Self-Organizing Maps on GPUs for clustering of molecular dynamics trajectories. Vincent Mallet, Michael Nilges, and Guillaume Bouvier. *Bioinformatics*, 2021
- 2020 ▷ OptiMol: Optimization of binding affinities in chemical space for drug discovery. Jacques Boitreaud\*, Vincent Mallet\*, Carlos Oliver, and Jérome Waldispuhl. *Journal of Chemical Information and Modeling*, 2020
- 2020 ▷ Augmented base pairing networks encode RNA-small molecule binding preferences. Carlos Oliver, Vincent Mallet, Roman Sarrazin Gendron, Vladimir Reinharz, William L Hamilton, Nicolas Moitessier, and Jérome Waldispuhl. *Nucleic Acids Research*, 2020

#### Conferences

2021 ▷ Reverse-complement equivariant networks for dna sequences. Vincent Mallet and Jean-Philippe Vert. Advances in Neural Information Processing Systems, 2021

### Preprints

- 2023 **Finding antibodies in cryo-em densities with crai**. Vincent Mallet, Chiara Rapisarda, Hervé Minoux, and Maks Ovsjanikov. *bioRxiv*, 2023. 10.1101/2023.09.27.559736
- 2023 ⊳ Somseq, displaying sequence similarity using self-organizing maps. Isaac Filella-Merce, Vincent Mallet, Eric Durand, Michael Nilges, Guillaume Bouvier, and Riccardo Pellarin. (In review, Bioinformatics), 2023

- 2023 ▷ Structure-based rna drug discovery with deep-learning assisted virtual screening. Juan Guillermo Carvajal Patiño, David Becerra, Vincent Mallet, David Hiraki, Nicolas Moitessier, Luis Fernando Niño, Carlos Oliver, and Jérôme Waldispühl. (In preparation), 2023
- 2023 Atomsurf: Surface representation for learning on protein structures. Vincent Mallet, Souhaib Attaiki, and Maks Ovsjanikov. *arXiv preprint arXiv:2309.16519*, 2023
- 2023 ▶ Approximate network motif mining via graph learning. Carlos Oliver, Dexiong Chen, Vincent Mallet, Pericles Philippopoulos, and Karsten Borgwardt. arXiv preprint arXiv:2206.01008, 2023
- 2022 **Edge-similarity-aware graph neural networks**. Vincent Mallet, Carlos G Oliver, and William L Hamilton. *arXiv preprint arXiv:2109.09432*, 2021
- 2019 ▷ Leveraging binding-site structure for drug discovery with pointcloud methods. Vincent Mallet, Carlos G Oliver, Nicolas Moitessier, and Jerome Waldispuhl. arXiv preprint arXiv:1905.12033, 2019

### **Talks**

- 2023 Modeling structural biology with geometric deep learning , Data Shapes Seminar, INRIA Paris
- 2023 **Geometric learning for drug design**, AstraZeneca drug discovery Seminar (online)
- 2023 RNAGlib: a platform for learning on RNA structure, MILA drug discovery Seminar
- 2023 Reverse-Complement Equivariant Networks for DNA Sequences , 20 min talk, Al in Structural Biology Workshop
- 2022 InDeep: learning for drug design on protein protein interactions, Workshop on Al in magnetic resonance
- 2022 Artificial Intelligence in prediction, determination and analysis of structures, EMBO workshop: Machine learning for structural biology
- 2022 Reverse-Complement Equivariant Networks for DNA Sequences, NeurlPS

## **Funding**

- 2022 ELISE Mobility Award
- 2019 PhD funding from the INCEPTION Program
- 2019 Granted AMX PhD funding
- 2018 Research Grant from the Computational Biology Group

### Publishing

Reviewer: Bioinformatics, NAR, NeurIPS, ICML Associate Editor: Frontiers in Bioinformatics

## Technical Skills

Libraries : Deep learning (PyTorch, DGL, Keras, TensorFlow, Sonnet, JAX, Haiku),

Bioinformatics (Chimerax, Biopython, RDKit, Pymol)

Developed packages: rnaglib, quicksom

## Languages

French: Native, English: Fluent (TOEFL 117), German, Italian:(Intermediate)

### Additional Interests

Scenography In charge of scenography for Phasm, an association organizing festivals and

concerts.

Sports Climbing, Running, Fencing, Skiing