

# Hugo Talibart

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## Profile

Computer scientist with 8+ years of expertise in algorithm design and implementation, with experience building and shipping software. Looking for a software engineering role where rigorous algorithmic thinking meets concrete development.

## Experience

**Postdoctoral Researcher** *Université Libre de Bruxelles, Brussels (Belgium)* 2023 – Present

- Designed ILP formulations for bias-controlled machine learning benchmark construction – applied to drug-target interaction prediction.
- Designed and trained a large-scale Transformer VAE for molecular generation – architecture design, training at scale. Released as a pip package.

🔗 [github.com/3BioCompBio/chembed](https://github.com/3BioCompBio/chembed)

- Collaborated on a sparse structure-informed probabilistic graphical model with C++ solver and Python interface, achieving orders of magnitude speedup over fully-connected models – applied to mutation effect prediction, state-of-the-art on standard benchmarks. Released as a pip package.

🔗 [github.com/3BioCompBio/StructureDCA](https://github.com/3BioCompBio/StructureDCA)

- Supervised two Master's students and an intern on benchmark design and deep learning topics.

**Postdoctoral Researcher** *Muséum National d'Histoire Naturelle, Sorbonne Université, Paris (France)* 2021 – 2023

- Extended my PhD's probabilistic graphical model alignment framework with a redesigned statistical inference pipeline – faster inference, improved robustness on shallow data.

**PhD Researcher** *INRIA, Université de Rennes, Rennes (France)* 2017 – 2021

- Introduced PPalin: first exact ILP-based alignment of probabilistic graphical models, solved in tractable time – applied to protein remote homology detection, outperforming HMM baselines.

## Education

**PhD in Computer Science** *INRIA, Université de Rennes 1, Rennes (France)* 2017 – 2021

**Master in Theoretical Computer Science** *Université de Rouen, Rouen (France)* 2016 – 2017

**Engineering Degree in Mathematics** *Institut National des Sciences Appliquées (INSA), Rouen (France)* 2012 – 2017

## Skills

**Languages** Python, C/C++, JavaScript, Bash

**ML & Data** PyTorch, NumPy, SciPy, Pandas, scikit-learn

**Tools** git, conda, docker/Singularity

**Environment** Linux, cluster computing (Slurm)

## Personal projects

**wikiddle.com** *Daily Wikipedia puzzle game*

- JavaScript frontend
- Python backend with FastAPI and SQL
- Self-hosted on a self-administered server
- 🔗 [github.com/htalibart/wikiddle](https://github.com/htalibart/wikiddle)

## Selected publications

- **Talibart H**, Gilis D. *Learning a chemistry-aware latent space for molecular encoding and generation with a large-scale Transformer Variational Autoencoder*. bioRxiv, 2025.
- Tshishyn M, **Talibart H**, Rooman M, Pucci F. *Structure-informed direct coupling analysis improves protein mutational landscape predictions*. bioRxiv, 2026.
- **Talibart H**, Coste F. *PPalign: optimal alignment of Potts models representing proteins with direct coupling information*. BMC Bioinformatics, 2021.

## Languages

**French:** Native

**English:** Fluent