## NICCOLO' DIPACE

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#### **EDUCATION**

# ÉCOLE NATIONALE SUPERIEURE DE CHIMIE DE PARIS-PSL UNIVERSITY Chemistry

Paris, France

Sept 2024 - Sept 2025

- · Double joint degree program, 2nd year, ICI (Integrative Chemistry & Innovation) Track
- · Relevant courses: Polymers, MOFs, MD simulations, DFT, Analitcal chemistry.
- · Thesis: DFT and TD-DFT on biocompatible, tunable, and reversible redox-responsive fluorescent metallocene-based probes

# UNIVERSITY OF FLORENCE

Florence, Italy

Chemistry Sciences
Sept 2023 – June 2024
Double joint degree program, 1st year, curriculum in Supramolecular, materials and nanosystems

- · Relevant courses: Nanoparticles, Computational chemistry, inorganic chemistry, nanosystems, supramolecular chemistry, materials.
- · Selected as 1st candidate for the Double Degree program

# UNIVERSITY OF FLORENCE Chemistry

Florence, Italy

Sept 2019 – Jun 2023

· Thesis: Molecular dynamics simulations of N-glycosylated proteins

#### **WORK EXPERIENCE**

# ÉCOLE NATIONALE SUPERIEURE DE CHIMIE DE PARIS-PSL UNIVERSITY Internship - Theoretical Investigation of Computational Protocols

Paris, France

Feb 2025 - Sept 2025

- · Focused on optimizing computational methods to accurately model the electronic properties
- · Performed simulations and analyzed the effects of redox reactions on the fluorescence properties of the probes.
- · Collaborated with Ilaria Ciofini to refine the protocol and improve computational accuracy.

### **UNIVERSITY OF FLORENCE**

Florence, Italy

### Internal Training – MD Studies of Protein Behavior with Different Water Models

March 2024 - May 2024

- $\cdot$  Replaced rigid water models with more flexible models to analyze the impact on protein structure and dynamics.
- · Evaluated protein conformational changes and stability by comparing results from simulations with different water models.

## UNIVERSITY OF FLORENCE

Florence, Italy

## Internship – Molecular Dynamics Simulations of N-glycosylated Proteins

Dic 2022 - May 2023

- $\cdot$  Conducted molecular dynamics simulations to study the behavior of N-glycosylated proteins.
- · Utilized GROMACS, VMD to model and simulate protein dynamics.
- Interpreted simulation results and contributed to the development of insights into protein function and folding mechanisms.

#### LEADERSHIP AND ACTIVITIES

## Waiter, Trattoria da Guido

Florence, Italy

Managed customer complaints in a professional manner, aiming for quick and efficient resolutions. Mai 2022 -- July 2024

Football coach: focus on team dynamics, leadership skills, and strategic game planning

Math and chemistry Tutor: Tutored undergraduates in core mathematical and chemistry concepts.

**Volunteering:** Developed communication and leadership skills by interacting with diverse groups and organizing workshops.

### SKILLS AND ADDITIONAL INFORMATION

**Languages:** English (Fully Proficient), Spanish (Fully Proficient), French (Intermediate – conversational and written), Italian (Native).

Programming and Software: Excel (Advanced), LaTeX (Proficient), Linux (Advanced), Igor (intermediate), ChemDraw

 $(Intermediate),\ Gaussian\ (Advanced),\ GROMACS\ (intermediate),\ Origin Lab\ (intermediate),\ VMD (Advanced).$ 

Interests and Sports: Football, Tennis, Running.