



Ramón L. PANADÉS-BARRUETA

Curriculum Vitae

Professional experience

- 2020–Present **Postdoctoral fellow**, [University of Twente](#), Computational Chemical Physics Group (CCP), Netherlands.
Project: Targeting Real chemical accuracy at the EXascale (TREX). European HPC Centre of Excellence (CoE)

Education

- 2017–2020 **PhD in Physics**, [University of Lille](#), Laboratoire de Physique des Lasers, Atomes et Molécules (PhLAM), France.
Thesis title: Full quantum simulations of the interaction between atmospheric molecules and model soot particles
Supervisor: Prof. Dr. Daniel PELÁEZ-RUIZ ([ISMO](#), Université Paris-Saclay)
- 2016–2017 **MSc in Physics (International Master 2 Atmospheric Environments)**, [University of Lille](#), Laboratoire de Physique des Lasers, Atomes et Molécules (PhLAM), France.
Thesis title: Towards a quantum dynamical description of the photodissociation of Cl_2 molecule adsorbed on ice
Supervisors: Prof. Dr. Daniel PELAEZ-RUIZ and Prof. Dr. Maurice MONNERVILLE
- 2011–2016 **BSc in Radiochemistry**, [University of Havana](#), Higher Institute of Technologies and Applied Sciences ([InSTEC](#)), Havana, Cuba.
Thesis title: Mean Potential Phase Space Theory study of the $Si(^3P) + OH(X^2\Pi) \rightarrow SiO(X^1\Sigma^+) + H(^2S)$ reaction
Supervisor: Dr. Alejandro RIVERO-SANTAMARÍA ([CMF](#), Universidad del País Vasco)

Publications

- 2020 Panadés-Barrueta, R. L. and Peláez, D. **Low-Rank Sum-of-Products Finite-Basis-Representation (SOP-FBR) of Potential Energy Surfaces**, [The Journal of Chemical Physics](#), *in review*
- 2019 Panadés-Barrueta, R. L., Martínez-Núñez, E., & Peláez, D. (2019). **Specific Reaction Parameter Multigrid POTFIT (SRP-MGPF): Automatic Generation of Sum-of-Products Form Potential Energy Surfaces for Quantum Dynamical Calculations**, [Frontiers in Chemistry](#), **7**, 576. Included in the book [Application of Optimization Algorithms in Chemistry](#)
- 2016 Panadés-Barrueta, R. L., Rubayo-Soneira, J., Monnerville, M., Larregaray, P., Dayou, F., & Rivero-Santamaría, A. (2016). **Mean Potential Phase Space Theory study of the $Si(^3P) + OH(X^2\Pi) \rightarrow SiO(X^1\Sigma^+) + H(^2S)$ reaction**, [Revista Cubana de Física](#), **33(2)**, 102–117.

Seminars and conferences

- August 2020 **On the automatic computation of global (intermolecular) potential energy surfaces for quantum dynamical simulations**, *Invited Speaker*.
[Symposium and Summer School on Physics of Ionized Gases](#)
Šabac, Serbia

February 2020 **Automatic computation of global (intermolecular) potential energy surfaces for (non) covalently bound systems**, *Contributed Talk*.

[Journée Chimie Théorique et Simulation Moléculaire IdF/Nord](#)
Chimie ParisTech. Paris, France

Teaching experience

2018–2019 **Laboratory of Thermodynamics**.

Place: IUT A de Lille (University of Lille), France No. hours: 64 Language: French

2013–2016 **Math Analysis and Linear Algebra**.

Place: InSTEC, Cuba No. hours: 72 Language: Spanish

Honors and Awards

2019 **Best Poster Prize** 10th International Meeting on Atomic and Molecular Physics and Chemistry (IMAMPC). Madrid, Spain.

2018 **Best Poster Prize** 6th High Dimensional Quantum Dynamics Workshop (HDQD). Lille, France.

2018 **PCCP Best Poster Prize** 9th International Meeting on Atomic and Molecular Physics and Chemistry (IMAMPC). Berlin, Germany.

2015 **ICPC Contestant** Caribbean Finals of the International Collegiate Programming Contest (ACM-ICPC). Havana, Cuba

2010 **IChO Contestant** Captain of the Cuba Team in the 42nd International Chemistry Olympiad ([IChO 2010](#)). Tokyo, Japan

Competitive research grants and fellowships

Oct.-Dec. **Research Grant German Academic Exchange Service (DAAD)**.

2019 Awarded by: Deutscher Akademischer Austausch Dienst
Place: Universität Heidelberg, Theoretical Chemistry Group
Supervisor: Prof. Dr. Oriol VENDRELL

2016–2017 **Labex CaPPA Fellowship**.

Awarded by: Laboratoire d'excellence CaPPA
Place: University of Lille

Research training

August 2019 **School EMIE-UP. Multiscale Dynamics in Molecular Systems**, *École de Physique des Houches. Haute-Savoie, France*.

June 2019 **3rd Mini-school on mathematics for theoretical chemistry and physics**, *Sorbonne Université, Pierre et Marie Curie. Paris, France*.

June 2018 **Bridging experiment and theory in precision spectroscopy (BETS) 4th MOLIM Training School**, *Nicolaus Copernicus University. Torun, Poland*.

January 2018 **Label of Theoretical Chemistry Île de France-Nord**, *Sorbonne Université, Pierre et Marie Curie. Paris, France*.

October 2017 **Quantum Dynamics with the Multi-Configuration Time-Dependent Hartree (MCTDH) method: future and perspectives**, *Université Paris-Saclay. Paris, France*.

Computer skills

OS UNIX, ([Arch](#)) Linux, Android

Languages Python (SciPy, NumPy, Matplotlib), C, FORTRAN, Bash, Lisp, Julia, \LaTeX

Software [MCTDH](#) (Developed the SOP-FBR and SRPTucker packages), MOPAC, MOLPRO, MOLCAS, Gaussian, [AutoMeKin](#), Inkscape, Git

Languages (CEFR)

Spanish Native speaker

English Proficient user (C1)

French Proficient user (C1)

German Basic user (A2)