PHD STUDENT



CONTACT



Address Rosario Argentina



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E-Mail jmlombardi@live.com

PERSONAL



Date of Birth 31/12/1993



Nationality Argentine

PLATFORMS



GitHub Link

LANGUAGES



Espanish Native



English Fluent



French Basic



German Basic (learning)

EDUCATION

02/2013 - 10/2017 Universidad Nacional de Rosario

Degree in Chemistry

Facultad De Ciencias Bioquímicas

Y Farmacéuticas

04/2018 - Ongoing expected completion date: 09/2023

Universidad Nacional de Rosario

Phd. in Physics

Facultad de Ciencias Exactas, ingeniería

y agrimensura

Topic: Theoretical study of electronic properties and reactivity of SAMs, MONCs and organic

molecules on metal surfaces.

Supervisor: Dr. Paula Natalia Abufager

Co-Supervisor: Dr. Heriberto Fabio Busnengo

PUBLICATIONS

2023 Structure and stability of metallo-porphyrin networks on Au(111)

J. M. Lombardi, D. Grumelli, R. Gutzler,

H. F. Busnengo y P. Abufager.

J. Phys. Chem. C 2023, 127, 13, 6569-6577

2020 Enhancing Hydrogen Evolution Activity of Au(111) in Alkaline Media Through Molecular Engineering of a 2D Polymer

P. Alexa, J. M. Lombardi, P. Abufager,

H. F. Busnengo, D. Grumelli, V. S. Vyas, F. Haase,

B Lotsch, R. Gutzer, K. Kern

Angewandte Chemie 59 8411-8415

2020 Three-way calibration using PARAFAC and MCR-ALS with previous synchronization of second-order chromatographic data through a new functional alignment of pure vectors for the quantification in the presence of retention time shifts in peak position and shape

S.Mazivila, J.M.Lombardi, R.Páscoa, S.Bortolato,

J.Leitão, J.Silva

Analytica Chimica Acta. 1146. 98-108

Functional data analysis, a new approach to aligning three-way liquid chromatographic with fluorescence detection data.

J.M.Lombardia, S.A.Bortolato,

Microchemical Journal 142, 219-228

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IN PROGRESS ARTICLES

In progress Catalysis of 2D metal-organic frameworks: porphyrins on Au(111)

J.M. Lombardi, D.Grumelli, R.Gutzler, H.F.Busnengo y P.Abufager

In progress Oxygen Reduction Reaction on gold-supported FePC and CoPC 2D networks:

effect of solvent co-deposited molecules

J. M. Lombardi, P. Abufager, H. F. Busnengo and D. Grumelli

In progress Functional data analysis, a comprehensive framework for processing non-quadrilinear

and low-selective data provided by four-way liquid chromatography analysis

M.Alcaraz, J.M. Lombardi, S.Bortolato

WORK EXPERICENCE

2022 - 2023 UNIVERSIDAD NACIONAL DE ROSARIO (UNR)

Teacher

Area: Physics

Department: Physical Chemistry

Facultad de Ciencias Bioquímicas y Farmacéuticas

2018 - 2022 UNIVERSIDAD NACIONAL DE ROSARIO (UNR)

Senior Teaching Assistant

Area: Physics

Department: Physical Chemistry

Facultad de Ciencias Bioquímicas y Farmacéuticas

2016 - 2018 UNIVERSIDAD NACIONAL DE ROSARIO (UNR)

Junior Teaching Assistant

Area: Physics

Department: Physical Chemistry

Facultad de Ciencias Bioquímicas y Farmacéuticas

2015 - 2017 UNIVERSIDAD NACIONAL DE ROSARIO (UNR)

Junior Teaching Assistant

Department: Organic Chemistry

Facultad de Ciencias Bioquímicas y Farmacéuticas

CONTRIBUTIONS PRESENTED AT INTERNATIONAL SCIENTIFIC AND TECHNOLOGICAL MEETINGS

2022 Reaction of 2D metal-organic frameworks: metal porphyrins on Au(111)

J.M. Lombardi, D. Grumelli, R. Gutzler, K. Kern, H.F. Busnengo, P. Abufager. Photo and ElectroCatalysis at the Atomic Scale (PECAS 2022), 20-23 Junio de 2022. San Sebastián, España. Comunicación oral

2020 Catalytic properties toward the Oxygen Reduction Reaction of 2D metal-organic frameworks: metalloporphyrins on Au(111)

J. M. Lombardi*, D. Grumelli, R.Gutzler, K. Kern, H. F. Busnengo and P. Abufager, CMD-2020-GEFES. Condensed Matter Divisions of the Spanish Royal Physics Society (RSEF-GEFES) and of the European Physical Society (EPS-CMD). 31st August-4th September.

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2020 Structure and electrocatalytic activity toward ORR of metallo-porphyrin networks

J. M. Lombardi, D. Grumelli, Gutzler, K. Kern, H.F. Busnengo, P.N. Abufager 2020 Express Conference on the Physics of materials and its application in Energy and environment 3th August - 6th August

2019 Structure and catalytic activity of metalloporphyrin networks

J. M. Lombardi, D. Grumelli, R.Gutzler, K. Kern, H. F. Busnengo and P. Abufager Gordon Research Conference on Dynamics at Surfaces (GRC) 28 de julio al 2 de Agosto, Newport, RI United States

2019 Structure and catalytic activity of metalloporphyrin networks

J. M. Lombardi, D. Grumelli, R.Gutzler, K. Kern, H. F. Busnengo and P. Abufager Gordon Research Seminar on Dynamics at Surfaces (GRS) 27-28 de Julio, Newport, RI United States

2018 On-surface transmetalation of metalloporphyrins self-assembled monolayers

J. M. Lombardi, D. Hötger, P. Abufager, C. Morchutt, P. Alexa, D. Grumelli, J.Dreiser, S. Stepanow, H. F. Busnengo, M. Etzkorn, R.Gutzler and K.Kern VI San Luis Conference on Surfaces, Interfaces and Catalysis, Santa Fe (Argentina), June, 2018

CONTRIBUTIONS PRESENTED AT ARGENTINIAN SCIENTIFIC AND TECHNOLOGICAL MEETINGS

Reacción de Reducción de Oxígeno en redes 2D de FePC/Au(111): efecto de moléculas de solvente co-depositadas.

J.M. Lombardi, H.F. Busnengo, P.Abufager, D.Grumelli NANO 2022, Río Cuarto, 9-11 Agosto 2022

Propiedades catalíticas frente a la Reacción de Reducción de Oxígeno de estructuras metal-orgánicas bidimensionales: porfirinas metálicas en Au(111)

J.M. Lombardi, H.F. Busnengo, D.Grumelli, P.Abufager NANO 2022, Río Cuarto, 9-11 August 2022

Award for Best e-Poster Presentation

Propiedades catalíticas frente a la Reacción de Reducción de Oxígeno de estructuras metal-orgánicas bidimensionales: porfirinas metálicas en Au(111).

J.M.Lombardi, H.F.Busnengo, D.Grumelli, P.Abufager.

IX Encuentro de Física y Química de Superficies y I Encuentro de Biología de Superficies, 26-28 octubre 2022, Santiago del Estero. Semi-plenary Conference

Estructura y actividad catalítica de redes de metaloporfirinas

J. M. Lombardi, D. Grumelli, Gutzler, K. Kern, H.F. Busnengo, P.N. Abufager. NANO 2019: XIX Encuentro de superficies y materiales nanoestructurados, 5-7 June 2019, BsAs, Argentina

Expositores XIII Jornada de Ciencia y Tecnología.

Luque, E.; Frattini, A.; Oliveros, M.; Batista da Silva, M. Lombardi J. M. (2019)

Construcción de equipamiento de bajo costo para el Laboratorio de Física de la FCByF de la UNR.

M.Oliveros, M.Batista de Silva, J.M.Lombardi, E.Luque, A.Frattini.

VI Jornadas de Física Aplicada a las Ciencias Biomédicas Rodolfo J. Rasia. 2019

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Restauracion de la multilinealidad en datos cromatograficos median un algoritmo basado en análisis funcional de datos.

J.M.Lombardi, M.Alcaraza, M.Montemurrob, S.A.Bortolato.

10° congreso de química analítica, (2019)

On-surface transmetalation of metalloporphyrins self-assembled monolayers.

J.M.Lombardi, D.Hötger, P.Abufager, C.Morchutt, P.Alexa, D.Grumelli, J.Dreiser, S.Stepanow, H.F.Busnengo, M.Etzkorn, R.Gutzler and K.Kern.

VI San Luis Conference on Surfaces, Interfaces and Catalysis, 2018, Santa Fe, Argentina

Cascada de clasificadores para la detección y cuantificación de parasitos en celulas infectadas con T.cruzi.

J.M. Lombardi.

Workshop on Scientific Programming Techniques, February 26th to March 9th, 2018, Universidad nacional de Quilmes (UNL). https://wtpc.github.io/

Desarrollo de un nuevo algoritmo basado en análisis funcional para restaurar la multilinealidad en datos cromatográficos.

J.M.Lombardi, S.Bortolato.

9th Analytical Chemistry Congress, November 7th to 10th, 2017, Rio Cuarto

FDA on chemometrics. VIII Workshop Wavelets and Information Theory.

J. M. Lombardi, S.Bortolato.

9th-11th August 2016, Universidad Nacional de la Plata(UNLP)

RECENT AWARDS

2022 Best Flash Poster selected by the Scientific Committee of NANO XX.

2021 BIO Hackathon 2nd place

Córdoba-Santa Fe. October 14th and 16th. 2021

2017 Nasa Space Apps Challenge - Winner

Category: best data analysis

SKILLS

Programming Skills (Most Relevant):

Python, Matlab, C, C++, Java, openGL, R, Fortran, Cuda Data science, Machine learning, Data mining

Libraries (Most Relevant):

Numpy, Pandas, Numba, openCV, openGL, modernGL, pyopenCl, TensorFlow, Theano, Keras, Sklearn, AMP, ASE

Tools (Most Relevant):

Visual studio, Sublime, Linux, VIM, VMD, Látex, Git, Blender, Photoshop, Ilustrator, Microsoft office, libre office, VASP, Gaussian, Hyperchem.