

Curriculum Vitae

PERSONAL DATA

Name: Miguel Ângelo dos Santos Machuqueiro
Place & Birth Date: Palmela, 27-12-1974
Nationality: Portuguese
Address: CQB and BioISI - Biosystems & Integrative Sciences Institute
Faculdade de Ciências, Universidade de Lisboa
Campo Grande, Edifício C8 (sala 8.5.52)
1749-016 Lisboa, Portugal
Phone: (+351) 217500112 Home: (+351) 265509539
Mobile: (+351) 967562285 Skype: miguel.machuqueiro
E-mail: machuque@ciencias.ulisboa.pt Alt. e-mail: miguel.machuqueiro@gmail.com
Group Web: <http://mms.rd.ciencias.ulisboa.pt>
Researcher ID: <http://www.researcherid.com/rid/C-8012-2011>
ORCID ID: <http://orcid.org/0000-0001-6923-8744>

ACADEMIC

2019- Assistant Researcher (CEEC-CI-2017) at the Chemistry and Biochemistry Department, Faculty of Sciences, University of Lisbon.
2014-2018 Invited Assistant Professor at the Chemistry and Biochemistry Department, Faculty of Sciences, University of Lisbon.
2009-2014 Assistant Researcher (Investigador Auxiliar do Programa Ciência2008) at the Chemistry and Biochemistry Center, Faculty of Sciences, University of Lisbon.
2003-2009 Pos-doctoral fellow at the Molecular Simulation and Protein Modelling groups, supervised by Dr. António Baptista and Prof. Cláudio Soares, at ITQB-NOVA, Oeiras, PT.
1999-2003 PhD from the University of Bern, Switzerland (Doctor of Philosophy; PhD; Dr. phil.-nat.), supervised by Prof. Tamis Darbre and Prof. Jean-Louis Reymond.
1992-1998 Diploma in Biochemistry, at the Faculty of Sciences University of Lisbon.

SPECIALIZATION DOMAINS

Chemistry and Biochemistry

Theoretical Chemistry and Biochemistry:

- Computational (Bio)chemistry
- Quantum Mechanics
- Molecular Dynamics
- Molecular Docking
- Molecular Mechanics
- Force Field Development
- Continuum Electrostatics
- Biological Membrane Biophysics

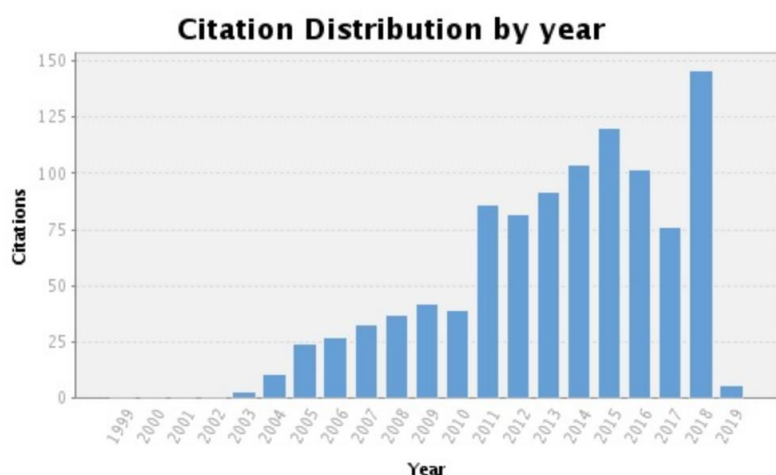
FUNDING: *Coordinator (PI)*:

- 2015- “CpHMD-L simulations of pHLIP peptides: design of new tumor-targeted drug delivery systems” (PTDC/QEQ-COM/5904/2014). Funding: **185.088 Euros**
- 2010-14 “Adding realism to the molecular modeling of lipidic membranes: inclusion of pH effects” (PTDC/QUI-BIQ/113721/2009). Funding: **150.000 Euros**

FUNDING: *Co-PI*:

- 2018-21 “Physical Basis of Disease: The case of Dialysis Related Amyloidosis” (PTDC/FIS-OUT/28210/2017) (Coordination: Patrícia Faísca, FCUL). Funding: **195.144,75 Euros**
- 2018-21 “Targeting multi-resistant tuberculosis with new potent isoniazid derivatives: an integrated medicinal chemistry approach” (PTDC/MED-QUI/29036/2017) (Coordination: Filomena Leitão, FCUL) Funding: **226.020,98 Euros**
- 2018-21 “Deal with PAINS: strategies to identify membrane modulators” (PTDC/BIA-BFS/28419/2017) (Coordination: Bruno L. Victor, FCUL) Funding: **235.111,50 Euros**

PUBLICATIONS STATS (ISI WEB OF SCIENCE)



Total Articles in Publication List: **50**

Articles With Citation Data: **49**

Sum of the Times Cited: **1030**

Average Citations per Article: **21.02**

h-index: **17**

Last Updated: **01/18/2019**
14:24 GMT

The number of publications may not be complete since
ISI WoS does not count recently accepted manuscripts

RECENT PUBLICATIONS:

- Dias, C., Pais, J., Nunes, R., Blázquez-Sánchez, M.T., Marquês, J.T., Almeida, A.F., Serra, P., Xavier, N.M., Vila-Viçosa, D., **Machuqueiro, M.**, Viana, A.S., Martins, A., Santos, M.S., Pelerito, A., Dias, R., Tenreiro, R., Oliveira, M.C., Contino, M., Colabufo, N.A., de Almeida, R.F.M., Rauter, A.P. (2018) "Sugar-Based Bactericides Targeting Phosphatidylethanolamine-Enriched Membranes", *Nat. Commun.*, 9, 4857. doi: [10.1038/s41467-018-06488-4](https://doi.org/10.1038/s41467-018-06488-4)
- Vila-Viçosa, D., Silva, T. F. D., Slaybaugh, G., Reshetnyak, Y. K., Andreev, O. A., **Machuqueiro, M.** (2018) "The membrane-induced pK_a shifts in wt-pHLIP and its L16H variant", *J. Chem. Theory Comput.*, 14, 3289-3297. doi: [10.1021/acs.jctc.8b00102](https://doi.org/10.1021/acs.jctc.8b00102)
- Silva, T. F. D., Vila-Viçosa, D., Reis, P. B. P. S., Victor, B. L., Diem, M., Oostenbrink, C., and **Machuqueiro, M.** (2018) "The impact of using single atomistic long range cutoff schemes with the GROMOS 54A7 force field", *J. Chem. Theory Comput.*, 14, 5823-5833. doi: [10.1021/acs.jctc.8b00758](https://doi.org/10.1021/acs.jctc.8b00758)