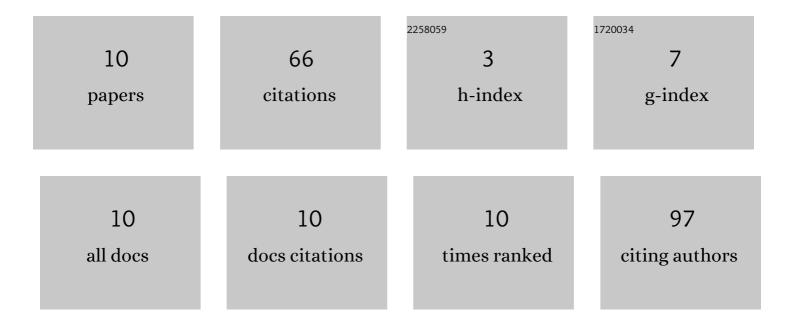
OtÃ;vio LuÃ-s de Santana

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Theoretical study of cooperative effects in the homo- and heteromeric hydrogen bond chains (HCN)nHF withn = 1, 2, and 3. International Journal of Quantum Chemistry, 2006, 106, 2714-2722.	2.0	29
2	Challenges encountered during development of Mn porphyrin-based, potent redox-active drug and superoxide dismutase mimic, MnTnBuOE-2-PyP5+, and its alkoxyalkyl analogues. Journal of Inorganic Biochemistry, 2017, 169, 50-60.	3.5	18
3	Green's function calculation of through-bond electronic coupling in donor–bridge–acceptor model systems. Chemical Physics Letters, 1999, 314, 508-515.	2.6	10
4	Theoretical calculations of the substituent effect on molecular properties of the RCNâ< HF hydrogen-bonded complexes with R=NH2, CH3O, CH3, OH, SH, H, Cl, F, CF3, CN and NO2. Vibrational Spectroscopy, 2009, 49, 133-141.	2.2	3
5	One-Electron Reduction Potentials: Calibration of Theoretical Protocols for Morita–Baylis–Hillman Nitroaromatic Compounds in Aprotic Media. Molecules, 2018, 23, 2129.	3.8	3
6	Electronic properties of the low-lying spin states of dimethylnitrosamine coordinated to Fe(III) heme models: An ab initio study. International Journal of Quantum Chemistry, 2014, 114, 508-520.	2.0	2
7	O ensino de reações orgânicas usando quÃmica computacional: I. reações de adição eletrofÃłica a alquenos. Quimica Nova, 2008, 31, 1243-1249.	0.3	1
8	Green's function calculation of through-bond electronic coupling in donor-bridge-acceptor model systems II: Importance factors applied to atomic sites. International Journal of Quantum Chemistry, 2005, 102, 265-274.	2.0	0
9	Diradicalar Character and Ring Stability of Mesoionic Heterocyclic Oxazoles and Thiazoles by Ab Initio Mono and Multi-Reference Methods. Molecules, 2020, 25, 4524.	3.8	0
10	A multireference configuration interaction study with singles and doubles of some mesoionic rings: reaction and activation free energies for the ringâ€opening reaction. International Journal of Quantum Chemistry, 2020, 120, e26391.	2.0	0