

# Ignacio Soteras

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1061798/publications.pdf>

Version: 2024-02-01

24  
papers

1,040  
citations

471509

17  
h-index

642732

23  
g-index

26  
all docs

26  
docs citations

26  
times ranked

1605  
citing authors

#	ARTICLE	IF	CITATIONS
1	Machine Learning applied to Wi-Fi fingerprinting: The experiences of the Ubiquim Challenge. , 2019, , .		6
2	Structural and energetic study of cation-π cation interactions in proteins. Physical Chemistry Chemical Physics, 2017, 19, 9849-9861.	2.8	19
3	Parametrization of halogen bonds in the CHARMM general force field: Improved treatment of ligand-protein interactions. Bioorganic and Medicinal Chemistry, 2016, 24, 4812-4825.	3.0	168
4	Unraveling phosphodiesterase surfaces. Identification of phosphodiesterase 7 allosteric modulation cavities. European Journal of Medicinal Chemistry, 2013, 70, 781-788.	5.5	5
5	Effect of Phosphodiesterase 7 (PDE7) Inhibitors in Experimental Autoimmune Encephalomyelitis Mice. Discovery of a New Chemically Diverse Family of Compounds. Journal of Medicinal Chemistry, 2012, 55, 3274-3284.	6.4	52
6	5-Imino-1,2,4-Thiadiazoles: First Small Molecules As Substrate Competitive Inhibitors of Glycogen Synthase Kinase 3. Journal of Medicinal Chemistry, 2012, 55, 1645-1661.	6.4	76
7	The boron-boron single bond in diborane(4) as a non-classical electron donor for hydrogen bonding. Physical Chemistry Chemical Physics, 2011, 13, 14026.	2.8	33
8	Exploring the Binding Sites of Glycogen Synthase Kinase 3. Identification and Characterization of Allosteric Modulation Cavities. Journal of Medicinal Chemistry, 2011, 54, 8461-8470.	6.4	91
9	Performance of the IEF-MST solvation continuum model in the SAMPL2 blind test prediction of hydration and tautomerization free energies. Journal of Computer-Aided Molecular Design, 2010, 24, 281-291.	2.9	24
10	Boron-Based Dipolar Multicomponent Reactions: Simple Generation of Substituted Aziridines, Oxazolidines and Pyrrolidines. Chemistry - A European Journal, 2010, 16, 7904-7915.	3.3	27
11	Polarizable Intermolecular Potentials for Water and Benzene Interacting with Halide and Metal Ions. Journal of Chemical Theory and Computation, 2009, 5, 3022-3031.	5.3	34
12	Performance of the IEF-MST Solvation Continuum Model in a Blind Test Prediction of Hydration Free Energies. Journal of Physical Chemistry B, 2009, 113, 9330-9334.	2.6	17
13	The impact of monovalent ion force field model in nucleic acids simulations. Physical Chemistry Chemical Physics, 2009, 11, 10596.	2.8	62
14	Induction effects in metal cation-benzene complexes. Physical Chemistry Chemical Physics, 2008, 10, 2616.	2.8	78
15	Structure-Directed Reversion in the $\pi$ -Facial Stereoselective Alkylation of Chiral Bicyclic Lactams. Journal of Organic Chemistry, 2008, 73, 7756-7763.	3.2	13
16	An ab initio strategy for handling induction phenomena in metal ion complexes. Molecular Physics, 2008, 106, 1685-1696.	1.7	3
17	Derivation of Distributed Models of Atomic Polarizability for Molecular Simulations. Journal of Chemical Theory and Computation, 2007, 3, 1901-1913.	5.3	41
18	Modeling Induction Phenomena in Intermolecular Interactions with an Ab Initio Force Field. Journal of Chemical Theory and Computation, 2007, 3, 1914-1926.	5.3	34

#	ARTICLE	IF	CITATIONS
19	Benzoderivatives of Nucleic Acid Bases as Modified DNA Building Blocks. Journal of Physical Chemistry A, 2006, 110, 510-518.	2.5	21
20	On the Origin of the Stereoselectivity in the Alkylation of Oxazolopiperidone Enolates. Journal of the American Chemical Society, 2006, 128, 6581-6588.	13.7	17
21	Extension of the MST model to the IEF formalism: HF and B3LYP parametrizations. Computational and Theoretical Chemistry, 2005, 727, 29-40.	1.5	79
22	MST Continuum Study of the Hydration Free Energies of Monovalent Ionic Species. Journal of Physical Chemistry B, 2005, 109, 3565-3574.	2.6	44
23	Group contributions to the solvation free energy from MST continuum calculations. Brazilian Journal of Physics, 2004, 34, 48-57.	1.4	7
24	Continuum solvation models: Dissecting the free energy of solvation. Physical Chemistry Chemical Physics, 2003, 5, 3827-3836.	2.8	89