

Sijie Luo

List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

11
papers

724
citations

9
h-index

12
g-index

12
ext. papers

821
ext. citations

5.5
avg, IF

3.81
L-index

#	Paper	IF	Citations
11	Multiconfiguration Pair-Density Functional Theory. <i>Journal of Chemical Theory and Computation</i> , 2014 , 10, 3669-80	6.4	246
10	Bulk Properties of Transition Metals: A Challenge for the Design of Universal Density Functionals. <i>Journal of Chemical Theory and Computation</i> , 2014 , 10, 3832-9	6.4	187
9	Validation of electronic structure methods for isomerization reactions of large organic molecules. <i>Physical Chemistry Chemical Physics</i> , 2011 , 13, 13683-9	3.6	67
8	Improved CO Adsorption Energies, Site Preferences, and Surface Formation Energies from a Meta-Generalized Gradient Approximation Exchange-Correlation Functional, M06-L. <i>Journal of Physical Chemistry Letters</i> , 2012 , 3, 2975-9	6.4	59
7	Density Functional Theory of Open-Shell Systems. The 3d-Series Transition-Metal Atoms and Their Cations. <i>Journal of Chemical Theory and Computation</i> , 2014 , 10, 102-21	6.4	55
6	How Evenly Can Approximate Density Functionals Treat the Different Multiplicities and Ionization States of 4d Transition Metal Atoms?. <i>Journal of Chemical Theory and Computation</i> , 2012 , 8, 4112-26	6.4	34
5	Noncollinear Spins Provide a Self-Consistent Treatment of the Low-Spin State of a Biomimetic Oxomanganese Synthetic Trimer Inspired by the Oxygen Evolving Complex of Photosystem II. <i>Journal of Physical Chemistry Letters</i> , 2011 , 2, 2629-2633	6.4	28
4	The Structure of Silica Surfaces Exposed to Atomic Oxygen. <i>Journal of Physical Chemistry C</i> , 2013 , 117, 9311-9321	3.8	22
3	Noncollinear Spin States for Density Functional Calculations of Open-Shell and Multi-Configurational Systems: Dissociation of MnO and NiO and Barrier Heights of O ₃ , BeH ₂ , and H ₄ . <i>Journal of Chemical Theory and Computation</i> , 2013 , 9, 5349-55	6.4	13
2	Computational Electrochemistry. Voltages of Lithium-Ion Battery Cathodes. <i>Journal of Physical Chemistry B</i> , 2016 , 120, 1437-9	3.4	5
1	Ligand-Mediated Ring \rightarrow Cube Transformation in a Catalytic Subnanocluster: Co ₄ O ₄ (MeCN) _n with n = 1-6. <i>Journal of Physical Chemistry Letters</i> , 2014 , 5, 2528-32	6.4	5