

William C Isley Iii

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.

13
papers

1,035
citations

11
h-index

13
g-index

13
ext. papers

1,166
ext. citations

8
avg. IF

3.54
L-index

#	Paper	IF	Citations
13	Destruction of chemical warfare agents using metal-organic frameworks. <i>Nature Materials</i> , 2015 , 14, 512-6	27	647
12	Design of a metal-organic framework with enhanced back bonding for separation of Ni^{2+} and Cu^{2+} . <i>Journal of the American Chemical Society</i> , 2014 , 136, 698-704	16.4	134
11	Proof for the concerted inversion mechanism in the trans \rightarrow cis isomerization of azobenzene using hydrogen bonding to induce isomer locking. <i>Journal of Organic Chemistry</i> , 2010 , 75, 4817-27	4.2	70
10	Predicting paramagnetic ^1H NMR chemical shifts and state-energy separations in spin-crossover host-guest systems. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 10620-8	3.6	31
9	Empirical and theoretical insights into the structural features and host-guest chemistry of M8L4 tube architectures. <i>Journal of the American Chemical Society</i> , 2014 , 136, 3972-80	16.4	28
8	Ab Initio Extension of the AMOEBA Polarizable Force Field to Fe(2.). <i>Journal of Chemical Theory and Computation</i> , 2013 , 9, 3062-71	6.4	26
7	Electrochemical Reduction of 2,4-Dinitrotoluene in Aprotic and pH-Buffered Media. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 13088-13097	3.8	25
6	Optimization and prediction of the electron-nuclear dipolar and scalar interaction in H and C liquid state dynamic nuclear polarization. <i>Chemical Science</i> , 2015 , 6, 6482-6495	9.4	21
5	Prediction of (^{19}F) NMR Chemical Shifts in Labeled Proteins: Computational Protocol and Case Study. <i>Molecular Pharmaceutics</i> , 2016 , 13, 2376-86	5.6	18
4	Quantum Chemical Methods for Modeling Covalent Modification of Biological Thiols. <i>Journal of Computational Chemistry</i> , 2020 , 41, 427-438	3.5	18
3	Molecular Modeling Characterization of a Conformationally Constrained Monolayer-Protected Gold Cluster. <i>Journal of Physical Chemistry C</i> , 2010 , 114, 16043-16050	3.8	11
2	Electric Potentials of Metastable Salt Clusters. <i>Journal of Physical Chemistry C</i> , 2019 , 123, 14010-14023	3.8	3
1	In situ microscopy across scales for the characterization of crystal growth mechanisms: the case of europium oxalate. <i>CrystEngComm</i> , 2018 , 20, 2822-2833	3.3	3