

Xin Wu

List of Publications by Year in descending order

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13
papers

1,077
citations

687363

13
h-index

996975

15
g-index

17
all docs

17
docs citations

17
times ranked

890
citing authors

#	ARTICLE	IF	CITATIONS
1	Big data analysis of <i>ab Initio</i> molecular integrals in the neglect of diatomic differential overlap approximation. <i>Journal of Computational Chemistry</i> , 2019, 40, 638-649.	3.3	10
2	Semiempirical Quantum-Chemical Methods with Orthogonalization and Dispersion Corrections. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 1743-1760.	5.3	45
3	Semiempirical Quantum-Chemical Orthogonalization-Corrected Methods: Theory, Implementation, and Parameters. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 1082-1096.	5.3	123
4	Semiempirical Quantum-Chemical Orthogonalization-Corrected Methods: Benchmarks for Ground-State Properties. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 1097-1120.	5.3	74
5	Specific Reaction Path Hamiltonian for Proton Transfer in Water: Reparameterized Semiempirical Models. <i>Journal of Chemical Theory and Computation</i> , 2013, 9, 2672-2686.	5.3	42
6	Semiempirical Quantum Chemical Calculations Accelerated on a Hybrid Multicore CPU-GPU Computing Platform. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 2272-2281.	5.3	44
7	Hybrid Quantum and Classical Simulations of the Dihydrofolate Reductase Catalyzed Hydride Transfer Reaction on an Accurate Semi-Empirical Potential Energy Surface. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 3420-3437.	5.3	45
8	Crystal Structures of Saturn-Like C ₅₀ Cl ₁₀ and Pineapple-Shaped C ₆₄ Cl ₄ : Geometric Implications of Double and Triple Pentagon-Fused Chlorofullerenes. <i>Angewandte Chemie - International Edition</i> , 2008, 47, 5340-5343.	13.8	116
9	Two Ih-symmetry-breaking C ₆₀ isomers stabilized by chlorination. <i>Nature Materials</i> , 2008, 7, 790-794.	27.5	114
10	An Entrant of Smaller Fullerene: C ₅₆ Captured by Chlorines and Aligned in Linear Chains. <i>Journal of the American Chemical Society</i> , 2008, 130, 15240-15241.	13.7	69
11	Dimetalloendofullerene U ₂ @C ₆₀ Has a U-U Multiple Bond Consisting of Sixfold One-Electron-Two-Center Bonds. <i>Journal of the American Chemical Society</i> , 2007, 129, 2171-2177.	13.7	95
12	Structures and Electronic Properties of M ₂ C ₂ @C ₇₈ (M = Ti, Zr, Hf): A Density Functional Theory Study. <i>Journal of Nanoscience and Nanotechnology</i> , 2007, 7, 1346-1352.	0.9	8
13	Isolation and Characterization of Sc ₂ C ₂ @C ₆₈ : A Metal-Carbide Endofullerene with a Non-IPR Carbon Cage. <i>Angewandte Chemie - International Edition</i> , 2006, 45, 2107-2111.	13.8	181