

Jia Chen

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

522
citations

1478505
6
h-index

1372567
10
g-index

10
all docs

10
docs citations

10
times ranked

459
citing authors

#	ARTICLE	IF	CITATIONS
1	Recent developments in the PySCF program package. <i>Journal of Chemical Physics</i> , 2020, 153, 024109.	3.0	388
2	Density functional plus dynamical mean-field theory of the spin-crossover molecule $\text{Fe}(\text{phen})_3^{2+}$. <i>Physical Review B</i> , 2015, 91, .		
3	Decoherence in Molecular Electron Spin Qubits: Insights from Quantum Many-Body Simulations. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 2074-2078.	4.6	32
4	Quantum-Inspired Algorithm for the Factorized Form of Unitary Coupled Cluster Theory. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 841-847.	5.3	21
5	Three Jahn-Teller States of Matter in Spin-Crossover System Mn(taa). <i>Physical Review Letters</i> , 2020, 124, 227201.	7.8	11
6	Intermolecular coupling and superconductivity in PbMo_6S_8 and other Chevrel phase compounds. <i>Physical Review Materials</i> , 2018, 2, .	2.4	9
7	Generalized quadrature for finite temperature Green's function methods. <i>Computer Physics Communications</i> , 2020, 253, 107178.	7.5	5
8	Long-Range Magnetic Exchange Pathways in Complex Clusters from First Principles. <i>Journal of Physical Chemistry C</i> , 2021, 125, 11124-11131.	3.1	4
9	Flexibility of the factorized form of the unitary coupled cluster Ansatz. <i>Journal of Chemical Physics</i> , 2022, 156, 044106.	3.0	3
10	Low-Depth Unitary Coupled Cluster Theory for Quantum Computation. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 2193-2198.	5.3	3