

Jianguo Yu

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/289276/publications.pdf>

Version: 2024-02-01

29
papers

1,622
citations

430874

18
h-index

526287

27
g-index

33
all docs

33
docs citations

33
times ranked

2796
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Reversible Sodium Ion Insertion in Single Crystalline Manganese Oxide Nanowires with Long Cycle Life. <i>Advanced Materials</i> , 2011, 23, 3155-3160. | 21.0 | 638 |
| 2 | Charge optimized many-body potential for the Si ⁺ •SiO ₂ system. <i>Physical Review B</i> , 2007, 75, . | 3.2 | 159 |
| 3 | Atomic-Resolution Visualization of Distinctive Chemical Mixing Behavior of Ni, Co, and Mn with Li in Layered Lithium Transition-Metal Oxide Cathode Materials. <i>Chemistry of Materials</i> , 2015, 27, 5393-5401. | 6.7 | 108 |
| 4 | First-principles study of defects and phase transition in UO ₂ . <i>Journal of Physics Condensed Matter</i> , 2009, 21, 435401. | 1.8 | 71 |
| 5 | Charge and Ion Transport in NiO and Aspects of Ni Oxidation from First Principles. <i>Journal of Physical Chemistry C</i> , 2012, 116, 1948-1954. | 3.1 | 62 |
| 6 | Coverage dependence and hydroperoxyl-mediated pathway of catalytic water formation on Pt (111) surface. <i>Journal of Chemical Physics</i> , 2006, 125, 054701. | 3.0 | 61 |
| 7 | Effects of Short-Range Attraction in Metal Epitaxial Growth. <i>Physical Review Letters</i> , 2002, 89, 286103. | 7.8 | 59 |
| 8 | Energetic recoils in UO ₂ simulated using five different potentials. <i>Journal of Chemical Physics</i> , 2009, 130, 174502. | 3.0 | 52 |
| 9 | Dynamical scaling behavior in two-dimensional ballistic deposition with shadowing. <i>Physical Review E</i> , 2002, 66, 021603. | 2.1 | 47 |
| 10 | Atomic-Scale Mechanisms of Enhanced Electrochemical Properties of Mo-Doped Co-Free Layered Oxide Cathodes for Lithium-Ion Batteries. <i>ACS Energy Letters</i> , 2019, 4, 2540-2546. | 17.4 | 40 |
| 11 | Unified interatomic potential for zircon, zirconia and silica systems. <i>Journal of Materials Chemistry</i> , 2009, 19, 3923. | 6.7 | 38 |
| 12 | Kinetic Monte Carlo Study of Ambipolar Lithium Ion and Electron Polaron Diffusion into Nanostructured TiO ₂ . <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2076-2081. | 4.6 | 38 |
| 13 | Homogeneous hydride formation path in H ₂ -Zr: Molecular dynamics simulations with the charge-optimized many-body potential. <i>Acta Materialia</i> , 2016, 111, 357-365. | 7.9 | 35 |
| 14 | Molecular-dynamics simulation of threshold displacement energies in zircon. <i>Nuclear Instruments & Methods in Physics Research B</i> , 2009, 267, 3431-3436. | 1.4 | 31 |
| 15 | Bubble formation and Kr distribution in Kr-irradiated UO ₂ . <i>Journal of Nuclear Materials</i> , 2015, 456, 125-132. | 2.7 | 29 |
| 16 | Scaling behavior of the surface in ballistic deposition. <i>Physical Review E</i> , 2002, 65, 060601. | 2.1 | 25 |
| 17 | Charge Localization and Transport in Lithiated Olivine Phosphate Materials. <i>Journal of Physical Chemistry C</i> , 2011, 115, 25001-25006. | 3.1 | 23 |
| 18 | Short-range attraction, surface currents, and mound formation in metal (111) epitaxial growth. <i>Physical Review B</i> , 2004, 69, . | 3.2 | 19 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 19 | Ab initio study of lithium transition metal fluorophosphate cathodes for rechargeable batteries. Journal of Materials Chemistry, 2011, 21, 12054. | 6.7 | 18 |
| 20 | Migration mechanisms of oxygen interstitial clusters in UO ₂ . Journal of Physics Condensed Matter, 2013, 25, 015003. | 1.8 | 18 |
| 21 | First-principles calculations of steering forces in epitaxial growth. Physical Review B, 2004, 69, . | 3.2 | 12 |
| 22 | Interatomic potential for the structure and energetics of tetrahedrally coordinated silica polymorphs. Physical Review B, 2007, 75, . | 3.2 | 10 |
| 23 | Oxygen transport in off-stoichiometric uranium dioxide mediated by defect clustering dynamics. Journal of Chemical Physics, 2015, 142, 094705. | 3.0 | 8 |
| 24 | Effects of short-range attraction in unstable metal (110) epitaxial growth: Molecular dynamics simulations for Cu, Ag, and Al. Physical Review B, 2006, 74, . | 3.2 | 6 |
| 25 | The impacts of charge transfer, localization, and metallicity on hydrogen retention and transport capacity. International Journal of Hydrogen Energy, 2022, 47, 20194-20204. | 7.1 | 4 |
| 26 | Near Surface Stoichiometry in UO ₂ : A Density Functional Theory Study. Journal of Chemistry, 2015, 2015, 1-8. | 1.9 | 2 |
| 27 | Influence of Alloying Elements and Effect of Stress on Anisotropic Hydrogen Diffusion in Zr-Based Alloys Predicted by Accelerated Kinetic Monte Carlo Simulations. Minerals, Metals and Materials Series, 2019, , 1815-1826. | 0.4 | 2 |
| 28 | Accident tolerant fuel rod failure under low stress: A case study of BWR under station blackout using Bison. Journal of Nuclear Materials, 2021, 553, 153037. | 2.7 | 1 |
| 29 | From Used Oxide Nuclear Fuel to Rechargeable Battery: A First-Principles Study. Materials Research Society Symposia Proceedings, 2013, 1541, 73201. | 0.1 | 0 |