

Yun-Jiang Wang

List of Publications by Year in descending order

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79
papers

2,051
citations

201385

27
h-index

288905

40
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79
all docs

79
docs citations

79
times ranked

1477
citing authors

#	ARTICLE	IF	CITATIONS
1	Atomistic interpretation of extra temperature and strain-rate sensitivity of heterogeneous dislocation nucleation in a multi-principal-element alloy. <i>International Journal of Plasticity</i> , 2022, 149, 103155.	4.1	18
2	Dynamic responses in shocked Cu-Zr nanoglasses with gradient microstructure. <i>International Journal of Plasticity</i> , 2022, 149, 103154.	4.1	15
3	Disentangling diffusion heterogeneity in high-entropy alloys. <i>Acta Materialia</i> , 2022, 224, 117527.	3.8	25
4	Correlation between vibrational anomalies and emergent anharmonicity of the local potential energy landscape in metallic glasses. <i>Physical Review B</i> , 2022, 105, .	1.1	12
5	Ergodic Structural Diversity Predicts Dynamics in Amorphous Materials. <i>Frontiers in Materials</i> , 2022, 9, .	1.2	0
6	Sluggish dynamics of homogeneous flow in high-entropy metallic glasses. <i>Scripta Materialia</i> , 2022, 214, 114673.	2.6	11
7	A hierarchically correlated flow defect model for metallic glass: Universal understanding of stress relaxation and creep. <i>International Journal of Plasticity</i> , 2022, 154, 103288.	4.1	29
8	Elastic criterion for shear-banding instability in amorphous solids. <i>Physical Review E</i> , 2022, 105, 045003.	0.8	8
9	Hidden spatiotemporal sequence in transition to shear band in amorphous solids. <i>Physical Review Research</i> , 2022, 4, .	1.3	10
10	Sluggish hydrogen diffusion and hydrogen decreasing stacking fault energy in a high-entropy alloy. <i>Materials Today Communications</i> , 2021, 26, 101902.	0.9	11
11	Inelastic deformation of metallic glasses under dynamic cyclic loading. <i>Scripta Materialia</i> , 2021, 194, 113675.	2.6	6
12	Grain boundary-mediated plasticity accommodating the cracking process in nanograined gold: In situ observations and simulations. <i>Scripta Materialia</i> , 2021, 194, 113693.	2.6	6
13	Synergistic strengthening mechanisms of rhenium in nickel-based single crystal superalloys. <i>Intermetallics</i> , 2021, 132, 107133.	1.8	15
14	Dynamic mechanical relaxation and thermal creep of high-entropy La ₃₀ Ce ₃₀ Ni ₁₀ Al ₂₀ Co ₁₀ bulk metallic glass. <i>Science China: Physics, Mechanics and Astronomy</i> , 2021, 64, 1.	2.0	37
15	Machine-learning integrated glassy defect from an intricate configurational-thermodynamic-dynamic space. <i>Physical Review B</i> , 2021, 104, .	1.1	15
16	Stress relaxation in high-entropy Pd ₂₀ Pt ₂₀ Cu ₂₀ Ni ₂₀ P ₂₀ metallic glass: Experiments, modeling and theory. <i>Mechanics of Materials</i> , 2021, 160, 103959.	1.7	5
17	Machine learning atomic-scale stiffness in metallic glass. <i>Extreme Mechanics Letters</i> , 2021, 48, 101446.	2.0	19
18	Microstructural effects on the dynamical relaxation of glasses and glass composites: A molecular dynamics study. <i>Acta Materialia</i> , 2021, 220, 117293.	3.8	9

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19	Dislocation nucleation and evolution at the ferrite-cementite interface under cyclic loadings. <i>Acta Materialia</i> , 2020, 186, 267-277.	3.8	30
20	“Self-sharpening” tungsten high-entropy alloy. <i>Acta Materialia</i> , 2020, 186, 257-266.	3.8	91
21	Ultrasonic plasticity of metallic glass near room temperature. <i>Applied Materials Today</i> , 2020, 21, 100866.	2.3	15
22	Complexity of plastic instability in amorphous solids: Insights from spatiotemporal evolution of vibrational modes. <i>European Physical Journal E</i> , 2020, 43, 56.	0.7	10
23	Unraveling strongly entropic effect on α -relaxation in metallic glass: Insights from enhanced atomistic samplings over experimentally relevant timescales. <i>Physical Review B</i> , 2020, 102, .	1.1	5
24	Unified perspective on structural heterogeneity of a LaCe-based metallic glass from versatile dynamic stimuli. <i>Intermetallics</i> , 2020, 125, 106922.	1.8	8
25	Statistical complexity of potential energy landscape as a dynamic signature of the glass transition. <i>Physical Review B</i> , 2020, 101, .	1.1	12
26	Atomistic structural mechanism for the glass transition: Entropic contribution. <i>Physical Review B</i> , 2020, 101, .	1.1	28
27	Novel atomic-scale mechanism of incipient plasticity in a chemically complex CrCoNi medium-entropy alloy associated with inhomogeneity in local chemical environment. <i>Acta Materialia</i> , 2020, 194, 283-294.	3.8	101
28	Ratchetting in Cold-Drawn Pearlitic Steel Wires. <i>Metallurgical and Materials Transactions A: Physical Metallurgy and Materials Science</i> , 2019, 50, 4561-4568.	1.1	8
29	Hierarchical-microstructure based modeling for plastic deformation of partial recrystallized copper. <i>Mechanics of Materials</i> , 2019, 139, 103207.	1.7	3
30	Revisiting the structure-property relationship of metallic glasses: Common spatial correlation revealed as a hidden rule. <i>Physical Review B</i> , 2019, 99, .	1.1	50
31	One-step annealing optimizes strength-ductility tradeoff in pearlitic steel wires. <i>Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing</i> , 2019, 757, 1-13.	2.6	25
32	Atomistic insights on the influence of pre-oxide shell layer and size on the compressive mechanical properties of nickel nanowires. <i>Journal of Applied Physics</i> , 2019, 125, .	1.1	4
33	Assessing the utility of structure in amorphous materials. <i>Journal of Chemical Physics</i> , 2019, 150, 114502.	1.2	34
34	Incorporating a soft ordered phase into an amorphous configuration enhances its uniform plastic deformation under shear. <i>AIP Advances</i> , 2019, 9, 015329.	0.6	1
35	Fast surface dynamics enabled cold joining of metallic glasses. <i>Science Advances</i> , 2019, 5, eaax7256.	4.7	87
36	Susceptibility of shear banding to chemical short-range order in metallic glasses. <i>Scripta Materialia</i> , 2019, 162, 141-145.	2.6	22

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37	Structural Parameter of Orientational Order to Predict the Boson Vibrational Anomaly in Glasses. <i>Physical Review Letters</i> , 2019, 122, 015501.	2.9	45
38	Enhancing strength without compromising ductility in copper by combining extrusion machining and heat treatment. <i>Journal of Materials Processing Technology</i> , 2019, 267, 52-60.	3.1	19
39	Investigation of high spin states in 133Cs. <i>European Physical Journal A</i> , 2018, 54, 1.	1.0	1
40	A free energy landscape perspective on the nature of collective diffusion in amorphous solids. <i>Acta Materialia</i> , 2018, 157, 165-173.	3.8	33
41	Oxyhydroxide of metallic nanowires in a molecular H ₂ O and H ₂ O ₂ environment and their effects on mechanical properties. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 17289-17303.	1.3	17
42	Atomic structure of the Fe/Fe ₃ C interface with the Isaichev orientation in pearlite. <i>Philosophical Magazine</i> , 2017, 97, 2375-2386.	0.7	22
43	Atomic theory of viscoelastic response and memory effects in metallic glasses. <i>Physical Review B</i> , 2017, 96, .	1.1	27
44	Strain gradient drives shear banding in metallic glasses. <i>Physical Review B</i> , 2017, 96, .	1.1	34
45	Universal structural softening in metallic glasses indicated by boson heat capacity peak. <i>Applied Physics Letters</i> , 2017, 111, .	1.5	15
46	Correlation between strain rate sensitivity and $\hat{\tau}$ relaxation of metallic glasses. <i>AIP Advances</i> , 2016, 6, 075022.	0.6	1
47	Effects of oxidation on tensile deformation of iron nanowires: Insights from reactive molecular dynamics simulations. <i>Journal of Applied Physics</i> , 2016, 120, .	1.1	27
48	Prediction of pressure-promoted thermal rejuvenation in metallic glasses. <i>Npj Computational Materials</i> , 2016, 2, .	3.5	67
49	Time-, stress-, and temperature-dependent deformation in nanostructured copper: Creep tests and simulations. <i>Journal of the Mechanics and Physics of Solids</i> , 2016, 94, 191-206.	2.3	54
50	Time, stress, and temperature-dependent deformation in nanostructured copper: Stress relaxation tests and simulations. <i>Acta Materialia</i> , 2016, 108, 252-263.	3.8	66
51	Mechanism transition and strong temperature dependence of dislocation nucleation from grain boundaries: An accelerated molecular dynamics study. <i>Physical Review B</i> , 2016, 94, .	1.1	32
52	Transition from stress-driven to thermally activated stress relaxation in metallic glasses. <i>Physical Review B</i> , 2016, 94, .	1.1	65
53	Understanding the serrated flow and Johari-Goldstein relaxation of metallic glasses. <i>Journal of Non-Crystalline Solids</i> , 2016, 444, 23-30.	1.5	17
54	Intrinsic structural defects on medium range in metallic glasses. <i>Intermetallics</i> , 2016, 75, 36-41.	1.8	17

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55	Direct atomic-scale evidence for shear–dilatation correlation in metallic glasses. <i>Scripta Materialia</i> , 2016, 112, 37-41.	2.6	28
56	Size-dependent plastic deformation and failure mechanisms of nanotwinned Ni3Al: Insights from an atomistic cracking model. <i>Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing</i> , 2016, 649, 449-460.	2.6	19
57	Universal enthalpy-entropy compensation rule for the deformation of metallic glasses. <i>Physical Review B</i> , 2015, 92, .	1.1	19
58	Publisher's Note: Universal enthalpy-entropy compensation rule for the deformation of metallic glasses [<i>Phys. Rev. B</i> 92, 174118 (2015)]. <i>Physical Review B</i> , 2015, 92, .	1.1	2
59	Thermal expansion accompanying the glass-liquid transition and crystallization. <i>AIP Advances</i> , 2015, 5, .	0.6	13
60	Characteristics of stress relaxation kinetics of La 60 Ni 15 Al 25 bulk metallic glass. <i>Acta Materialia</i> , 2015, 98, 43-50.	3.8	89
61	Bridging shear transformation zone to the atomic structure of amorphous solids. <i>Journal of Non-Crystalline Solids</i> , 2015, 410, 100-105.	1.5	5
62	Studying the elastic properties of nanocrystalline copper using a model of randomly packed uniform grains. <i>Computational Materials Science</i> , 2013, 79, 56-62.	1.4	34
63	Atomistic understanding of diffusion kinetics in nanocrystals from molecular dynamics simulations. <i>Physical Review B</i> , 2013, 88, .	1.1	17
64	Entropic effect on creep in nanocrystalline metals. <i>Acta Materialia</i> , 2013, 61, 3866-3871.	3.8	28
65	Size-dependent transition of deformation mechanism, and nonlinear elasticity in Ni3Al nanowires. <i>Applied Physics Letters</i> , 2013, 102, .	1.5	19
66	Atomistic Design of High Strength Crystalline-Amorphous Nanocomposites. <i>Materials Transactions</i> , 2013, 54, 1592-1596.	0.4	9
67	First Report of Lily Blight and Wilt Caused by <i>Fusarium tricinctum</i> in China. <i>Plant Disease</i> , 2013, 97, 993-993.	0.7	19
68	Grain Size Dependence of Creep in Nanocrystalline Copper by Molecular Dynamics. <i>Materials Transactions</i> , 2012, 53, 156-160.	0.4	35
69	Effect of water stress on leaf photosynthesis, chlorophyll content, and growth of oriental lily. <i>Russian Journal of Plant Physiology</i> , 2011, 58, 844-850.	0.5	56
70	Transition of creep mechanism in nanocrystalline metals. <i>Physical Review B</i> , 2011, 84, .	1.1	51
71	CO adsorption on small Au _n (n = 1 – 7) clusters supported on a reduced rutile TiO ₂ (110) surface: a first-principles study. <i>Chinese Physics B</i> , 2011, 20, 036801.	0.7	6
72	A comparison of the ideal strength between L12Co3(Al,W) and Ni3Al under tension and shear from first-principles calculations. <i>Applied Physics Letters</i> , 2009, 94, .	1.5	72

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73	Mechanical properties and electronic structure of superhard diamondlike BC5: A first-principles study. <i>Journal of Applied Physics</i> , 2009, 106, .	1.1	14
74	Influence of alloying elements on the elastic properties of ternary and quaternary nickel-base superalloys. <i>Philosophical Magazine</i> , 2009, 89, 2935-2947.	0.7	22
75	Influence of the alloying element Re on the ideal tensile and shear strength of $\hat{1}^3\hat{1}^2$ -Ni3Al. <i>Scripta Materialia</i> , 2009, 61, 197-200.	2.6	51
76	Mechanical and electronic properties of 5d transition metal diborides MB2 (M=Re, W, Os, Ru). <i>Journal of Applied Physics</i> , 2009, 105, .	1.1	32
77	Effect of Alloying Elements on the Elastic Properties of $\hat{1}^3$ -Ni and $\hat{1}^3$ -Ni3Al from First-principles Calculations. <i>Materials Research Society Symposia Proceedings</i> , 2009, 1224, 1.	0.1	10
78	The alloying mechanisms of Re, Ru in the quaternary Ni-based superalloys interface: A first principles calculation. <i>Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing</i> , 2008, 490, 242-249.	2.6	42
79	A first-principles survey of the partitioning behaviors of alloying elements on $\hat{1}^3/\hat{1}^2$ interface. <i>Journal of Applied Physics</i> , 2008, 104, 013109.	1.1	15