

Ronald E Miller

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76
papers

6,215
citations

33
h-index

78
g-index

78
ext. papers

6,714
ext. citations

3.8
avg, IF

6.03
L-index

#	Paper	IF	Citations
76	Size-dependent elastic properties of nanosized structural elements. <i>Nanotechnology</i> , 2000 , 11, 139-147	3.4	1435
75	An adaptive finite element approach to atomic-scale mechanics: the quasicontinuum method. <i>Journal of the Mechanics and Physics of Solids</i> , 1999 , 47, 611-642	5	479
74	Atomistic/continuum coupling in computational materials science. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2003 , 11, R33-R68	2	402
73	The Quasicontinuum Method: Overview, applications and current directions. <i>Journal of Computer-Aided Materials Design</i> , 2002 , 9, 203-239		326
72	A unified framework and performance benchmark of fourteen multiscale atomistic/continuum coupling methods. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2009 , 17, 053001	2	266
71	Quasicontinuum Models of Interfacial Structure and Deformation. <i>Physical Review Letters</i> , 1998 , 80, 742-745	7.45	262
70	Multiscale plasticity modeling: coupled atomistics and discrete dislocation mechanics. <i>Journal of the Mechanics and Physics of Solids</i> , 2004 , 52, 755-787	5	222
69	Nanoindentation and incipient plasticity. <i>Journal of Materials Research</i> , 1999 , 14, 2233-2250	2.5	214
68	Quasicontinuum simulation of fracture at the atomic scale. <i>Modelling and Simulation in Materials Science and Engineering</i> , 1998 , 6, 607-638	2	186
67	Modeling Materials: Continuum, Atomistic and Multiscale Techniques 2011 ,		180
66	A continuum plasticity model for the constitutive and indentation behaviour of foamed metals. <i>International Journal of Mechanical Sciences</i> , 2000 , 42, 729-754	5.5	174
65	Coupled atomistic and discrete dislocation plasticity. <i>Physical Review Letters</i> , 2002 , 89, 025501	7.4	165
64	Failure of sandwich beams with metallic foam cores. <i>International Journal of Solids and Structures</i> , 2001 , 38, 4901-4920	3.1	156
63	Finite-temperature quasicontinuum: molecular dynamics without all the atoms. <i>Physical Review Letters</i> , 2005 , 95, 060202	7.4	139
62	Quasicontinuum models of fracture and plasticity. <i>Engineering Fracture Mechanics</i> , 1998 , 61, 427-444	4.2	124
61	A coupled atomistic/continuum model of defects in solids. <i>Journal of the Mechanics and Physics of Solids</i> , 2002 , 50, 2085-2106	5	121
60	Modelling grain-boundary resistance in intergranular dislocation slip transmission. <i>Philosophical Magazine A: Physics of Condensed Matter, Structure, Defects and Mechanical Properties</i> , 2002 , 82, 2511-2527		96

59	Atomic-scale simulations of nanoindentation-induced plasticity in copper crystals with nanometer-sized nickel coatings. <i>Acta Materialia</i> , 2006 , 54, 33-45	8.4	94
58	A coupled atomistics and discrete dislocation plasticity simulation of nanoindentation into single crystal thin films. <i>Acta Materialia</i> , 2004 , 52, 271-284	8.4	94
57	The potential of atomistic simulations and the knowledgebase of interatomic models. <i>Jom</i> , 2011 , 63, 17-17	2.1	91
56	A finite-temperature dynamic coupled atomistic/discrete dislocation method. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2005 , 13, 1101-1118	2	90
55	Deformation characteristics and stress-strain response of nanotwinned copper via molecular dynamics simulation. <i>Acta Materialia</i> , 2009 , 57, 4364-4373	8.4	78
54	On the nonlocal nature of dislocation nucleation during nanoindentation. <i>Journal of the Mechanics and Physics of Solids</i> , 2008 , 56, 1203-1223	5	68
53	Coupled Atomistic/Discrete Dislocation Simulations of Nanoindentation at Finite Temperature. <i>Journal of Engineering Materials and Technology, Transactions of the ASME</i> , 2005 , 127, 358-368	1.8	58
52	Long-Range, Entangled Carbon Nanotube Networks in Polycarbonate. <i>Advanced Functional Materials</i> , 2003 , 13, 868-872	15.6	57
51	Multiscale simulation of material removal processes at the nanoscale. <i>Journal of the Mechanics and Physics of Solids</i> , 2007 , 55, 2384-2405	5	55
50	A stress-gradient based criterion for dislocation nucleation in crystals. <i>Journal of the Mechanics and Physics of Solids</i> , 2004 , 52, 1507-1525	5	52
49	New Perspectives in Plasticity Theory: Dislocation Nucleation, Waves, and Partial Continuity of Plastic Strain Rate. <i>Mathematics and Mechanics of Solids</i> , 2008 , 13, 292-315	2.3	51
48	Atomistic simulation of nanoindentation into copper multilayers. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2005 , 13, 1089-1099	2	41
47	Finite-Temperature Quasi-Continuum. <i>Applied Mechanics Reviews</i> , 2013 , 65,	8.6	40
46	A non-local formulation of the peierls dislocation model. <i>Journal of the Mechanics and Physics of Solids</i> , 1998 , 46, 1845-1867	5	40
45	Hybrid Continuum Mechanics and Atomistic Methods for Simulating Materials Deformation and Failure. <i>MRS Bulletin</i> , 2007 , 32, 920-926	3.2	38
44	A molecular dynamics study of twin width, grain size and temperature effects on the toughness of 2D-columnar nanotwinned copper. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2009 , 17, 055009	2	33
43	Critical analysis of local constitutive models for slip and decohesion. <i>Philosophical Magazine A: Physics of Condensed Matter, Structure, Defects and Mechanical Properties</i> , 1996 , 73, 803-827		30
42	Multiscale modeling of crack initiation and propagation at the nanoscale. <i>Journal of the Mechanics and Physics of Solids</i> , 2016 , 88, 35-49	5	26

41	Crystallographic Texture and Volume Fraction of β and β' Phases in Zr-2.5Nb Pressure Tube Material During Heating and Cooling. <i>Metallurgical and Materials Transactions A: Physical Metallurgy and Materials Science</i> , 2012 , 43, 806-821	2.3	21
40	An Energy Balance Criterion for Nanoindentation-Induced Single and Multiple Dislocation Events. <i>Journal of Applied Mechanics, Transactions ASME</i> , 2006 , 73, 327-334	2.7	21
39	Direct Coupling of Atomistic and Continuum Mechanics in Computational Materials Science. <i>International Journal for Multiscale Computational Engineering</i> , 2003 , 1, 16	2.4	18
38	Grain boundary motion assisted via radiation cascades in bcc Fe. <i>Physical Review B</i> , 2008 , 78,	3.3	17
37	Continuum Mechanics and Thermodynamics: From Fundamental Concepts to Governing Equations 2011 ,		13
36	Molecular dynamics at constant Cauchy stress. <i>Journal of Chemical Physics</i> , 2016 , 144, 184107	3.9	12
35	A perspective on atomistic-continuum multiscale modeling. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2017 , 25, 071004	2	11
34	Superior Dynamic Penetration Resistance of Nanoscale Multilayer Polymer/Metal Films. <i>Journal of Applied Mechanics, Transactions ASME</i> , 2020 , 87,	2.7	10
33	Multiscale modeling of solids at the nanoscale: dynamic approach. <i>Canadian Journal of Physics</i> , 2008 , 86, 391-400	1.1	9
32	Energy absorption mechanisms of nanoscopic multilayer structures under ballistic impact loading. <i>Computational Materials Science</i> , 2021 , 195, 110504	3.2	9
31	Physical properties of liquid hexane and derived polar by-products of hexane autoxidation: molecular dynamics calculations using the TraPPE-UA force field. <i>Molecular Simulation</i> , 2013 , 39, 882-894	2.4	8
30	Synthesis and characterization of partially silane-terminated polyurethanes reinforced with acid-treated halloysite nanotubes for transparent armour systems. <i>Scientific Reports</i> , 2020 , 10, 13805	4.9	7
29	Density Functional Theory Rate Calculation of Hydrogen Abstraction Reactions of N-Phenyl- β -naphthylamine Antioxidants. <i>Industrial & Engineering Chemistry Research</i> , 2018 , 57, 876-880	3.9	6
28	Umbrella spherical integration: a stable meshless method for non-linear solids. <i>International Journal for Numerical Methods in Engineering</i> , 2007 , 69, 2807-2847	2.4	6
27	Molecular dynamics study of the mechanical behaviour of ultrathin polymer/metal multilayers under extreme dynamic conditions. <i>Computational Materials Science</i> , 2020 , 184, 109951	3.2	6
26	Molecular dynamics study of the penetration resistance of multilayer polymer/ceramic nanocomposites under supersonic projectile impacts. <i>Extreme Mechanics Letters</i> , 2021 , 44, 101238	3.9	6
25	Crack Behaviour at Bi-Crystal Interfaces: A Mixed Atomistic and Continuum Approach. <i>Materials Research Society Symposia Proceedings</i> , 2000 , 653, 1		5
24	Molecular Dynamics Simulations of Shock Propagation and Spallation in Amorphous Polymers. <i>Journal of Applied Mechanics, Transactions ASME</i> , 2021 , 88,	2.7	5

23	Multiscale approach for determining hydrogen diffusivity in zirconium. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2018 , 26, 085002	2	5
22	First principles study of hydrogen in lead zirconate titanate. <i>Smart Materials and Structures</i> , 2019 , 28, 034002	3-4	4
21	A Continuum Plasticity Model for the Constitutive Behaviour of Foamed Metals. <i>Materials Research Society Symposia Proceedings</i> , 1998 , 521, 39		4
20	Effect of single initial overload and mean load on the low-cycle fatigue life of normalized 300 M alloy steel. <i>International Journal of Fatigue</i> , 2020 , 130, 105273	5	4
19	The Theory and Implementation of the Quasicontinuum Method 2005 , 663-682		4
18	Benchmarking, validation and reproducibility of concurrent multiscale methods are still needed. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2017 , 25, 071001	2	3
17	Multiscale modeling of ductile crystals at the nanoscale subjected to cyclic indentation. <i>Acta Materialia</i> , 2008 , 56, 2799-2809	8.4	3
16	A multi-state modified embedded atom method potential for titanium. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2017 , 25, 015010	2	2
15	A Differential Scanning Calorimetry (DSC) Study of Phase Changes in an As-Received Zr-2.5Nb Pressure Tube Material during Continuous Heating and Cooling. <i>Materials Science Forum</i> , 2012 , 706-709, 853-858	0.4	2
14	Experimental observations of void growth in the Zr ₂ .5Nb pressure tube alloy. <i>Journal of Nuclear Materials</i> , 2005 , 341, 231-234	3-3	2
13	Microstructural evidence of the toughening mechanisms of polyurethane reinforced with halloysite nanotubes under high strain-rate tensile loading. <i>Scientific Reports</i> , 2021 , 11, 13161	4-9	2
12	Molecular-level investigation on the spallation of polyurea. <i>MRS Communications</i> , 2021 , 11, 532-538	2.7	2
11	Crack Behaviour at Bi-Crystal Interfaces: A Mixed Atomistic and Continuum Approach. <i>Materials Research Society Symposia Proceedings</i> , 2000 , 653,		1
10	Effect of a Halloysite-polyurethane nanocomposite interlayer on the ballistic performance of laminate transparent armour. <i>Composites Part C: Open Access</i> , 2022 , 7, 100231	1.6	1
9	Quantum and classical molecular dynamics simulations of shocked polyurea and polyurethane. <i>Computational Materials Science</i> , 2022 , 203, 111166	3.2	1
8	A study of nano-indentation using coupled atomistic and discrete dislocation (CADD) modeling 2003 , 455-459		1
7	Finite Temperature Coupled Atomistic/Continuum Discrete Dislocation Dynamics Simulation of Nanoindentation 2006 , 225-234		1
6	Molecular dynamics study on the shock induced spallation of polyethylene. <i>Journal of Applied Physics</i> , 2022 , 131, 025102	2.5	0

- 5 Molecular dynamics simulation of pull-out Halloysite nanotube from polyurethane matrix. *Advances in Mechanical Engineering*, **2021**, 13, 168781402110446 1.2 0
- 4 Transiting the molecular potential energy surface along low energy pathways: the TRREAT algorithm. *Journal of Computational Chemistry*, **2013**, 34, 2502-13 3.5
- 3 Coupled atomistic/continuum modelling of plasticity in materials **2007**, 189-219
- 2 Denuded zones in zirconium pressure vessels: oxygen's role examined via multi-scale diffusion model. *Modelling and Simulation in Materials Science and Engineering*, **2020**, 28, 065005 2
- 1 Mesoscopic Length Scales for Deformed Nanostructures **2006**, 263-275