David J Srolovitz

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488 25,655 84 140 h-index g-index citations papers 27,618 7.06 504 5.9 L-index avg, IF ext. citations ext. papers

#	Paper	IF	Citations
488	High-Rate, Gas-Phase Growth of MoS2 Nested Inorganic Fullerenes and Nanotubes. <i>Science</i> , 1995 , 267, 222-5	33.3	1045
487	Development of new interatomic potentials appropriate for crystalline and liquid iron. <i>Philosophical Magazine</i> , 2003 , 83, 3977-3994	1.6	941
486	On the stability of surfaces of stressed solids. <i>Acta Metallurgica</i> , 1989 , 37, 621-625		888
485	Computer simulation of grain growth Kinetics. Acta Metallurgica, 1984, 32, 783-791		876
484	Development of an interatomic potential for phosphorus impurities in '-iron. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S2629-S2642	1.8	442
483	Computer simulation of grain growth II. Grain size distribution, topology, and local dynamics. <i>Acta Metallurgica</i> , 1984 , 32, 793-802		415
482	Computer simulation of normal grain growth in three dimensions. <i>The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties</i> , 1989 , 59, 293-329		333
481	Simulation and theory of abnormal grain growthlinisotropic grain boundary energies and mobilities. <i>Acta Metallurgica</i> , 1989 , 37, 1227-1240		296
480	Crystal-melt interfacial free energies in hcp metals: A molecular dynamics study of Mg. <i>Physical Review B</i> , 2006 , 73,	3.3	281
479	Capillary instabilities in thin films. I. Energetics. <i>Journal of Applied Physics</i> , 1986 , 60, 247-254	2.5	272
478	Computer simulation of grain growth. Abnormal grain growth. Acta Metallurgica, 1985, 33, 2233-2247		261
477	Structural defects in amorphous solids Statistical analysis of a computer model. <i>Philosophical Magazine A: Physics of Condensed Matter, Structure, Defects and Mechanical Properties</i> , 1981 , 44, 847-86	6	256
476	Computer simulation of grain growth-III. Influence of a particle dispersion. <i>Acta Metallurgica</i> , 1984 , 32, 1429-1438		255
475	Physical Origins of Intrinsic Stresses in Volmer-Weber Thin Films. MRS Bulletin, 2002, 27, 19-25	3.2	249
474	Kinetics of ordering in two dimensions. II. Quenched systems. <i>Physical Review B</i> , 1983 , 28, 2705-2716	3.3	231
473	Finite-temperature defect properties from free-energy minimization. <i>Physical Review Letters</i> , 1989 , 63, 624-627	7.4	227
472	Cracklike surface instabilities in stressed solids. <i>Physical Review Letters</i> , 1993 , 71, 1593-1596	7.4	225

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471	Computer simulation of grain growth I V. Anisotropic grain boundary energies. <i>Acta Metallurgica</i> , 1985 , 33, 509-520		216	
470	Computer simulation of recrystallization[] Homogeneous nucleation and growth. <i>Acta Metallurgica</i> , 1986 , 34, 1833-1845		215	
469	An atomistic study of deformation of amorphous metals. <i>Acta Metallurgica</i> , 1983 , 31, 335-352		214	
468	The von Neumann relation generalized to coarsening of three-dimensional microstructures. <i>Nature</i> , 2007 , 446, 1053-5	50.4	213	
467	Capillary instabilities in thin films. II. Kinetics. <i>Journal of Applied Physics</i> , 1986 , 60, 255-260	2.5	201	
466	Engineering the shape and structure of materials by fractal cut. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014 , 111, 17390-5	11.5	196	
465	Surface stress model for intrinsic stresses in thin films. <i>Journal of Materials Research</i> , 2000 , 15, 2468-247	4 .5	187	
464	Local structural fluctuations in amorphous and liquid metals: a simple theory of the glass transition. <i>Journal of Physics F: Metal Physics</i> , 1982 , 12, 2141-2163		184	
463	Radial distribution function and structural relaxation in amorphous solids. <i>Physical Review B</i> , 1981 , 24, 6936-6944	3.3	183	
462	Measurement of the cleavage energy of graphite. <i>Nature Communications</i> , 2015 , 6, 7853	17.4	175	
461	Dislocation E win interaction mechanisms for ultrahigh strength and ductility in nanotwinned metals. <i>Acta Materialia</i> , 2009 , 57, 4508-4518	3.4	160	
460	Effects of lattice anisotropy and temperature on domain growth in the two-dimensional Potts model. <i>Physical Review A</i> , 1991 , 43, 2662-2668	2.6	158	
459	Computer simulation on surfaces and [001] symmetric tilt grain boundaries in Ni, Al, and Ni3Al. Journal of Materials Research, 1989, 4, 62-77	2.5	155	
458	Grain-boundary kinetics: A unified approach. <i>Progress in Materials Science</i> , 2018 , 98, 386-476	1 2.2	154	
457	Domain-growth kinetics for the Q-state Potts model in two and three dimensions. <i>Physical Review B</i> , 1988 , 38, 4752-4760	3.3	153	
456	Elastic fracture in random materials. <i>Physical Review B</i> , 1988 , 37, 5500-5507	3.3	151	
455	Diffusionally modified dislocation-particle elastic interactions. <i>Acta Metallurgica</i> , 1984 , 32, 1079-1088		151	
454	Oscillatory surface relaxations in Ni, Al, and their ordered alloys. <i>Physical Review Letters</i> , 1986 , 57, 1308-	⊁3 ₄ 11	151	

453	Simultaneous grain boundary migration and grain rotation. <i>Acta Materialia</i> , 2006 , 54, 1707-1719	8.4	148
452	Computer simulation of recrystallization in non-uniformly deformed metals. <i>Acta Metallurgica</i> , 1989 , 37, 627-639		145
451	Computer simulation of recrystallization II. Heterogeneous nucleation and growth. <i>Acta Metallurgica</i> , 1988 , 36, 2115-2128		144
450	Grain boundaries exhibit the dynamics of glass-forming liquids. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009 , 106, 7735-40	11.5	141
449	Size-dependent deformation of nanocrystalline Pt nanopillars. <i>Nano Letters</i> , 2012 , 12, 6385-92	11.5	137
448	Columnar growth in thin films. <i>Physical Review Letters</i> , 1988 , 60, 424-427	7.4	137
447	The Thermodynamics and Kinetics of film agglomeration. <i>Jom</i> , 1995 , 47, 31-36	2.1	136
446	Metal-ceramic adhesion and the Harris functional. <i>Physical Review Letters</i> , 1994 , 72, 4021-4024	7.4	135
445	Dislocation distributions in two dimensions. <i>Scripta Metallurgica</i> , 1989 , 23, 1347-1352		133
444	Kinetics of the Q-State Potts Model in Two Dimensions. <i>Physical Review Letters</i> , 1983 , 50, 263-266	7.4	129
443	Void formation during film growth: A molecular dynamics simulation study. <i>Journal of Applied Physics</i> , 1996 , 79, 1448-1457	2.5	127
442	Oxygen Diffusion in Yttria-Stabilized Zirconia: A New Simulation Model. <i>Journal of the American Ceramic Society</i> , 2005 , 87, 1821-1830	3.8	125
441	Inhibition of grain growth by second phase particles: Three dimensional Monte Carlo computer simulations. <i>Scripta Metallurgica</i> , 1989 , 23, 753-758		121
440	Misorientation dependence of intrinsic grain boundary mobility: simulation and experiment. <i>Acta Materialia</i> , 1999 , 47, 3901-3914	8.4	120
439	A kinetic Monte Carlo method for the atomic-scale simulation of chemical vapor deposition: Application to diamond. <i>Journal of Applied Physics</i> , 1997 , 82, 6293-6300	2.5	117
438	Theory of metal@eramic adhesion. <i>Acta Metallurgica Et Materialia</i> , 1995 , 43, 2721-2730		115
437	Twisted Bilayer Graphene: Moir with a Twist. <i>Nano Letters</i> , 2016 , 16, 5923-7	11.5	112
436	Size effect in compression of single-crystal gold microparticles. <i>Acta Materialia</i> , 2011 , 59, 5202-5215	8.4	111

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435	Brittle fracture in polycrystalline microstructures with the extended finite element method. <i>International Journal for Numerical Methods in Engineering</i> , 2003 , 56, 2015-2037	2.4	111
434	Boundary Mobility and Energy Anisotropy Effects on Microstructural Evolution During Grain Growth. <i>Journal of Materials Science</i> , 2002 , 10, 201-216		110
433	Microstructural simulation of dynamic recrystallization. Acta Metallurgica Et Materialia, 1992, 40, 43-55		110
432	Morphology of nested fullerenes. <i>Physical Review Letters</i> , 1995 , 74, 1779-1782	7.4	109
431	Crosshatched surface morphology in strained III-V semiconductor films. <i>Journal of Applied Physics</i> , 1990 , 67, 4093-4098	2.5	109
430	Stress relaxation and misfit dislocation nucleation in the growth of misfitting films: A molecular dynamics simulation study. <i>Journal of Applied Physics</i> , 1998 , 83, 217-227	2.5	105
429	Computer Simulation of Final-Stage Sintering: I, Model Kinetics, and Microstructure. <i>Journal of the American Ceramic Society</i> , 1990 , 73, 2857-2864	3.8	102
428	Computer simulation of the elastically driven migration of a flat grain boundary. <i>Acta Materialia</i> , 2004 , 52, 2569-2576	8.4	100
427	Thermodynamic properties of metastable Ag-Cu alloys. <i>Journal of Applied Physics</i> , 1993 , 74, 3144-3149	2.5	99
426	Analytical and numerical modeling of columnar evolution in thin films. <i>Journal of Vacuum Science and Technology A: Vacuum, Surfaces and Films</i> , 1988 , 6, 2371-2380	2.9	99
425	Computer simulation of grain boundaries in Ni3Al: The effect of grain boundary composition. <i>Scripta Metallurgica</i> , 1986 , 20, 1389-1394		99
424	Systematic prediction of kinetically limited crystal growth morphologies. <i>Physical Review Letters</i> , 2005 , 95, 155503	7.4	98
423	A two-dimensional molecular dynamics simulation of thin film growth by oblique deposition. Journal of Applied Physics, 1996 , 80, 5682-5690	2.5	98
422	Impurity effects on domain-growth kinetics. I. Ising model. <i>Physical Review B</i> , 1985 , 32, 3014-3020	3.3	97
421	Atomistic simulation of the deformation of gold nanopillars. Acta Materialia, 2007, 55, 2085-2099	8.4	96
420	Kinetic Monte Carlo Simulation of Chemical Vapor Deposition. <i>Annual Review of Materials Research</i> , 2002 , 32, 297-319	12.8	96
419	Grain growth phenomena in films: A Monte Carlo approach. <i>Journal of Vacuum Science and Technology A: Vacuum, Surfaces and Films</i> , 1986 , 4, 2925-2931	2.9	96
418	First-principles study of graphene edge properties and flake shapes. <i>Physical Review B</i> , 2010 , 81,	3.3	95

417	Curvature driven grain boundary migration in aluminum: molecular dynamics simulations. <i>Acta Materialia</i> , 2005 , 53, 79-86	8.4	95
416	Self-interstitials in V and Mo. <i>Physical Review B</i> , 2002 , 66,	3.3	94
415	Grain growth in two dimensions. Scripta Metallurgica, 1983, 17, 241-246		94
414	Effect of Fe segregation on the migration of a non-symmetric B tilt grain boundary in Al. <i>Journal of Materials Research</i> , 2005 , 20, 208-218	2.5	93
413	Kinetics of buckling of a compressed film on a viscous substrate. <i>Applied Physics Letters</i> , 2001 , 78, 2482-	-254.84	92
412	On the volume fraction dependence of particle limited grain growth. Scripta Metallurgica, 1987, 21, 675	5-679	92
411	Nanowire failure: long = brittle and short = ductile. <i>Nano Letters</i> , 2012 , 12, 910-4	11.5	91
410	Texture development mechanisms in ion beam assisted deposition. <i>Journal of Applied Physics</i> , 1998 , 84, 5261-5269	2.5	90
409	Nonlinear geometric effects in mechanical bistable morphing structures. <i>Physical Review Letters</i> , 2012 , 109, 114302	7.4	89
408	Molecular dynamics simulation of triple junction migration. <i>Acta Materialia</i> , 2002 , 50, 1405-1420	8.4	89
407	Simulation of faceted film growth in two-dimensions: microstructure, morphology and texture. <i>Acta Materialia</i> , 1999 , 47, 2269-2281	8.4	89
406	Mechanical behavior and interface design of MoSi2-based alloys and composites. <i>Materials Science</i> & <i>A: Structural Materials: Properties, Microstructure and Processing</i> , 1992 , 155, 147-158	5.3	86
405	Nanoindentation size effect in single-crystal nanoparticles and thin films: A comparative experimental and simulation study. <i>Acta Materialia</i> , 2011 , 59, 2309-2321	8.4	85
404	Reconciling grain growth and shear-coupled grain boundary migration. <i>Nature Communications</i> , 2017 , 8, 1764	17.4	83
403	Crystal-melt interfacial free energies in metals: fcc versus bcc. <i>Physical Review B</i> , 2004 , 69,	3.3	83
402	A level set method for dislocation dynamics. <i>Acta Materialia</i> , 2003 , 51, 5499-5518	8.4	82
401	Tunable helical ribbons. <i>Applied Physics Letters</i> , 2011 , 98, 011906	3.4	81
400	First-principles calculation of the thermodynamics of InxGa1\(\text{Inx} \) alloys: Effect of lattice vibrations. <i>Physical Review B</i> , 2006 , 73,	3.3	81

399	Atomistic Simulation of Curvature Driven Grain Boundary Migration. <i>Journal of Materials Science</i> , 1998 , 6, 41-58		80
398	Grain-boundary metastability and its statistical properties. <i>Acta Materialia</i> , 2016 , 104, 259-273	8.4	78
397	Simulation of faceted film growth in three dimensions: microstructure, morphology and texture. <i>Acta Materialia</i> , 2005 , 53, 1191-1204	8.4	78
396	Brittle fracture in materials with random defects. <i>Physical Review B</i> , 1989 , 39, 9273-9281	3.3	77
395	Effects of particle size on inhibited grain growth. Scripta Metallurgica Et Materialia, 1990, 24, 101-106		77
394	Atomic-scale analysis of liquid-gallium embrittlement of aluminum grain boundaries. <i>Acta Materialia</i> , 2014 , 73, 312-325	8.4	75
393	Defect interactions on solid surfaces. Surface Science, 1993, 284, 211-221	1.8	75
392	van der Waals bilayer energetics: Generalized stacking-fault energy of graphene, boron nitride, and graphene/boron nitride bilayers. <i>Physical Review B</i> , 2015 , 92,	3.3	74
391	Thin film compressive stresses due to adatom insertion into grain boundaries. <i>Physical Review Letters</i> , 2007 , 99, 036102	7.4	74
390	Level set simulations of dislocation-particle bypass mechanisms. <i>Acta Materialia</i> , 2004 , 52, 1745-1760	8.4	73
389	Diffusional relaxation of the dislocation-inclusion repulsion. <i>Philosophical Magazine A: Physics of Condensed Matter, Structure, Defects and Mechanical Properties</i> , 1983 , 48, 795-809		73
388	Grain growth in three dimensions: A lattice model. <i>Scripta Metallurgica</i> , 1985 , 19, 225-230		73
387	Onset of plasticity in gold nanopillar compression. <i>Nano Letters</i> , 2007 , 7, 101-7	11.5	71
386	Deformation mechanisms, length scales and optimizing the mechanical properties of nanotwinned metals. <i>Acta Materialia</i> , 2011 , 59, 6890-6900	8.4	70
385	Surface morphology evolution in stressed solids: Surface diffusion controlled crack initiation. <i>Journal of the Mechanics and Physics of Solids</i> , 1994 , 42, 1551-1574	5	69
384	Mechanisms of failure in nanoscale metallic glass. <i>Nano Letters</i> , 2014 , 14, 5858-64	11.5	68
383	Abnormal grain growth induced by sub-boundary-enhanced solid-state wetting: Analysis by phase-field model simulations. <i>Acta Materialia</i> , 2009 , 57, 838-845	8.4	68
382	Computer simulation of recrystallizationIII. Influence of a dispersion of fine particles. <i>Acta Metallurgica Et Materialia</i> , 1992 , 40, 3475-3495		68

381	Strain engineering of 2D semiconductors and graphene: from strain fields to band-structure tuning and photonic applications. <i>Light: Science and Applications</i> , 2020 , 9, 190	16.7	68
380	A Monte Carlo-finite element model for strain energy controlled microstructural evolution: Bafting[In superalloys. <i>Acta Metallurgica</i> , 1989 , 37, 641-650		67
379	Topological framework for local structure analysis in condensed matter. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, E5769-76	11.5	66
378	Structure and evolution of quenched Ising clusters. <i>Physical Review B</i> , 1984 , 30, 5150-5155	3.3	66
377	Microstructural evolution in two-dimensional two-phase polycrystals. <i>Acta Metallurgica Et Materialia</i> , 1993 , 41, 1119-1136		65
376	Etching effects during the chemical vapor deposition of (100) diamond. <i>Journal of Chemical Physics</i> , 1999 , 111, 4291-4299	3.9	64
375	Impurity effects on grain boundary migration. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2002 , 10, R79-R109	2	63
374	Shadowing effects on the microstructure of obliquely deposited films. <i>Journal of Applied Physics</i> , 2002 , 91, 1963-1972	2.5	63
373	Local structure and topology of a model amorphous metal. <i>Journal of Physics F: Metal Physics</i> , 1981 , 11, 2209-2219		62
372	Phase field approach for simulating solid-state dewetting problems. <i>Acta Materialia</i> , 2012 , 60, 5578-55	98.4	61
371	Interatomic potential for vanadium suitable for radiation damage simulations. <i>Journal of Applied Physics</i> , 2003 , 93, 3328-3335	2.5	61
370	Grain boundary migration: misorientation dependence. <i>Current Opinion in Solid State and Materials Science</i> , 2001 , 5, 9-14	12	61
369	Nanostructure and surface effects on yield in Cu nanowires. <i>Acta Materialia</i> , 2013 , 61, 1831-1842	8.4	60
368	A more accurate three-dimensional grain growth algorithm. <i>Acta Materialia</i> , 2011 , 59, 6837-6847	8.4	60
367	Dislocation climb effects on particle bypass mechanisms. <i>Philosophical Magazine</i> , 2006 , 86, 3937-3957	1.6	60
366	Abnormal grain growth in three dimensions. <i>Scripta Metallurgica Et Materialia</i> , 1990 , 24, 661-665		59
365	The inverse hallpetch relation in nanocrystalline metals: A discrete dislocation dynamics analysis. Journal of the Mechanics and Physics of Solids, 2016 , 88, 252-266	5	58
364	Dislocation motion in the presence of diffusing solutes: a computer simulation study. <i>Acta Materialia</i> , 2000 , 48, 2163-2175	8.4	58

363	Phase separation during film growth. Journal of Applied Physics, 1992, 72, 442-446	2.5	58
362	Molecular dynamics simulation of single asperity contact. <i>Acta Materialia</i> , 2004 , 52, 3983-3996	8.4	57
361	Thermal conductivity of crystalline quartz from classical simulations. <i>Physical Review B</i> , 2004 , 70,	3.3	57
360	Adhesion in NiAl-Cr from first principles. <i>Physical Review B</i> , 1996 , 53, 13883-13890	3.3	56
359	Kinetics of Domain Growth: Universality of Kinetic Exponents. <i>Physical Review Letters</i> , 1984 , 52, 1321-1	3 ≱ .4₄	56
358	Effect of strain on the stacking fault energy of copper: A first-principles study. <i>Physical Review B</i> , 2013 , 88,	3.3	55
357	Large-scale molecular dynamics simulations of wear in diamond-like carbon at the nanoscale. <i>Applied Physics Letters</i> , 2013 , 103, 073118	3.4	55
356	Computer Simulation of Final-Stage Sintering: II, Influence of Initial Pore Size. <i>Journal of the American Ceramic Society</i> , 1990 , 73, 2865-2872	3.8	54
355	Characterization of atomic motion governing grain boundary migration. <i>Physical Review B</i> , 2006 , 74,	3.3	53
354	Mechanism of texture development in ion-beam-assisted deposition. <i>Applied Physics Letters</i> , 1999 , 75, 584-586	3.4	53
353	Mobility of low-angle grain boundaries in pure metals. <i>Philosophical Magazine</i> , 2010 , 90, 3107-3128	1.6	52
352	Adhesion effects in material transfer in mechanical contacts. <i>Acta Materialia</i> , 2006 , 54, 5305-5312	8.4	52
351	Effects of boundary inclination and boundary type on shear-driven grain boundary migration. <i>Philosophical Magazine</i> , 2008 , 88, 243-256	1.6	51
350	Grain boundary energy and grain growth in Al films: Comparison of experiments and simulations. <i>Scripta Materialia</i> , 2006 , 54, 1059-1063	5.6	51
349	Effect of material properties on liquid metal embrittlement in the Al G a system. <i>Acta Materialia</i> , 2009 , 57, 1546-1553	8.4	50
348	Stress distributions in growing oxide films. <i>Acta Materialia</i> , 2003 , 51, 2171-2190	8.4	50
347	The mechanism of texture formation during film growth: The roles of preferential sputtering and shadowing. <i>Applied Physics Letters</i> , 1996 , 69, 3007-3009	3.4	50
346	A new method for the simulation of alloys: Application to interfacial segregation. <i>Acta Metallurgica Et Materialia</i> , 1991 , 39, 3071-3082		50

345	Elastic field of a surface step: Atomistic simulations and anisotropic elastic theory. <i>Physical Review B</i> , 1996 , 53, 11120-11127	3.3	49
344	Clock-model description of incommensurate ferroelectric films and of nematic-liquid-crystal films. <i>Physical Review B</i> , 1986 , 34, 1815-1819	3.3	49
343	Edge dislocation-circular inclusion interactions at elevated temperatures. <i>Acta Metallurgica</i> , 1983 , 31, 2151-2159		49
342	Influence of flexoelectric coupling on domain patterns in ferroelectrics. <i>Physical Review B</i> , 2014 , 89,	3.3	48
341	Microstructure versus flaw: mechanisms of failure and strength in nanostructures. <i>Nano Letters</i> , 2013 , 13, 5703-9	11.5	48
340	Extended ensemble molecular dynamics method for constant strain rate uniaxial deformation of polymer systems. <i>Journal of Chemical Physics</i> , 1997 , 107, 4396-4407	3.9	48
339	Simulation of the interaction between Fe impurities and point defects in V. <i>Physical Review B</i> , 2007 , 76,	3.3	48
338	Monte Carlo simulation of phase separation during thin-film codeposition. <i>Journal of Applied Physics</i> , 1993 , 74, 1707-1715	2.5	47
337	Elastic step interactions on vicinal surfaces of fcc metals. Surface Science, 1994, 317, 221-234	1.8	47
336	Polycrystal deformation in a discrete dislocation dynamics framework. <i>Acta Materialia</i> , 2014 , 75, 92-10.	5 8.4	46
335	Origins of growth stresses in amorphous semiconductor thin films. <i>Physical Review Letters</i> , 2003 , 91, 096101	7.4	46
334	The effect of randomness on the strength of high-entropy alloys. <i>Acta Materialia</i> , 2019 , 166, 424-434	8.4	46
333	Stress and morphology evolution during island growth. <i>Physical Review Letters</i> , 2006 , 96, 186103	7.4	45
332	Metal / ceramic adhesion: a first principles study of MgO/Al and MgO/Ag. <i>Journal of Adhesion Science and Technology</i> , 1994 , 8, 837-851	2	45
331	Effects of diffusing impurities on domain growth in the Ising model. <i>Physical Review B</i> , 1987 , 35, 6902-6	59319	45
330	Structure and energy of (111) low-angle twist boundaries in Al, Cu and Ni. <i>Acta Materialia</i> , 2013 , 61, 13	278.1433	7 ₄₄
329	Twinning in thin films Elastic analysis. <i>Acta Materialia</i> , 1996 , 44, 4085-4096	8.4	44
328	Local stress calculation in simulations of multicomponent systems. <i>Journal of Computational Physics</i> , 2009 , 228, 8467-8479	4.1	43

327	Comparison of molecular dynamics simulation methods for the study of grain boundary migration. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2013 , 21, 045017	2	42	
326	Atomic-scale simulations of chemical vapor deposition on flat and vicinal diamond substrates. Journal of Crystal Growth, 1998, 194, 353-368	1.6	42	
325	A regular solution model for impurity drag on a migrating grain boundary. <i>Acta Materialia</i> , 2001 , 49, 589	9-55.247	42	
324	Impurity effects on adhesion: Nb, C, O, B, and S at a Mo/MoSi2 interface. <i>Physical Review B</i> , 1993 , 47, 13615-13625	3.3	42	
323	Machine learning determination of atomic dynamics at grain boundaries. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018 , 115, 10943-10947	11.5	42	
322	Complete topology of cells, grains, and bubbles in three-dimensional microstructures. <i>Physical Review Letters</i> , 2012 , 109, 095505	7.4	41	
321	MICROSTRUCTURAL STABILITY OF STRESSED LAMELLAR AND FIBER COMPOSITES. <i>Acta Materialia</i> , 1997 , 45, 2715-2733	8.4	41	
320	First-principles study of wurtzite InN (0001) and (0001) surfaces. <i>Physical Review B</i> , 2006 , 74,	3.3	41	
319	Simulation and analysis of the migration mechanism of B tilt grain boundaries in an fcc metal. <i>Acta Materialia</i> , 2006 , 54, 623-633	8.4	41	
318	Thermodynamics of solid and liquid embedded-atom-method metals: A variational study. <i>Journal of Chemical Physics</i> , 1991 , 94, 5090-5097	3.9	41	
317	Dislocation generation in the two-dimensional Frenkel-Kontorova model at high stresses. <i>Physical Review Letters</i> , 1986 , 57, 2702-2705	7.4	41	
316	Misfit effects in adhesion calculations. <i>Modelling and Simulation in Materials Science and Engineering</i> , 1998 , 6, 153-164	2	40	
315	Phase separation during co-deposition of Allie thin films. <i>Journal of Materials Research</i> , 1992 , 7, 653-66	6 2.5	40	
314	Impurity effects on domain-growth kinetics. II. Potts model. <i>Physical Review B</i> , 1985 , 32, 3021-3025	3.3	40	
313	Geometric and topological properties of the canonical grain-growth microstructure. <i>Physical Review E</i> , 2015 , 92, 063308	2.4	39	
312	Morphologies of diamond films from atomic-scale simulations of chemical vapor deposition. <i>Diamond and Related Materials</i> , 1997 , 6, 1198-1206	3.5	39	
311	Morphology evolution during the growth of strained-layer superlattices. <i>Physical Review B</i> , 2000 , 62, 8397-8409	3.3	39	
310	First-principles study of the 🖽 2O3 (0001)/Cu(111) interface. <i>Journal of Materials Science</i> , 1996 , 3, 289		39	

309	Three-dimensional formulation of dislocation climb. <i>Journal of the Mechanics and Physics of Solids</i> , 2015 , 83, 319-337	5	38
308	Manipulating ferroelectric domains in nanostructures under electron beams. <i>Physical Review Letters</i> , 2013 , 111, 165702	7.4	38
307	Microstructural Mechanics Model of Anisotropic-Thermal-Expansion-Induced Microcracking. <i>Journal of the American Ceramic Society</i> , 1994 , 77, 1123-1138	3.8	38
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