Miroslav Cerny

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	The Impact of Vibrational Entropy on the Segregation of Cu to Antiphase Boundaries in Fe3Al. Magnetochemistry, 2021, 7, 108.	2.4	3
2	A novel multiscale approach to brittle fracture of nano/microâ€sized components. Fatigue and Fracture of Engineering Materials and Structures, 2020, 43, 1630-1645.	3.4	9
3	Impact of Antiphase Boundaries on Structural, Magnetic and Vibrational Properties of Fe3Al. Materials, 2020, 13, 4884.	2.9	6
4	Extraordinary Response of H-Charged and H-Free Coherent Grain Boundaries in Nickel to Multiaxial Loading. Crystals, 2020, 10, 590.	2.2	4
5	On the effect of supercell size and strain localization in computational tensile tests. Modelling and Simulation in Materials Science and Engineering, 2020, 28, 065011.	2.0	4
6	Ab initio study of the theoretical strength and magnetism of the Feâ^'Pd, Feâ^'Pt and Feâ^'Cu nanocomposites. Journal of Magnetism and Magnetic Materials, 2019, 469, 100-107.	2.3	3
7	Ab initio aided strain gradient elasticity theory in prediction of nanocomponent fracture. Mechanics of Materials, 2019, 136, 103074.	3.2	9
8	Correlating structural and mechanical properties of AlN/TiN superlattice films. Scripta Materialia, 2019, 165, 159-163.	5.2	29
9	Atomistic approaches to cleavage of interfaces. Modelling and Simulation in Materials Science and Engineering, 2019, 27, 035007.	2.0	24
10	Strength of FePd/MgO and FePt/MgO interfaces from first principles. Modelling and Simulation in Materials Science and Engineering, 2018, 26, 035002.	2.0	3
11	Energetics of NiTi allotropes under uniaxial compression. Scripta Materialia, 2017, 133, 92-95.	5.2	4
12	The origin of lattice instability in bcc tungsten under triaxial loading. Philosophical Magazine, 2017, 97, 2971-2984.	1.6	9
13	Interface-induced electronic structure toughening of nitride superlattices. Surface and Coatings Technology, 2017, 325, 410-416.	4.8	34
14	Stress Coupling Effect on Ideal Shear Strength: Tungsten as a Case Study. Advances in Materials Science and Engineering, 2016, 2016, 1-5.	1.8	1
15	Ab initio tensile tests of grain boundaries in the fcc crystals of Ni and Co with segregated sp-impurities. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 2016, 669, 218-225.	5.6	29
16	Multiaxial stress–strain response and displacive transformations in NiTi alloy from first principles. Acta Materialia, 2016, 109, 223-229.	7.9	6
17	Mechanical stability of Ni and Ir under hydrostatic and uniaxial loading. Modelling and Simulation in Materials Science and Engineering, 2015, 23, 055010.	2.0	14
18	Ab initio calculations of mechanical properties: Methods and applications. Progress in Materials Science, 2015, 73, 127-158.	32.8	114

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#	Article	IF	CITATIONS
19	Modeling Load-displacement Curve and Pop-in Effect in Nanoindentation Tests. , 2014, 3, 1111-1116.		2
20	On the effect of deformation twins on stability of B19′ structure in NiTi martensite. Computational Materials Science, 2014, 87, 107-111.	3.0	11
21	Stability and strength of covalent crystals under uniaxial and triaxial loading from first principles. Journal of Physics Condensed Matter, 2013, 25, 035401.	1.8	11
22	Shear instabilities in perfect bcc crystals during simulated tensile tests. Physical Review B, 2013, 87, .	3.2	17
23	Dynamic stability of fcc crystals under isotropic loading from first principles. Journal of Physics Condensed Matter, 2012, 24, 215403.	1.8	13
24	Strength of bcc crystals under combined shear and axial loading from first principles. Computational Materials Science, 2012, 55, 337-343.	3.0	12
25	The theoretical strength of fcc crystals under multiaxial loading. Computational Materials Science, 2011, 50, 2257-2261.	3.0	10
26	Can twinning stabilize B19′ structure in NiTi martensite?. Intermetallics, 2011, 19, 1567-1572.	3.9	21
27	The theoretical shear strength of fcc crystals under superimposed triaxial stress. Acta Materialia, 2010, 58, 3117-3123.	7.9	19
28	Ideal tensile strength of cubic crystals under superimposed transverse biaxial stresses from first principles. Physical Review B, 2010, 82, .	3.2	51
29	Influence of superimposed normal stress on shear strength of perfect bcc crystals. Computational Materials Science, 2010, 47, 907-910.	3.0	14
30	The theoretical tensile strength of fcc crystals predicted from shear strength calculations. Journal of Physics Condensed Matter, 2009, 21, 145406.	1.8	29
31	Influence of comp ound twinning on Young's moduli in NiTi martensite. , 2009, , .		2
32	Tensile Strength of Perfect Cubic Crystals under Superimposed Transverse Plain Stress. Materials Science Forum, 2008, 567-568, 73-76.	0.3	6
33	Elastic properties of B19' structure of NiTi alloy under uniaxial and hydrostatic loading from first principles. Strength of Materials, 2008, 40, 12-15.	0.5	9
34	Multiscale modelling of nanoindentation test in copper crystal. Engineering Fracture Mechanics, 2008, 75, 3755-3762.	4.3	9
35	Influence of normal stress on theoretical shear strength of fcc metals. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 2008, 483-484, 692-694.	5.6	25
36	Influence of normal stress on shear strength along ã€^111〉{110} and ã€^111〉{112} deformation paths ir alloy from first principles. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 2008, 481-482, 247-249.	1 B2 NiTi 5.6	7

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#	ARTICLE	IF	CITATIONS
37	xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si17.gif" overflow="scroll"> <mml:mrow><mml:mo stretchy="false">ã€^<mml:mn>1</mml:mn><mml:mspace <br="" width="0.12em">/><mml:mn>1</mml:mn>122</mml:mspace></mml:mo </mml:mrow>	3.0	14
38	stretchy="lalse">ā6‰ <mml:mo stretchy="lalse">{</mml:mo> <mml:mo><mml:mspace Effect of normal stress on the ideal shear strength in covalent crystals. Physical Review B, 2008, 77, .</mml:mspace </mml:mo>	3.2	52
39	Onset of Microplasticity in Copper Crystal during Nanoindentation. Key Engineering Materials, 2007, 348-349, 801-804.	0.4	2
40	Influence of superimposed biaxial stress on the tensile strength of perfect crystals from first principles. Physical Review B, 2007, 76, .	3.2	46
41	Elastic stability of magnetic crystals under isotropic compression and tension. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 2007, 462, 432-435.	5.6	16
42	Higher-energy Structures and Stability of Cu and Al Crystals Along Displacive Transformation Paths. Journal of Computer-Aided Materials Design, 2006, 12, 161-173.	0.7	8
43	Elasticity and Stability of Fe-P Ordered Systems from First Principles. Materials Science Forum, 2005, 482, 135-138.	0.3	8
44	Theoretical Strength of Metals and Intermetallics from First Principles. Materials Science Forum, 2005, 482, 33-38.	0.3	10
45	Calculations of theoretical strength: State of the art and history. Journal of Computer-Aided Materials Design, 2004, 11, 1-28.	0.7	101
46	Ab initio analysis of theoretical isotropic strength and elasticity of nickel aluminide compounds. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 2004, 387-389, 923-925.	5.6	11
47	Ab initiocalculations of ideal tensile strength and mechanical stability in copper. Journal of Physics Condensed Matter, 2004, 16, 1045-1052.	1.8	61
48	Stability of fcc crystals under hydrostatic loading. Journal of Alloys and Compounds, 2004, 378, 159-162.	5.5	15
49	Ab initiocalculations of elastic and magnetic properties of Fe, Co, Ni, and Cr crystals under isotropic deformation. Physical Review B, 2003, 67, .	3.2	74
50	Ab initio calcuation of ideal strength for cubic crystals under three-axial tension. European Physical Journal D, 1999, 49, 1495-1501.	0.4	20
51	Stabilizing Effect of (100) Compound Twinning in NiTi Martensite. Key Engineering Materials, 0, 465, 81-84.	0.4	1
52	First Principles Study of Ideal Composites Reinforced by Coherent Nano-Fibres. Key Engineering Materials, 0, 465, 73-76.	0.4	3
53	The [100] Compressive Strength of Perfect Cubic Crystals under Superimposed Biaxial Stresses. Key Engineering Materials, 0, 465, 183-186.	0.4	3
54	Dynamic Stability of Ni FCC Crystal under Isotropic Tension. Key Engineering Materials, 0, 592-593, 47-50.	0.4	0