

Josef Kudrnovsky

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271
papers

8,143
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45
h-index

82
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272
ext. papers

8,586
ext. citations

3.1
avg. IF

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L-index

#	Paper	IF	Citations
271	First-principles theory of dilute magnetic semiconductors. <i>Reviews of Modern Physics</i> , 2010 , 82, 1633-1690	10.5	855
270	Room-temperature antiferromagnetic memory resistor. <i>Nature Materials</i> , 2014 , 13, 367-74	27	435
269	Ab initio calculations of exchange interactions, spin-wave stiffness constants, and Curie temperatures of Fe, Co, and Ni. <i>Physical Review B</i> , 2001 , 64,	3.3	392
268	Electronic Structure of Disordered Alloys, Surfaces and Interfaces 1997 ,		329
267	Magnetic percolation in diluted magnetic semiconductors. <i>Physical Review Letters</i> , 2004 , 93, 137202	7.4	250
266	Exchange interactions in III-V and group-IV diluted magnetic semiconductors. <i>Physical Review B</i> , 2004 , 69,	3.3	248
265	Electronic structure of random alloys by the linear band-structure methods. <i>Physical Review B</i> , 1990 , 41, 7515-7528	3.3	165
264	Electronic and nuclear chemical reactivity. <i>Journal of Chemical Physics</i> , 1994 , 101, 8988-8997	3.9	150
263	Magnetism without magnetic impurities in ZrO ₂ oxide. <i>Applied Physics Letters</i> , 2008 , 92, 212503	3.4	124
262	Effects of resonant interface states on tunneling magnetoresistance. <i>Physical Review B</i> , 2002 , 65,	3.3	121
261	Electronic structures and magnetic moments of Fe _{3+y} Si _{1-y} and Fe _{3-x} V _x Si alloys with DO ₃ -derived structure. <i>Physical Review B</i> , 1991 , 43, 5924-5933	3.3	116
260	Exchange interactions and Curie temperatures in Ni ₂ MnSb alloys: First-principles study. <i>Physical Review B</i> , 2006 , 73,	3.3	103
259	Exchange interactions, spin waves, and transition temperatures in itinerant magnets. <i>Philosophical Magazine</i> , 2006 , 86, 1713-1752	1.6	101
258	Reactivity kernels, the normal modes of chemical reactivity, and the hardness and softness spectra. <i>Journal of Chemical Physics</i> , 1995 , 103, 3543-3551	3.9	99
257	Calculating the Curie temperature reliably in diluted III-V ferromagnetic semiconductors. <i>Europhysics Letters</i> , 2005 , 69, 812-818	1.6	98
256	Interface resistance of disordered magnetic multilayers. <i>Physical Review B</i> , 2001 , 63,	3.3	96
255	Oscillatory curie temperature of two-dimensional ferromagnets. <i>Physical Review Letters</i> , 2000 , 85, 5424-7.4	7.4	94

254	On-site Coulomb interaction and the magnetism of (GaMn)N and (GaMn)As. <i>Physical Review B</i> , 2004 , 69,	3.3	91
253	Canonical description of electron states in random alloys. <i>Physical Review B</i> , 1987 , 35, 2487-2489	3.3	91
252	Disordered magnetic multilayers: Electron transport within the coherent potential approximation. <i>Physical Review B</i> , 2006 , 73,	3.3	90
251	Ab initio theory of perpendicular magnetotransport in metallic multilayers. <i>Physical Review B</i> , 2000 , 62, 15084-15095	3.3	85
250	Mn-stabilized zirconia: from imitation diamonds to a new potential high-Tc ferromagnetic spintronics material. <i>Physical Review Letters</i> , 2007 , 98, 016101	7.4	84
249	Interatomic electron transport by semiempirical and ab initio tight-binding approaches. <i>Physical Review B</i> , 2002 , 65,	3.3	83
248	Itinerant magnetism of disordered Fe-Co and Ni-Cu alloys in two and three dimensions. <i>Physical Review B</i> , 1994 , 49, 3352-3362	3.3	77
247	Interlayer magnetic coupling: Effect of interface roughness. <i>Physical Review B</i> , 1996 , 53, 5125-5128	3.3	74
246	Dilute moment n-type ferromagnetic semiconductor Li(Zn,Mn)As. <i>Physical Review Letters</i> , 2007 , 98, 067202	7.4	69
245	Lattice constant in diluted magnetic semiconductors (Ga,Mn)As. <i>Physical Review B</i> , 2003 , 67,	3.3	68
244	On the calculation of the surface Green function by the tight-binding linear muffin-tin orbital method. <i>Journal of Physics Condensed Matter</i> , 1989 , 1, 9893-9897	1.8	67
243	Ab initio theory of exchange interactions and the Curie temperature of bulk Gd. <i>Journal of Physics Condensed Matter</i> , 2003 , 15, 2771-2782	1.8	66
242	Magnetic impurities and materials design for semiconductor spintronics. <i>Physica B: Condensed Matter</i> , 2003 , 340-342, 863-869	2.8	66
241	Ferromagnetism in diluted magnetic semiconductors: A comparison between ab initio mean-field, RPA, and Monte Carlo treatments. <i>Physical Review B</i> , 2003 , 68,	3.3	65
240	Coulomb correlation effects on the electronic structure of III-V diluted magnetic semiconductors. <i>Physical Review B</i> , 2004 , 69,	3.3	62
239	Magnetic properties and disorder effects in diluted magnetic semiconductors. <i>Physical Review B</i> , 2005 , 72,	3.3	55
238	Application of ab initio and CALPHAD thermodynamics to Mo-Ta-W alloys. <i>Physical Review B</i> , 2005 , 71,	3.3	53
237	Potential, core-level, and d band shifts at transition-metal surfaces. <i>Physical Review B</i> , 1996 , 54, 8892-8898	3.3	52

236	Interlayer exchange coupling: The effect of substitutional disorder. <i>Physical Review Letters</i> , 1996 , 76, 4254-4257	7.4	52
235	First-principles study of surface segregation in Cu-Ni alloys. <i>Physical Review B</i> , 1993 , 48, 2704-2710	3.3	52
234	Orbital symmetry, reactivity, and transition metal surface chemistry. <i>Physical Review Letters</i> , 1994 , 72, 3222-3225	7.4	52
233	Magnetic anisotropy energy of disordered tetragonal Fe-Co systems from ab initio alloy theory. <i>Physical Review B</i> , 2012 , 86,	3.3	51
232	Self-consistent Green's-function method for random overlayers. <i>Physical Review B</i> , 1992 , 46, 4222-4228	3.3	51
231	Ab initio theory of galvanomagnetic phenomena in ferromagnetic metals and disordered alloys. <i>Physical Review B</i> , 2012 , 86,	3.3	49
230	Effective interatomic interactions in inhomogeneous semi-infinite systems. <i>Physical Review B</i> , 1992 , 45, 14328-14334	3.3	49
229	Compensation, interstitial defects, and ferromagnetism in diluted ferromagnetic semiconductors. <i>Physical Review B</i> , 2005 , 72,	3.3	48
228	Theory of Oscillatory Exchange Coupling in Fe/(V,Cr) and Fe/(Cr,Mn). <i>Physical Review Letters</i> , 1995 , 74, 4063-4066	7.4	47
227	Origins of surface alloy formation: Cu(001)c(2 x 2)-Pd as a case study. <i>Physical Review Letters</i> , 1992 , 69, 308-311	7.4	46
226	Physical properties of FeRh alloys: The antiferromagnetic to ferromagnetic transition. <i>Physical Review B</i> , 2015 , 91,	3.3	45
225	Electronic, magnetic, and transport properties and magnetic phase transition in quaternary (Cu,Ni)MnSb Heusler alloys. <i>Physical Review B</i> , 2008 , 78,	3.3	45
224	Self-consistent Green's-function method for surfaces of random alloys. <i>Physical Review B</i> , 1993 , 47, 16525-16534	3.3	44
223	Ferromagnetism of imperfect ultrathin Ru and Rh films on a Ag(001) substrate. <i>Physical Review Letters</i> , 1995 , 74, 2551-2554	7.4	43
222	Exchange coupling in transition-metal ferromagnets. <i>Physical Review B</i> , 2000 , 62, 5293-5296	3.3	42
221	Magnetic properties of fcc Ni-based transition metal alloys. <i>Physical Review B</i> , 2008 , 77,	3.3	41
220	The Auger spectra of metals: effect of electron correlations in partially filled narrow bands. <i>Journal of Physics F: Metal Physics</i> , 1984 , 14, 2443-2453		41
219	Anomalous Hall effect in stoichiometric Heusler alloys with native disorder: A first-principles study. <i>Physical Review B</i> , 2013 , 88,	3.3	40

218	Spin-disorder resistivity of ferromagnetic metals from first principles: The disordered-local-moment approach. <i>Physical Review B</i> , 2012 , 86,	3-3	39
217	First-principles theoretical studies of half-metallic ferromagnetism in CrTe. <i>Physical Review B</i> , 2010 , 82,	3-3	39
216	Mn-doped Ga(As,P) and (Al,Ga)As ferromagnetic semiconductors: Electronic structure calculations. <i>Physical Review B</i> , 2007 , 75,	3-3	39
215	Magnetic coupling of interfaces: A surface-Green's-function approach. <i>Physical Review B</i> , 1994 , 50, 16105-16108	3-3	39
214	First-principles study of the electronic structure and exchange interactions in bcc europium. <i>Physical Review B</i> , 2003 , 68,	3-3	38
213	Comparative study of the electronic structure of ordered, partially ordered, and disordered phases of the Cu ₃ Au alloy. <i>Physical Review B</i> , 1991 , 43, 4613-4621	3-3	37
212	Approximate treatment of charge selfconsistency and lattice relaxations in random transition metal alloys: Application to CuPd system. <i>Solid State Communications</i> , 1989 , 70, 577-580	1.6	37
211	Can correlated substitution enhance the Curie temperature in diluted magnetic semiconductors?. <i>Applied Physics Letters</i> , 2004 , 85, 4941-4943	3-4	36
210	First-principles study of stability and local order in substitutional Ta-W alloys. <i>Physical Review B</i> , 2001 , 64,	3-3	36
209	Electronic structure of semiinfinite crystals with substitutional disorder in surface layer. <i>Surface Science</i> , 1977 , 64, 411-424	1.8	34
208	Microscopic analysis of the valence band and impurity band theories of (Ga,Mn)As. <i>Physical Review Letters</i> , 2010 , 105, 227202	7-4	33
207	Electronic structure and magnetic properties of random alloys: Fully relativistic spin-polarized linear muffin-tin-orbital method. <i>Physical Review B</i> , 1996 , 54, 1610-1621	3-3	33
206	Substrate-induced antiferromagnetism of a Fe monolayer on the Ir(001) surface. <i>Physical Review B</i> , 2009 , 80,	3-3	32
205	Theoretical study of ordering in Fe-Al alloys based on a density-functional generalized-perturbation method. <i>Physical Review B</i> , 1997 , 55, 8184-8193	3-3	32
204	Curie temperatures of fcc and bcc nickel and permalloy: Supercell and Green's function methods. <i>Physical Review B</i> , 2008 , 77,	3-3	32
203	Interlayer magnetic coupling: The torque method. <i>Physical Review B</i> , 1996 , 53, 15036-15044	3-3	32
202	Electronic properties of surfaces of disordered alloys. <i>Physical Review B</i> , 1991 , 44, 6410-6415	3-3	32
201	Fermi sea term in the relativistic linear muffin-tin-orbital transport theory for random alloys. <i>Physical Review B</i> , 2014 , 89,	3-3	30

200	Relativistic electronic structure of random alloys and their surfaces by linear band-structure methods. <i>Physical Review B</i> , 1994 , 50, 7903-7914	3.3	30
199	Stability and ordering properties of fcc alloys based on Rh, Ir, Pd, and Pt. <i>Physical Review B</i> , 2006 , 74,	3.3	29
198	Magnetic nature of (100) surfaces of bcc RuV, RhV, and PdV binary alloys. <i>Physical Review B</i> , 1998 , 57, R11065-R11068	3.3	29
197	Interlayer magnetic coupling: Effect of alloying in the spacer. <i>Physical Review B</i> , 1996 , 54, R3738-R3741	3.3	29
196	Adlayer Core-Level Shifts of Random Metal Overlayers on Transition-Metal Substrates. <i>Physical Review Letters</i> , 1997 , 78, 1807-1810	7.4	28
195	Disorder effects in diluted ferromagnetic semiconductors. <i>Physical Review B</i> , 2003 , 68,	3.3	26
194	Electronic structure of fluorite-type compounds and mixed crystals. <i>Physical Review B</i> , 1991 , 43, 12597-12606	3.3	26
193	Magnetism of mixed quaternary Heusler alloys: (Ni,T) ₂ MnSn (T=Cu,Pd) as a case study. <i>Physical Review B</i> , 2010 , 82,	3.3	25
192	Exchange interactions and Curie temperatures in Cr-based alloys in the zinc blende structure: Volume- and composition-dependence from first-principles calculations. <i>Physical Review B</i> , 2010 , 81,	3.3	25
191	Theory of surface segregation in metallic alloys: The generalized perturbation method. <i>Computational Materials Science</i> , 1999 , 15, 144-168	3.2	25
190	Magnetism-induced ordering in two and three dimensions. <i>Physical Review B</i> , 1994 , 50, 9603-9606	3.3	25
189	Electronic structure of substitutionally disordered metal alloys: Single-site approximation for canonical d bands. <i>Physical Review B</i> , 1985 , 31, 6424-6434	3.3	25
188	Nonlocal torque operators in ab initio theory of the Gilbert damping in random ferromagnetic alloys. <i>Physical Review B</i> , 2015 , 92,	3.3	24
187	Pressure dependence of Curie temperature and resistivity in complex Heusler alloys. <i>Physical Review B</i> , 2011 , 84,	3.3	24
186	Relativistic LMTO method for systems of light elements. <i>Philosophical Magazine</i> , 2008 , 88, 2787-2798	1.6	24
185	Ab Initio Study of Curie Temperatures of Diluted Magnetic Semiconductors. <i>Journal of Superconductivity and Novel Magnetism</i> , 2003 , 16, 119-122		24
184	Dynamical correlations in multiorbital Hubbard models: fluctuation exchange approximations. <i>Journal of Physics Condensed Matter</i> , 2005 , 17, 61-74	1.8	24
183	Electron transport in magnetic multilayers: Effect of disorder. <i>Physical Review B</i> , 2002 , 65,	3.3	24

182	Interlayer exchange coupling: Effect of the cap. <i>Physical Review B</i> , 1997 , 56, 8919-8927	3.3	23
181	Importance of complex band structure and resonant states for tunneling. <i>Journal of Magnetism and Magnetic Materials</i> , 2002 , 240, 108-113	2.8	23
180	Electronic structure of random semiconductor alloys by the tight-binding linear muffin-tin orbital method. <i>Physical Review B</i> , 1989 , 40, 10029-10032	3.3	23
179	Effect of disorder on the electronic structure of palladium. <i>Physical Review B</i> , 1990 , 41, 7988-7998	3.3	22
178	Magnetism without magnetic impurities in oxides ZrO ₂ and TiO ₂ . <i>Philosophical Magazine</i> , 2008 , 88, 2755-2764	2.7	21
177	Half-metallicity and magnetism of GeTe doped with transition metals V, Cr, and Mn: A theoretical study from the viewpoint of application in spintronics. <i>Journal of Applied Physics</i> , 2012 , 112, 053902	2.5	20
176	Electronic structure of ordered and disordered CuxPd1-x alloys via the linear-muffin-tin-orbitals method. <i>Physical Review B</i> , 1992 , 45, 8272-8282	3.3	20
175	Transport properties of iron at Earth's core conditions: The effect of spin disorder. <i>Physical Review B</i> , 2017 , 96,	3.3	19
174	Temperature dependence of the interlayer exchange coupling in magnetic multilayers: An ab initio approach. <i>Physical Review B</i> , 1999 , 60, 9588-9595	3.3	19
173	Ab initio theory of surface segregation: Self-consistent determination of the concentration profile. <i>Physical Review B</i> , 1996 , 54, 8202-8212	3.3	19
172	Amplitude and Phase of the Oscillatory Exchange Coupling between Fe-Co-Ni Alloy Layers across a Cu Spacer Layer. <i>Physical Review Letters</i> , 1997 , 78, 358-361	7.4	18
171	Residual resistivity of diluted III-V magnetic semiconductors. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5607-S5614	1.8	18
170	Frustration and long-range behavior of the exchange interactions in AuFe spin-glass alloys. <i>Physical Review B</i> , 2004 , 70,	3.3	18
169	Interface reflectivities and quantum-well states in magnetic multilayers. <i>Physical Review B</i> , 1998 , 58, 13721-13733	3.3	18
168	Ab initio theory of exchange interactions in itinerant magnets. <i>Physica Status Solidi (B): Basic Research</i> , 2003 , 236, 318-324	1.3	17
167	Theory of the Auger Spectra of Narrow-Band Metals with Impurities. <i>Physica Status Solidi (B): Basic Research</i> , 1981 , 108, 683-692	1.3	17
166	Electronic structure and magnetism of diluted magnetic semiconductors. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5481-S5489	1.8	16
165	Magnetoresistance in domain walls: effect of randomness. <i>Surface Science</i> , 2001 , 482-485, 1107-1112	1.8	16

164	Electronic structure of disordered overlayers on metal substrates. <i>Physical Review B</i> , 1991 , 44, 4068-4073,	3.3	16
163	Interlayer exchange coupling through ordered and disordered alloy spacers. <i>Journal of Magnetism and Magnetic Materials</i> , 1997 , 165, 128-133	2.8	15
162	Electronic structure of ordered and disordered Pd ₃ Fe. <i>Journal of Magnetism and Magnetic Materials</i> , 1990 , 87, 97-105	2.8	15
161	Effects of atomic and magnetic order on electronic transport in Pd-rich Pd-Fe alloys. <i>Physical Review B</i> , 2011 , 84,	3.3	14
160	Dynamical electron correlations in weakly interacting systems: TB-LMTO approach to metals and random alloys. <i>Physical Review B</i> , 1999 , 60, 15664-15673	3.3	14
159	Unified approach to electronic, thermodynamical, and transport properties of Fe ₃ Si and Fe ₃ Al alloys. <i>Physical Review B</i> , 2014 , 90,	3.3	13
158	Physical properties of the tetragonal CuMnAs: A first-principles study. <i>Physical Review B</i> , 2017 , 96,	3.3	13
157	Noncollinear magnetic ordering in compressed FePd ₃ ordered alloy: A first principles study. <i>Physical Review B</i> , 2012 , 86,	3.3	13
156	Electronic and transport properties of the Mn-doped topological insulator Bi ₂ Te ₃ : A first-principles study. <i>Physical Review B</i> , 2016 , 93,	3.3	12
155	First-principles study of spin-disorder resistivity of heavy rare-earth metals: Gd _m series. <i>Physical Review B</i> , 2012 , 85,	3.3	12
154	Electronic structure and transport properties of CrAs ₂ /TaAs ₂ /rAs trilayers from first principles theory. <i>Physical Review B</i> , 2004 , 70,	3.3	12
153	On the orientational dependence of giant magnetoresistance. <i>European Physical Journal B</i> , 1999 , 9, 245-250	2.5	12
152	Phase diagram of the Cu-Pd surface alloy: A first-principles calculation. <i>Physical Review B</i> , 1995 , 51, 17910-17915	3.3	12
151	Electronic structure in random alloys: Cooperation of structural and chemical disorders. <i>Solid State Communications</i> , 1986 , 58, 67-70	1.6	12
150	Theory of chemisorption. <i>European Physical Journal D</i> , 1985 , 35, 1017-1032		12
149	Correlated Doping in Semiconductors: the Role of Donors in III-V Diluted Magnetic Semiconductors. <i>Acta Physica Polonica A</i> , 2002 , 102, 673-678	0.6	12
148	Effective magnetic Hamiltonians from first principles. <i>EPJ Web of Conferences</i> , 2013 , 40, 11001	0.3	11
147	Origin of the negative giant magnetoresistance effect in Co _{1-x} Cr _x /Cu/Co (111) trilayers. <i>Physical Review B</i> , 2004 , 69,	3.3	11

146	Ordering and segregation in XPt (X=V, Cu, and Au) random alloys. <i>Physical Review B</i> , 2001 , 64,	3.3	11
145	New type of oscillatory exchange coupling induced by ordering in the magnetic layers. <i>Physical Review Letters</i> , 1996 , 76, 3834-3837	7.4	11
144	Galvanomagnetic properties of partially ordered L10 FePt alloys. <i>Physical Review B</i> , 2014 , 89,	3.3	10
143	Electronic structure of random Ag-Pd and Ag-vacancy overlayers on an fcc Pd(001) substrate. <i>Physical Review B</i> , 1993 , 48, 1870-1876	3.3	10
142	Electronic structure of random non-isocoric transition metal alloys. <i>Solid State Communications</i> , 1988 , 65, 613-616	1.6	10
141	Compositional Dependence of the Formation Energies of Substitutional and Interstitial Mn in Partially Compensated (Ga,Mn)As. <i>Acta Physica Polonica A</i> , 2004 , 105, 637-644	0.6	10
140	Relation of Curie temperature and conductivity: (Ga,Mn)As alloy as a case study. <i>Applied Physics Letters</i> , 2007 , 91, 102509	3.4	9
139	Exchange interactions and critical temperatures in diluted magnetic semiconductors. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5571-S5578	1.8	9
138	Ab-initio calculations of the electronic and atomic structure of solids and their surfaces. <i>Computer Physics Communications</i> , 1996 , 97, 111-123	4.2	9
137	The Electronic Structure of Palladium-Noble Metal Alloys. <i>Physica Status Solidi (B): Basic Research</i> , 1988 , 148, K23-K27	1.3	9
136	Theory of auger spectra from disordered alloys the effect of electron correlations in filled narrow bands. <i>Physica Status Solidi (B): Basic Research</i> , 1982 , 114, 627-635	1.3	9
135	Electrical conductivity of electrons in a model binary disordered alloy with long range order. <i>European Physical Journal D</i> , 1977 , 27, 71-87		9
134	Electronic and transport properties of a new quaternary Heusler alloy CoMnFeSi. <i>Physical Review B</i> , 2018 , 97,	3.3	8
133	Influence of oxygen and hydrogen adsorption on the magnetic structure of an ultrathin iron film on an Ir(001) surface. <i>Physical Review B</i> , 2013 , 88,	3.3	8
132	Magnetic phase stability of monolayers: Fe on a TaxW1-x(001) random alloy as a case study. <i>Physical Review B</i> , 2010 , 81,	3.3	8
131	Pressure effect on magnetic moments in ordered Ni3Mn and disordered Ni100-x Mn x alloys: ab initio calculation and experiment. <i>High Pressure Research</i> , 2011 , 31, 116-120	1.6	8
130	Chemical ordering and composition fluctuations at the (001) surface of the Fe64Ni36 Invar alloy. <i>Physical Review B</i> , 2006 , 74,	3.3	8
129	The CPP transport in metallic magnetic multilayers. <i>Surface Science</i> , 2000 , 454-456, 918-924	1.8	8

128	Reformulation of the Korringa - Kohn - Rostoker coherent potential approximation for the treatment of space-filling cell potentials and charge-transfer effects. <i>Journal of Physics Condensed Matter</i> , 1996 , 8, 7869-7881	1.8	8
127	Calculation of equilibrium lattice parameters and the heat of mixing for the system Au/Pd by the relativistic Korringa-Kohn-Rostoker coherent-potential-approximation method. <i>Physical Review B</i> , 1993 , 48, 7866-7871	3.3	8
126	Disordered Alloys and Their Surfaces: The Coherent Potential Approximation 1999 , 349-378		8
125	Half-metallicity, magnetism and electrical resistivity of Sn _{1-x} Mn _x Te alloys in rock salt and zinc blende structures. <i>Journal of Magnetism and Magnetic Materials</i> , 2015 , 375, 15-25	2.8	7
124	Magnetism, half-metallicity and electrical transport properties of V- and Cr-doped semiconductor SnTe: A theoretical study. <i>Journal of Applied Physics</i> , 2013 , 114, 213704	2.5	7
123	Charge-transfer effects in disordered alloys: the test case of Al - Li alloys. <i>Journal of Physics Condensed Matter</i> , 1996 , 8, 7883-7898	1.8	7
122	Phase stability and ordering in diluted magnetic IIIV semiconductors. <i>Philosophical Magazine</i> , 2004 , 84, 1889-1905	1.6	7
121	Interference, resonances, and bound states at the Pd(001) and Rh(001) surfaces. <i>Physical Review B</i> , 1994 , 50, 11142-11145	3.3	7
120	Electronic theory of surface segregation in transition metal alloys. <i>Surface Science</i> , 1994 , 307-309, 821-825		7
119	Theory of angle-resolved photoemission from alloys: the normal spectra from copper-rich alloys. <i>Journal of Physics F: Metal Physics</i> , 1986 , 16, 943-959		7
118	The Two Band Models in the Theory of Disordered Substitutional Alloys. <i>Physica Status Solidi (B): Basic Research</i> , 1975 , 70, 759-766	1.3	7
117	Temperature-dependent resistivity and anomalous Hall effect in NiMnSb from first principles. <i>Physical Review B</i> , 2019 , 99,	3.3	6
116	Band mapping of the weakly off-stoichiometric Heusler alloy Ni _{49.7} Mn _{29.1} Ga _{21.2} in the austenitic phase. <i>Physical Review B</i> , 2015 , 91,	3.3	6
115	Exchange interactions and correlations in Heusler alloys. <i>Journal of Magnetism and Magnetic Materials</i> , 2007 , 310, 1654-1656	2.8	6
114	Exchange interactions and Curie temperatures in diluted magnetic semiconductor. <i>Journal of Magnetism and Magnetic Materials</i> , 2004 , 272-276, 1983-1984	2.8	6
113	Ab-initio theory of the CPP-magnetoconductance. <i>European Physical Journal D</i> , 1999 , 49, 1583-1589		6
112	Ordering tendencies in Fe-Al alloys in magnetic and non-magnetic models. <i>Materials Science and Engineering B: Solid-State Materials for Advanced Technology</i> , 1996 , 37, 237-241	3.1	6
111	Pressure dependence of electronic densities of states and superconducting transition temperatures in NiZr glasses. <i>Physical Review B</i> , 1991 , 43, 110-118	3.3	6

110	First-principles calculations of electronic structure in random hcp alloys: A Ru-Re example. <i>Physical Review B</i> , 1990 , 41, 10459-10462	3.3	6
109	Electronic structure of sputter-deposited alloy films: application to the Fe-Cu-Ag system. <i>Journal of Physics Condensed Matter</i> , 1990 , 2, 6847-6851	1.8	6
108	Second-order perturbation treatment of correlations in transition metal alloys. <i>Physica Status Solidi (B): Basic Research</i> , 1983 , 116, 119-128	1.3	6
107	Electronic structure of nin-isocoric transition metal alloys: Cu_xRh_{1-x} . <i>Solid State Communications</i> , 1985 , 54, 981-984	1.6	6
106	Correlation Effects on Adatoms: The Self-Consistent T-Matrix Approximation. <i>Physica Status Solidi (B): Basic Research</i> , 1980 , 97, K57-K60	1.3	6
105	Defect-induced magnetic structure of CuMnSb. <i>Physical Review B</i> , 2016 , 94,	3.3	6
104	The disordered local moment approach to the spin-disorder resistivity of metallic ferromagnets. <i>EPJ Web of Conferences</i> , 2013 , 40, 12001	0.3	5
103	Ordering effects in diluted magnetic semiconductors. <i>Phase Transitions</i> , 2007 , 80, 333-350	1.3	5
102	Ab initio theory of transport in FeRh-based natural magnetic multilayers. <i>Journal of Magnetism and Magnetic Materials</i> , 2002 , 240, 162-164	2.8	5
101	Aspects of magnetotunnelling drawn from ab-initio-type calculations. <i>The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties</i> , 2002 , 82, 1027-1045		5
100	Electronic states in mixed pseudobinary (Pb,Sr)S crystals. <i>Physical Review B</i> , 1991 , 43, 9758-9762	3.3	5
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