

Andrés Vega

List of Publications by Year in descending order

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172
papers

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174
all docs

174
docs citations

174
times ranked

1972
citing authors

#	ARTICLE	IF	CITATIONS
1	Structure and magnetism of cobalt clusters. Physical Review B, 2003, 67, .	1.1	128
2	Systematicab initiostudy of the electronic and magnetic properties of different pure and mixed iron systems. Physical Review B, 2000, 61, 13639-13646.	1.1	98
3	Structure and magnetism of small rhodium clusters. Physical Review B, 2002, 66, .	1.1	96
4	Calculatedsp-electron andspd-hybridization effects on the magnetic properties of smallFeNclusters. Physical Review B, 1993, 47, 4742-4746.	1.1	89
5	Magnetic moments of Ni clusters. Physical Review B, 1998, 57, 12469-12475.	1.1	73
6	Spin polarization at the Fe/V interface. Physical Review B, 1993, 48, 985-992.	1.1	69
7	Ti-decorated zigzag graphene nanoribbons for hydrogen storage. A van der Waals-corrected density-functional study. International Journal of Hydrogen Energy, 2015, 40, 4960-4968.	3.8	65
8	Structural and magnetic properties of X12Y (X, Y=Fe, Co, Ni, Ru, Rh, Pd, and Pt) nanoalloys. Journal of Chemical Physics, 2010, 132, .	1.2	61
9	Comparative of ab initio ρ -study of the structural, electronic, and magnetic trends of isoelectronic late transition metal clusters. Physical Review B, 2008, 78, .	1.1	59
10	Antiferromagnetic interlayer coupling in Fe/V and Fe/Cr. Journal of Applied Physics, 1991, 69, 4544-4546.	1.1	56
11	Study of the Structural and Electronic Properties of Rh _N and Ru _N Clusters ($N \leq 20$) within the Density Functional Theory. Journal of Physical Chemistry A, 2009, 113, 13483-13491.	1.1	53
12	Theoretical investigation of free-standing CoPd nanoclusters as a function of cluster size and stoichiometry in the Pd-rich phase: Geometry, chemical order, magnetism, and metallic behavior. Physical Review B, 2006, 74, .	1.1	52
13	Geometrical effects on the magnetism of small Ni clusters. Physical Review B, 1997, 55, 13279-13282.	1.1	48
14	Optical properties of large band gap \hat{I}^2 -In ₂ S ₃ compounds obtained by physical vapour deposition. Optical Materials, 2005, 27, 647-653.	1.7	48
15	Average magnetization and local magnetic moments of FeNclusters($N < 230$). Physical Review B, 1999, 60, 434-439.	1.1	43
16	Induced spin polarization inV:FenVmsuperlattices and thin V films on Fe substrates. Physical Review B, 1999, 59, 14510-14515.	1.1	40
17	Stability, magnetic behavior, and chemical order of M_n clusters. Physical Review B, 2009, 79, .	1.1	40
18	Possibility of various magnetic configurations in the Cr (Fe) monolayer deposited on vicinal surfaces of Fe (Cr). Physical Review B, 1995, 51, 11546-11554.	1.1	39

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19	Topological antiferromagnetism at Cr surfaces and interfaces. <i>Physical Review B</i> , 1994, 49, 12797-12800.	1.1	36
20	Li-decorated Pmmn8 phase of borophene for hydrogen storage. A van der Waals corrected density-functional theory study. <i>International Journal of Hydrogen Energy</i> , 2019, 44, 1021-1033.	3.8	35
21	All-electron and pseudopotential study of the spin-polarization of the V(001) surface: LDA versus GGA. <i>Physical Review B</i> , 2001, 63, .	1.1	34
22	Magnetic and electronic properties of substitutional FeN cluster impurities in Cr: Transition from antiferromagnetic to ferromagnetic FeN. <i>Physical Review B</i> , 1994, 50, 3899-3906.	1.1	32
23	Noncollinear magnetism of thin Cr films deposited on a stepped Fe (001) surface. <i>Physical Review B</i> , 2003, 68, .	1.1	32
24	What will freestanding borophene nanoribbons look like? An analysis of their possible structures, magnetism and transport properties. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 1054-1061.	1.3	32
25	Noncollinear Fe spin structure in (Sm-Co)/Fe exchange-spring bilayers: Layer-resolved $\langle \mathbf{m}_i \rangle$ and $\langle \mathbf{m}_i^2 \rangle$ spectroscopy and electronic structure calculations. <i>Physical Review B</i> , 2012, 85, .	1.1	31
26	Electronic quasiparticle structure of ferromagnetic bcc iron. <i>European Physical Journal B</i> , 1995, 96, 357-372.	0.6	30
27	Magnetic-Order Transition in Thin Fe Overlayers on Cr: Role of the Interfacial Roughness. <i>Europhysics Letters</i> , 1995, 31, 561-566.	0.7	29
28	Magnetism of Co nanoparticles supported on the Cu(111) substrate: Size and environment dependence. <i>Physical Review B</i> , 1997, 55, 445-451.	1.1	29
29	Stability, structural, and magnetic phase diagrams of ternary ferromagnetic 3d-transition-metal clusters with five and six atoms. <i>Journal of Chemical Physics</i> , 2011, 134, 054101.	1.2	29
30	Antiferromagnetic interlayer coupling in Fe/c-SiFe/Fe sandwiches and multilayers. <i>Physical Review B</i> , 2001, 65, .	1.1	28
31	Hydrogen Interaction in Pd-Pt Alloy Nanoparticles. <i>Journal of Physical Chemistry C</i> , 2012, 116, 126-133.	1.5	28
32	Origin of dead magnetic Fe overlayers on V(110). <i>Physical Review B</i> , 2001, 64, .	1.1	27
33	Magnetic properties of Pd atomic clusters from different theoretical approaches. <i>European Physical Journal D</i> , 2007, 44, 125-131.	0.6	26
34	C(2 Å ⁻²) ferrimagnetic order versus P(1 Å ⁻¹) ferromagnetic order in V, Cr, Mn monolayers on Fe(001). <i>Thin Solid Films</i> , 1996, 275, 103-105.	0.8	25
35	Magnetic behavior of monoatomic Co wires on Pd(110). <i>Physical Review B</i> , 2000, 61, 6848-6853.	1.1	25
36	Ab Initio Study of the Adsorption of NO on the Rh6+Cluster. <i>Journal of Physical Chemistry A</i> , 2011, 115, 8350-8360.	1.1	25

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37	Al enhances the H storage capacity of graphene at nanoribbon borders but not at central sites: A study using nonlocal van der Waals density functionals. <i>Physical Review B</i> , 2012, 85, .	1.1	25
38	Structure, fragmentation patterns, and magnetic properties of small cobalt oxide clusters. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 21732-21741.	1.3	25
39	Self-consistent theory of overlap interactions in the tight-binding method. <i>Physical Review B</i> , 1993, 47, 12995-12998.	1.1	24
40	Structural and magnetic properties of CoRh nanoparticles. <i>Physical Review B</i> , 2004, 70, .	1.1	24
41	A density-functional study of the structures, binding energies and magnetic moments of the clusters Mo_N ($N = 2-13$), $Mo_{12}Fe$, $Mo_{12}Co$ and $Mo_{12}Ni$. <i>Nanotechnology</i> , 2008, 19, 145704.	1.3	24
42	Spin configuration in a frustrated ferromagnetic/antiferromagnetic thin-film system. <i>Nanotechnology</i> , 2007, 18, 235702.	1.3	23
43	Structural and zero-point vibrational effects on the electric dipole moments and static dipole polarizabilities of sodium clusters. <i>Physical Review B</i> , 2011, 84, .	1.1	23
44	Insulating or Metallic: Coexistence of Different Electronic Phases in Zinc Clusters. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 2111-2115.	7.2	23
45	Molecular hydrogen uptake by zigzag graphene nanoribbons doped with early 3d transition-metal atoms. <i>International Journal of Hydrogen Energy</i> , 2013, 38, 8872-8880.	3.8	22
46	Tight-binding study of the ionization of iron clusters. <i>Physical Review B</i> , 1996, 54, 3003-3006.	1.1	21
47	A density-functional study of the possibility of noncollinear magnetism in small Mn clusters using SIESTA and the generalized gradient approximation to exchange and correlation. <i>Journal of Chemical Physics</i> , 2008, 128, 114315.	1.2	21
48	Finite-temperature ferromagnetism of f.c.c. cobalt. <i>Physica Status Solidi (B): Basic Research</i> , 1996, 193, 177-187.	0.7	20
49	Tailoring the spin density waves in Fe/Cr multilayers by selective inclusion of Sn, V and Mn. <i>Surface Science</i> , 2009, 603, 117-124.	0.8	20
50	Magnetic magic numbers are not magic for clusters embedded in noble metals. <i>Physical Review B</i> , 2002, 66, .	1.1	19
51	Density functional study of ternary $Fe_xCo_yNi_z$ ($x+y+z=7$) clusters. <i>Theoretical Chemistry Accounts</i> , 2013, 132, 1.	0.5	19
52	Nonmetal-metal transition in Ni clusters. <i>Solid State Communications</i> , 1997, 104, 635-639.	0.9	18
53	Competition between two- and three-dimensional growth of Co clusters deposited on Cu(001): Influence on the magnetic properties. <i>Physical Review B</i> , 2001, 63, .	1.1	18
54	Metallic behavior of Pd atomic clusters. <i>Nanotechnology</i> , 2007, 18, 365706.	1.3	18

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55	Magnetic properties of semi-infinite systems of Cr(Fe) on a Fe(Cr) (001) surface. <i>Surface Science</i> , 1991, 251-252, 51-54.	0.8	17
56	Magnetism of Vicinal Surfaces of Vanadium. <i>Europhysics Letters</i> , 1994, 27, 165-170.	0.7	17
57	Magnetism in small Pd clusters. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2004, 332, 107-114.	0.9	17
58	A new magnetic superatom: Cr@Zn ₁₇ . <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 28033-28043.	1.3	17
59	Structure and magnetic properties of small Fe clusters supported on the Ni(001) surface. <i>Physical Review B</i> , 2005, 71, .	1.1	16
60	Ab initio study of lithium decoration of popgraphene and hydrogen storage capacity of the hybrid nanostructure. <i>International Journal of Hydrogen Energy</i> , 2021, 46, 15724-15737.	3.8	16
61	Stepped Fe(001) surface magnetism. <i>Journal of Magnetism and Magnetic Materials</i> , 1992, 104-107, 1687-1688.	1.0	15
62	Optical conductivity in substoichiometric titanium carbides. <i>Physical Review B</i> , 1998, 58, 3507-3510.	1.1	15
63	Collinear versus noncollinear magnetic order in Pd atomic clusters: Ab initio calculations. <i>Physical Review B</i> , 2006, 74, .	1.1	15
64	Antiferromagnetic-like coupling in the cationic iron cluster of thirteen atoms. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 14458.	1.3	15
65	Are zinc clusters really amorphous? A detailed protocol for locating global minimum structures of clusters. <i>Nanoscale</i> , 2018, 10, 19162-19181.	2.8	15
66	Hydrogen storage capacity of Li-decorated borophene and pristine graphene slit pores: A combined ab initio and quantum-thermodynamic study. <i>Applied Surface Science</i> , 2021, 562, 150019.	3.1	15
67	Magnetic moments of. <i>European Physical Journal D</i> , 1999, 6, 235.	0.6	15
68	The magnetization reversal process in spin spring magnets. <i>Nanotechnology</i> , 2008, 19, 315401.	1.3	14
69	$\text{Mo}_{n-4}\text{Cu}_n$ Structural transition and electronic structure of interest in spintronics. <i>Physical Review B</i> , 2009, 79, 114401.	1.1	14
70	Structural, Electronic, and Magnetic Properties Of Co _n Cu _m Nanoalloys ($m + n = 12$) from First Principles Calculations. <i>Journal of Physical Chemistry A</i> , 2012, 116, 9353-9360.	1.1	14
71	Antiferromagnetic versus ferromagnetic orders in V, Cr and Mn overlayers on Fe(103). <i>Journal of Magnetism and Magnetic Materials</i> , 1995, 148, 177-178.	1.0	13
72	Ab initio optical conductivity in LaMO ₃ (M=Ti, Cu). <i>Physical Review B</i> , 1996, 54, 11271-11275.	1.1	13

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73	Tight-binding and evolutionary search approach for nanoscale CoRh alloys. <i>Physica B: Condensed Matter</i> , 2005, 370, 200-214.	1.3	13
74	Twining effects in the magnetism of small Pd clusters. <i>Solid State Communications</i> , 2005, 133, 573-578.	0.9	13
75	Engineering the magnetic structure of Fe clusters by Mn alloying. <i>Nanotechnology</i> , 2008, 19, 245701.	1.3	13
76	New structural and electronic properties of (TiO ₂) ₁₀ . <i>Journal of Chemical Physics</i> , 2016, 144, 234312.	1.2	13
77	Spin-polarized transport in hydrogen-passivated graphene and silicene nanoribbons with magnetic transition-metal substituents. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 22606-22616.	1.3	13
78	Unusual effect of interatomic interactions on magnetism: Rh adatoms on the Ag(001) surface. <i>Physical Review B</i> , 1998, 57, R14020-R14023.	1.1	12
79	Structural and magnetic properties of Fe clusters at the Al (001) surface: Early transition from paramagnetic to ferromagnetic Fe. <i>Physical Review B</i> , 2004, 69, .	1.1	12
80	Multiple collinear magnetic arrangements in thin Mn films supported on Fe(001). Antiferromagnetic versus ferromagnetic behavior. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2005, 337, 469-472.	0.9	12
81	Impact of dimerization and stretching on the transport properties of molybdenum atomic wires. <i>Nanotechnology</i> , 2010, 21, 095205.	1.3	12
82	Hydrogen insertion in Pd core/Pt shell cubo-octahedral nanoparticles. <i>Physical Review B</i> , 2011, 83, .	1.1	12
83	A new family of star-like icosahedral structures for small cobalt clusters. <i>Chemical Physics</i> , 2013, 415, 106-111.	0.9	12
84	Structure, fragmentation patterns, and magnetic properties of small nickel oxide clusters. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 3366-3383.	1.3	12
85	Density functional study of the structural, electronic, and magnetic properties of Mo _n and Mo _n S (n) Tj ETQq1 1 0,784314 rgBT /Ove	0,8	12
86	Density functional study of the optical response of FeN and CoN nitrides with zinc-blend and rock-salt structures. <i>Journal of Physics and Chemistry of Solids</i> , 2017, 100, 148-153.	1.9	12
87	Magnetic behaviour of selected geometries of Pd clusters: icosahedral versus fcc structures. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2004, 330, 126-130.	0.9	11
88	Magnetic structure of cobalt clusters. <i>Journal of Alloys and Compounds</i> , 2004, 369, 93-96.	2.8	11
89	Properties Design: Prediction and Experimental Validation of the Luminescence Properties of a New Eu^{II}-Based Phosphor. <i>Chemistry - A European Journal</i> , 2018, 24, 16276-16281.	1.7	11
90	Density Functional Study of Charge Transfer at the Graphene/Ionic Liquid Interface. <i>Journal of Physical Chemistry C</i> , 2018, 122, 15070-15077.	1.5	11

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91	Vanadium: From cluster to semi-infinite crystal. <i>Journal of Applied Physics</i> , 1993, 73, 6207-6209.	1.1	10
92	Magnetic moments of Ni monolayers and small ground-state Ni clusters at the Al (001) surface. <i>Physical Review B</i> , 2000, 62, 11104-11108.	1.1	10
93	Magnetic Cooperative Effects in Small Ni-Ru Clusters. <i>Journal of Physical Chemistry A</i> , 2011, 115, 13950-13955.	1.1	10
94	Relation between structural patterns and magnetism in small iron oxide clusters: reentrance of the magnetic moment at high oxidation ratios. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 246-272.	1.3	10
95	Theoretical study of the gap evolution of In ₂ X ₃ (X=O, S, Se, Te) with lattice compression. <i>Optical Materials</i> , 2001, 17, 497-499.	1.7	9
96	Electronic structure calculations of low-dimensional transition metals. <i>Handbook of Magnetic Materials</i> , 2003, 15, 199-288.	0.6	9
97	Density Functional Study of the Structures and Electronic Properties of Nitrogen-Doped Ni _n Clusters, $n = 1-10$. <i>Journal of Physical Chemistry A</i> , 2011, 115, 13997-14005.	1.1	9
98	On the electric dipole moments of small sodium clusters from different theoretical approaches. <i>Chemical Physics</i> , 2012, 399, 252-257.	0.9	9
99	Calculation of electronic properties of the (100) surface of Fe including overlap interactions. <i>Surface Science</i> , 1991, 251-252, 55-58.	0.8	8
100	Interplay between structure and magnetism in hydride iron-vanadium systems. <i>Physical Review B</i> , 2008, 78, .	1.1	8
101	Possibility of collinear magnetic order in frustrated free-standing Fe_{2n} . <i>Physical Review B</i> , 2008, 77, .	1.1	8
102	Structure and electronic properties of molybdenum monatomic wires encapsulated in carbon nanotubes. <i>Journal of Physics Condensed Matter</i> , 2011, 23, 265302.	0.7	8
103	Theoretical study of Al _n V ⁺ clusters and their interaction with Ar. <i>Journal of Chemical Physics</i> , 2013, 139, 214305.	1.2	8
104	Spin-orbit effects on the structural, homotop, and magnetic configurations of small pure and Fe-doped Pt clusters. <i>Journal of Nanoparticle Research</i> , 2014, 16, 1.	0.8	8
105	Borophene vs. graphene interfaces: Tuning the electric double layer in ionic liquids. <i>Journal of Molecular Liquids</i> , 2020, 303, 112647.	2.3	8
106	Defects and Magnetic Properties: The Cr/Fe(001) Interfaces. <i>Materials Research Society Symposia Proceedings</i> , 1995, 384, 247.	0.1	7
107	Spin polarization at Fe/Cr interfaces. <i>Journal of Applied Physics</i> , 1997, 81, 4347-4349.	1.1	7
108	Is the vanadium(001) surface magnetic? Pseudopotential toward all-electron calculations. <i>International Journal of Quantum Chemistry</i> , 2003, 91, 230-233.	1.0	7

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109	Noncollinear magnetism caused by intermixing at ferromagnetic/antiferromagnetic interfaces. <i>Physical Review B</i> , 2006, 74, .	1.1	7
110	A density-functional study of the structures and electronic properties of neutral, anionic, and endohedrally doped In _x P _x clusters. <i>Journal of Chemical Physics</i> , 2009, 131, 074504.	1.2	7
111	Structural and Electronic Properties of TM _n [(BN) ₃ H ₆] ₆ Complexes with TM = Co (<i>n</i> = 1–3) and with TM = Fe, Ni, Ru, Rh, Pd (<i>n</i> = 1–3). <i>Journal of Physical Chemistry A</i> , 2014, 118, 2976-2983.	1.1	7
112	Zn ₁₇ Superatom Cage Doped with 3d Transition-Metal (TM) Impurities (TM = Sc, Ti, V, Cr, Mn,). <i>J. Phys. Chem. C</i> , 2015, 119, 11200-11209.	1.5	7
113	Onset of C(2Å–2) ferrimagnetic order in Cr islands deposited on Fe(001) as a function of island size. <i>Journal of Applied Physics</i> , 1996, 79, 5834.	1.1	6
114	Spin-polarization of thin Mn films on Fe(107). <i>Journal of Magnetism and Magnetic Materials</i> , 1996, 156, 199-201.	1.0	6
115	Interlayer exchange coupling in Fe/SiFe/Fe sandwiches. <i>Surface Science</i> , 2001, 482-485, 994-997.	0.8	6
116	Magnetic behavior of Pd nanoclusters. <i>Physica B: Condensed Matter</i> , 2004, 354, 271-277.	1.3	6
117	Electronic structure and transport properties of monatomic Fe chains in a vacuum and anchored to a graphene nanoribbon. <i>Journal of Physics Condensed Matter</i> , 2012, 24, 455304.	0.7	6
118	Structural, Electronic, and Magnetic Properties of Iron Disulfide Fe ₂ S ₂ ± ⁰ Å [±] (<i>n</i> = 1–6) Clusters. <i>Journal of Physical Chemistry A</i> , 2017, 121, 3768-3780.	1.1	6
119	Nanoscale reactivity of Zn x Mg 20 [±] x investigated by structural and electronic indicators. <i>Corrosion Science</i> , 2017, 124, 35-45.	3.0	6
120	Magnetic properties of Co islands grown on Cu(111). <i>Surface Science</i> , 1996, 352-354, 902-906.	0.8	5
121	Magnetism in Rh clusters under hydrostatic deformations. <i>European Physical Journal D</i> , 2003, 23, 343-349.	0.6	5
122	Electronic structure investigation of the exchange-spring behavior during the magnetic reversal process. <i>Physical Review B</i> , 2008, 77, .	1.1	5
123	Ab initio study of hydrogen insertion in ultrathin transition metal doped V films: Structural and electronic properties. <i>Physical Review B</i> , 2010, 81, .	1.1	5
124	Charge and spin transport properties of Mo ₂ X ₂ nanoribbons. <i>Journal of Physics Condensed Matter</i> , 2014, 26, 165302.	0.7	5
125	Spin-dependent electronic conduction along zigzag graphene nanoribbons bearing adsorbed Ni and Fe nanostructures. <i>Journal of Physics Condensed Matter</i> , 2014, 26, 165302.	0.7	5
126	Structural, Vibrational, and Magnetic Properties of FeCoO ₂ ± ⁰ (<i>n</i> = 1–6) Bimetallic Oxide Clusters. <i>Journal of Physical Chemistry C</i> , 2015, 119, 11200-11209.	1.5	5

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127	Structural properties, magnetism and reactivity of $\text{Ni}_{13}\text{Fe}_x$ nanoclusters. <i>Journal of Applied Physics</i> , 2011, 110, 124305.	1.0	5
128	Why are Zn-rich Zn-Mg nanoalloys optimal protective coatings against corrosion? A first-principles study of the initial stages of the oxidation process. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 24685-24698.	1.3	5
129	Geometrical and chemical environment effects on the magnetism of stepped surfaces of V and V over Fe. <i>European Physical Journal D</i> , 1993, 43, 1045-1050.	0.4	4
130	Antiferromagnetic versus ferromagnetic coupling in Fe/Cr(107) and Cr/Fe(107). <i>Journal of Applied Physics</i> , 1994, 76, 6989-6991.	1.1	4
131	Electronic structure and optical response of L-CePd ₅ . <i>Physical Review B</i> , 1995, 51, 4823-4829.	1.1	4
132	Magnetic moments in Ni clusters with deformations. <i>Solid State Communications</i> , 2001, 117, 477-482.	0.9	4
133	Effects of the structural deformations on the magnetism of Rh ₆ and Rh ₁₃ clusters. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2003, 318, 473-479.	0.9	4
134	Magnetism in segregated bimetallic CoRh nanoclusters. <i>Physica B: Condensed Matter</i> , 2004, 354, 278-281.	1.3	4
135	Theoretical study of the charge transfer in supported transition metal microclusters. <i>European Physical Journal D</i> , 2005, 34, 51-54.	0.6	4
136	Improvement of hydrogen uptake in iron and vanadium matrices by doping with 3d atomic impurities. <i>Journal of Alloys and Compounds</i> , 2012, 545, 19-27.	2.8	4
137	Impact of S doping on the structural, electronic and magnetic properties of Cr _n (n = 2-6) clusters. <i>European Physical Journal D</i> , 2017, 71, 1.	0.6	4
138	Predicting photon cascade emission in Pr ³⁺ -doped fluorides. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 15503-15511.	1.3	4
139	Structural and Electronic Rearrangements in Fe ₂ S ₂ , Fe ₃ S ₄ , and Fe ₄ S ₄ Atomic Clusters under the Attack of NO, CO, and O ₂ . <i>Journal of Physical Chemistry A</i> , 2019, 123, 10919-10929.	1.1	4
140	Incorporating charge transfer effects into a metallic empirical potential for accurate structure determination in (ZnMg) _N nanoalloys. <i>Nanoscale</i> , 2020, 12, 20432-20448.	2.8	4
141	A neural network potential for searching the atomic structures of pure and mixed nanoparticles. Application to ZnMg nanoalloys with an eye on their anticorrosive properties. <i>Acta Materialia</i> , 2021, 220, 117341.	3.8	4
142	Magnetic properties of very thin V films adsorbed on Fe substrate. <i>Journal of Magnetism and Magnetic Materials</i> , 1993, 121, 177-179.	1.0	3
143	Optical absorption and interband transitions in CePd ₇ . <i>Physical Review B</i> , 1996, 53, 6881-6884.	1.1	3
144	Temperature-dependent magnetic behavior of ideal and stepped Fe surfaces. <i>Physical Review B</i> , 2001, 63, .	1.1	3

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145	The structure of an Fe monolayer on the Ni (111) surface. A density-functional study using the generalized gradient approximation. Solid State Communications, 2006, 137, 129-131.	0.9	3
146	Magnetic coupling in the cluster Fe ₂ Mn ₄ : A fully unconstrained density-functional study. Solid State Communications, 2006, 140, 480-482.	0.9	3
147	Response of Mn overlayers on Fe to external magnetic fields: Electronic structure calculations. Surface Science, 2009, 603, 2537-2543.	0.8	3
148	Spin currents and filtering behavior in zigzag graphene nanoribbons with adsorbed molybdenum chains. Journal of Physics Condensed Matter, 2015, 27, 135301.	0.7	3
149	Small Ni clusters at the (110) and (111) surfaces of Al: structures and lack of magnetic moment. Surface Science, 2001, 482-485, 976-980.	0.8	2
150	Morphology and magnetism of Fe monolayers and small Fe clusters (n= 2-19) supported on the Ni(111) surface. Nanotechnology, 2007, 18, 055701.	1.3	2
151	Proximity effects on the spin density waves in X/Cr(001) multilayers (X=Sn, V, and Mn). Thin Solid Films, 2011, 519, 2037-2042.	0.8	2
152	Magnetization reversal process at atomic scale in systems with itinerant electrons. Journal of Physics Condensed Matter, 2012, 24, 176002.	0.7	2
153	Structural, magnetic, and vibrational properties of stoichiometric clusters of $\langle \text{scp} \rangle \text{C} \langle / \text{scp} \rangle \langle \text{scp} \rangle \text{N} \langle / \text{scp} \rangle$. International Journal of Quantum Chemistry, 2015, 115, 523-528.	1.0	2
154	Structure, fragmentation patterns, and electronic properties of small indium oxide clusters. Theoretical Chemistry Accounts, 2018, 137, 1.	0.5	2
155	Tuning the Magnetic Moment of Small Late 3d-Transition-Metal Oxide Clusters by Selectively Mixing the Transition-Metal Constituents. Nanomaterials, 2020, 10, 1814.	1.9	2
156	Stepped Fe(100) and V/Fe(100) Magnetism. Materials Research Society Symposia Proceedings, 1991, 231, 323.	0.1	1
157	Magnetic properties of small 3-d transition metal clusters: Role of the sp-electrons and spd-hybridization. Scripta Materialia, 1993, 3, 359-363.	0.5	1
158	Calculation of the magnetic properties of FeN clusters embedded in 3d transition-metal matrices. Computational Materials Science, 1994, 2, 463-467.	1.4	1
159	Electronic, magnetic, and optical properties of. Journal of Physics Condensed Matter, 1997, 9, 6267-6277.	0.7	1
160	Competition between CoCr hybridization and low-coordination effects in the magnetism of a Co monolayer deposited on Cr substrates. Journal of Magnetism and Magnetic Materials, 1997, 165, 265-267.	1.0	1
161	Deformation Effects in the Magnetic Moments of Ni Clusters. , 2001, , 77-85.		1
162	Magnetism of the Fe ₉ nanocluster supported on Ni(001). Computational Materials Science, 2006, 35, 307-310.	1.4	1

#	ARTICLE	IF	CITATIONS
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