

Vladimir Sokolovskiy

List of Publications by Year in descending order

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157
papers

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163
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citing authors

#	ARTICLE	IF	CITATIONS
1	c and structural properties of Co-Ni- Z xmins:mml= http://www.w3.org/1998/Math/MathML display="inline" id="d1e1016"> $\text{Z} \approx 1.0784314 \text{ rgBT /Overclock 10 Tf 50 747 Td}$	2.3	2
2	Electronic and Vibrational Properties of Fe _x NiAl and Co _x NiAl Full Heusler Alloys: A First-Principles Comparison. <i>IEEE Transactions on Magnetics</i> , 2022, 58, 1-5.	2.1	0
3	Magnetocaloric effect and magnetic phase diagram of Ni-Mn-Ga Heusler alloy in steady and pulsed magnetic fields. <i>Journal of Alloys and Compounds</i> , 2022, 904, 164051.	5.5	14
4	Impact of local arrangement of Fe and Ni on the phase stability and magnetocrystalline anisotropy in Fe-Ni-Al Heusler alloys. <i>Physical Review Materials</i> , 2022, 6, .	2.4	6
5	Structural and magnetic properties of Fe _x Al alloys: Ab initio studies. <i>Journal of Magnetism and Magnetic Materials</i> , 2022, 557, 169437.	2.3	3
6	Segregation tendency and properties of FeRh _{1-x} Pt alloys. <i>Journal of Magnetism and Magnetic Materials</i> , 2022, 556, 169403.	2.3	1
7	Exchange Correlation Effects in Modulated Martensitic Structures of the Mn ₂ NiGa Alloy. <i>Physics of Metals and Metallography</i> , 2022, 123, 375-380.	1.0	3
8	Review of Modern Theoretical Approaches for Study of Magnetocaloric Materials. <i>Physics of Metals and Metallography</i> , 2022, 123, 319-374.	1.0	12
9	Study of Heat Transfer Processes in a System Containing Fe _x Rh Microwires. <i>Physics of Metals and Metallography</i> , 2022, 123, 381-385.	1.0	0
10	Magnetocaloric Effect in Metals and Alloys. <i>Physics of Metals and Metallography</i> , 2022, 123, 315-318.	1.0	13
11	Prediction of a Heusler alloy with switchable metal-to-half-metal behavior. <i>Physical Review B</i> , 2021, 103, .	3.2	8
12	A Ternary Map of Ni _x Mn _{1-x} Ga Heusler Alloys from Ab Initio Calculations. <i>Metals</i> , 2021, 11, 973.	2.3	4
13	Phase transitions in Fe-(23-x)Ga alloys: Experimental results and modeling. <i>Journal of Alloys and Compounds</i> , 2021, 885, 160917.	5.5	3
14	Design of a Stable Heusler Alloy with Switchable Metal-to-Half-Metal Transition at Finite Temperature. <i>Advanced Theory and Simulations</i> , 2021, 4, 2100311.	2.8	6
15	FIRST-PRINCIPLES STUDIES OF THE PHASE TRANSITIONS IN Fe-Si ALLOYS. <i>Bulletin of the South Ural State University Series Mathematics Mechanics Physics</i> , 2021, 13, 52-58.	0.2	0
16	Statistical model for the martensitic transformation simulation in Heusler alloys. <i>Physica B: Condensed Matter</i> , 2020, 578, 411874.	2.7	2
17	Structural, magnetic and electronic properties of FeRh _x Pd _{1-x} compounds: Ab initio study. <i>Physica B: Condensed Matter</i> , 2020, 578, 411882.	2.7	2
18	First principles study of structural and magnetic properties in Fe _{100-x} Ge alloys. <i>Physica B: Condensed Matter</i> , 2020, 580, 411934.	2.7	4

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19	Phase transitions in Fe ₃ Al-based alloys: <i>ab initio</i> study. <i>Phase Transitions</i> , 2020, 93, 43-53.	1.3	1
20	Electronic structure beyond the generalized gradient approximation for $\text{Ni}_{2+x}\text{Mn}_{1-x}\text{Ga}$. <i>Physical Review B</i> , 2020, 102, .		
21	Superconducting and antiferromagnetic properties of dual-phase V ₃ Ga. <i>Applied Physics Letters</i> , 2020, 117, 062401.	3.3	5
22	Magnetocaloric properties of $\text{Ni}_{2+x}\text{Mn}_{1-x}\text{Ga}$ with coupled magnetostructural phase transition. <i>Journal of Applied Physics</i> , 2020, 127, .	2.5	9
23	Ab Initio Studies of Phase Transformations in Fe _{100-x} Si _x . <i>Physics of the Solid State</i> , 2020, 62, 739-743.	0.6	2
24	Exchange-correlation corrections for electronic properties of half-metallic Co ₂ FeSi and nonmagnetic semiconductor CoFeTiAl. <i>Journal of Applied Physics</i> , 2020, 127, .	2.5	10
25	Electronic and Magnetic Properties of DyFe ₄ Ge ₂ Alloys near a Phase Transition. <i>Physics of the Solid State</i> , 2020, 62, 931-936.	0.6	0
26	A Study of the Structure and Magnetic Properties of FeRh _{1-x} Ir _x (x = 0.5-1) Alloys by First-Principles Methods. <i>Physics of the Solid State</i> , 2020, 62, 963-967.	0.6	2
27	First-Principles Study of the Structure and Properties of Fe ₃ Pd and Fe-Pd-Rh Alloys. <i>Shape Memory and Superelasticity</i> , 2020, 6, 61-66.	2.2	2
28	Effects of magnetic and structural phase transitions on the normal and anomalous Hall effects in Ni-Mn-In-B Heusler alloys. <i>Physical Review B</i> , 2020, 101, .	3.2	24
29	Coulomb correlation in noncollinear antiferromagnetic $\text{Mn}_{1\pm x}\text{Ga}$. <i>Physical Review B</i> , 2020, 101, .	3.2	27
30	Prediction of giant magnetocaloric effect in Ni ₄₀ Co ₁₀ Mn ₃₆ Al ₁₄ Heusler alloys: An insight from ab initio and Monte Carlo calculations. <i>Journal of Applied Physics</i> , 2020, 127, 163901.	2.5	8
31	Phase Transformations in Ni(Co)-Mn(Cr,C)-(In,Sn) Alloys: An Ab Initio Study. <i>Physics of Metals and Metallography</i> , 2020, 121, 202-209.	1.0	4
32	Theoretical Approach to Investigation of the Magnetic and Magnetocaloric Properties of Heusler Ni-Mn-Ga Alloys. <i>Physics of the Solid State</i> , 2020, 62, 785-792.	0.6	4
33	Magnetocaloric and Shape Memory Effects in the Mn ₂ NiGa Heusler Alloy. <i>Physics of the Solid State</i> , 2020, 62, 815-820.	0.6	8
34	VOLUME MAGNETOSTRICTION OF FE-GA ALLOYS: CALCULATION FROM FIRST PRINCIPLES. <i>Bulletin of the South Ural State University Series Mathematics Mechanics Physics</i> , 2020, 12, 57-62.	0.2	0
35	Structural and Elastic Properties of Fe-Ge Alloys: <i>ab initio</i> studies. <i>Bulletin of the South Ural State University Series Mathematics Mechanics Physics</i> , 2020, 12, 49-56.	0.2	0
36	Phase Diagram of Fe-Al Alloys: A Study from First Principles. <i>Bulletin of the Russian Academy of Sciences: Physics</i> , 2019, 83, 844-846.	0.6	3

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37	Correlation effects in the ground state of Ni-(Co)-Mn-Sn Heusler compounds. <i>MRS Advances</i> , 2019, 4, 441-446.	0.9	3
38	Monte Carlo simulations of hysteresis effects at the martensitic transformation. <i>Physica B: Condensed Matter</i> , 2019, 575, 411692.	2.7	2
39	Correlation effects on ground-state properties of ternary Heusler alloys: First-principles study. <i>Physical Review B</i> , 2019, 99, .	3.2	28
40	Phase diagram of magnetostrictive Fe-Ga alloys: insights from theory and experiment. <i>Phase Transitions</i> , 2019, 92, 101-116.	1.3	33
41	Magnetostriction of Fe _{100-x} Gax alloys from first principles calculations. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 476, 120-123.	2.3	7
42	Investigation of electronic, magnetic and structural properties of the Fe _{1-x} MnxRh. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 476, 325-328.	2.3	6
43	First-principles study of Ni-Co-Mn-Sn alloys with regular and inverse Heusler structure. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 476, 546-550.	2.3	7
44	Ground state and magnetic properties of the Cr-doped Ni-Mn-(Ga, Ge, In, Sn) alloys: Insights from ab initio study. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 470, 123-126.	2.3	7
45	Modeling of the structural and magnetic properties of Fe-Rh-(Z) (Z = Mn, Pt) alloys by first principles methods. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 470, 69-72.	2.3	7
46	Ternary diagrams of magnetic properties of Ni-Mn-Ga Heusler alloys from ab initio and Monte Carlo studies. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 470, 64-68.	2.3	10
47	Peculiarities of phonons in Ni-Mn-Ga alloys: Ab initio studies. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 470, 73-76.	2.3	2
48	Magnetic properties of Fe _{100-x} Ga : Ab initio and Monte Carlo study. <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 470, 118-122.	2.3	8
49	Segregation tendency of Heusler alloys. <i>Physical Review Materials</i> , 2019, 3, .	2.4	12
50	Theoretical study of heat transfer processes in Heusler-type magnetic microwires. <i>Letters on Materials</i> , 2019, 9, 395-399.	0.7	3
51	Ab Initio Calculation of Vacancy Formation Energy in Antiperovskite Mn ₃ GaC. <i>Bulletin of the South Ural State University Series Mathematics Mechanics Physics</i> , 2019, 11, 58-64.	0.2	0
52	FIRST-PRINCIPLES INVESTIGATIONS OF REFERENCE STATES OF Co ₂ CrIn HEUSLER ALLOYS. <i>Bulletin of the South Ural State University Series Mathematics Mechanics Physics</i> , 2019, 11, 59-66.	0.2	0
53	Modelling of Rhombohedral Magnetostriction in Fe-Ga Alloys. <i>Bulletin of the South Ural State University, Series: Mathematical Modelling, Programming and Computer Software</i> , 2019, 12, 158-165.	0.4	1
54	Probing Structural and Magnetic Instabilities and Hysteresis in Heuslers by Density Functional Theory Calculations (<i>Phys. Status Solidi B</i> 2/2018). <i>Physica Status Solidi (B): Basic Research</i> , 2018, 255, 1870108.	1.5	2

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55	Phenomenological analysis of thermal hysteresis in Ni-Mn-Ga Heusler alloys. <i>Phase Transitions</i> , 2018, 91, 469-476.	1.3	1
56	Structural, magnetic and thermodynamic properties of Mn ₃ -X-C (X = Ga, Sn) compounds: ab initio study. <i>Physica B: Condensed Matter</i> , 2018, 549, 94-97.	2.7	1
57	Probing Structural and Magnetic Instabilities and Hysteresis in Heuslers by Density Functional Theory Calculations. <i>Physica Status Solidi (B): Basic Research</i> , 2018, 255, 1700296.	1.5	11
58	Magnetocaloric effect in Ni-Co-Mn-(Sn, Al) Heusler alloys: Theoretical study. <i>Journal of Magnetism and Magnetic Materials</i> , 2018, 459, 295-300.	2.3	9
59	Magnetic states of Ni ₂ MnZ and Ni ₂ CrZ (Z = Al, As, Bi, Ga, Ge, In, P, Pb, Sb, Si, Sn, Tl) Heusler alloys. <i>Journal of Magnetism and Magnetic Materials</i> , 2018, 459, 78-83.	2.3	11
60	Monte Carlo Simulations of Thermal Hysteresis in Ni-Mn-Based Heusler Alloys. <i>Physica Status Solidi (B): Basic Research</i> , 2018, 255, 1700265.	1.5	3
61	The effect of antisite disorder on magnetic and magnetocaloric properties of Ni-Co-Mn-In alloys: ab initio and Monte Carlo studies., , 2018, , .	0	
62	Effect of disorder on magnetic properties and martensitic transformation of Co-doped Ni-Mn-Al Heusler alloy. <i>Intermetallics</i> , 2018, 102, 132-139.	3.9	12
63	Investigation of structural and magnetic properties of Fe-Rh-(Z) (Z = Co, Pt) alloys by first principles method. <i>EPJ Web of Conferences</i> , 2018, 185, 05005.	0.3	1
64	The Effect of Anti-Site Disorder on Structural and Magnetic Properties of Ni-Co-Mn-In Alloys: <italic>Ab Initio</italic> and Monte Carlo Studies. <i>IEEE Transactions on Magnetics</i> , 2018, 54, 1-5.	2.1	5
65	Ab Initio Study of the Structural, Magnetic, Electronic, and Thermodynamic Properties of Pd ₂ MnZ (Z =) T _j ETQq1 1 0.784314 gBT /Overlock 1		
66	First-Principles Study of the Structure and Magnetic Properties of Fe ₈ Rh ₈ xZx (Z = Mn, Pt, Co; x = 1,) T _j ETQq0 0.0 rgBT /Overlock 1		
67	Properties and Decomposition of Heusler Alloys. <i>Energy Technology</i> , 2018, 6, 1478-1490.	3.8	24
68	Ab initio study of magnetic and structural properties of Fe-Ga alloys. <i>EPJ Web of Conferences</i> , 2018, 185, 04013.	0.3	9
69	Structural and magnetic properties of heusler alloys Pd ₂ MnZ (Z=Ga, Ge, As): AB INITIO study. <i>EPJ Web of Conferences</i> , 2018, 185, 05007.	0.3	3
70	Complex investigations of phase diagram of Ni-Pt-Mn-Ga Heusler alloys. <i>Letters on Materials</i> , 2018, 8, 21-26.	0.7	4
71	First-principles and Monte Carlo studies of the Ni ₂ (Mn,Cr)Ga Heusler alloys electronic and magnetic properties. <i>Materials Research Express</i> , 2017, 4, 026105.	1.6	10
72	Magnetic properties and martensitic transformation of Ni-Mn-Ge Heusler alloys from first-principles and Monte Carlo studies. <i>Journal Physics D: Applied Physics</i> , 2017, 50, 195001.	2.8	8

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73	Ab initio calculations of structural and magnetic properties of Ni-Co-Mn-Cr-Sn supercell. <i>Intermetallics</i> , 2017, 87, 55-60.	3.9	11
74	Complex investigation of structural and magnetic properties of the Ni-Mn-(Ga, Ge) alloys within ab initio approach. <i>Materials Today: Proceedings</i> , 2017, 4, 4616-4620.	1.8	0
75	Effect of structural disorder on the ground state properties of Co ₂ CrAl Heusler alloy. <i>Physica B: Condensed Matter</i> , 2017, 519, 82-89.	2.7	16
76	Ab initio study of the composite phase diagram of Ni-Mn-Ga shape memory alloys. <i>Journal of Experimental and Theoretical Physics</i> , 2017, 125, 104-110.	0.9	5
77	Structural, magnetic and electronic properties of Ni-Mn-Ga-Cr Heusler alloys: ab initio and Monte Carlo studies. <i>Materials Today: Proceedings</i> , 2017, 4, 4621-4625.	1.8	1
78	First principles study of the structural and magnetic properties of Fe(Rh, Pd) and Fe(Rh, Ni) alloys. <i>Materials Today: Proceedings</i> , 2017, 4, 4642-4646.	1.8	11
79	Large change of magnetic moment in Ni₁₃Co₃Mn₁₃Sn₃ and Ni₁₃Co₃Mn₁₃Sn₂Al₁ Heusler alloys at martensitic transitions: investigation from first principles., 2017, ,	0	
80	Modeling of heat transfer processes in Ni ₂ MnIn magnetic wires. <i>Physica Status Solidi (A) Applications and Materials Science</i> , 2016, 213, 390-398.	1.8	6
81	Electronic and magnetic properties of the Co ₂ -based Heusler compounds under pressure: first-principles and Monte Carlo studies. <i>Journal Physics D: Applied Physics</i> , 2016, 49, 355004.	2.8	41
82	Magnetic, thermal, and electrical properties of an Ni _{45.37} Mn _{40.91} In _{13.72} Heusler alloy. <i>Journal of Experimental and Theoretical Physics</i> , 2016, 122, 874-882.	0.9	18
83	< i>Ab Initio</i> Investigations of Structural and Magnetic Properties of Cr-Doped Ni-Co-Mn-Sn Heusler Alloys. <i>Materials Science Forum</i> , 2016, 845, 134-137.	0.3	3
84	First principles study of the structural properties of Ni _{1.75} Co _{0.25} Mn _{1.5-x} Cr _x In _{0.5} Heusler alloy. <i>MATEC Web of Conferences</i> , 2015, 33, 02002.	0.2	1
85	Magnetic states of C-doped Ni _{43.75} Co _{6.25} Mn _{37.5} In _{12.5} Heusler alloys. <i>MATEC Web of Conferences</i> , 2015, 33, 05001.	0.2	1
86	Ab initio calculations of structural and magnetic properties of Ni-Co-Mn-Cr-Sn alloys. <i>MATEC Web of Conferences</i> , 2015, 33, 05003.	0.2	0
87	First-principles and Monte Carlo studies of C-doped Ni ₄₅ Co ₅ Mn ₃₇ In ₁₃ Heusler alloys. <i>MATEC Web of Conferences</i> , 2015, 33, 05004.	0.2	1
88	Compositional trends in Ni-Mn-Ga Heusler alloys: first-principles approach. <i>MATEC Web of Conferences</i> , 2015, 33, 05005.	0.2	4
89	Achieving large magnetocaloric effects in Co- and Cr-substituted Heusler alloys: Predictions from first-principles and Monte Carlo studies. <i>Physical Review B</i> , 2015, 91, .	3.2	36
90	Large magnetocaloric effects in magnetic intermetallics: First-principles and Monte Carlo studies. <i>MATEC Web of Conferences</i> , 2015, 33, 02001.	0.2	2

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91	Martensitic transformation in shape memory crystal with defects: Monte Carlo simulations and Landau theory. <i>Physica Status Solidi (B): Basic Research</i> , 2015, 252, 2309-2316.	1.5	2
92	First Principles and Monte Carlo Calculations of Structural and Magnetic Properties of $\text{Fe}_{\text{x}}\text{Ni}_{2-\text{x}}\text{Mn}_{1+\text{y}}\text{Al}_{1-\text{y}}$ Heusler Alloys. <i>MATEC Web of Conferences</i> , 2015, 33, 05002.	0.2	0
93	Magnetic States of the $\text{Ni}_{1.75}\text{Co}_{0.25}\text{Mn}_{1.25}\text{Cr}_{0.25}\text{In}_{0.5}$ Heusler Alloy. <i>IEEE Transactions on Magnetics</i> , 2015, 51, 1-4.		
94	First-Principles Calculations of Magnetic Properties of Cr-Doped $\text{Ni}_{45}\text{Co}_5\text{Mn}_{37}\text{In}_{13}$ Heusler Alloys. <i>IEEE Transactions on Magnetics</i> , 2015, 51, 1-4.	2.1	5
95	The metamagnetic behavior and giant inverse magnetocaloric effect in $\text{Ni}_{\text{x}}\text{Co}_{\text{x}}\text{Mn}_{\text{x}}(\text{Ga}, \text{In}, \text{Sn})$ Heusler alloys. <i>Journal of Magnetism and Magnetic Materials</i> , 2015, 385, 193-197.	2.3	22
96	First Principles Investigation of Magnetic Properties of Fe-Ni-Mn-Al Heusler Alloys. <i>Physics Procedia</i> , 2015, 75, 1427-1434.	1.2	7
97	First Principles Investigations of Structural and Magnetic Properties of Fe-Ni-Mn-Al Heusler Alloys. <i>Solid State Phenomena</i> , 2015, 233-234, 187-191.	0.3	3
98	Predictions of a Large Magnetocaloric Effect in Co- and Cr-Substituted Heusler Alloys Using First-Principles and Monte Carlo Approaches. <i>Physics Procedia</i> , 2015, 75, 1381-1388.	1.2	3
99	First principles investigation of structural and magnetic properties of $\text{Ni}_{\text{x}}\text{Co}_{\text{x}}\text{Mn}_{\text{x}}\text{In}$ Heusler alloys. <i>Journal Physics D: Applied Physics</i> , 2015, 48, 164005.	2.8	18
100	First-principles calculations of magnetic properties of Cr-doped $\text{Ni}_{45}\text{Co}_5\text{Mn}_{37}\text{In}_{13}$ Heusler alloys. , 2015, , .		0
101	Monte Carlo Study of the Polycrystalline Ni_2MnGa Heusler Alloy. <i>Solid State Phenomena</i> , 2015, 233-234, 251-254.	0.3	1
102	The magnetic states of the $\text{Ni}_{1.75}\text{Co}_{0.25}\text{Mn}_{1.25}\text{Cr}_{0.25}\text{In}_{0.5}$ Heusler alloy. , 2015, , .		
103	Structural and Magnetic Properties of Mn_2NiZ ($Z = \text{Ga}, \text{In}, \text{Sn}, \text{Sb}$) Heusler Alloys from <i>Ab Initio</i> Calculations. <i>Solid State Phenomena</i> , 2015, 233-234, 229-232.	0.3	8
104	First-principles study of the structural and magnetic properties of the Ni 45 Co 5 Mn 39 Sn 11 Heusler alloy. <i>Journal of Magnetism and Magnetic Materials</i> , 2015, 383, 180-185.	2.3	7
105	Calculation of Electronic Structure and Field Induced Magnetic Collapse in Ferroic Materials. , 2015, , 405-408.		0
106	Monte Carlo and first-principles approaches for single crystal and polycrystalline Ni_2MnGa Heusler alloys. <i>Journal Physics D: Applied Physics</i> , 2014, 47, 425002.	2.8	20
107	Theoretical prediction of the spin glass behavior in the low-temperature phase of $\text{Ni}_2\text{Mn}_{1.36}\text{In}_{0.64}$ Heusler alloy. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2014, 11, 1110-1115.	0.8	1
108	Investigation of structural and magnetic properties of Heusler $\text{Fe}_{2+\text{x}}\text{Mn}_{1-\text{x}}\text{Al}$ alloys by first principles method. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2014, 11, 979-983.	0.8	3

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109	First-principles calculation of the instability leading to giant inverse magnetocaloric effects. <i>Physical Review B</i> , 2014, 89, .	3.2	73
110	Interacting magnetic cluster-spin glasses and strain glasses in Ni-Mn based Heusler structured intermetallics. <i>Physica Status Solidi (B): Basic Research</i> , 2014, 251, 2135-2148.	1.5	37
111	Tuning magnetic exchange interactions to enhance magnetocaloric effect in Ni ₅₀ Mn ₃₄ In ₁₆ Heusler alloy: Monte Carlo and ab initio studies. <i>International Journal of Refrigeration</i> , 2014, 37, 273-280.	3.4	14
112	Ab Initio and Monte Carlo Approaches For the Magnetocaloric Effect in Co- and In-Doped Ni-Mn-Ga Heusler Alloys. <i>Entropy</i> , 2014, 16, 4992-5019.	2.2	40
113	Magnetocaloric and magnetic properties of Ni ₂ Mn _{1-x} Cu _x Ga Heusler alloys: An insight from the direct measurements and ab initio and Monte Carlo calculations. <i>Journal of Applied Physics</i> , 2013, 114, .	2.5	30
114	Complex magnetic ordering as a driving mechanism of multifunctional properties of Heusler alloys from first principles. <i>European Physical Journal B</i> , 2013, 86, 1.	1.5	88
115	Publisher's Note: First-principles investigation of chemical and structural disorder in magnetic Ni ₂ Mn _{1+x} Sn _{1-x} Heusler alloys [Phys. Rev. B86, 134418 (2012)]. <i>Physical Review B</i> , 2013, 87, .	3.2	3
116	Quaternary Ni-Mn-In-Y Heusler alloys: a way to achieve materials with better magnetocaloric properties?. <i>Journal Physics D: Applied Physics</i> , 2013, 46, 305003.	2.8	24
117	Optimization of smart Heusler alloys from first principles. <i>Journal of Alloys and Compounds</i> , 2013, 577, S107-S112.	5.5	46
118	Theoretical treatment and direct measurements of magnetocaloric effect in Ni _{2.19} xFe _x Mn _{0.81} Ga Heusler alloys. <i>Journal of Magnetism and Magnetic Materials</i> , 2013, 343, 6-12.	2.3	14
119	Ab initio study of magnetic properties of Fe-Mn-Al Heusler alloys. <i>Materials Research Society Symposia Proceedings</i> , 2013, 1581, 1.	0.1	1
120	The Supercell Scaling Investigation of Magnetic Properties in Ni-Mn-X (X=Ga, In, Sn, Sb) Heusler Alloys by Means of First-principles Methods. <i>Materials Research Society Symposia Proceedings</i> , 2013, 1581, 1.	0.1	0
121	Ab initio investigation of the structural and magnetic properties of Ni-Pt-Mn-Ga alloys. <i>Materials Research Society Symposia Proceedings</i> , 2013, 1581, 1.	0.1	1
122	Interaction of phase transformation and magneto- and elastocaloric properties of Heusler alloys. <i>Materials Research Society Symposia Proceedings</i> , 2013, 1581, 1.	0.1	1
123	Theoretical Study of Magnetic Properties and Twin Boundary Motion in Heusler Ni-Mn-X Shape Memory Alloys Using First Principles and Monte Carlo Method. <i>Advances in Science and Technology</i> , 2012, 78, 7-12.	0.2	1
124	First-principles investigations of caloric effects in ferroic materials. , 2012, , .		6
125	Phase Diagrams of Conventional and Inverse Functional Magnetic Heusler Alloys: New Theoretical and Experimental Investigations. <i>Springer Series in Materials Science</i> , 2012, , 19-47.	0.6	6
126	Monte Carlo Study of the Magnetic and Magnetocaloric Properties of La _{1-x} Ca _x MnO ₃ (x = 0.33 and 0.5). <i>Solid State Phenomena</i> , 2012, 190, 347-350.	0.3	7

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145	Modelling the phase diagram of magnetic shape memory Heusler alloys. Journal Physics D: Applied Physics, 2006, 39, 865-889.	2.8	306
146	Phase transitions in $Ni_{2+x}Mn_{1-x}Ga$ with a high Ni excess. Physical Review B, 2005, 72, .	3.2	176
147	Structural and magnetic phase transitions in shape-memory alloys $Ni_{2+x}Mn_{1-x}Ga$. Physical Review B, 1999, 59, 1113-1120.	3.2	401
148	Monte-Carlo Calculation of the Magnetocaloric Effect in Ni-Mn-Ga Alloys. Solid State Phenomena, 0, 152-153, 493-496.	0.3	1
149	Theoretical Modeling of Magnetocaloric Effect in Heusler Ni-Mn-In Alloy by Monte Carlo Study. Materials Science Forum, 0, 635, 137-142.	0.3	7
150	Theoretical Study of Twin Boundary Motion in Heusler Ni-Mn-Ga Alloys Using Monte Carlo Method. Solid State Phenomena, 0, 190, 327-330.	0.3	0
151	Theoretical Study of Magnetic Properties and Multiple Twin Boundary Motion in Heusler Ni-Mn-X Shape Memory Alloys Using First Principles and Monte Carlo Method. Materials Science Forum, 0, 738-739, 461-467.	0.3	0
152	<i>Ab Initio</i> Study of Magnetic Properties and Phase Diagram of Ni-Mn-Ga Heusler Alloys. Materials Science Forum, 0, 738-739, 473-477.	0.3	3
153	First Principles Calculations of Magnetic Exchange Parameters of Fe-Mn-Al Heusler Alloys. Solid State Phenomena, 0, 215, 131-136.	0.3	3
154	Novel Achievements in the Research Field of Multifunctional Shape Memory Ni-Mn-In and Ni-Mn-In-Z Heusler Alloys. Materials Science Foundations, 0, 81-82, 38-76.	0.2	15
155	First Principles Study of the Structural and Magnetic Properties of Cr-Doped $Ni_{1.75}Co_{0.25}Mn_{1.5}In_{0.5}$ Heusler Alloys. Materials Science Forum, 0, 845, 138-141.	0.2	6
156	Ternary Diagrams of Ni-Mn-Ga from First Principles. Materials Science Forum, 0, 845, 130-133.	0.3	1
157	Density of States of Co- and Cr-Doped $Ni_{2.0}Mn_{1.5}Sn_{0.5}$ Heusler Alloys. Materials Science Forum, 0, 845, 162-165.	0.3	0