

Ferenc Tasnadi

List of Publications by Year in descending order

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2,722
citations

147566

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76
all docs

76
docs citations

76
times ranked

2461
citing authors

#	ARTICLE	IF	CITATIONS
1	Predicting elastic properties of hard-coating alloys using ab-initio and machine learning methods. Npj Computational Materials, 2022, 8, .	3.5	16
2	The effect of strain and pressure on the electron-phonon coupling and superconductivity in MgB ₂ —Benchmark of theoretical methodologies and outlook for nanostructure design. Journal of Applied Physics, 2022, 131, 063902.	1.1	1
3	Materials synthesis at terapascal static pressures. Nature, 2022, 605, 274-278.	13.7	35
4	High-Pressure Synthesis of Dirac Materials: Layered van der Waals Bonded BeN_4 Polymorph. Physical Review Letters, 2021, 126, 175501.	2.9	90
5	Temperature-dependent elastic properties of binary and multicomponent high-entropy refractory carbides. Materials and Design, 2021, 204, 109634.	3.3	26
6	Finite-temperature interplay of structural stability, chemical complexity, and elastic properties of bcc multicomponent alloys from ab initio trained machine-learning potentials. Physical Review Materials, 2021, 5, .	0.9	9
7	Realization of an Ideal Cairo Tessellation in Nickel Diazenide NiN_2 : High-Pressure Route to Pentagonal 2D Materials. ACS Nano, 2021, 15, 13539-13546.	7.3	55
8	Efficient prediction of elastic properties of $\text{Ti}_0.5\text{Al}_0.5\text{N}$ at elevated temperature using machine learning interatomic potential. Thin Solid Films, 2021, 737, 138927.	0.8	4
9	Accurate prediction of high-temperature elastic constants of $\text{Ti}_0.5\text{Al}_0.5\text{N}$ random alloy. Thin Solid Films, 2021, 735, 138872.	0.8	4
10	Thermodynamic and electronic properties of ReN_2 polymorphs at high pressure. Physical Review B, 2021, 104, .	1.1	1
11	Achieving low elastic moduli of bcc Ti-V alloys in vicinity of mechanical instability. AIP Advances, 2020, 10, 105322.	0.6	3
12	Temperature dependence of the Kohn anomaly in bcc Nb from first-principles self-consistent phonon calculations. Physical Review B, 2020, 101, .	1.1	11
13	High-Pressure Synthesis of Metal-Inorganic Frameworks Hf_4N_{20} , WN_8 , and Os_5N_{28} with Polymeric Nitrogen Linkers. Angewandte Chemie - International Edition, 2020, 59, 10321-10326.	7.2	36
14	High-Pressure Synthesis of Metal-Inorganic Frameworks Hf_4N_{20} , WN_8 , and Os_5N_{28} with Polymeric Nitrogen Linkers. Angewandte Chemie, 2020, 132, 10407-10412.	1.6	8
15	High-Pressure Synthesis of Metal-Inorganic Frameworks Hf_4N_{20} , WN_8 , and Os_5N_{28} with Polymeric Nitrogen Linkers (Angew. Chem.) Tj ETQq1 1 0.784314	1.6	8
16	Elinvar effect in $\hat{\Gamma}^2$ -Ti simulated by on-the-fly trained moment tensor potential. New Journal of Physics, 2020, 22, 113005.	1.2	20
17	Strengthening, transformation toughening, and fracture dynamics of rocksalt-structure TiO_2 thin films. Physical Review Applied, 2020, 12, 041101.	1.2	20
18	High-pressure synthesis of ultraincompressible hard rhenium nitride pernitride $\text{Re}_2(\text{N}_2)(\text{N})_2$ stable at ambient conditions. Nature Communications, 2019, 10, 2994.	5.8	65

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19	Thermal expansion of quaternary nitride coatings. Journal of Physics Condensed Matter, 2018, 30, 135901.	0.7	5
20	High temperature thermodynamics of spinodal decomposition in arc deposited Ti _x Nb _y Al _z N coatings. Materials and Design, 2018, 150, 165-170.	3.3	7
21	Assessing the SCAN functional for itinerant electron ferromagnets. Physical Review B, 2018, 98, .	1.1	64
22	High-Pressure Synthesis of a Nitrogen-Rich Inclusion Compound ReN ₈ with Conjugated Polymeric Nitrogen Chains. Angewandte Chemie - International Edition, 2018, 57, 9048-9053.	7.2	70
23	High-Pressure Synthesis of a Nitrogen-Rich Inclusion Compound ReN ₈ with Conjugated Polymeric Nitrogen Chains. Angewandte Chemie, 2018, 130, 9186-9191.	1.6	16
24	Fe-N system at high pressure reveals a compound featuring polymeric nitrogen chains. Nature Communications, 2018, 9, 2756.	5.8	153
25	Non-equilibrium vacancy formation energies in metastable alloys – A case study of Ti _{0.5} Al _{0.5} N. Materials and Design, 2017, 114, 484-493.	3.3	13
26	Systematic ab initio investigation of the elastic modulus in quaternary transition metal nitride alloys and their coherent multilayers. Acta Materialia, 2017, 127, 124-132.	3.8	44
27	Exploring the high entropy alloy concept in (AlTiVNbCr)N. Thin Solid Films, 2017, 636, 346-352.	0.8	27
28	Topological transitions of the Fermi surface of osmium under pressure: an LDA+DMFT study. New Journal of Physics, 2017, 19, 033020.	1.2	10
29	Strong piezoelectric response in stable TiZnN ₂ , ZrZnN ₂ , and HfZnN ₂ found by ab initio high-throughput approach. Journal of Applied Physics, 2016, 120, .	1.1	17
30	Effects of configurational disorder on the elastic properties of icosahedral boron-rich alloys based on B ₆ O, B ₁₃ C ₂ , and B ₄ C, and their mixing thermodynamics. Journal of Chemical Physics, 2016, 144, 134503.	1.2	18
31	Growth and thermal stability of TiN/ZrAlN: Effect of internal interfaces. Acta Materialia, 2016, 121, 396-406.	3.8	44
32	Coherency effects on the mixing thermodynamics of cubic Ti _{1-x} N _x TiN(001) multilayers. Physical Review B, 2016, 93, .	1.1	10
33	Ab initio calculations and experimental study of piezoelectric YIn _{1-x} N thin films deposited using reactive magnetron sputter epitaxy. Acta Materialia, 2016, 105, 199-206.	3.8	20
34	N and Ti adatom dynamics on stoichiometric polar TiN(111) surfaces. Surface Science, 2016, 649, 72-79.	0.8	32
35	Large piezoelectric response of quaternary wurtzite nitride alloys and its physical origin from first principles. Physical Review B, 2015, 92, .	1.1	41
36	Temperature-dependent elastic properties of Ti _{1-x} Al _x N alloys. Applied Physics Letters, 2015, 107, .	1.5	46

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37	Stabilization of wurtzite $\text{Sc}_{0.4}\text{Al}_{0.6}\text{N}$ in pseudomorphic epitaxial $\text{ScAl}_x\text{N}/\text{InAl}_x\text{N}$ superlattices. <i>Acta Materialia</i> , 2015, 94, 101-110.	3.8	19
38	Special quasirandom structure method in application for advanced properties of alloys: A study on $\text{Ti}_{0.5}\text{Al}_{0.5}\text{N}$ and $\text{TiN}/\text{Ti}_{0.5}\text{Al}_{0.5}\text{N}$ multilayer. <i>Computational Materials Science</i> , 2015, 103, 194-199.	1.4	9
39	High temperature phase decomposition in $\text{Ti}_x\text{Zr}_y\text{Al}_z\text{N}$. <i>AIP Advances</i> , 2014, 4, .	0.6	13
40	Macroscopic elastic properties of textured ZrN-AlN polycrystalline aggregates: From <i>ab initio</i> calculations to grain-scale interactions. <i>Physical Review B</i> , 2014, 90, .	1.1	34
41	Anomalous epitaxial stability of (001) interfaces in ZrN/SiN_x multilayers. <i>APL Materials</i> , 2014, 2, 046106.	2.2	10
42	Effect of Al substitution on Ti, Al, and N adatom dynamics on $\text{TiN}(001)$, (011), and (111) surfaces. <i>Surface Science</i> , 2014, 630, 28-40.	0.8	37
43	Importance of Correlation Effects in hcp Iron Revealed by a Pressure-Induced Electronic Topological Transition. <i>Physical Review Letters</i> , 2013, 110, 117206.	2.9	58
44	Temperature dependence of TiN elastic constants from <i>ab initio</i> molecular dynamics simulations. <i>Physical Review B</i> , 2013, 87, .	1.1	78
45	Systematic theoretical search for alloys with increased thermal stability for advanced hard coatings applications. <i>New Journal of Physics</i> , 2013, 15, 095010.	1.2	15
46	Volume matching condition to establish the enhanced piezoelectricity in ternary $(\text{Sc},\text{Y})_{0.5}(\text{Al},\text{Ga},\text{In})_{0.5}$ alloys. <i>Physical Review B</i> , 2013, 87, .	1.1	53
47	High pressure and high temperature stabilization of cubic AlN in $\text{Ti}_{0.6}\text{Al}_{0.4}\text{N}$. <i>Journal of Applied Physics</i> , 2013, 113, .	1.1	34
48	<i>Ab initio</i> elastic tensor of cubic $\text{TiAl}_{0.5}\text{N}$ alloys: Dependence of elastic constants on size and shape of the supercell model and their convergence. <i>Physical Review B</i> , 2013, 87, .	1.1	125
49	Elastic constants, composition and piezoelectric polarization in $\text{InAl}_{0.5}\text{N}$ surfaces from first principles. <i>Physical Review B</i> , 2013, 87, .	1.1	33
50	From <i>ab initio</i> calculations to experimental implications for the applicability of Vegard's rule. First-principles study of the $\text{SiN}/\text{TiN}(001)$ interface. <i>Physical Review B</i> , 2012, 85, .	1.1	31
51	From <i>ab initio</i> calculations to experimental implications for the applicability of Vegard's rule. First-principles study of the $\text{SiN}/\text{TiN}(001)$ interface. <i>Physical Review B</i> , 2012, 85, .	1.1	17
52	$\text{Y}_x\text{Al}_{1-x}\text{N}$ thin films. <i>Journal Physics D: Applied Physics</i> , 2012, 45, 422001.	1.3	42
53	Density functional investigation of rhombohedral stacks of graphene: Topological surface states, nonlinear dielectric response, and bulk limit. <i>Physical Review B</i> , 2011, 84, .	1.1	43
54	Phase Stability and Elasticity of TiAlN . <i>Materials</i> , 2011, 4, 1599-1618.	1.3	80

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55	Improving thermal stability of hard coating films via a concept of multicomponent alloying. Applied Physics Letters, 2011, 99, .	1.5	95
56	Significant elastic anisotropy in $Ti_{1-x}Al_xN$ alloys. Applied Physics Letters, 2010, 97, .	1.5	107
57	Origin of the Anomalous Piezoelectric Response in Wurtzite $Sc_xAl_{1-x}N$ Alloys. Physical Review Letters, 2010, 104, 137601.	2.9	305
58	Increased electromechanical coupling in $Sc_xAl_{1-x}N$. Applied Physics Letters, 2010, 97, .	1.5	149
59	Elastic properties and electrostructural correlations in ternary scandium-based cubic inverse perovskites: A first-principles study. Physical Review B, 2009, 79, .	1.1	87
60	Stability of the ternary perovskites Sc_3		

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73	Local self-interaction-free approximate exchange-correlation potentials in the variational density functional theory for individual excited states. <i>Chemical Physics Letters</i> , 2002, 366, 496-503.	1.2	5