Vladan StevanoviÄ

List of Publications by Year
in descending order

50	4,859	172457	197818
papers	citations	29 h-index	49 g-index

51	51	51	7563
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Predicting Oxygen Off-Stoichiometry and Hydrogen Incorporation in Complex Perovskite Oxides. Chemistry of Materials, 2022, 34, 510-518.	6.7	7
2	Role of disorder in the synthesis of metastable zinc zirconium nitrides. Physical Review Materials, 2022, 6, .	2.4	14
3	Triple ionicâ€“electronic conducting oxides for next-generation electrochemical devices. Nature Materials, 2021, 20, 301-313.	27.5	160
4	Computationally Guided Discovery of Axis-Dependent Conduction Polarity in NaSnAs Crystals. Chemistry of Materials, 2021, 33, 946-951.	6.7	13
5	Computational Fermi level engineering and doping-type conversion of Mg:Ga2O3 via three-step synthesis process. Journal of Applied Physics, 2021, 129, .	2.5	14
6	Predicting energy and stability of known and hypothetical crystals using graph neural network. Patterns, 2021, 2, 100361.	5.9	16
7	Metastable materials discovery in the age of large-scale computation. Applied Physics Reviews, 2021, 8, .	11.3	16
8	Devil is in the Defects: Electronic Conductivity in Solid Electrolytes. Chemistry of Materials, 2021, 33, 7484-7498.	6.7	49
9	Theoretical Insights for Improving the Schottky-Barrier Height at the <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline" overflow="scroll"><mml:msub><mml:mi>Ga</mml:mi><mml:mn>2</mml:mn></mml:msub><mml:msub><mml:mrow><mml:mi mathvariant="normal">O</mml:mi></mml:mrow><mml:mn>3</mml:mn></mml:msub><mml:mrow><mml:mo>/</mml:mo><mml:mi>P Interface. Physical Review Applied, 2021, 16, .	3.8	5
10	The glassy solid as a statistical ensemble of crystalline microstates. Npj Computational Materials, 2020, 6, .	8.7	14
11	Minimization of Atomic Displacements as a Guiding Principle of the Martensitic Phase Transformation. Physical Review Letters, 2020, 125, 125502.	7.8	8
12	Matching crystal structures atom-to-atom. Journal of Chemical Physics, 2020, 152, 074106.	3.0	16
13	A simple chemical guide for finding novel n-type dopable Zintl pnictide thermoelectric materials. Journal of Materials Chemistry A, 2019, 7, 19385-19395.	10.3	29
14	New frontiers for the materials genome initiative. Npj Computational Materials, 2019, 5, .	8.7	312
15	The importance of phase equilibrium for doping efficiency: iodine doped PbTe. Materials Horizons, 2019, 6, 1444-1453.	12.2	42
16	Polycrystalline <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mi>ZrTe</mml:mi></mml:mrow><mml:mrow><mml:mn>5</mml:mn></mml:mrow></mml:msub></mml:mrow></mml:math> Parametrized as a Narrow-Band-Gap Semiconductor for Thermoelectric Performance. Physical Review Applied, 2018, 9, .	3.8	26
17	Origin of Pronounced Nonlinear Band Gap Behavior in Leadâ€“Tin Hybrid Perovskite Alloys. Chemistry of Materials, 2018, 30, 3920-3928.	6.7	166
18	Enhanced Piezoelectric Response of AlN via CrN Alloying. Physical Review Applied, 2018, 9, .	3.8	57

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19	Experimental and computational phase boundary mapping of Co₄Sn₆Te₆. Journal of Materials Chemistry A, 2018, 6, 24175-24185.	10.3	26
20	Physical descriptor for the Gibbs energy of inorganic crystalline solids and temperature-dependent materials chemistry. Nature Communications, 2018, 9, 4168.	12.8	152
21	Structural and Chemical Features Giving Rise to Defect Tolerance of Binary Semiconductors. Chemistry of Materials, 2018, 30, 5583-5592.	6.7	36
22	Predicting kinetics of polymorphic transformations from structure mapping and coordination analysis. Physical Review Materials, 2018, 2, .	2.4	13
23	Metastable rocksalt ZnO is <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mi>p</mml:mi></mml:math> -type dopable. Physical Review Materials, 2018, 2, .	2.4	15
24	Capturing Anharmonicity in a Lattice Thermal Conductivity Model for High-Throughput Predictions. Chemistry of Materials, 2017, 29, 2494-2501.	6.7	88
25	A computational framework for automation of point defect calculations. Computational Materials Science, 2017, 130, 1-9.	3.0	131
26	Thermoelectric Performance and Defect Chemistry in n-Type Zintl KGaSb₄. Chemistry of Materials, 2017, 29, 4523-4534.	6.7	59
27	Searching for â€œDefect-Tolerantâ€•Photovoltaic Materials: Combined Theoretical and Experimental Screening. Chemistry of Materials, 2017, 29, 4667-4674.	6.7	275
28	Raman spectroscopy and x-ray diffraction of <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:msup><mml:mrow><mml:mi mathvariant="italic">sp</mml:mi></mml:mrow><mml:mn>3</mml:mn></mml:msup><mml:mspace width="4pt" /><mml:mi>CaC</mml:mi><mml:msub><mml:mi mathvariant="normal">O</mml:mi><mml:mn>3</mml:mn></mml:msub></mml:mrow></mml:math> at lower mantle pressures. Physical Review B, 2017, 96, .	3.2	54
29	Large Area Atomically Flat Surfaces via Exfoliation of Bulk Bi₂Se₃ Single Crystals. Chemistry of Materials, 2017, 29, 8472-8477.	6.7	8
30	SnO as a potential oxide thermoelectric candidate. Journal of Materials Chemistry C, 2017, 5, 8854-8861.	5.5	72
31	Polymorphism in elemental silicon: Probabilistic interpretation of the realizability of metastable structures. Physical Review B, 2017, 96, .	3.2	14
32	Tuning the piezoelectric and mechanical properties of the AlN system via alloying with YN and BN. Journal of Applied Physics, 2017, 122, .	2.5	49
33	Computational identification of promising thermoelectric materials among known quasi-2D binary compounds. Journal of Materials Chemistry A, 2016, 4, 11110-11116.	10.3	55
34	Defect Tolerance in Methylammonium Lead Triiodide Perovskite. ACS Energy Letters, 2016, 1, 360-366.	17.4	500
35	Synthesis of a mixed-valent tin nitride and considerations of its possible crystal structures. Journal of Chemical Physics, 2016, 144, 144201.	3.0	29
36	Energy conversion properties of ZnSiP<inf>2</inf>, a lattice-matched material for silicon-based tandem photovoltaics. , 2016, , .		2

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37	Sampling Polymorphs of Ionic Solids using Random Superlattices. Physical Review Letters, 2016, 116, 075503.	7.8	39
38	Thermoelectricity in transition metal compounds: the role of spin disorder. Physical Chemistry Chemical Physics, 2016, 18, 31777-31786.	2.8	27
39	Discovery of a ternary pseudobrookite phase in the earth-abundant Tiâ€“Znâ€“O system. Dalton Transactions, 2016, 45, 1572-1581.	3.3	6
40	Identifying defect-tolerant semiconductors with high minority-carrier lifetimes: beyond hybrid lead halide perovskites. MRS Communications, 2015, 5, 265-275.	1.8	662
41	Hybrid Organicâ€“Inorganic Perovskites (HOIPs): Opportunities and Challenges. Advanced Materials, 2015, 27, 5102-5112.	21.0	372
42	Semiconducting properties of spinel tin nitride and other IV₃N₄ polymorphs. Journal of Materials Chemistry C, 2015, 3, 1389-1396.	5.5	49
43	Control of the Electrical Properties in Spinel Oxides by Manipulating the Cation Disorder. Advanced Functional Materials, 2014, 24, 610-618.	14.9	109
44	Non-equilibrium deposition of phase pure Cu2O thin films at reduced growth temperature. APL Materials, 2014, 2, .	5.1	55
45	Defect Tolerant Semiconductors for Solar Energy Conversion. Journal of Physical Chemistry Letters, 2014, 5, 1117-1125.	4.6	304
46	Convergence of density and hybrid functional defect calculations for compound semiconductors. Physical Review B, 2013, 88, .	3.2	94
47	Correcting density functional theory for accurate predictions of compound enthalpies of formation: Fitted elemental-phase reference energies. Physical Review B, 2012, 85, .	3.2	454
48	Large insulating gap in topological insulators induced by negative spin-orbit splitting. Physical Review B, 2012, 86, .	3.2	26
49	Universal Electrostatic Origin of Cation Ordering in A₂BO₄ Spinel Oxides. Journal of the American Chemical Society, 2011, 133, 11649-11654.	13.7	71
50	Simple Point-Ion Electrostatic Model Explains the Cation Distribution in Spinel Oxides. Physical Review Letters, 2010, 105, 075501.	7.8	48