

# Alexander Urban

## List of Publications by Year in descending order

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Version: 2024-02-01

39  
papers

5,183  
citations

236612

25  
h-index

315357

38  
g-index

39  
all docs

39  
docs citations

39  
times ranked

5489  
citing authors



#	ARTICLE	IF	CITATIONS
19	Influence of Inversion on Mg Mobility and Electrochemistry in Spinel. <i>Chemistry of Materials</i> , 2017, 29, 7918-7930.	3.2	75
20	First-Principles Simulation of the (Li <sup>+</sup> Ni <sup>+</sup> Vacancy)O Phase Diagram and Its Relevance for the Surface Phases in Ni-Rich Li-Ion Cathode Materials. <i>Chemistry of Materials</i> , 2017, 29, 7840-7851.	3.2	79
21	Efficient and accurate machine-learning interpolation of atomic energies in compositions with many species. <i>Physical Review B</i> , 2017, 96, .	1.1	228
22	Construction of ground-state preserving sparse lattice models for predictive materials simulations. <i>Npj Computational Materials</i> , 2017, 3, .	3.5	15
23	Computational Design and Preparation of Cation-Disordered Oxides for High-Energy-Density Li-Ion Batteries. <i>Advanced Energy Materials</i> , 2016, 6, 1600488.	10.2	93
24	The Effect of Cation Disorder on the Average Li Intercalation Voltage of Transition-Metal Oxides. <i>Chemistry of Materials</i> , 2016, 28, 3659-3665.	3.2	62
25	Lithium Batteries: Computational Design and Preparation of Cation-Disordered Oxides for High-Energy-Density Li-Ion Batteries ( <i>Adv. Energy Mater.</i> 15/2016). <i>Advanced Energy Materials</i> , 2016, 6, .	10.2	0
26	Understanding the Effect of Cation Disorder on the Voltage Profile of Lithium Transition-Metal Oxides. <i>Chemistry of Materials</i> , 2016, 28, 5373-5383.	3.2	79
27	Computational understanding of Li-ion batteries. <i>Npj Computational Materials</i> , 2016, 2, .	3.5	411
28	Finding and proving the exact ground state of a generalized Ising model by convex optimization and MAX-SAT. <i>Physical Review B</i> , 2016, 94, .	1.1	25
29	The structural and chemical origin of the oxygen redox activity in layered and cation-disordered Li-excess cathode materials. <i>Nature Chemistry</i> , 2016, 8, 692-697.	6.6	1,022
30	An implementation of artificial neural-network potentials for atomistic materials simulations: Performance for TiO <sub>2</sub> . <i>Computational Materials Science</i> , 2016, 114, 135-150.	1.4	377
31	Calibrating transition-metal energy levels and oxygen bands in first-principles calculations: Accurate prediction of redox potentials and charge transfer in lithium transition-metal oxides. <i>Physical Review B</i> , 2015, 92, .	1.1	126
32	The Intercalation Phase Diagram of Mg in V <sub>2</sub> O <sub>5</sub> from First-Principles. <i>Chemistry of Materials</i> , 2015, 27, 3733-3742.	3.2	130
33	A disordered rock-salt Li-excess cathode material with high capacity and substantial oxygen redox activity: Li <sub>1.25</sub> Nb <sub>0.25</sub> Mn <sub>0.5</sub> O <sub>2</sub> . <i>Electrochemistry Communications</i> , 2015, 60, 70-73.	2.3	145
34	Designing New Lithium-Excess Cathode Materials from Percolation Theory: Nanohighways in Li <sub>x</sub> Ni <sub>2</sub> Sb <sub>x</sub> O <sub>2</sub> . <i>Nano Letters</i> , 2015, 15, 596-602.	4.5	54
35	Electrodes: The Configurational Space of Rocksalt-Type Oxides for High-Capacity Lithium Battery Electrodes ( <i>Adv. Energy Mater.</i> 13/2014). <i>Advanced Energy Materials</i> , 2014, 4, n/a-n/a.	10.2	3
36	The Configurational Space of Rocksalt-Type Oxides for High-Capacity Lithium Battery Electrodes. <i>Advanced Energy Materials</i> , 2014, 4, 1400478.	10.2	256

#	ARTICLE	IF	CITATIONS
37	Unlocking the Potential of Cation-Disordered Oxides for Rechargeable Lithium Batteries. Science, 2014, 343, 519-522.	6.0	943
38	Growth of One-Dimensional Pd Nanowires on the Terraces of a Reduced SnO <sub>2</sub> (101) Surface. Physical Review Letters, 2007, 98, 186102.	2.9	16
39	Tuning the Reactivity of Oxide Surfaces by Charge-Accepting Adsorbates. Angewandte Chemie - International Edition, 2007, 46, 7315-7318.	7.2	53