

Alexander Urban

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Stacking-Fault Enhanced Oxygen Redox in Li_2MnO_3 . <i>Advanced Energy Materials</i> , 2022, 12, .	19.5	17
2	Data-driven approach to parameterize $\langle \text{mml:math} \text{xmlns:mml="http://www.w3.org/1998/Math/MathML"} \rangle \langle \text{mml:mi} \text{SCAN} \langle \text{mml:mi} \rangle \langle \text{mml:mo} \rangle + \langle \text{mml:mo} \rangle \langle \text{mml:mi} \rangle \text{U} \langle \text{mml:mi} \rangle \langle \text{mml:math} \text{xmlns:mml="http://www.w3.org/1998/Math/MathML"} \rangle$ for an accurate description of $\langle \text{mml:math} \text{xmlns:mml="http://www.w3.org/1998/Math/MathML"} \rangle \langle \text{mml:mrow} \rangle \langle \text{mml:mn} \rangle 3 \langle \text{mml:mn} \rangle \langle \text{mml:mi} \rangle \text{d} \langle \text{mml:mi} \rangle \langle \text{mml:mrow} \rangle \langle \text{mml:math} \text{xmlns:mml="http://www.w3.org/1998/Math/MathML"} \rangle$ transition metal oxide thermochemistry. <i>Physical Review Materials</i> , 2022, 6, .	2.4	6
3	Understanding the Onset of Surface Degradation in LiNiO_2 Cathodes. <i>ACS Applied Energy Materials</i> , 2022, 5, 5730-5741.	5.1	10
4	Accelerated Atomistic Modeling of Solid-State Battery Materials With Machine Learning. <i>Frontiers in Energy Research</i> , 2021, 9, .	2.3	25
5	Strategies for the construction of machine-learning potentials for accurate and efficient atomic-scale simulations. <i>Machine Learning: Science and Technology</i> , 2021, 2, 031001.	5.0	42
6	Anisotropic to Isotropic Transition in Monolayer Group-IV Tellurides. <i>Materials</i> , 2021, 14, 4495.	2.9	4
7	Realizing continuous cation order-to-disorder tuning in a class of high-energy spinel-type Li-ion cathodes. <i>Matter</i> , 2021, 4, 3897-3916.	10.0	32
8	Augmenting zero-Kelvin quantum mechanics with machine learning for the prediction of chemical reactions at high temperatures. <i>Nature Communications</i> , 2021, 12, 7012.	12.8	10
9	Effect of fluorination and Li-excess on the Li migration barrier in Mn-based cathode materials. <i>Journal of Materials Chemistry A</i> , 2020, 8, 19965-19974.	10.3	20
10	Potential and pH Dependence of the Buried Interface of Membrane-Coated Electrocatalysts. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 52125-52135.	8.0	2
11	Ultrahigh power and energy density in partially ordered lithium-ion cathode materials. <i>Nature Energy</i> , 2020, 5, 213-221.	39.5	158
12	Understanding the Origin of Higher Capacity for Ni-Based Disordered Rock-Salt Cathodes. <i>Chemistry of Materials</i> , 2020, 32, 3447-3461.	6.7	16
13	Effect of Fluorination on Lithium Transport and Short-Range Order in Disordered Rock-Salt Type Lithium-Ion Battery Cathodes. <i>Advanced Energy Materials</i> , 2020, 10, 1903240.	19.5	83
14	Efficient training of ANN potentials by including atomic forces via Taylor expansion and application to water and a transition-metal oxide. <i>Npj Computational Materials</i> , 2020, 6, .	8.7	40
15	Hidden structural and chemical order controls lithium transport in cation-disordered oxides for rechargeable batteries. <i>Nature Communications</i> , 2019, 10, 592.	12.8	162
16	Constructing first-principles phase diagrams of amorphous Li_xSi using machine-learning-assisted sampling with an evolutionary algorithm. <i>Journal of Chemical Physics</i> , 2018, 148, 241711.	3.0	121
17	Stoichiometric Layered Potassium Transition Metal Oxide for Rechargeable Potassium Batteries. <i>Chemistry of Materials</i> , 2018, 30, 6532-6539.	6.7	108
18	Electronic-Structure Origin of Cation Disorder in Transition-Metal Oxides. <i>Physical Review Letters</i> , 2017, 119, 176402.	7.8	135

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19	Influence of Inversion on Mg Mobility and Electrochemistry in Spinel. <i>Chemistry of Materials</i> , 2017, 29, 7918-7930.	6.7	75
20	First-Principles Simulation of the (Li ⁺ Ni ⁺ 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.	6.7	79
21	Efficient and accurate machine-learning interpolation of atomic energies in compositions with many species. <i>Physical Review B</i> , 2017, 96, .	3.2	228
22	Construction of ground-state preserving sparse lattice models for predictive materials simulations. <i>Npj Computational Materials</i> , 2017, 3, .	8.7	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.	19.5	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.	6.7	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, .	19.5	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.	6.7	79
27	Computational understanding of Li-ion batteries. <i>Npj Computational Materials</i> , 2016, 2, .	8.7	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, .	3.2	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.	13.6	1,022
30	An implementation of artificial neural-network potentials for atomistic materials simulations: Performance for TiO ₂ . <i>Computational Materials Science</i> , 2016, 114, 135-150.	3.0	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, .	3.2	126
32	The Intercalation Phase Diagram of Mg in V ₂ O ₅ from First-Principles. <i>Chemistry of Materials</i> , 2015, 27, 3733-3742.	6.7	130
33	A disordered rock-salt Li-excess cathode material with high capacity and substantial oxygen redox activity: Li _{1.25} Nb _{0.25} Mn _{0.5} O ₂ . <i>Electrochemistry Communications</i> , 2015, 60, 70-73.	4.7	145
34	Designing New Lithium-Excess Cathode Materials from Percolation Theory: Nanohighways in Li _x Ni ₂ Sb _x O ₂ . <i>Nano Letters</i> , 2015, 15, 596-602.	9.1	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.	19.5	3
36	The Configurational Space of Rocksalt-Type Oxides for High-Capacity Lithium Battery Electrodes. <i>Advanced Energy Materials</i> , 2014, 4, 1400478.	19.5	256

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37	Unlocking the Potential of Cation-Disordered Oxides for Rechargeable Lithium Batteries. Science, 2014, 343, 519-522.	12.6	943
38	Growth of One-Dimensional Pd Nanowires on the Terraces of a Reduced SnO ₂ (101) Surface. Physical Review Letters, 2007, 98, 186102.	7.8	16
39	Tuning the Reactivity of Oxide Surfaces by Charge-Accepting Adsorbates. Angewandte Chemie - International Edition, 2007, 46, 7315-7318.	13.8	53