

Itai Leven

List of Publications by Year in descending order

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Version: 2024-02-01

23
papers

1,029
citations

567281

15
h-index

752698

20
g-index

24
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docs citations

24
times ranked

1207
citing authors

#	ARTICLE	IF	CITATIONS
1	Robust Superlubricity in Graphene/h-BN Heterojunctions. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 115-120.	4.6	184
2	Interlayer Potential for Graphene/h-BN Heterostructures. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 2896-2905.	5.3	107
3	Coherent commensurate electronic states at the interface between misoriented graphene layers. <i>Nature Nanotechnology</i> , 2016, 11, 752-757.	31.5	107
4	Can electric fields drive chemistry for an aqueous microdroplet?. <i>Nature Communications</i> , 2022, 13, 280.	12.8	102
5	Ultrahigh Torsional Stiffness and Strength of Boron Nitride Nanotubes. <i>Nano Letters</i> , 2012, 12, 6347-6352.	9.1	72
6	Inter-layer potential for hexagonal boron nitride. <i>Journal of Chemical Physics</i> , 2014, 140, 104106.	3.0	72
7	Interlayer Potential for Homogeneous Graphene and Hexagonal Boron Nitride Systems: Reparametrization for Many-Body Dispersion Effects. <i>Journal of Physical Chemistry C</i> , 2017, 121, 22826-22835.	3.1	61
8	Development of an Advanced Force Field for Water Using Variational Energy Decomposition Analysis. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 5001-5013.	5.3	49
9	Multiwalled nanotube faceting unravelled. <i>Nature Nanotechnology</i> , 2016, 11, 1082-1086.	31.5	47
10	NewtonNet: a Newtonian message passing network for deep learning of interatomic potentials and forces. , 2022, 1, 333-343.		42
11	Recent Advances for Improving the Accuracy, Transferability, and Efficiency of Reactive Force Fields. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 3237-3251.	5.3	41
12	Quantifying the two-state facilitated diffusion model of protein-DNA interactions. <i>Nucleic Acids Research</i> , 2019, 47, 5530-5538.	14.5	25
13	C-GeM: Coarse-Grained Electron Model for Predicting the Electrostatic Potential in Molecules. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 6820-6826.	4.6	20
14	A Reactive Force Field with Coarse-Grained Electrons for Liquid Water. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 9240-9247.	4.6	18
15	Smallest Archimedean Screw: Facet Dynamics and Friction in Multiwalled Nanotubes. <i>Nano Letters</i> , 2017, 17, 5321-5328.	9.1	16
16	Inertial extended-Lagrangian scheme for solving charge equilibration models. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 18652-18659.	2.8	16
17	Nanotube Motion on Layered Materials: A Registry Perspective. <i>Journal of Physical Chemistry C</i> , 2016, 120, 4466-4470.	3.1	14
18	Proton Traffic Jam: Effect of Nanoconfinement and Acid Concentration on Proton Hopping Mechanism. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 25419-25427.	13.8	14

#	ARTICLE	IF	CITATIONS
19	Protein C-GeM: A Coarse-Grained Electron Model for Fast and Accurate Protein Electrostatics Prediction. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 4357-4369.	5.4	9
20	A benchmark dataset for Hydrogen Combustion. <i>Scientific Data</i> , 2022, 9, 215.	5.3	6
21	Stochastic Constrained Extended System Dynamics for Solving Charge Equilibration Models. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 5991-5998.	5.3	5
22	Proton Traffic Jam: Effect of Nanoconfinement and Acid Concentration on Proton Hopping Mechanism. <i>Angewandte Chemie</i> , 0, , .	2.0	2
23	Rücktitelbild: Proton Traffic Jam: Effect of Nanoconfinement and Acid Concentration on Proton Hopping Mechanism (<i>Angew. Chem.</i> 48/2021). <i>Angewandte Chemie</i> , 2021, 133, 25788-25788.	2.0	0