

# Arun Kumar Mannodi-Kanakkithodi

## List of Publications by Year in descending order

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Version: 2024-02-01

38  
papers

2,542  
citations

394421

19  
h-index

377865

34  
g-index

40  
all docs

40  
docs citations

40  
times ranked

2882  
citing authors

#	ARTICLE	IF	CITATIONS
1	Machine learning in materials informatics: recent applications and prospects. Npj Computational Materials, 2017, 3, .	8.7	1,013
2	Machine Learning Strategy for Accelerated Design of Polymer Dielectrics. Scientific Reports, 2016, 6, 20952.	3.3	279
3	Rational Co-Design of Polymer Dielectrics for Energy Storage. Advanced Materials, 2016, 28, 6277-6291.	21.0	149
4	Scoping the polymer genome: A roadmap for rational polymer dielectrics design and beyond. Materials Today, 2018, 21, 785-796.	14.2	143
5	A polymer dataset for accelerated property prediction and design. Scientific Data, 2016, 3, 160012.	5.3	139
6	Accelerated materials property predictions and design using motif-based fingerprints. Physical Review B, 2015, 92, .	3.2	136
7	Heterogeneous mesoporous manganese/cobalt oxide catalysts for selective oxidation of 5-hydroxymethylfurfural to 2,5-diformylfuran. Chemical Communications, 2017, 53, 11751-11754.	4.1	65
8	Poly(dimethyltin glutarate) as a Prospective Material for High Dielectric Applications. Advanced Materials, 2015, 27, 346-351.	21.0	64
9	Rational Design of Organotin Polyesters. Macromolecules, 2015, 48, 2422-2428.	4.8	54
10	Mining Materials Design Rules from Data: The Example of Polymer Dielectrics. Chemistry of Materials, 2017, 29, 9001-9010.	6.7	48
11	Critical assessment of regression-based machine learning methods for polymer dielectrics. Computational Materials Science, 2016, 125, 123-135.	3.0	45
12	Comprehensive Computational Study of Partial Lead Substitution in Methylammonium Lead Bromide. Chemistry of Materials, 2019, 31, 3599-3612.	6.7	37
13	Optimization of Organotin Polymers for Dielectric Applications. ACS Applied Materials & Interfaces, 2016, 8, 21270-21277.	8.0	33
14	Effect of selenium and chlorine co-passivation in polycrystalline CdSeTe devices. Applied Physics Letters, 2019, 115, .	3.3	33
15	Machine-learned impurity level prediction for semiconductors: the example of Cd-based chalcogenides. Npj Computational Materials, 2020, 6, .	8.7	32
16	Multi-objective optimization techniques to design the Pareto front of organic dielectric polymers. Computational Materials Science, 2016, 125, 92-99.	3.0	31
17	A rational co-design approach to the creation of new dielectric polymers with high energy density. IEEE Transactions on Dielectrics and Electrical Insulation, 2017, 24, 732-743.	2.9	26
18	Data-driven design of novel halide perovskite alloys. Energy and Environmental Science, 2022, 15, 1930-1949.	30.8	26

#	ARTICLE	IF	CITATIONS
19	Infrared-pump electronic-probe of methylammonium lead iodide reveals electronically decoupled organic and inorganic sublattices. <i>Nature Communications</i> , 2019, 10, 482.	12.8	25
20	Universal machine learning framework for defect predictions in zinc blende semiconductors. <i>Patterns</i> , 2022, 3, 100450.	5.9	22
21	Defect Energetics in Pseudo-Cubic Mixed Halide Lead Perovskites from First-Principles. <i>Journal of Physical Chemistry C</i> , 2020, 124, 16729-16738.	3.1	19
22	Computational Data-Driven Materials Discovery. <i>Trends in Chemistry</i> , 2021, 3, 79-82.	8.5	19
23	Materials Data Infrastructure and Materials Informatics. , 2018, , 193-225.		18
24	Defective Ultrathin ZnIn <sub>2</sub> S <sub>4</sub> for Photoreductive Deuteration of Carbonyls Using D <sub>2</sub> O as the Deuterium Source. <i>Advanced Science</i> , 2022, 9, e2103408.	11.2	15
25	Critical role of morphology on the dielectric constant of semicrystalline polyolefins. <i>Journal of Chemical Physics</i> , 2016, 144, 234905.	3.0	14
26	Charge Transfer Dynamics of Phase-Segregated Halide Perovskites: CH <sub>3</sub> NH <sub>3</sub> PbCl <sub>3</sub> and CH <sub>3</sub> NH <sub>3</sub> PbI <sub>3</sub> or (CH <sub>3</sub> HNH <sub>3</sub> ) <sub>2</sub> (CH <sub>3</sub> NH <sub>3</sub> ) <sub>n</sub> Pb <sub>i</sub> Mixtures. <i>ACS Applied Materials &amp; Interfaces</i> , 2019, 11, 9583-9593.	8.0	14
27	Accelerated screening of functional atomic impurities in halide perovskites using high-throughput computations and machine learning. <i>Journal of Materials Science</i> , 2022, 57, 10736-10754.	3.7	10
28	A material genome approach towards exploration of Zn and Cd coordination complex polyester as dielectrics: Design, synthesis and characterization. <i>Polymer</i> , 2018, 159, 95-105.	3.8	7
29	Machine learning for impurity charge-state transition levels in semiconductors from elemental properties using multi-fidelity datasets. <i>Journal of Chemical Physics</i> , 2022, 156, 114110.	3.0	5
30	Spectroscopic Comparison of Thermal Transport at Organic-Inorganic and Organic-Hybrid Interfaces Using CsPbBr <sub>3</sub> and FAPbBr <sub>3</sub> (FA = Formamidinium) Perovskite Nanocrystals. <i>Nano Letters</i> , 2019, 19, 8155-8160.	9.1	4
31	Deep learning the properties of inorganic perovskites. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2022, 30, 034005.	2.0	4
32	First-principles Study of Intrinsic and Extrinsic Point Defects in Lead-Based Hybrid Perovskites. , 2018, , .		3
33	Efficient CdTe photovoltaics by co-passivating grain boundaries. , 2018, , .		2
34	First principles modeling of polymer encapsulant degradation in Si photovoltaic modules. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 10357-10364.	2.8	2
35	Computational design of passivants for CdTe grain boundaries. <i>Solar Energy Materials and Solar Cells</i> , 2021, 232, 111279.	6.2	2
36	A first principles investigation of ternary and quaternary II-VI zincblende semiconductor alloys. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2022, 30, 044001.	2.0	1

#	ARTICLE	IF	CITATIONS
37	Machine learning defect properties in Cd-based chalcogenides. , 2019, , .		0
38	A Rational Co-Design Approach for Next Generation Dielectric Materials with the Transition Metal Containing Coordination Polymers. ECS Meeting Abstracts, 2017, , .	0.0	0