

N M Anoop Krishnan

List of Publications by Year in descending order

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Version: 2024-02-01

106
papers

2,634
citations

185998

28
h-index

243296

44
g-index

108
all docs

108
docs citations

108
times ranked

1764
citing authors

#	ARTICLE	IF	CITATIONS
1	Extracting processing and testing parameters from materials science literature for improved property prediction of glasses. <i>Chemical Engineering and Processing: Process Intensification</i> , 2022, 180, 108607.	1.8	11
2	Predicting the near field underwater explosion response of coated composite cylinders using multiscale simulations, experiments, and machine learning. <i>Composite Structures</i> , 2022, 283, 115157.	3.1	7
3	Elucidating the auxetic behavior of cementitious cellular composites using finite element analysis and interpretable machine learning. <i>Materials and Design</i> , 2022, 213, 110341.	3.3	18
4	Interpreting the optical properties of oxide glasses with machine learning and Shapely additive explanations. <i>Journal of the American Ceramic Society</i> , 2022, 105, 4046-4057.	1.9	17
5	Prediction of concrete strengths enabled by missing data imputation and interpretable machine learning. <i>Cement and Concrete Composites</i> , 2022, 128, 104414.	4.6	55
6	Elucidating the influence of structure and Ag ⁺ -Na ⁺ ion-exchange on crack-resistance and ionic conductivity of Na ₃ Al _{1.8} Si _{1.6} P _{1.8} O ₁₂ glass electrolyte. <i>Acta Materialia</i> , 2022, 227, 117745.	3.8	6
7	Fracture response of wollastonite fiber-reinforced cementitious composites: Evaluation using micro-indentation and finite element simulation. <i>Ceramics International</i> , 2022, , .	2.3	4
8	Graphene Oxide Tribofilms Enhance the Scratch Resistance of Silica Glasses. <i>ACS Applied Nano Materials</i> , 2022, 5, 4812-4822.	2.4	4
9	MatSciBERT: A materials domain language model for text mining and information extraction. <i>Npj Computational Materials</i> , 2022, 8, .	3.5	50
10	Reactive molecular simulation of shockwave propagation in calcium silicate hydrate gels. <i>Journal of Non-Crystalline Solids</i> , 2022, 590, 121677.	1.5	1
11	Natural language processing-guided meta-analysis and structure factor database extraction from glass literature. <i>Journal of Non-Crystalline Solids: X</i> , 2022, 15, 100103.	0.5	5
12	Strain sensing efficiency of hierarchical nano-engineered smart twill-weave composites: Evaluations using multiscale numerical simulations. <i>Composite Structures</i> , 2021, 255, 112905.	3.1	7
13	Realizing cool and warm white-LEDs based on color controllable (Sr,Ba) ₂ Al ₃ O ₆ :Eu ²⁺ phosphors obtained via a microwave-assisted diffusion method. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 15245-15256.	1.3	10
14	Scalable Gaussian processes for predicting the optical, physical, thermal, and mechanical properties of inorganic glasses with large datasets. <i>Materials Advances</i> , 2021, 2, 477-487.	2.6	28
15	Rigidity theory of glass: Determining the onset temperature of topological constraints by molecular dynamics. <i>Journal of Non-Crystalline Solids</i> , 2021, 554, 120614.	1.5	5
16	Artificial intelligence and machine learning in glass science and technology: 21 challenges for the 21 st century. <i>International Journal of Applied Glass Science</i> , 2021, 12, 277-292.	1.0	28
17	Modeling the nanoindentation response of silicate glasses by peridynamic simulations. <i>Journal of the American Ceramic Society</i> , 2021, 104, 3531-3544.	1.9	10
18	Machine learning-aided cost prediction and optimization in construction operations. <i>Engineering, Construction and Architectural Management</i> , 2021, , .	1.8	5

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19	Analytical model of the network topology and rigidity of calcium aluminosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2021, 104, 3947-3962.	1.9	14
20	Towards understanding the scratchability in functional glasses. <i>Ceramics International</i> , 2021, 47, 20821-20843.	2.3	9
21	Effect of irradiation on the atomic structure of borosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2021, 104, 6194-6206.	1.9	6
22	Looking through glass: Knowledge discovery from materials science literature using natural language processing. <i>Patterns</i> , 2021, 2, 100290.	3.1	25
23	Disorder-induced expansion of silicate minerals arises from the breakage of weak topological constraints. <i>Journal of Non-Crystalline Solids</i> , 2021, 564, 120846.	1.5	5
24	Ionic Conductivity of $\text{Na}_3\text{Al}_2\text{P}_3\text{O}_{12}$ Glass Electrolytes—Role of Charge Compensators. <i>Inorganic Chemistry</i> , 2021, 60, 12893-12905.	1.9	20
25	Finite Element-Based Numerical Simulations to Evaluate the Influence of Wollastonite Microfibers on the Dynamic Compressive Behavior of Cementitious Composites. <i>Materials</i> , 2021, 14, 4435.	1.3	2
26	Integrating multiscale numerical simulations with machine learning to predict the strain sensing efficiency of nano-engineered smart cementitious composites. <i>Materials and Design</i> , 2021, 209, 109995.	3.3	11
27	Role of steric repulsions on the precipitation kinetics and the structure of calcium-silicate-hydrate gels. <i>Soft Matter</i> , 2021, 17, 8902-8914.	1.2	3
28	Irradiation-induced brittle-to-ductile transition in α -quartz. <i>Journal of the American Ceramic Society</i> , 2020, 103, 3962-3970.	1.9	10
29	Dynamics of confined water and its interplay with alkali cations in sodium aluminosilicate hydrate gel: insights from reactive force field molecular dynamics. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 23707-23724.	1.3	10
30	Fracture toughness of fly ash-based geopolymer gels: Evaluations using nanoindentation experiment and molecular dynamics simulation. <i>Construction and Building Materials</i> , 2020, 262, 120797.	3.2	22
31	Elucidating the constitutive relationship of calcium-silicate-hydrate gel using high throughput reactive molecular simulations and machine learning. <i>Scientific Reports</i> , 2020, 10, 21336.	1.6	12
32	An adaptive, interacting, cluster-based model for predicting the transmission dynamics of COVID-19. <i>Heliyon</i> , 2020, 6, e05722.	1.4	12
33	Drift Response Evaluation of Buckling-Restrained Braced Frames (BRBFs) under Sequential Seismic Disturbances. <i>IOP Conference Series: Materials Science and Engineering</i> , 2020, 936, 012040.	0.3	0
34	A Peridynamics-Based Micromechanical Modeling Approach for Random Heterogeneous Structural Materials. <i>Materials</i> , 2020, 13, 1298.	1.3	11
35	Cooling rate effects on the structure of 45S5 bioglass: Insights from experiments and simulations. <i>Journal of Non-Crystalline Solids</i> , 2020, 534, 119952.	1.5	31
36	Understanding the role of post-indentation recovery on the hardness of glasses: Case of silica, borate, and borosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2020, 534, 119955.	1.5	21

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37	Class Transition and Crystallization in Hexagonal Boron Nitride: Crucial Role of Orientational Order. <i>Advanced Theory and Simulations</i> , 2020, 3, 1900174.	1.3	1
38	Fracture toughness of sodium aluminosilicate hydrate (NASH) gels: Insights from molecular dynamics simulations. <i>Journal of Applied Physics</i> , 2020, 127, .	1.1	10
39	On the equivalence of vapor-deposited and melt-quenched glasses. <i>Journal of Chemical Physics</i> , 2020, 152, 164504.	1.2	7
40	Deep learning aided rational design of oxide glasses. <i>Materials Horizons</i> , 2020, 7, 1819-1827.	6.4	54
41	Topological optimization of cementitious binders: Advances and challenges. <i>Cement and Concrete Composites</i> , 2019, 101, 5-14.	4.6	22
42	Redox Sensitive Self-Assembling Dipeptide for Sustained Intracellular Drug Delivery. <i>Bioconjugate Chemistry</i> , 2019, 30, 2458-2468.	1.8	19
43	Predicting Young's modulus of oxide glasses with sparse datasets using machine learning. <i>Journal of Non-Crystalline Solids</i> , 2019, 524, 119643.	1.5	58
44	Realistic atomic structure of fly ash-based geopolymer gels: Insights from molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2019, 151, .	1.2	27
45	Predicting the dissolution kinetics of silicate glasses by topology-informed machine learning. <i>Npj Materials Degradation</i> , 2019, 3, .	2.6	59
46	Glass Fracture Upon Ballistic Impact: New Insights From Peridynamics Simulations. <i>Frontiers in Materials</i> , 2019, 6, .	1.2	17
47	Dynamic compressive behavior of metallic particulate-reinforced cementitious composites: SHPB experiments and numerical simulations. <i>Construction and Building Materials</i> , 2019, 227, 116668.	3.2	17
48	Densityâ€“stiffness scaling in minerals upon disordering: Irradiation vs. vitrification. <i>Acta Materialia</i> , 2019, 166, 611-617.	3.8	23
49	Simulating the Fracture of Notched Mortar Beams through Extended Finite-Element Method and Peridynamics. <i>Journal of Engineering Mechanics - ASCE</i> , 2019, 145, 04019049.	1.6	13
50	Predicting the Youngâ€™s Modulus of Silicate Glasses using High-Throughput Molecular Dynamics Simulations and Machine Learning. <i>Scientific Reports</i> , 2019, 9, 8739.	1.6	86
51	Atomic picture of structural relaxation in silicate glasses. <i>Applied Physics Letters</i> , 2019, 114, .	1.5	26
52	Experimental and Numerical Investigation of Fracture Behavior of Particle-Reinforced Alkali-Activated Slag Mortars. <i>Journal of Materials in Civil Engineering</i> , 2019, 31, 04019043.	1.3	11
53	Structural percolation controls the precipitation kinetics of colloidal calciumâ€“silicateâ€“hydrate gels. <i>Journal Physics D: Applied Physics</i> , 2019, 52, 315301.	1.3	9
54	Spatial damage sensing ability of metallic particulate-reinforced cementitious composites: Insights from electrical resistance tomography. <i>Materials and Design</i> , 2019, 175, 107817.	3.3	14

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55	Evidence of a two-dimensional glass transition in graphene: Insights from molecular simulations. <i>Scientific Reports</i> , 2019, 9, 4517.	1.6	19
56	Influence of microencapsulated phase change materials (PCMs) on the chloride ion diffusivity of concretes exposed to Freeze-thaw cycles: Insights from multiscale numerical simulations. <i>Construction and Building Materials</i> , 2019, 212, 317-328.	3.2	16
57	Long-term creep deformations in colloidal calcium silicate hydrate gels by accelerated aging simulations. <i>Journal of Colloid and Interface Science</i> , 2019, 542, 339-346.	5.0	19
58	Elucidating the formation of Al–NBO bonds, Al–O–Al linkages and clusters in alkaline-earth aluminosilicate glasses based on molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 23966-23977.	1.3	20
59	The effect of irradiation on the atomic structure and chemical durability of calcite and dolomite. <i>Npj Materials Degradation</i> , 2019, 3, .	2.6	17
60	Fracture response of metallic particulate-reinforced cementitious composites: Insights from experiments and multiscale numerical simulations. <i>Cement and Concrete Composites</i> , 2019, 97, 154-165.	4.6	18
61	Microstructure-guided numerical simulation to evaluate the influence of phase change materials (PCMs) on the freeze-thaw response of concrete pavements. <i>Construction and Building Materials</i> , 2019, 201, 246-256.	3.2	41
62	A microstructure-guided numerical approach to evaluate strain sensing and damage detection ability of random heterogeneous self-sensing structural materials. <i>Computational Materials Science</i> , 2019, 156, 195-205.	1.4	21
63	Effects of polydispersity and disorder on the mechanical properties of hydrated silicate gels. <i>Journal of the Mechanics and Physics of Solids</i> , 2019, 122, 555-565.	2.3	35
64	Predicting the dissolution kinetics of silicate glasses using machine learning. <i>Journal of Non-Crystalline Solids</i> , 2018, 487, 37-45.	1.5	100
65	The hydrophilic-to-hydrophobic transition in glassy silica is driven by the atomic topology of its surface. <i>Journal of Chemical Physics</i> , 2018, 148, 074503.	1.2	35
66	Crack Healing in Cementitious Mortars Using Enzyme-Induced Carbonate Precipitation: Quantification Based on Fracture Response. <i>Journal of Materials in Civil Engineering</i> , 2018, 30, .	1.3	53
67	A new transferable interatomic potential for molecular dynamics simulations of borosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2018, 498, 294-304.	1.5	121
68	Hardness of silicate glasses: Atomic-scale origin of the mixed modifier effect. <i>Journal of Non-Crystalline Solids</i> , 2018, 489, 16-21.	1.5	31
69	Microstructure-guided numerical simulations to predict the thermal performance of a hierarchical cement-based composite material. <i>Cement and Concrete Composites</i> , 2018, 87, 20-28.	4.6	23
70	Effect of irradiation on silicate aggregates' density and stiffness. <i>Journal of Nuclear Materials</i> , 2018, 512, 126-136.	1.3	21
71	New insights into the atomic structure of amorphous TiO ₂ using tight-binding molecular dynamics. <i>Journal of Chemical Physics</i> , 2018, 149, 094501.	1.2	11
72	Role of Electrochemical Surface Potential and Irradiation on Garnet-Type Almandine's Dissolution Kinetics. <i>Journal of Physical Chemistry C</i> , 2018, 122, 17268-17277.	1.5	15

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73	Elucidating the influences of compliant microscale inclusions on the fracture behavior of cementitious composites. <i>Cement and Concrete Composites</i> , 2018, 94, 13-23.	4.6	8
74	Effect of nanoscale phase separation on the fracture behavior of glasses: Toward tough, yet transparent glasses. <i>Physical Review Materials</i> , 2018, 2, .	0.9	21
75	Buckling analysis of cylindrical thin-shells using strain gradient elasticity theory. <i>Meccanica</i> , 2017, 52, 1369-1379.	1.2	12
76	Irradiation- vs. vitrification-induced disordering: The case of α -quartz and glassy silica. <i>Journal of Chemical Physics</i> , 2017, 146, 204502.	1.2	35
77	Irradiation-induced topological transition in SiO ₂ : Structural signature of networks' rigidity. <i>Journal of Non-Crystalline Solids</i> , 2017, 463, 25-30.	1.5	43
78	Ion exchange strengthening and thermal expansion of glasses: Common origin and critical role of network connectivity. <i>Journal of Non-Crystalline Solids</i> , 2017, 455, 70-74.	1.5	36
79	Cooling rate effects in sodium silicate glasses: Bridging the gap between molecular dynamics simulations and experiments. <i>Journal of Chemical Physics</i> , 2017, 147, 074501.	1.2	107
80	Effects of Irradiation on Albite's Chemical Durability. <i>Journal of Physical Chemistry A</i> , 2017, 121, 7835-7845.	1.1	37
81	Revealing the Effect of Irradiation on Cement Hydrates: Evidence of a Topological Self-Organization. <i>ACS Applied Materials & Interfaces</i> , 2017, 9, 32377-32385.	4.0	40
82	Topological Control on the Structural Relaxation of Atomic Networks under Stress. <i>Physical Review Letters</i> , 2017, 119, 035502.	2.9	51
83	Enthalpy Landscape Dictates the Irradiation-Induced Disordering of Quartz. <i>Physical Review X</i> , 2017, 7, .	2.8	27
84	Coaxial Boron-Nitride/Carbon Nanotubes as a Potential Replacement for Double-Walled Carbon Nanotubes for High Strain Applications. <i>Journal of Nanoscience and Nanotechnology</i> , 2017, 17, 5252-5260.	0.9	3
85	Mechanics of Metal-Nanocomposites at Multiple Length Scales: Case of Al-BNNT. <i>Journal of Nanomechanics & Micromechanics</i> , 2017, 7, 04017014.	1.4	0
86	Irradiation-driven amorphous-to-glassy transition in quartz: The crucial role of the medium-range order in crystallization. <i>Physical Review Materials</i> , 2017, 1, .	0.9	27
87	Confined Water in Layered Silicates: The Origin of Anomalous Thermal Expansion Behavior in Calcium-Silicate-Hydrates. <i>ACS Applied Materials & Interfaces</i> , 2016, 8, 35621-35627.	4.0	43
88	A microstructure-guided constitutive modeling approach for random heterogeneous materials: Application to structural binders. <i>Computational Materials Science</i> , 2016, 119, 52-64.	1.4	31
89	Finite element-based micromechanical modeling of the influence of phase properties on the elastic response of cementitious mortars. <i>Construction and Building Materials</i> , 2016, 127, 153-166.	3.2	18
90	The influence of microencapsulated phase change material (PCM) characteristics on the microstructure and strength of cementitious composites: Experiments and finite element simulations. <i>Cement and Concrete Composites</i> , 2016, 73, 29-41.	4.6	128

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91	Elucidating the Crack Resistance of Alkali-Activated Slag Mortars Using Coupled Fracture Tests and Image Correlation. Journal of the American Ceramic Society, 2016, 99, 273-280.	1.9	6
92	Strain energy and process zone based fracture characterization of a novel iron carbonate binding material. Engineering Fracture Mechanics, 2016, 156, 1-15.	2.0	16
93	Micromechanical Modeling for Material Design of Durable Infrastructural Materials: The Influence of Aggregate and Matrix Modification on Elastic Behavior of Mortars. , 2016, , .		5
94	Flexural fracture response of a novel iron carbonate matrix " Glass fiber composite and its comparison to Portland cement-based composites. Construction and Building Materials, 2015, 93, 360-370.	3.2	22
95	Crack propagation and strain localization in metallic particulate-reinforced cementitious mortars. Materials & Design, 2015, 79, 15-25.	5.1	28
96	Fracture process zone and tensile behavior of blended binders containing limestone powder. Cement and Concrete Research, 2015, 73, 51-62.	4.6	36
97	Effective properties of a fly ash geopolymer: Synergistic application of X-ray synchrotron tomography, nanoindentation, and homogenization models. Cement and Concrete Research, 2015, 78, 252-262.	4.6	107
98	Pore- and micro-structural characterization of a novel structural binder based on iron carbonation. Materials Characterization, 2014, 98, 168-179.	1.9	25
99	Defect induced plasticity and failure mechanism of boron nitride nanotubes under tension. Journal of Applied Physics, 2014, 116, .	1.1	14
100	The fracture response of blended formulations containing limestone powder: Evaluations using two-parameter fracture model and digital image correlation. Cement and Concrete Composites, 2014, 53, 316-326.	4.6	54
101	Synthesis and Properties of a Novel Structural Binder Utilizing the Chemistry of Iron Carbonation. ACS Applied Materials & Interfaces, 2014, 6, 8295-8304.	4.0	39
102	Chirality dependent elastic properties of single-walled boron nitride nanotubes under uniaxial and torsional loading. Journal of Applied Physics, 2014, 115, .	1.1	23
103	A novel method for studying the buckling of nanotubes considering geometrical imperfections. Applied Physics A: Materials Science and Processing, 2014, 117, 945-953.	1.1	2
104	Stochastic buckling analysis of carbon nanotubes. , 2014, , 833-836.		0
105	The profiles of first and second SARS-CoV-2 waves in the top ten COVID-19 affected countries. Journal of Global Health Reports, 0, 5, .	1.0	5
106	Quantifying the Densification and Shear Flow under Indentation Deformation in Borosilicate Glasses. International Journal of Applied Glass Science, 0, , .	1.0	0