

Morten Mattrup Smedskjr

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

193
papers

4,202
citations

35
h-index

53
g-index

201
ext. papers

5,095
ext. citations

5
avg, IF

6.02
L-index

#	Paper	IF	Citations
193	StatMechGlass: Python based software for composition structure prediction in oxide glasses using statistical mechanics. <i>SoftwareX</i> , 2022 , 17, 100913	2.7	0
192	Statistical mechanical model for the formation of octahedral silicon in phosphosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2022 , 105, 1031	3.8	0
191	Indentation deformation and cracking behavior of hydrated aluminoborate glasses. <i>Journal of the American Ceramic Society</i> , 2022 , 105, 1039	3.8	0
190	Impact of network topology on the thermal and mechanical properties of lithium germanate glasses. <i>Journal of the American Ceramic Society</i> , 2022 , 105, 977	3.8	0
189	Topological Constraint Theory of Glass: Counting Constraints by Molecular Dynamics Simulations 2022 , 123-148		
188	Oxide glasses under pressure: Recent insights from experiments and simulations. <i>Journal of Applied Physics</i> , 2022 , 131, 170901	2.5	0
187	Mechanical Properties of Oxide Glasses. <i>Reviews in Mineralogy and Geochemistry</i> , 2022 , 87, 229-281	7.1	1
186	Irradiation-induced toughening of calcium aluminoborosilicate glasses. <i>Materials Today Communications</i> , 2022 , 31, 103649	2.5	0
185	Advancing the Mechanical Performance of Glasses: Perspectives and Challenges. <i>Advanced Materials</i> , 2021 , e2109029	24	8
184	Flexible inorganic-organic hybrids with dual inorganic components. <i>Materials Today Chemistry</i> , 2021 , 22, 100584	6.2	2
183	Predicting Fracture Propensity in Amorphous Alumina from Its Static Structure Using Machine Learning. <i>ACS Nano</i> , 2021 ,	16.7	5
182	Modeling the nanoindentation response of silicate glasses by peridynamic simulations. <i>Journal of the American Ceramic Society</i> , 2021 , 104, 3531-3544	3.8	2
181	Bond Switching in Densified Oxide Glass Enables Record-High Fracture Toughness. <i>ACS Applied Materials & Interfaces</i> , 2021 , 13, 17753-17765	9.5	9
180	Thermal conductivity of densified borosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2021 , 557, 120649	4.9	3
179	Interatomic potential parameterization using particle swarm optimization: Case study of glassy silica. <i>Journal of Chemical Physics</i> , 2021 , 154, 134505	3.9	1
178	Analytical model of the network topology and rigidity of calcium aluminosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2021 , 104, 3947-3962	3.8	4
177	Predicting the early-stage creep dynamics of gels from their static structure by machine learning. <i>Acta Materialia</i> , 2021 , 210, 116817	8.4	8

176	Indentation Response of Calcium Aluminoborosilicate Glasses Subjected to Humid Aging and Hot Compression. <i>Materials</i> , 2021 , 14,	3.5	1
175	Mechanics, Ionics, and Optics of Metal-Organic Framework and Coordination Polymer Glasses. <i>Nano Letters</i> , 2021 , 21, 6382-6390	11.5	6
174	Volume relaxation in a borosilicate glass hot compressed by three different methods. <i>Journal of the American Ceramic Society</i> , 2021 , 104, 816-823	3.8	1
173	Decoupling of indentation modulus and hardness in silicate glasses: Evidence of a shear- to densification-dominated transition. <i>Journal of Non-Crystalline Solids</i> , 2021 , 553, 120518	3.9	1
172	Structural densification of lithium phosphoaluminoborate glasses. <i>Journal of the American Ceramic Society</i> , 2021 , 104, 1345-1359	3.8	0
171	Bond switching is responsible for nanoductility in zeolitic imidazolate framework glasses. <i>Dalton Transactions</i> , 2021 , 50, 6126-6132	4.3	7
170	Deformation mechanism of a metal-organic framework glass under indentation. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 16923-16931	3.6	1
169	Structural control of self-healing silica-poly(tetrahydropyran)-poly(ϵ -caprolactone) hybrids. <i>Journal of Materials Chemistry B</i> , 2021 , 9, 4400-4410	7.3	2
168	Rigidity theory of glass: Determining the onset temperature of topological constraints by molecular dynamics. <i>Journal of Non-Crystalline Solids</i> , 2021 , 554, 120614	3.9	3
167	A glass act. <i>Nature Chemistry</i> , 2021 , 13, 723-724	17.6	2
166	Mechanical properties of hydrated cesium-lithium aluminoborate glasses. <i>Physical Review Materials</i> , 2021 , 5,	3.2	1
165	Radiation effects on structure and mechanical properties of borosilicate glasses. <i>Journal of Nuclear Materials</i> , 2021 , 552, 153025	3.3	3
164	Toughening of soda-lime-silica glass by nanoscale phase separation: Molecular dynamics study. <i>Physical Review Materials</i> , 2021 , 5,	3.2	1
163	Revealing the medium-range structure of glassy silica using force-enhanced atomic refinement. <i>Journal of Non-Crystalline Solids</i> , 2021 , 573, 121138	3.9	1
162	Vibrational disorder and densification-induced homogenization of local elasticity in silicate glasses.. <i>Scientific Reports</i> , 2021 , 11, 24454	4.9	
161	Mixed Alkali Effect in Silicate Glass Structure: Viewpoint of Si Nuclear Magnetic Resonance and Statistical Mechanics. <i>Journal of Physical Chemistry B</i> , 2020 , 124, 10292-10299	3.4	6
160	Bauchy et al. Reply. <i>Physical Review Letters</i> , 2020 , 124, 199602	7.4	
159	Topological model of alkali germanate glasses and exploration of the germanate anomaly. <i>Journal of the American Ceramic Society</i> , 2020 , 103, 4224-4233	3.8	4

158	New insights into the structure of sodium silicate glasses by force-enhanced atomic refinement. <i>Journal of Non-Crystalline Solids</i> , 2020 , 536, 120006	3.9	10
157	Metal-Organic Framework Glasses Possess Higher Thermal Conductivity than Their Crystalline Counterparts. <i>ACS Applied Materials & Interfaces</i> , 2020 , 12, 18893-18903	9.5	21
156	Structure Dependence of Poisson's Ratio in Cesium Silicate and Borate Glasses. <i>Materials</i> , 2020 , 13,	3.5	1
155	On the relation between fracture toughness and crack resistance in oxide glasses. <i>Journal of Non-Crystalline Solids</i> , 2020 , 534, 119946	3.9	14
154	Cooling rate effects on the structure of 45S5 bioglass: Insights from experiments and simulations. <i>Journal of Non-Crystalline Solids</i> , 2020 , 534, 119952	3.9	15
153	Observation of indentation-induced shear bands in a metal-organic framework glass. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020 , 117, 10149-10154	11.5	17
152	On the equivalence of vapor-deposited and melt-quenched glasses. <i>Journal of Chemical Physics</i> , 2020 , 152, 164504	3.9	5
151	Achieving ultrahigh crack resistance in glass through humid aging. <i>Physical Review Materials</i> , 2020 , 4,	3.2	4
150	Fracture toughness of a metal-organic framework glass. <i>Nature Communications</i> , 2020 , 11, 2593	17.4	31
149	Indentation cracking and deformation mechanism of sodium aluminoborosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2020 , 103, 1656-1665	3.8	6
148	Composition and pressure effects on the structure, elastic properties and hardness of aluminoborosilicate glass. <i>Journal of Non-Crystalline Solids</i> , 2020 , 530, 119797	3.9	15
147	Heat conduction in oxide glasses: Balancing diffusons and propagons by network rigidity. <i>Applied Physics Letters</i> , 2020 , 117, 031901	3.4	7
146	Predicting Cation Interactions in Alkali Aluminoborate Glasses using Statistical Mechanics. <i>Journal of Non-Crystalline Solids</i> , 2020 , 544, 120099	3.9	3
145	Relaxation behavior of densified sodium aluminoborate glass. <i>Acta Materialia</i> , 2020 , 198, 153-167	8.4	1
144	Atomic structure of hot compressed borosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2020 , 103, 6215-6225	3.8	5
143	Revealing hidden medium-range order in amorphous materials using topological data analysis. <i>Science Advances</i> , 2020 , 6,	14.3	14
142	Competitive effects of free volume, rigidity, and self-adaptivity on indentation response of silicoaluminoborate glasses. <i>Journal of the American Ceramic Society</i> , 2020 , 103, 944-954	3.8	4
141	Predicting the dissolution kinetics of silicate glasses by topology-informed machine learning. <i>Npj Materials Degradation</i> , 2019 , 3,	5.7	32

140	Modifier clustering and avoidance principle in borosilicate glasses: A molecular dynamics study. <i>Journal of Chemical Physics</i> , 2019 , 150, 044502	3.9	11
139	Structure, properties, and fabrication of calcium aluminate-based glasses. <i>International Journal of Applied Glass Science</i> , 2019 , 10, 488-501	1.8	4
138	Predicting the Young's Modulus of Silicate Glasses using High-Throughput Molecular Dynamics Simulations and Machine Learning. <i>Scientific Reports</i> , 2019 , 9, 8739	4.9	49
137	Mechanical property optimization of a zinc borate glass by lanthanum doping. <i>Journal of Non-Crystalline Solids</i> , 2019 , 520, 119461	3.9	8
136	Atomic picture of structural relaxation in silicate glasses. <i>Applied Physics Letters</i> , 2019 , 114, 233703	3.4	16
135	Quantifying the internal stress in over-constrained glasses by molecular dynamics simulations. <i>Journal of Non-Crystalline Solids: X</i> , 2019 , 1, 100013	2.5	8
134	Permanent Densification of Calcium Aluminophosphate Glasses. <i>Frontiers in Materials</i> , 2019 , 6,	4	5
133	Prediction of the Young's modulus of silicate glasses by topological constraint theory. <i>Journal of Non-Crystalline Solids</i> , 2019 , 514, 15-19	3.9	23
132	Structural dependence of chemical durability in modified aluminoborate glasses. <i>Journal of the American Ceramic Society</i> , 2019 , 102, 1157-1168	3.8	17
131	Breaking the Limit of Micro-Ductility in Oxide Glasses. <i>Advanced Science</i> , 2019 , 6, 1901281	13.6	24
130	Statistical Mechanical Model of Topological Fluctuations and the Intermediate Phase in Binary Phosphate Glasses. <i>Journal of Physical Chemistry B</i> , 2019 , 123, 7640-7648	3.4	3
129	Revisiting the Dependence of Poisson's Ratio on Liquid Fragility and Atomic Packing Density in Oxide Glasses. <i>Materials</i> , 2019 , 12,	3.5	17
128	Predicting Composition-Structure Relations in Alkali Borosilicate Glasses Using Statistical Mechanics. <i>Frontiers in Materials</i> , 2019 , 6,	4	8
127	Boron anomaly in the thermal conductivity of lithium borate glasses. <i>Physical Review Materials</i> , 2019 , 3,	3.2	7
126	Liquid fragility determination of oxide glass-formers using temperature-modulated DSC. <i>International Journal of Applied Glass Science</i> , 2019 , 10, 321-329	1.8	4
125	Luminescence behaviour of Eu ³⁺ in hot-compressed silicate glasses. <i>Journal of Non-Crystalline Solids: X</i> , 2019 , 4, 100041	2.5	2
124	The role of the network-modifier's field-strength in the chemical durability of aluminoborate glasses. <i>Journal of Non-Crystalline Solids</i> , 2019 , 505, 279-285	3.9	17
123	Elasticity, hardness, and fracture toughness of sodium aluminoborosilicate glasses. <i>Journal of the American Ceramic Society</i> , 2019 , 102, 4520-4537	3.8	13

122	Statistical Mechanical Modeling of Borate Glass Structure and Topology: Prediction of Superstructural Units and Glass Transition Temperature. <i>Journal of Physical Chemistry B</i> , 2019 , 123, 12062-1213	3.4	22
121	Indentation deformation in oxide glasses: Quantification, structural changes, and relation to cracking. <i>Journal of Non-Crystalline Solids: X</i> , 2019 , 1, 100007	2.5	34
120	Predicting the dissolution kinetics of silicate glasses using machine learning. <i>Journal of Non-Crystalline Solids</i> , 2018 , 487, 37-45	3.9	63
119	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	3.9	26
118	Time and humidity dependence of indentation cracking in aluminosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2018 , 491, 64-70	3.9	14
117	Parametric study of temperature-modulated differential scanning calorimetry for high-temperature oxide glasses with varying fragility. <i>Journal of Non-Crystalline Solids</i> , 2018 , 484, 84-94	3.9	3
116	Structural stability of NaPON glass upon heating in air and nitrogen. <i>Journal of Non-Crystalline Solids</i> , 2018 , 482, 137-146	3.9	5
115	Pressure-induced structural changes in titanophosphate glasses studied by neutron and X-ray total scattering analyses. <i>Journal of Non-Crystalline Solids</i> , 2018 , 483, 50-59	3.9	10
114	A new transferable interatomic potential for molecular dynamics simulations of borosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2018 , 498, 294-304	3.9	76
113	Hardness of silicate glasses: Atomic-scale origin of the mixed modifier effect. <i>Journal of Non-Crystalline Solids</i> , 2018 , 489, 16-21	3.9	25
112	Competitive effects of modifier charge and size on mechanical and chemical resistance of aluminoborate glasses. <i>Journal of Non-Crystalline Solids</i> , 2018 , 499, 264-271	3.9	5
111	Predicting Q-Speciation in Binary Phosphate Glasses Using Statistical Mechanics. <i>Journal of Physical Chemistry B</i> , 2018 , 122, 7609-7615	3.4	11
110	Structural impact of nitrogen incorporation on properties of alkali germanophosphate glasses. <i>Journal of the American Ceramic Society</i> , 2018 , 101, 5004-5019	3.8	4
109	Combining high hardness and crack resistance in mixed network glasses through high-temperature densification. <i>Physical Review Materials</i> , 2018 , 2,	3.2	7
108	Effect of nanoscale phase separation on the fracture behavior of glasses: Toward tough, yet transparent glasses. <i>Physical Review Materials</i> , 2018 , 2,	3.2	11
107	Predictive model for the composition dependence of glassy dynamics. <i>Journal of the American Ceramic Society</i> , 2018 , 101, 1169-1179	3.8	13
106	Nano-phase separation and structural ordering in silica-rich mixed network former glasses. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 15707-15717	3.6	9
105	Deformation and cracking behavior of La ₂ O ₃ -doped oxide glasses with high Poisson's ratio. <i>Journal of Non-Crystalline Solids</i> , 2018 , 494, 86-93	3.9	7

104	Structural Compromise between High Hardness and Crack Resistance in Aluminoborate Glasses. <i>Journal of Physical Chemistry B</i> , 2018 , 122, 6287-6295	3.4	17
103	Structural origin of high crack resistance in sodium aluminoborate glasses. <i>Journal of Non-Crystalline Solids</i> , 2017 , 460, 54-65	3.9	53
102	Correlating the Network Topology of Oxide Glasses with their Chemical Durability. <i>Journal of Physical Chemistry B</i> , 2017 , 121, 1139-1147	3.4	41
101	Fragility and configurational heat capacity of calcium aluminosilicate glass-forming liquids. <i>Journal of Non-Crystalline Solids</i> , 2017 , 461, 24-34	3.9	17
100	Topological engineering of glasses using temperature-dependent constraints. <i>MRS Bulletin</i> , 2017 , 42, 29-33	3.2	17
99	Mixed alkali silicophosphate oxynitride glasses: Structure-property relations. <i>Journal of Non-Crystalline Solids</i> , 2017 , 462, 51-64	3.9	12
98	Photoelastic response of permanently densified oxide glasses. <i>Optical Materials</i> , 2017 , 67, 155-161	3.3	4
97	Pressure-driven structural depolymerization of zinc phosphate glass. <i>Journal of Non-Crystalline Solids</i> , 2017 , 469, 31-38	3.9	10
96	Structure of MgO/CaO sodium aluminosilicate glasses: Raman spectroscopy study. <i>Journal of Non-Crystalline Solids</i> , 2017 , 470, 145-151	3.9	24
95	Discovery of Ultra-Crack-Resistant Oxide Glasses with Adaptive Networks. <i>Chemistry of Materials</i> , 2017 , 29, 5865-5876	9.6	77
94	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	3.9	25
93	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	3.9	83
92	Thermometer Effect: Origin of the Mixed Alkali Effect in Glass Relaxation. <i>Physical Review Letters</i> , 2017 , 119, 095501	7.4	41
91	Dissolution Kinetics of Hot Compressed Oxide Glasses. <i>Journal of Physical Chemistry B</i> , 2017 , 121, 9063-9072	3.4	24
90	Accessing Forbidden Glass Regimes through High-Pressure Sub-T Annealing. <i>Scientific Reports</i> , 2017 , 7, 46631	4.9	6
89	Topological Origin of the Network Dilation Anomaly in Ion-Exchanged Glasses. <i>Physical Review Applied</i> , 2017 , 8,	4.3	13
88	Network Glasses Under Pressure: Permanent Densification in Modifier-Free Al ₂ O ₃ B ₂ O ₃ B ₂ O ₅ BiO ₂ Systems. <i>Physical Review Applied</i> , 2017 , 7,	4.3	21
87	Pressure-Induced Densification of Oxide Glasses at the Glass Transition. <i>Frontiers in Materials</i> , 2017 , 4,	4	45

86	Modifier field strength effects on densification behavior and mechanical properties of alkali aluminoborate glasses. <i>Physical Review Materials</i> , 2017 , 1,	3.2	22
85	Fracture toughness anomalies: Viewpoint of topological constraint theory. <i>Acta Materialia</i> , 2016 , 121, 234-239	8.4	62
84	Crucial effect of angular flexibility on the fracture toughness and nano-ductility of aluminosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2016 , 454, 46-51	3.9	18
83	Accelerating the Design of Functional Glasses through Modeling. <i>Chemistry of Materials</i> , 2016 , 28, 4267-4277	9.7	136
82	Thermal history dependence of indentation induced densification in an aluminosilicate glass. <i>Journal of Non-Crystalline Solids</i> , 2016 , 445-446, 34-39	3.9	11
81	Linking Equilibrium and Nonequilibrium Dynamics in Glass-Forming Systems. <i>Journal of Physical Chemistry B</i> , 2016 , 120, 3226-31	3.4	21
80	Universal behavior of changes in elastic moduli of hot compressed oxide glasses. <i>Chemical Physics Letters</i> , 2016 , 651, 88-91	2.5	21
79	Effects of Thermal and Pressure Histories on the Chemical Strengthening of Sodium Aluminosilicate Glass. <i>Frontiers in Materials</i> , 2016 , 3,	4	11
78	Impact of nitridation of metaphosphate glasses on liquid fragility. <i>Journal of Non-Crystalline Solids</i> , 2016 , 441, 22-28	3.9	20
77	A medium range order structural connection to the configurational heat capacity of borate-silicate mixed glasses. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 10887-95	3.6	14
76	Structure and mechanical properties of compressed sodium aluminosilicate glasses: Role of non-bridging oxygens. <i>Journal of Non-Crystalline Solids</i> , 2016 , 441, 49-57	3.9	71
75	Raman spectroscopy study of pressure-induced structural changes in sodium borate glass. <i>Journal of Non-Crystalline Solids</i> , 2016 , 443, 130-135	3.9	12
74	Pressure-induced structural transformations in phosphorus oxynitride glasses. <i>Journal of Non-Crystalline Solids</i> , 2016 , 452, 153-160	3.9	6
73	Volume and structural relaxation in compressed sodium borate glass. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 29879-29891	3.6	19
72	Confocal depth-resolved micro-X-ray absorption spectroscopy study of chemically strengthened boroaluminosilicate glasses. <i>RSC Advances</i> , 2016 , 6, 24060-24065	3.7	3
71	Structure-property relations in calcium aluminate glasses containing different divalent cations and SiO ₂ . <i>Journal of Non-Crystalline Solids</i> , 2015 , 427, 160-165	3.9	18
70	Role of elastic deformation in determining the mixed alkaline earth effect of hardness in silicate glasses. <i>Journal of Applied Physics</i> , 2015 , 117, 034903	2.5	8
69	Hardness of oxynitride glasses: topological origin. <i>Journal of Physical Chemistry B</i> , 2015 , 119, 4109-15	3.4	22

68	Structure-topology-property correlations of sodium phosphosilicate glasses. <i>Journal of Chemical Physics</i> , 2015 , 143, 064510	3.9	35
67	Indentation deformation mechanism of isostatically compressed mixed alkali aluminosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2015 , 426, 175-183	3.9	42
66	Temperature-dependent densification of sodium borosilicate glass. <i>RSC Advances</i> , 2015 , 5, 78845-78851	3.7	22
65	Unique effects of thermal and pressure histories on glass hardness: Structural and topological origin. <i>Journal of Chemical Physics</i> , 2015 , 143, 164505	3.9	43
64	Cation Diffusivity and the Mixed Network Former Effect in Borosilicate Glasses. <i>Journal of Physical Chemistry B</i> , 2015 , 119, 7106-15	3.4	14
63	Effect of divalent cations and SiO ₂ on the crystallization behavior of calcium aluminate glasses. <i>Journal of Non-Crystalline Solids</i> , 2015 , 413, 20-23	3.9	6
62	Sub-critical crack growth in silicate glasses: Role of network topology. <i>Applied Physics Letters</i> , 2015 , 107, 141901	3.4	19
61	Mixed alkaline earth effect in the compressibility of aluminosilicate glasses. <i>Journal of Chemical Physics</i> , 2014 , 140, 054511	3.9	42
60	Composition-Structure-Property Relations of Compressed Borosilicate Glasses. <i>Physical Review Applied</i> , 2014 , 2,	4.3	38
59	Principles of Pyrex [®] glass chemistry: structure-property relationships. <i>Applied Physics A: Materials Science and Processing</i> , 2014 , 116, 491-504	2.6	32
58	On the origin of the mixed alkali effect on indentation in silicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2014 , 406, 22-26	3.9	28
57	Pressure-induced changes in interdiffusivity and compressive stress in chemically strengthened glass. <i>ACS Applied Materials & Interfaces</i> , 2014 , 6, 10436-44	9.5	19
56	Alkali diffusivity in alkaline earth sodium boroaluminosilicate glasses. <i>Solid State Ionics</i> , 2014 , 263, 95-98	3.3	9
55	Statistical mechanics of glass. <i>Journal of Non-Crystalline Solids</i> , 2014 , 396-397, 41-53	3.9	72
54	Irreversibility of pressure induced boron speciation change in glass. <i>Scientific Reports</i> , 2014 , 4, 3770	4.9	46
53	Topological Model for Boroaluminosilicate Glass Hardness. <i>Frontiers in Materials</i> , 2014 , 1,	4	36
52	Indentation size effect and the plastic compressibility of glass. <i>Applied Physics Letters</i> , 2014 , 104, 251906	3.4	16
51	Hardness and incipient plasticity in silicate glasses: Origin of the mixed modifier effect. <i>Applied Physics Letters</i> , 2014 , 104, 051913	3.4	41

50	Effect of Na ₂ CO ₃ as foaming agent on dynamics and structure of foam glass melts. <i>Journal of Non-Crystalline Solids</i> , 2014 , 400, 1-5	3.9	32
49	Non-conservation of the total alkali concentration in ion-exchanged glass. <i>Journal of Non-Crystalline Solids</i> , 2014 , 387, 71-75	3.9	7
48	Structure and properties of sodium aluminosilicate glasses from molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2013 , 139, 044507	3.9	96
47	Compositional control of the photoelastic response of silicate glasses. <i>Optical Materials</i> , 2013 , 35, 2435-2439	3.9	12
46	Relaxation kinetics of the mechanical properties of an aluminosilicate glass. <i>Journal of Non-Crystalline Solids</i> , 2013 , 362, 40-46	3.9	19
45	Viscosity and Fragility of Alkaline-Earth Sodium Boroaluminosilicate Liquids. <i>Journal of the American Ceramic Society</i> , 2013 , 96, 2831-2838	3.8	10
44	Liquidus surface of MgO-CaO-Al ₂ O ₃ -SiO ₂ glass-forming systems. <i>Journal of Non-Crystalline Solids</i> , 2013 , 363, 39-45	3.9	16
43	Are the dynamics of a glass embedded in its elastic properties?. <i>Journal of Chemical Physics</i> , 2013 , 138, 12A501	3.9	14
42	Impact of ZnO on the structure and properties of sodium aluminosilicate glasses: Comparison with alkaline earth oxides. <i>Journal of Non-Crystalline Solids</i> , 2013 , 381, 58-64	3.9	23
41	Mixed alkaline earth effect in sodium aluminosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2013 , 369, 61-68	3.9	62
40	Elastic and micromechanical properties of isostatically compressed soda-borate glasses. <i>Journal of Non-Crystalline Solids</i> , 2013 , 364, 44-52	3.9	50
39	Microscopic Origins of Compositional Trends in Aluminosilicate Glass Properties. <i>Journal of the American Ceramic Society</i> , 2013 , 96, 1436-1443	3.8	29
38	Environmental effects on fatigue of alkaline earth aluminosilicate glass with varying fictive temperature. <i>Journal of Non-Crystalline Solids</i> , 2013 , 379, 161-168	3.9	15
37	Liquidus Temperature of SrO-Al ₂ O ₃ -SiO ₂ Glass-Forming Compositions. <i>International Journal of Applied Glass Science</i> , 2013 , 4, 225-230	1.8	6
36	Topological Model for the Viscosity of Multicomponent Glass-Forming Liquids. <i>International Journal of Applied Glass Science</i> , 2013 , 4, 408-413	1.8	42
35	Minimalist landscape model of glass relaxation. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2012 , 391, 3446-3459	3.3	17
34	Influence of aluminum speciation on the stability of aluminosilicate glasses against crystallization. <i>Applied Physics Letters</i> , 2012 , 101, 041906	3.4	28
33	Aging in chalcogenide glasses: Origin and consequences. <i>Journal of Non-Crystalline Solids</i> , 2012 , 358, 129-132	3.3	11

32	Glass-forming ability of soda lime borate liquids. <i>Journal of Non-Crystalline Solids</i> , 2012 , 358, 658-665	3.9	19
31	Relationship between viscous dynamics and the configurational thermal expansion coefficient of glass-forming liquids. <i>Journal of Non-Crystalline Solids</i> , 2012 , 358, 648-651	3.9	4
30	Composition-structure-property relationships in boroaluminosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2012 , 358, 993-1002	3.9	76
29	Surface-luminescence from thermally reduced bismuth-doped sodium aluminosilicate glasses. <i>Journal of Non-Crystalline Solids</i> , 2012 , 358, 3193-3199	3.9	13
28	Distinguishability of particles in glass-forming systems. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2012 , 391, 5392-5403	3.3	11
27	Unified physics of stretched exponential relaxation and Weibull fracture statistics. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2012 , 391, 6121-6127	3.3	38
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