

Thomas M Truskett

List of Publications by Year in descending order

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178
papers

8,124
citations

57631

44
h-index

56606

83
g-index

186
all docs

186
docs citations

186
times ranked

6251
citing authors

#	ARTICLE	IF	CITATIONS
1	Is Random Close Packing of Spheres Well Defined?. <i>Physical Review Letters</i> , 2000, 84, 2064-2067.	2.9	1,173
2	Modeling Water, the Hydrophobic Effect, and Ion Solvation. <i>Annual Review of Biophysics and Biomolecular Structure</i> , 2005, 34, 173-199.	18.3	363
3	Hydrophobic hydration from small to large lengthscales: Understanding and manipulating the crossover. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 9475-9480.	3.3	268
4	Towards a quantification of disorder in materials: Distinguishing equilibrium and glassy sphere packings. <i>Physical Review E</i> , 2000, 62, 993-1001.	0.8	258
5	Layering and Position-Dependent Diffusive Dynamics of Confined Fluids. <i>Physical Review Letters</i> , 2008, 100, 145901.	2.9	161
6	A single-bond approach to orientation-dependent interactions and its implications for liquid water. <i>Journal of Chemical Physics</i> , 1999, 111, 2647-2656.	1.2	157
7	Excess-entropy-based anomalies for a waterlike fluid. <i>Journal of Chemical Physics</i> , 2006, 125, 244502.	1.2	148
8	Structural precursor to freezing in the hard-disk and hard-sphere systems. <i>Physical Review E</i> , 1998, 58, 3083-3088.	0.8	147
9	Thermodynamic implications of confinement for a waterlike fluid. <i>Journal of Chemical Physics</i> , 2001, 114, 2401-2418.	1.2	143
10	Thermodynamics Predicts How Confinement Modifies the Dynamics of the Equilibrium Hard-Sphere Fluid. <i>Physical Review Letters</i> , 2006, 96, 177804.	2.9	133
11	The Kauzmann Paradox Revisited. <i>Journal of Physical Chemistry B</i> , 2001, 105, 11809-11816.	1.2	131
12	A simple molecular thermodynamic theory of hydrophobic hydration. <i>Journal of Chemical Physics</i> , 2002, 116, 2907-2921.	1.2	118
13	Effects of lengthscales and attractions on the collapse of hydrophobic polymers in water. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007, 104, 733-738.	3.3	113
14	Free volume in the hard sphere liquid. <i>Molecular Physics</i> , 1998, 95, 289-297.	0.8	112
15	Relationship between thermodynamics and dynamics of supercooled liquids. <i>Journal of Chemical Physics</i> , 2006, 125, 076102.	1.2	106
16	Concentrated Dispersions of Equilibrium Protein Nanoclusters That Reversibly Dissociate into Active Monomers. <i>ACS Nano</i> , 2012, 6, 1357-1369.	7.3	104
17	The Equation of State of an Energy Landscape. <i>Journal of Physical Chemistry B</i> , 1999, 103, 7390-7397.	1.2	103
18	Anomalous structure and dynamics of the Gaussian-core fluid. <i>Physical Review E</i> , 2009, 79, 031203.	0.8	100

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19	Steric stabilization of nanoparticles with grafted low molecular weight ligands in highly concentrated brines including divalent ions. <i>Soft Matter</i> , 2016, 12, 2025-2039.	1.2	99
20	Quantitative Link between Single-Particle Dynamics and Static Structure of Supercooled Liquids. <i>Journal of Physical Chemistry B</i> , 2006, 110, 18147-18150.	1.2	95
21	Relationships between Self-Diffusivity, Packing Fraction, and Excess Entropy in Simple Bulk and Confined Fluids. <i>Journal of Physical Chemistry B</i> , 2007, 111, 10054-10063.	1.2	94
22	Synergistic Formation and Stabilization of Oil-in-Water Emulsions by a Weakly Interacting Mixture of Zwitterionic Surfactant and Silica Nanoparticles. <i>Langmuir</i> , 2014, 30, 984-994.	1.6	90
23	Generalized Rosenfeld scalings for tracer diffusivities in not-so-simple fluids: Mixtures and soft particles. <i>Physical Review E</i> , 2009, 80, 061205.	0.8	79
24	Charged Gold Nanoparticles with Essentially Zero Serum Protein Adsorption in Undiluted Fetal Bovine Serum. <i>Journal of the American Chemical Society</i> , 2013, 135, 7799-7802.	6.6	79
25	Modified Montmorillonite Clay Microparticles for Stable Oil-in-Seawater Emulsions. <i>ACS Applied Materials & Interfaces</i> , 2014, 6, 11502-11513.	4.0	78
26	Inverse methods for material design. <i>AIChE Journal</i> , 2014, 60, 2732-2740.	1.8	77
27	Inverse design of simple pairwise interactions with low-coordinated 3D lattice ground states. <i>Soft Matter</i> , 2013, 9, 3866.	1.2	74
28	Structural Ensemble of an Intrinsically Disordered Polypeptide. <i>Journal of Physical Chemistry B</i> , 2013, 117, 118-124.	1.2	72
29	Does confining the hard-sphere fluid between hard walls change its average properties?. <i>Journal of Chemical Physics</i> , 2007, 126, 244708.	1.2	65
30	Predicting How Nanoconfinement Changes the Relaxation Time of a Supercooled Liquid. <i>Physical Review Letters</i> , 2013, 111, 235901.	2.9	65
31	Inverse methods for design of soft materials. <i>Journal of Chemical Physics</i> , 2020, 152, 140902.	1.2	63
32	On the Use of Excess Entropy Scaling to Describe the Dynamic Properties of Water. <i>Journal of Physical Chemistry B</i> , 2010, 114, 10558-10566.	1.2	61
33	Tuning Density Profiles and Mobility of Inhomogeneous Fluids. <i>Physical Review Letters</i> , 2008, 100, 106001.	2.9	60
34	How short-range attractions impact the structural order, self-diffusivity, and viscosity of a fluid. <i>Journal of Chemical Physics</i> , 2007, 127, 044502.	1.2	59
35	Theory of supercooled liquids and glasses: Energy landscape and statistical geometry perspectives. <i>Advances in Chemical Engineering</i> , 2001, 28, 21-79.	0.5	58
36	A Simple Statistical Mechanical Model of Water. <i>Journal of Physical Chemistry B</i> , 2002, 106, 11829-11842.	1.2	58

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37	Predicting water's phase diagram and liquid-state anomalies. <i>Journal of Chemical Physics</i> , 2002, 117, 5101-5104.	1.2	54
38	Communication: Inverse design for self-assembly via on-the-fly optimization. <i>Journal of Chemical Physics</i> , 2016, 145, .	1.2	52
39	Nonequilibrium hard-disk packings with controlled orientational order. <i>Journal of Chemical Physics</i> , 2000, 113, 4844.	1.2	51
40	Communication: Generalizing Rosenfeld's excess-entropy scaling to predict long-time diffusivity in dense fluids of Brownian particles: From hard to ultrasoft interactions. <i>Journal of Chemical Physics</i> , 2011, 134, 081101.	1.2	51
41	Equilibrium Gold Nanoclusters Quenched with Biodegradable Polymers. <i>ACS Nano</i> , 2013, 7, 239-251.	7.3	51
42	Contrasting the Influence of Cationic Amino Acids on the Viscosity and Stability of a Highly Concentrated Monoclonal Antibody. <i>Pharmaceutical Research</i> , 2017, 34, 193-207.	1.7	50
43	Indocyanine Green J Aggregates in Polymersomes for Near-Infrared Photoacoustic Imaging. <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 46437-46450.	4.0	46
44	Linking Semiconductor Nanocrystals into Gel Networks through Al ³⁺ -Inorganic Bridges. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 14840-14844.	7.2	45
45	Probabilistic inverse design for self-assembling materials. <i>Journal of Chemical Physics</i> , 2017, 146, .	1.2	44
46	Available states and available space: static properties that predict self-diffusivity of confined fluids. <i>Journal of Statistical Mechanics: Theory and Experiment</i> , 2009, 2009, P04006.	0.9	43
47	Reversible Self-Assembly of Glutathione-Coated Gold Nanoparticle Clusters via pH-Tunable Interactions. <i>Langmuir</i> , 2017, 33, 12244-12253.	1.6	43
48	Liquid Structure via Cavity Size Distributions. <i>Journal of Physical Chemistry B</i> , 2000, 104, 12028-12034.	1.2	41
49	Structural anomalies of fluids: Origins in second and higher coordination shells. <i>Physical Review E</i> , 2008, 77, 041201.	0.8	41
50	Assembly of nothing: equilibrium fluids with designed structured porosity. <i>Soft Matter</i> , 2016, 12, 2663-2667.	1.2	41
51	Coarse-Grained Strategy for Modeling Protein Stability in Concentrated Solutions. <i>Biophysical Journal</i> , 2005, 89, 2372-2384.	0.2	40
52	On the Use of Excess Entropy Scaling To Describe Single-Molecule and Collective Dynamic Properties of Hydrocarbon Isomer Fluids. <i>Journal of Physical Chemistry B</i> , 2010, 114, 16487-16493.	1.2	40
53	High concentration tangential flow ultrafiltration of stable monoclonal antibody solutions with low viscosities. <i>Journal of Membrane Science</i> , 2016, 508, 113-126.	4.1	40
54	The subtleties of water in small spaces. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2003, 100, 10139-10140.	3.3	38

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55	Evidence that Amorphous Water below 160 K Is Not a Fragile Liquid. <i>Journal of Physical Chemistry B</i> , 2006, 110, 11033-11036.	1.2	38
56	Improving Viscosity and Stability of a Highly Concentrated Monoclonal Antibody Solution with Concentrated Proline. <i>Pharmaceutical Research</i> , 2018, 35, 133.	1.7	38
57	On the formation of equilibrium gels via a macroscopic bond limitation. <i>Journal of Chemical Physics</i> , 2016, 145, 074906.	1.2	37
58	Heteropolymer Collapse Theory for Protein Folding in the Pressure-Temperature Plane. <i>Biophysical Journal</i> , 2006, 91, 2427-2435.	0.2	36
59	Origin and detection of microstructural clustering in fluids with spatial-range competitive interactions. <i>Physical Review E</i> , 2015, 91, 042312.	0.8	36
60	Unsupervised machine learning for detection of phase transitions in off-lattice systems. I. Foundations. <i>Journal of Chemical Physics</i> , 2018, 149, 194109.	1.2	36
61	Coarse-Grained Strategy for Modeling Protein Stability in Concentrated Solutions. II: Phase Behavior. <i>Biophysical Journal</i> , 2006, 90, 1949-1960.	0.2	34
62	Transport in Amorphous Solid Water Films: Implications for Self-Diffusivity. <i>Journal of Physical Chemistry B</i> , 2006, 110, 17987-17997.	1.2	34
63	Nanocrystal superlattices that exhibit improved order on heating: an example of inverse melting?. <i>Faraday Discussions</i> , 2015, 181, 181-192.	1.6	34
64	Excess entropy scaling of dynamic quantities for fluids of dumbbell-shaped particles. <i>Journal of Chemical Physics</i> , 2010, 133, 104506.	1.2	33
65	Orientationally Ordered Silicon Nanocrystal Cuboctahedra in Superlattices. <i>Nano Letters</i> , 2016, 16, 7814-7821.	4.5	33
66	Confinement, entropy, and single-particle dynamics of equilibrium hard-sphere mixtures. <i>Journal of Chemical Physics</i> , 2007, 127, 154513.	1.2	32
67	Dimensionality and Design of Isotropic Interactions that Stabilize Honeycomb, Square, Simple Cubic, and Diamond Lattices. <i>Physical Review X</i> , 2014, 4, .	2.8	32
68	Equilibrium cluster fluids: pair interactions via inverse design. <i>Soft Matter</i> , 2015, 11, 9342-9354.	1.2	32
69	Designing convex repulsive pair potentials that favor assembly of kagome and snub square lattices. <i>Journal of Chemical Physics</i> , 2016, 145, 054901.	1.2	32
70	Gelation of plasmonic metal oxide nanocrystals by polymer-induced depletion attractions. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, 8925-8930.	3.3	32
71	Protein-Protein Interactions of Highly Concentrated Monoclonal Antibody Solutions via Static Light Scattering and Influence on the Viscosity. <i>Journal of Physical Chemistry B</i> , 2019, 123, 739-755.	1.2	32
72	Composition and concentration anomalies for structure and dynamics of Gaussian-core mixtures. <i>Journal of Chemical Physics</i> , 2009, 131, 161101.	1.2	31

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73	High Interfacial Activity of Polymers Grafted through Functionalized Iron Oxide Nanoparticle Clusters. <i>Langmuir</i> , 2014, 30, 10188-10196.	1.6	31
74	Fluids with competing interactions. I. Decoding the structure factor to detect and characterize self-limited clustering. <i>Journal of Chemical Physics</i> , 2016, 145, .	1.2	31
75	Density fluctuations in many-body systems. <i>Physical Review E</i> , 1998, 58, 7369-7380.	0.8	30
76	Configurational Properties and Corresponding States in Simple Fluids and Water. <i>Journal of Physical Chemistry B</i> , 1999, 103, 5106-5116.	1.2	30
77	Polymer Conductivity through Particle Connectivity. <i>Chemistry of Materials</i> , 2009, 21, 1948-1954.	3.2	30
78	Mapping between long-time molecular and Brownian dynamics. <i>Soft Matter</i> , 2011, 7, 9859.	1.2	30
79	Impact of surface roughness on diffusion of confined fluids. <i>Journal of Chemical Physics</i> , 2011, 135, 154502.	1.2	30
80	Tunable equilibrium nanocluster dispersions at high protein concentrations. <i>Soft Matter</i> , 2013, 9, 1766-1771.	1.2	30
81	Viscosity Reduction of a Concentrated Monoclonal Antibody with Arginine-HCl and Arginine-Glutamate. <i>Industrial & Engineering Chemistry Research</i> , 2016, 55, 11225-11234.	1.8	30
82	Ideal glass transitions in thin films: An energy landscape perspective. <i>Journal of Chemical Physics</i> , 2003, 119, 1897-1900.	1.2	29
83	Interactions and design rules for assembly of porous colloidal mesophases. <i>Soft Matter</i> , 2017, 13, 1335-1343.	1.2	29
84	Structure and phase behavior of polymer-linked colloidal gels. <i>Journal of Chemical Physics</i> , 2019, 151, 124901.	1.2	28
85	Charge Shielding Prevents Aggregation of Supercharged GFP Variants at High Protein Concentration. <i>Molecular Pharmaceutics</i> , 2017, 14, 3269-3280.	2.3	27
86	Inverse design of multicomponent assemblies. <i>Journal of Chemical Physics</i> , 2018, 148, 104509.	1.2	27
87	X-ray Scattering and Coarse-Grained Simulations for Clustering and Interactions of Monoclonal Antibodies at High Concentrations. <i>Journal of Physical Chemistry B</i> , 2019, 123, 5274-5290.	1.2	27
88	Assembly of Linked Nanocrystal Colloids by Reversible Covalent Bonds. <i>Chemistry of Materials</i> , 2020, 32, 10235-10245.	3.2	27
89	Assembling Inorganic Nanocrystal Gels. <i>Nano Letters</i> , 2022, 22, 1457-1466.	4.5	27
90	Coarse-Grained Molecular Dynamics Simulations for Understanding the Impact of Short-Range Anisotropic Attractions on Structure and Viscosity of Concentrated Monoclonal Antibody Solutions. <i>Molecular Pharmaceutics</i> , 2020, 17, 1748-1756.	2.3	26

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91	Colloidal Nanocrystal Gels from Thermodynamic Principles. <i>Accounts of Chemical Research</i> , 2021, 54, 798-807.	7.6	26
92	Iso-g(2)Processes in Equilibrium Statistical Mechanics. <i>Journal of Physical Chemistry B</i> , 2001, 105, 6592-6597.	1.2	25
93	Coarse-Grained Strategy for Modeling Protein Stability in Concentrated Solutions. III: Directional Protein Interactions. <i>Biophysical Journal</i> , 2007, 92, 4316-4324.	0.2	25
94	Communication: From close-packed to topologically close-packed: Formation of Laves phases in moderately polydisperse hard-sphere mixtures. <i>Journal of Chemical Physics</i> , 2018, 148, 191101.	1.2	25
95	Free volume in the hard sphere liquid. <i>Molecular Physics</i> , 1998, 95, 289-297.	0.8	25
96	Attractions, Water Structure, and Thermodynamics of Hydrophobic Polymer Collapse. <i>Journal of Physical Chemistry B</i> , 2008, 112, 13193-13196.	1.2	23
97	Unsupervised machine learning for detection of phase transitions in off-lattice systems. II. Applications. <i>Journal of Chemical Physics</i> , 2018, 149, 194110.	1.2	23
98	A simple analytical model of water. <i>Biophysical Chemistry</i> , 2003, 105, 449-459.	1.5	22
99	Using available volume to predict fluid diffusivity in random media. <i>Physical Review E</i> , 2006, 74, 040102.	0.8	22
100	Excess-entropy scaling of dynamics for a confined fluid of dumbbell-shaped particles. <i>Physical Review E</i> , 2010, 82, 041201.	0.8	22
101	Formation of Small Gold Nanoparticle Chains with High NIR Extinction through Bridging with Calcium Ions. <i>Langmuir</i> , 2016, 32, 1127-1138.	1.6	21
102	Using Energy Landscapes To Predict the Properties of Thin Films. <i>Journal of Physical Chemistry B</i> , 2004, 108, 19769-19779.	1.2	20
103	Inverse Design of Self-Assembling Frank-Kasper Phases and Insights Into Emergent Quasicrystals. <i>Journal of Physical Chemistry B</i> , 2018, 122, 5547-5556.	1.2	20
104	Assembly of particle strings via isotropic potentials. <i>Journal of Chemical Physics</i> , 2019, 150, 124903.	1.2	20
105	The statistical geometry of voids in liquids. <i>Fluid Phase Equilibria</i> , 1999, 158-160, 549-556.	1.4	19
106	Protein-Protein Interactions, Clustering, and Rheology for Bovine IgG up to High Concentrations Characterized by Small Angle X-Ray Scattering and Molecular Dynamics Simulations. <i>Journal of Pharmaceutical Sciences</i> , 2020, 109, 696-708.	1.6	19
107	Designing pairwise interactions that stabilize open crystals: Truncated square and truncated hexagonal lattices. <i>Journal of Chemical Physics</i> , 2017, 146, 144501.	1.2	17
108	Triangle Distribution and Equation of State for Classical Rigid Disks. <i>Journal of Statistical Physics</i> , 2000, 100, 49-72.	0.5	16

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109	Implications of the effective one-component analysis of pair correlations in colloidal fluids with polydispersity. <i>Journal of Chemical Physics</i> , 2011, 135, 124513.	1.2	16
110	Quenched Assembly of NIR-Active Gold Nanoclusters Capped with Strongly Bound Ligands by Tuning Particle Charge via pH and Salinity. <i>Journal of Physical Chemistry C</i> , 2014, 118, 14291-14298.	1.5	16
111	Enhancing Stability and Reducing Viscosity of a Monoclonal Antibody With Cosolutes by Weakening Protein-Protein Interactions. <i>Journal of Pharmaceutical Sciences</i> , 2019, 108, 2517-2526.	1.6	16
112	Universal Gelation of Metal Oxide Nanocrystals via Depletion Attractions. <i>Nano Letters</i> , 2020, 20, 4007-4013.	4.5	16
113	Model for the free-volume distributions of equilibrium fluids. <i>Journal of Chemical Physics</i> , 2006, 124, 214502.	1.2	15
114	Structural signatures of mobility on intermediate time scales in a supercooled fluid. <i>Journal of Chemical Physics</i> , 2010, 132, .	1.2	15
115	Putting the squeeze on cavities in liquids: Quantifying pressure effects on solvation using simulations and scaled-particle theory. <i>Journal of Chemical Physics</i> , 2011, 134, 014507.	1.2	15
116	Connection between Thermodynamics and Dynamics of Simple Fluids in Highly Attractive Pores. <i>Langmuir</i> , 2013, 29, 14527-14535.	1.6	15
117	Breadth versus depth: Interactions that stabilize particle assemblies to changes in density or temperature. <i>Journal of Chemical Physics</i> , 2016, 144, 084502.	1.2	15
118	Relating Collective Diffusion, Protein-Protein Interactions, and Viscosity of Highly Concentrated Monoclonal Antibodies through Dynamic Light Scattering. <i>Industrial & Engineering Chemistry Research</i> , 2019, 58, 22456-22471.	1.8	15
119	Effects of linker flexibility on phase behavior and structure of linked colloidal gels. <i>Journal of Chemical Physics</i> , 2021, 154, 074901.	1.2	15
120	Residual multiparticle entropy does not generally change sign near freezing. <i>Journal of Chemical Physics</i> , 2008, 128, 161101.	1.2	14
121	Enhancing tracer diffusivity by tuning interparticle interactions and coordination shell structure. <i>Soft Matter</i> , 2012, 8, 4083-4089.	1.2	14
122	Communication: Phase behavior of materials with isotropic interactions designed by inverse strategies to favor diamond and simple cubic lattice ground states. <i>Journal of Chemical Physics</i> , 2013, 139, 141102.	1.2	14
123	Insights Into Crowding Effects on Protein Stability From a Coarse-Grained Model. <i>Journal of Biomechanical Engineering</i> , 2009, 131, 071002.	0.6	13
124	Antibody nanoparticle dispersions formed with mixtures of crowding molecules retain activity and In Vivo bioavailability. <i>Journal of Pharmaceutical Sciences</i> , 2012, 101, 3763-3778.	1.6	13
125	Structure, Thermodynamics, and Position-Dependent Diffusivity in Fluids with Sinusoidal Density Variations. <i>Langmuir</i> , 2014, 30, 8247-8252.	1.6	13
126	Cooling Dodecanethiol-Capped 2 nm Diameter Gold Nanocrystal Superlattices below Room Temperature Induces a Reversible Order-Disorder Structure Transition. <i>Journal of Physical Chemistry C</i> , 2016, 120, 27682-27687.	1.5	13

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127	Fluids with competing interactions. II. Validating a free energy model for equilibrium cluster size. <i>Journal of Chemical Physics</i> , 2016, 145, .	1.2	13
128	Self-diffusion of a highly concentrated monoclonal antibody by fluorescence correlation spectroscopy: insight into protein-protein interactions and self-association. <i>Soft Matter</i> , 2019, 15, 6660-6676.	1.2	13
129	Prediction and Optimization of Ion Transport Characteristics in Nanoparticle-Based Electrolytes Using Convolutional Neural Networks. <i>Journal of Physical Chemistry B</i> , 2021, 125, 4838-4849.	1.2	13
130	Machine Learning-Assisted Design of Material Properties. <i>Annual Review of Chemical and Biomolecular Engineering</i> , 2022, 13, 235-254.	3.3	13
131	Connection Between Thermodynamics and Dynamics of Simple Fluids in Pores: Impact of Fluid-Fluid Interaction Range and Fluid-Solid Interaction Strength. <i>Journal of Physical Chemistry C</i> , 2017, 121, 16316-16327.	1.5	12
132	Design of two-dimensional particle assemblies using isotropic pair interactions with an attractive well. <i>AIP Advances</i> , 2017, 7, .	0.6	12
133	Quantized bounding volume hierarchies for neighbor search in molecular simulations on graphics processing units. <i>Computational Materials Science</i> , 2019, 164, 139-146.	1.4	12
134	Colorimetric quantification of linking in thermoreversible nanocrystal gel assemblies. <i>Science Advances</i> , 2022, 8, eabm7364.	4.7	12
135	Note: Position-dependent and pair diffusivity profiles from steady-state solutions of color reaction-counterdiffusion problems. <i>Journal of Chemical Physics</i> , 2014, 141, 046101.	1.2	11
136	Communication: Local structure-mobility relationships of confined fluids reverse upon supercooling. <i>Journal of Chemical Physics</i> , 2015, 142, 161102.	1.2	11
137	Breakup of Oil Jets into Droplets in Seawater with Environmentally Benign Nanoparticle and Surfactant Dispersants. <i>Industrial & Engineering Chemistry Research</i> , 2015, 54, 4243-4251.	1.8	10
138	Transport Mechanisms Underlying Ionic Conductivity in Nanoparticle-Based Single-Ion Electrolytes. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 6970-6975.	2.1	10
139	Connecting Solute Diffusion to Morphology in Triblock Copolymer Membranes. <i>Macromolecules</i> , 2020, 53, 2336-2343.	2.2	10
140	Wertheim's thermodynamic perturbation theory with double-bond association and its application to colloid-linker mixtures. <i>Journal of Chemical Physics</i> , 2021, 154, 024905.	1.2	10
141	Influence of morphology of colloidal nanoparticle gels on ion transport and rheology. <i>Journal of Chemical Physics</i> , 2019, 150, 214903.	1.2	9
142	Phase diagram for two-dimensional layer of soft particles. <i>Soft Matter</i> , 2019, 15, 4162-4169.	1.2	9
143	The role of pressure in inverse design for assembly. <i>Journal of Chemical Physics</i> , 2019, 151, 104104.	1.2	9
144	Influence of pore morphology on the diffusion of water in triblock copolymer membranes. <i>Journal of Chemical Physics</i> , 2020, 152, 014904.	1.2	9

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145	Free Volumes and the Anomalous Self-Diffusivity of Attractive Colloids. <i>Journal of Physical Chemistry B</i> , 2006, 110, 5166-5169.	1.2	8
146	Dual nature of magnetic nanoparticle dispersions enables control over short-range attraction and long-range repulsion interactions. <i>Communications Chemistry</i> , 2022, 5, .	2.0	8
147	Effect of Dilute Nitric Acid on Crystallization and Fracture of Amorphous Solid Water Films. <i>Journal of Physical Chemistry C</i> , 2007, 111, 10438-10447.	1.5	7
148	Shear-rate-dependent structural order and viscosity of a fluid with short-range attractions. <i>Physical Review E</i> , 2008, 78, 010201.	0.8	7
149	Gold nanoparticles with high densities of small protuberances on nanocluster cores with strong NIR extinction. <i>RSC Advances</i> , 2015, 5, 104674-104687.	1.7	7
150	Equation of state of the rigid disk fluid from its triangle distribution. <i>Journal of Chemical Physics</i> , 2000, 113, 10186-10190.	1.2	6
151	Volatile Diffusional Character of Cytoplasm. <i>Journal of Physical Chemistry B</i> , 2006, 110, 25606-25607.	1.2	6
152	RELATIONSHIP BETWEEN SHEAR VISCOSITY AND STRUCTURE OF A MODEL COLLOIDAL SUSPENSION. <i>Chemical Engineering Communications</i> , 2009, 197, 63-75.	1.5	6
153	Graphoepitaxy for Pattern Multiplication of Nanoparticle Monolayers. <i>Physical Review Letters</i> , 2014, 113, 085503.	2.9	6
154	Tuning structure and mobility of solvation shells surrounding tracer additives. <i>Journal of Chemical Physics</i> , 2015, 142, 124501.	1.2	6
155	Impact of solvent granularity and layering on tracer hydrodynamics in confinement. <i>Soft Matter</i> , 2016, 12, 9561-9574.	1.2	6
156	Response to "Comment on "A simple molecular thermodynamic theory of hydrophobic hydration" [J. Chem. Phys. 119, 10448 (2003)]. <i>Journal of Chemical Physics</i> , 2003, 119, 10450-10451.	1.2	5
157	Communication: Fine discretization of pair interactions and an approximate analytical strategy for predicting equilibrium behavior of complex fluids. <i>Journal of Chemical Physics</i> , 2013, 139, 161102.	1.2	5
158	Effects of protein engineering and rational mutagenesis on crystal lattice of single chain antibody fragments. <i>Proteins: Structure, Function and Bioinformatics</i> , 2014, 82, 1884-1895.	1.5	5
159	A Conversation with John McKetta. <i>Annual Review of Chemical and Biomolecular Engineering</i> , 2017, 8, 1-11.	3.3	5
160	Control of Primary Particle Spacing in Gold Nanoparticle Clusters for Both High NIR Extinction and Full Reversibility. <i>Langmuir</i> , 2017, 33, 3413-3426.	1.6	5
161	Cross-stream migration of a Brownian droplet in a polymer solution under Poiseuille flow. <i>Soft Matter</i> , 2019, 15, 3168-3178.	1.2	5
162	Response to "Comment on "Residual multiparticle entropy does not generally change sign near freezing" [J. Chem. Phys. 130, 037101 (2009)]. <i>Journal of Chemical Physics</i> , 2009, 130, 037102.	1.2	4

#	ARTICLE	IF	CITATIONS
163	How Local and Average Particle Diffusivities of Inhomogeneous Fluids Depend on Microscopic Dynamics. <i>Journal of Physical Chemistry B</i> , 2015, 119, 9103-9113.	1.2	4
164	Position-Dependent Dynamics Explain Pore-Averaged Diffusion in Strongly Attractive Adsorptive Systems. <i>Langmuir</i> , 2017, 33, 13955-13963.	1.6	4
165	Geometric model of crack-templated networks for transparent conductive films. <i>Applied Physics Letters</i> , 2022, 120, .	1.5	4
166	Comment on "Observations on an equation of state for water confined in narrow slit-pores". <i>J. Chem. Phys.</i> 116, 2565 (2002)]. <i>Journal of Chemical Physics</i> , 2002, 117, 8162-8163.	1.2	3
167	Predicting the structure of fluids with piecewise constant interactions: Comparing the accuracy of five efficient integral equation theories. <i>Physical Review E</i> , 2015, 91, 043307.	0.8	3
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170	Intrinsic vulnerabilities to mechanical failure in nanoscale films. <i>Mechanics of Materials</i> , 2006, 38, 924-932.	1.7	2
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178	Stability of force-driven shear flows in nonequilibrium molecular simulations with periodic boundaries. <i>Journal of Chemical Physics</i> , 2020, 152, 214113.	1.2	0