

Zhen-Gang Wang

List of Publications by Year in descending order

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105
papers

4,021
citations

94433

37
h-index

133252

59
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all docs

105
docs citations

105
times ranked

3269
citing authors

#	ARTICLE	IF	CITATIONS
1	Thermodynamics of Electrolyte Solutions Near Charged Surfaces: Constant Surface Charge vs. Constant Surface Potential. <i>Journal of Chemical Physics</i> , 2022, 156, 174704.	3.0	2
2	A coarse-grained model of room-temperature ionic liquids between metal electrodes: a molecular dynamics study. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 11573-11584.	2.8	3
3	Mechanisms of Flow-Induced Polymer Translocation. <i>Macromolecules</i> , 2022, 55, 3602-3612.	4.8	9
4	Local-Average Free Volume Correlates with Dynamics in Glass Formers. <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 3957-3964.	4.6	6
5	Supernatant Phase in Polyelectrolyte Complex Coacervation: Cluster Formation, Binodal, and Nucleation. <i>Macromolecules</i> , 2022, 55, 3910-3923.	4.8	12
6	Complexation between Oppositely Charged Polyelectrolytes in Dilute Solution: Effects of Charge Asymmetry. <i>Macromolecules</i> , 2022, 55, 3898-3909.	4.8	15
7	Surface Charge Density in Electrical Double Layer Capacitors with Nanoscale Cathode-Anode Separation. <i>Journal of Physical Chemistry B</i> , 2021, 125, 625-636.	2.6	20
8	Like dissolves like: A first-principles theory for predicting liquid miscibility and mixture dielectric constant. <i>Science Advances</i> , 2021, 7, .	10.3	47
9	Nonelectrostatic Adsorption of Polyelectrolytes and Mediated Interactions between Solid Surfaces. <i>Langmuir</i> , 2021, 37, 5483-5493.	3.5	8
10	Image-charge effects on ion adsorption near aqueous interfaces. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	7.1	36
11	Preferential Ion Adsorption in Blue Energy Applications. <i>ACS Sustainable Chemistry and Engineering</i> , 2021, 9, 9230-9239.	6.7	7
12	Effects of Confinement and Ion Adsorption in Ionic Liquid Supercapacitors with Nanoporous Electrodes. <i>ACS Nano</i> , 2021, 15, 11724-11733.	14.6	24
13	Salt-Induced Liquid-Liquid Phase Separation: Combined Experimental and Theoretical Investigation of Water-Acetonitrile-Salt Mixtures. <i>Journal of the American Chemical Society</i> , 2021, 143, 773-784.	13.7	35
14	Shear Banding in Entangled Polymers: Stress Plateau, Banding Location, and Lever Rule. <i>ACS Macro Letters</i> , 2021, 10, 1517-1523.	4.8	8
15	Interfacial Structure and Tension of Polyelectrolyte Complex Coacervates. <i>Macromolecules</i> , 2021, 54, 10994-11007.	4.8	15
16	Coil-to-Globule Transition in Polymeric Solvents. <i>Macromolecules</i> , 2021, 54, 10984-10993.	4.8	8
17	Electrostatic Correlations and Temperature-Dependent Dielectric Constant Can Model LCST in Polyelectrolyte Complex Coacervation. <i>Macromolecules</i> , 2021, 54, 11326-11337.	4.8	20
18	Ion transport in small-molecule and polymer electrolytes. <i>Journal of Chemical Physics</i> , 2020, 153, 100903.	3.0	53

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19	Effects of Surface Transition and Adsorption on Ionic Liquid Capacitors. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 1767-1772.	4.6	15
20	Microscopic origins of the swim pressure and the anomalous surface tension of active matter. <i>Physical Review E</i> , 2020, 101, 012604.	2.1	37
21	Revisiting the $\hat{\tau}$ Point. <i>Macromolecules</i> , 2020, 53, 10409-10420.	4.8	12
22	Two-step relaxation and the breakdown of the Stokes-Einstein relation in glass-forming liquids. <i>Physical Review E</i> , 2019, 100, 052607.	2.1	10
23	Nonlinear Rheological Behaviors in Polymer Melts after Step Shear. <i>Macromolecules</i> , 2019, 52, 4103-4110.	4.8	8
24	Food Polyelectrolytes Compress the Colonic Mucus Hydrogel by a Donnan Mechanism. <i>Biomacromolecules</i> , 2019, 20, 2675-2683.	5.4	11
25	On the origin of oscillatory interactions between surfaces mediated by polyelectrolyte solution. <i>Journal of Chemical Physics</i> , 2019, 151, 214901.	3.0	12
26	Swimming to Stability: Structural and Dynamical Control via Active Doping. <i>ACS Nano</i> , 2019, 13, 560-572.	14.6	27
27	Nonphysical Behavior in Several Statistical Mechanically Based Equations of State. <i>Industrial & Engineering Chemistry Research</i> , 2019, 58, 1382-1395.	3.7	8
28	Polyelectrolyte Chain Structure and Solution Phase Behavior. <i>Macromolecules</i> , 2018, 51, 1706-1717.	4.8	60
29	Improved local lattice Monte Carlo simulation for charged systems. <i>Journal of Chemical Physics</i> , 2018, 148, 114105.	3.0	5
30	Mechanisms of Diffusion in Associative Polymer Networks: Evidence for Chain Hopping. <i>Journal of the American Chemical Society</i> , 2018, 140, 14185-14194.	13.7	30
31	Statistical field theory for polar fluids. <i>Journal of Chemical Physics</i> , 2018, 149, 124108.	3.0	9
32	Polyelectrolyte complex coacervation: Effects of concentration asymmetry. <i>Journal of Chemical Physics</i> , 2018, 149, 163303.	3.0	71
33	Salt Partitioning in Complex Coacervation of Symmetric Polyelectrolytes. <i>Macromolecules</i> , 2018, 51, 5586-5593.	4.8	83
34	Globally Suppressed Dynamics in Ion-Doped Polymers. <i>ACS Macro Letters</i> , 2018, 7, 734-738.	4.8	20
35	Density functional theory for charged fluids. <i>Soft Matter</i> , 2018, 14, 5878-5887.	2.7	28
36	A priori determination of the region of the three physical volume root loci in the Perturbed-Chain SAFT EOS. <i>Fluid Phase Equilibria</i> , 2017, 434, 152-166.	2.5	8

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37	Analysis and Control of Chain Mobility in Protein Hydrogels. <i>Journal of the American Chemical Society</i> , 2017, 139, 3796-3804.	13.7	33
38	Electrostatic correlations and the polyelectrolyte self energy. <i>Journal of Chemical Physics</i> , 2017, 146, 084901.	3.0	69
39	Accurate Determination of Ion Polarizabilities in Aqueous Solutions. <i>Journal of Physical Chemistry B</i> , 2017, 121, 6416-6424.	2.6	42
40	Designing Polymer Electrolytes for Safe and High Capacity Rechargeable Lithium Batteries. <i>Accounts of Chemical Research</i> , 2017, 50, 590-593.	15.6	149
41	Variational Methods in Statistical Thermodynamics—A Pedagogical Introduction. <i>Molecular Modeling and Simulation</i> , 2017, , 1-29.	0.2	3
42	Shear-Induced Heterogeneity in Associating Polymer Gels: Role of Network Structure and Dilatancy. <i>Physical Review Letters</i> , 2017, 119, 117801.	7.8	16
43	<i>50th Anniversary Perspective</i>: Polymer Conformation—A Pedagogical Review. <i>Macromolecules</i> , 2017, 50, 9073-9114.	4.8	113
44	Salting-Out and Salting-In of Polyelectrolyte Solutions: A Liquid-State Theory Study. <i>Macromolecules</i> , 2016, 49, 9720-9730.	4.8	63
45	Inhomogeneous screening near the dielectric interface. <i>Journal of Chemical Physics</i> , 2016, 144, 134902.	3.0	16
46	Molecular-Based Theory for Electron-Transfer Reorganization Energy in Solvent Mixtures. <i>Journal of Physical Chemistry B</i> , 2016, 120, 6373-6382.	2.6	3
47	Effects of Ion-Induced Cross-Linking on the Phase Behavior in Salt-Doped Polymer Blends. <i>Macromolecules</i> , 2016, 49, 425-431.	4.8	35
48	The scaling behavior of the second virial coefficient of linear and ring polymer. <i>Science China Chemistry</i> , 2016, 59, 619-623.	8.2	10
49	A molecularly based theory for electron transfer reorganization energy. <i>Journal of Chemical Physics</i> , 2015, 143, 224502.	3.0	4
50	On the theoretical description of weakly charged surfaces. <i>Journal of Chemical Physics</i> , 2015, 142, 104705.	3.0	36
51	Celebrating Soft Matter's 10th Anniversary: Chain configuration and rate-dependent mechanical properties in transient networks. <i>Soft Matter</i> , 2015, 11, 2085-2096.	2.7	32
52	Density-Functional Theory for Mixtures of AB Random Copolymer and CO ₂ . <i>Macromolecules</i> , 2015, 48, 6035-6046.	4.8	10
53	Combined Theoretical and Experimental Study of Refractive Indices of Water—Acetonitrile—Salt Systems. <i>Journal of Physical Chemistry B</i> , 2015, 119, 10701-10709.	2.6	60
54	Systematic Computational and Experimental Investigation of Lithium-Ion Transport Mechanisms in Polyester-Based Polymer Electrolytes. <i>ACS Central Science</i> , 2015, 1, 198-205.	11.3	162

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55	Chemically Specific Dynamic Bond Percolation Model for Ion Transport in Polymer Electrolytes. <i>Macromolecules</i> , 2015, 48, 7346-7358.	4.8	77
56	An efficient dissipative particle dynamics-based algorithm for simulating electrolyte solutions. <i>Journal of Chemical Physics</i> , 2015, 142, 024103.	3.0	16
57	Influence of Topology on the Free Energy and Metric Properties of an Ideal Ring Polymer Confined in a Slit. <i>Macromolecules</i> , 2015, 48, 8675-8680.	4.8	10
58	Continuous Self-Energy of Ions at the Dielectric Interface. <i>Physical Review Letters</i> , 2014, 112, 136101.	7.8	33
59	Origin of Stress Overshoot during Startup Shear of Entangled Polymer Melts. <i>ACS Macro Letters</i> , 2014, 3, 569-573.	4.8	41
60	Phase Behavior of a Block Copolymer/Salt Mixture through the Order-to-Disorder Transition. <i>Macromolecules</i> , 2014, 47, 2666-2673.	4.8	50
61	Thermodynamics of Salt-Doped Block Copolymers. <i>ACS Macro Letters</i> , 2014, 3, 708-711.	4.8	46
62	Theory of Polymer Chains in Poor Solvent: Single-Chain Structure, Solution Thermodynamics, and $\hat{\Gamma}$ Point. <i>Macromolecules</i> , 2014, 47, 4094-4102.	4.8	47
63	Effects of dielectric inhomogeneity in polyelectrolyte solutions. <i>Soft Matter</i> , 2013, 9, 5686.	2.7	24
64	First-Order Disordered-to-Lamellar Phase Transition in Lithium Salt-Doped Block Copolymers. <i>ACS Macro Letters</i> , 2013, 2, 478-481.	4.8	57
65	Evolution of Chain Conformation and Entanglements during Startup Shear. <i>ACS Macro Letters</i> , 2013, 2, 561-565.	4.8	22
66	Discontinuous Bubble Nucleation Due to a Metastable Condensation Transition in Polymer-CO ₂ Mixtures. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 1639-1643.	4.6	13
67	Effects of image charges on double layer structure and forces. <i>Journal of Chemical Physics</i> , 2013, 139, 124702.	3.0	54
68	Ion Solvation in Liquid Mixtures: Effects of Solvent Reorganization. <i>Physical Review Letters</i> , 2012, 109, 257802.	7.8	57
69	Minimum free energy paths for a nanoparticle crossing the lipid membrane. <i>Soft Matter</i> , 2012, 8, 12066.	2.7	21
70	Theory of Polymers in Poor Solvent: Phase Equilibrium and Nucleation Behavior. <i>Macromolecules</i> , 2012, 45, 6266-6271.	4.8	28
71	Salt-doped block copolymers: ion distribution, domain spacing and effective $\hat{\Gamma}$ parameter. <i>Soft Matter</i> , 2012, 8, 9356.	2.7	113
72	Effects of ion solvation on phase equilibrium and interfacial tension of liquid mixtures. <i>Journal of Chemical Physics</i> , 2011, 135, 014707.	3.0	35

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73	Thermodynamics of Ion-Containing Polymer Blends and Block Copolymers. <i>Physical Review Letters</i> , 2011, 107, 198301.	7.8	129
74	Thermodynamic basis for the genome to capsid charge relationship in viral encapsidation. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2011, 108, 16986-16991.	7.1	56
75	Coil-to-globule transition by dissipative particle dynamics simulation. <i>Journal of Chemical Physics</i> , 2011, 134, 244904.	3.0	52
76	Monte Carlo simulation of a single ring among linear chains: Structural and dynamic heterogeneity. <i>Journal of Chemical Physics</i> , 2010, 133, 064901.	3.0	39
77	Fluctuation in electrolyte solutions: The self energy. <i>Physical Review E</i> , 2010, 81, 021501.	2.1	170
78	A simple model for the anomalous intrinsic viscosity of dendrimers. <i>Soft Matter</i> , 2010, 6, 2619.	2.7	26
79	Theory of Side-Chain Liquid Crystal Polymers: Bulk Behavior and Chain Conformation. <i>Macromolecules</i> , 2010, 43, 10096-10106.	4.8	22
80	Thermodynamic Properties of Block Copolymer Electrolytes Containing Imidazolium and Lithium Salts. <i>Macromolecules</i> , 2010, 43, 8282-8289.	4.8	131
81	Electrostatic Regulation of Genome Packaging in Human Hepatitis B Virus. <i>Biophysical Journal</i> , 2009, 96, 3065-3073.	0.5	27
82	Effects of Ion Solvation on the Miscibility of Binary Polymer Blends. <i>Journal of Physical Chemistry B</i> , 2008, 112, 16205-16213.	2.6	96
83	VARIATIONAL ELECTROSTATICS FOR CHARGE SOLVATION. <i>Journal of Theoretical and Computational Chemistry</i> , 2008, 07, 397-419.	1.8	15
84	Nucleation of membrane adhesions. <i>Physical Review E</i> , 2008, 77, 021906.	2.1	23
85	Metastable cluster intermediates in the condensation of charged macromolecule solutions. <i>Journal of Chemical Physics</i> , 2007, 127, 084912.	3.0	51
86	Polymer-Tethered Ligand-Receptor Interactions between Surfaces II. <i>Langmuir</i> , 2007, 23, 13024-13039.	3.5	18
87	Polymer-tethered ligand-receptor interactions between surfaces. <i>Journal of Polymer Science, Part B: Polymer Physics</i> , 2006, 44, 2621-2637.	2.1	12
88	Challenges and opportunities in polymer theory. <i>Journal of Polymer Science, Part B: Polymer Physics</i> , 2006, 44, 3445-3447.	2.1	1
89	End-to-end distance vector distribution with fixed end orientations for the wormlike chain model. <i>Physical Review E</i> , 2005, 72, 041802.	2.1	56
90	Nature of Disordered Micelles in Sphere-Forming Block Copolymer Melts. <i>Macromolecules</i> , 2005, 38, 1979-1988.	4.8	47

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91	DNA Packaging in Bacteriophage: Is Twist Important?. <i>Biophysical Journal</i> , 2005, 88, 3912-3923.	0.5	98
92	Exact Results for a Semiflexible Polymer Chain in an Aligning Field. <i>Macromolecules</i> , 2004, 37, 5814-5823.	4.8	73
93	Dynamics of Water near a Protein Surface. <i>Journal of Physical Chemistry B</i> , 2003, 107, 13218-13228.	2.6	100
94	Semiflexible polymer solutions. I. Phase behavior and single-chain statistics. <i>Journal of Chemical Physics</i> , 2003, 119, 13113-13128.	3.0	42
95	Nucleation of stable cylinders from a metastable lamellar phase in a diblock copolymer melt. <i>Journal of Chemical Physics</i> , 2003, 118, 10293-10305.	3.0	45
96	Nucleation in binary polymer blends: Effects of adding diblock copolymers. <i>Journal of Chemical Physics</i> , 2003, 118, 8997-9006.	3.0	11
97	Nucleation in binary polymer blends: A self-consistent field study. <i>Journal of Chemical Physics</i> , 2002, 116, 2289-2300.	3.0	67
98	Concentration fluctuation in binary polymer blends: ξ parameter, spinodal and Ginzburg criterion. <i>Journal of Chemical Physics</i> , 2002, 117, 481-500.	3.0	111
99	Computationally focusing the directed evolution of proteins. <i>Journal of Cellular Biochemistry</i> , 2001, 84, 58-63.	2.6	29
100	On the direct evaluation of the equilibrium distribution of clusters by simulation. II. <i>Journal of Chemical Physics</i> , 2001, 115, 6898-6906.	3.0	6
101	Transient instability upon temperature quench in weakly ordered block copolymers. <i>Journal of Chemical Physics</i> , 1999, 111, 10681-10688.	3.0	16
102	On the direct evaluation of the equilibrium distribution of clusters by simulation. <i>Journal of Chemical Physics</i> , 1999, 111, 9958-9964.	3.0	19
103	Interfacial Curvature in Graft and Diblock Copolymers and Implications for Long-Range Order in Cylindrical Morphologies. <i>Macromolecules</i> , 1997, 30, 6771-6782.	4.8	31
104	Chain Dimensions in Amorphous Polymer Melts. <i>Macromolecules</i> , 1995, 28, 570-576.	4.8	19
105	Chiral Symmetry Breaking and Pattern Formation in Two-Dimensional Films. <i>Materials Research Society Symposia Proceedings</i> , 1992, 292, 235.	0.1	1