Martin Krger

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234 8,557 47 82 g-index

260 9,583 4.7 6.69 ext. papers ext. citations avg, IF L-index

#	Paper	IF	Citations
234	Poly(N-isopropylacrylamide) Phase Diagrams: Fifty Years of Research. <i>Angewandte Chemie - International Edition</i> , 2015 , 54, 15342-67	16.4	557
233	Shortest multiple disconnected path for the analysis of entanglements in two- and three-dimensional polymeric systems. <i>Computer Physics Communications</i> , 2005 , 168, 209-232	4.2	310
232	Simple models for complex nonequilibrium fluids. <i>Physics Reports</i> , 2004 , 390, 453-551	27.7	238
231	Topological analysis of polymeric melts: chain-length effects and fast-converging estimators for entanglement length. <i>Physical Review E</i> , 2009 , 80, 031803	2.4	214
230	Shape effect in cellular uptake of PEGylated nanoparticles: comparison between sphere, rod, cube and disk. <i>Nanoscale</i> , 2015 , 7, 16631-46	7.7	204
229	Rheology and structural changes of polymer melts via nonequilibrium molecular dynamics. <i>Journal of Rheology</i> , 1993 , 37, 1057-1079	4.1	196
228	Rheological evidence for a dynamical crossover in polymer melts via nonequilibrium molecular dynamics. <i>Physical Review Letters</i> , 2000 , 85, 1128-31	7.4	190
227	Crossover from the Rouse to the Entangled Polymer Melt Regime: Signals from Long, Detailed Atomistic Molecular Dynamics Simulations, Supported by Rheological Experiments. <i>Macromolecules</i> , 2003 , 36, 1376-1387	5.5	179
226	Primitive Path Networks Generated by Annealing and Geometrical Methods: Insights into Differences. <i>Macromolecules</i> , 2007 , 40, 2897-2903	5.5	160
225	Challenges in Multiscale Modeling of Polymer Dynamics. <i>Polymers</i> , 2013 , 5, 751-832	4.5	143
224	Endocytosis of PEGylated nanoparticles accompanied by structural and free energy changes of the grafted polyethylene glycol. <i>Biomaterials</i> , 2014 , 35, 8467-78	15.6	142
223	Nanoparticle effect on the dynamics of polymer chains and their entanglement network. <i>Physical Review Letters</i> , 2012 , 109, 118001	7.4	141
222	Primitive Path Identification and Entanglement Statistics in Polymer Melts: Results from Direct Topological Analysis on Atomistic Polyethylene Models. <i>Macromolecules</i> , 2006 , 39, 4207-4216	5.5	137
221	Flow Effects on Melt Structure and Entanglement Network of Linear Polymers: Results from a Nonequilibrium Molecular Dynamics Simulation Study of a Polyethylene Melt in Steady Shear. <i>Macromolecules</i> , 2010 , 43, 6886-6902	5.5	128
220	Combined molecular algorithms for the generation, equilibration and topological analysis of entangled polymers: methodology and performance. <i>International Journal of Molecular Sciences</i> , 2009 , 10, 5054-89	6.3	124
219	Tuning polymer thickness: synthesis and scaling theory of homologous series of dendronized polymers. <i>Journal of the American Chemical Society</i> , 2009 , 131, 11841-54	16.4	121
218	Effect of charge, hydrophobicity, and sequence of nucleoporins on the translocation of model particles through the nuclear pore complex. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013 , 110, 3363-8	11.5	113

(2014-1997)

217	Polymer Melts under Uniaxial Elongational Flow: Stress (Dptical Behavior from Experiments and Nonequilibrium Molecular Dynamics Computer Simulations. <i>Macromolecules</i> , 1997 , 30, 526-539	5.5	106
216	The largest synthetic structure with molecular precision: towards a molecular object. <i>Angewandte Chemie - International Edition</i> , 2011 , 50, 737-40	16.4	102
215	A thermodynamically admissible reptation model for fast flows of entangled polymers. II. Model predictions for shear and extensional flows. <i>Journal of Rheology</i> , 2000 , 44, 1293-1317	4.1	100
214	Viscoelastic flows studied by smoothed particle dynamics. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2002 , 105, 35-51	2.7	98
213	Modeling of Polymer Structure and Conformations in Polymer Nanocomposites from Atomistic to Mesoscale: A Review. <i>Polymer Reviews</i> , 2016 , 56, 385-428	14	97
212	A predictive multiscale computational framework for viscoelastic properties of linear polymers. <i>Polymer</i> , 2012 , 53, 5935-5952	3.9	91
211	Nanoparticle Geometrical Effect on Structure, Dynamics and Anisotropic Viscosity of Polyethylene Nanocomposites. <i>Macromolecules</i> , 2012 , 45, 2099-2112	5.5	87
21 0	Quantifying chain reptation in entangled polymer melts: topological and dynamical mapping of atomistic simulation results onto the tube model. <i>Journal of Chemical Physics</i> , 2010 , 132, 124904	3.9	87
209	Structure and Dynamics of Dilute Polymer Solutions under Shear Flow via Nonequilibrium Molecular Dynamics. <i>Macromolecules</i> , 1999 , 32, 5660-5672	5.5	84
208	Direct observation of the dynamics of semiflexible polymers in shear flow. <i>Physical Review Letters</i> , 2013 , 110, 108302	7.4	83
207	Influence of Nanorod Inclusions on Structure and Primitive Path Network of Polymer Nanocomposites at Equilibrium and Under Deformation. <i>Macromolecules</i> , 2011 , 44, 1034-1045	5.5	83
206	A theoretical evaluation of the effects of carbon nanotube entanglement and bundling on the structural and mechanical properties of buckypaper. <i>Carbon</i> , 2012 , 50, 1793-1806	10.4	81
205	Morphology control of hairy nanopores. ACS Nano, 2011, 5, 4737-47	16.7	80
204	Collapse of Thermoresponsive Brushes and the Tuning of Protein Adsorption. <i>Macromolecules</i> , 2011 , 44, 6986-7005	5.5	79
203	From Dendrimers to Dendronized Polymers and Forests: Scaling Theory and its Limitations. <i>Macromolecules</i> , 2010 , 43, 6213-6224	5.5	76
202	Adsorption of core-shell nanoparticles at liquid I quid interfaces. Soft Matter, 2011, 7, 7663	3.6	75
201	Molecular simulation guided constitutive modeling on finite strain viscoelasticity of elastomers. Journal of the Mechanics and Physics of Solids, 2016 , 88, 204-226	5	67
200	Dynamic structure of unentangled polymer chains in the vicinity of non-attractive nanoparticles. <i>Soft Matter</i> , 2014 , 10, 1723-37	3.6	67

199	Wormlike micelles under shear flow: A microscopic model studied by nonequilibrium-molecular-dynamics computer simulations. <i>Physical Review E</i> , 1996 , 53, 2531-2536	2.4	66
198	Colloid-Brush Interactions: The Effect of Solvent Quality. <i>Macromolecules</i> , 2011 , 44, 3622-3638	5.5	61
197	Simple, admissible, and accurate approximants of the inverse Langevin and Brillouin functions, relevant for strong polymer deformations and flows. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2015 , 223, 77-87	2.7	58
196	Variance reduced Brownian simulation of a bead-spring chain under steady shear flow considering hydrodynamic interaction effects. <i>Journal of Chemical Physics</i> , 2000 , 113, 4767-4773	3.9	58
195	Dendronized Polymers: Molecular Objects between Conventional Linear Polymers and Colloidal Particles. <i>ACS Macro Letters</i> , 2014 , 3, 991-998	6.6	56
194	Structure, dimensions, and entanglement statistics of long linear polyethylene chains. <i>Journal of Physical Chemistry B</i> , 2009 , 113, 442-55	3.4	56
193	Chaotic orientational behavior of a nematic liquid crystal subjected to a steady shear flow. <i>Physical Review E</i> , 2002 , 66, 040702	2.4	55
192	Ternary protein adsorption onto brushes: strong versus weak. <i>Langmuir</i> , 2009 , 25, 11621-34	4	54
191	Chaotic and regular shear-induced orientational dynamics of nematic liquid crystals. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2002 , 315, 537-568	3.3	51
190	Molecular dynamics simulations of polymer crystallization under confinement: Entanglement effect. <i>Polymer</i> , 2017 , 109, 71-84	3.9	49
190 189		3.9 15.6	49 49
	effect. <i>Polymer</i> , 2017 , 109, 71-84 Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive		
189	effect. <i>Polymer</i> , 2017 , 109, 71-84 Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive polymer brushes. <i>Biomaterials</i> , 2012 , 33, 4975-87 Rheological and entanglement characteristics of linear-chain polyethylene liquids in planar Couette	15.6	49
189 188	Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive polymer brushes. <i>Biomaterials</i> , 2012 , 33, 4975-87 Rheological and entanglement characteristics of linear-chain polyethylene liquids in planar Couette and planar elongational flows. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008 , 152, 168-183 Systematic time-scale-bridging molecular dynamics applied to flowing polymer melts. <i>Physical</i>	15.6 2.7	49
189 188 187	Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive polymer brushes. <i>Biomaterials</i> , 2012 , 33, 4975-87 Rheological and entanglement characteristics of linear-chain polyethylene liquids in planar Couette and planar elongational flows. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008 , 152, 168-183 Systematic time-scale-bridging molecular dynamics applied to flowing polymer melts. <i>Physical Review E</i> , 2009 , 79, 011802 Molecular dynamics of model liquid crystals composed of semiflexible molecules. <i>Physical Review E</i> ,	15.6 2.7 2.4	49 48 47
189 188 187 186	Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive polymer brushes. <i>Biomaterials</i> , 2012 , 33, 4975-87 Rheological and entanglement characteristics of linear-chain polyethylene liquids in planar Couette and planar elongational flows. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008 , 152, 168-183 Systematic time-scale-bridging molecular dynamics applied to flowing polymer melts. <i>Physical Review E</i> , 2009 , 79, 011802 Molecular dynamics of model liquid crystals composed of semiflexible molecules. <i>Physical Review E</i> , 1996 , 54, 5178-5186 Aggregation of polyethylene glycol polymers suppresses receptor-mediated endocytosis of	15.6 2.7 2.4	49 48 47 47
189 188 187 186	Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive polymer brushes. <i>Biomaterials</i> , 2012 , 33, 4975-87 Rheological and entanglement characteristics of linear-chain polyethylene liquids in planar Couette and planar elongational flows. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008 , 152, 168-183 Systematic time-scale-bridging molecular dynamics applied to flowing polymer melts. <i>Physical Review E</i> , 2009 , 79, 011802 Molecular dynamics of model liquid crystals composed of semiflexible molecules. <i>Physical Review E</i> , 1996 , 54, 5178-5186 Aggregation of polyethylene glycol polymers suppresses receptor-mediated endocytosis of PEGylated liposomes. <i>Nanoscale</i> , 2018 , 10, 4545-4560	15.6 2.7 2.4 2.4 7.7	49 48 47 47 46

(1997-2008)

181	Universal scaling, entanglements, and knots of model chain molecules. <i>Physical Review Letters</i> , 2008 , 101, 265702	7.4	44
180	Consistent closure schemes for statistical models of anisotropic fluids. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008 , 149, 40-55	2.7	44
179	Magnetoviscosity of semidilute ferrofluids and the role of dipolar interactions: comparison of molecular simulations and dynamical mean-field theory. <i>Physical Review E</i> , 2005 , 71, 031205	2.4	44
178	Viscoelasticity of polymeric melts and concentrated solutions. The effect of flow-induced alignment of chain ends. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1993 , 195, 336-353	3.3	44
177	Kinetics of gene derepression by ERK signaling. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013 , 110, 10330-5	11.5	43
176	Thermomechanical properties of the WCAllennard-Jones model system in its fluid and solid states. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1998 , 250, 58-82	3.3	43
175	Height and Width of Adsorbed Dendronized Polymers: Electron and Atomic Force Microscopy of Homologous Series. <i>Macromolecules</i> , 2011 , 44, 6785-6792	5.5	41
174	Rotation and Deformation of a Finitely Extendable Flexible Polymer Molecule in a Steady Shear Flow. <i>Macromolecules</i> , 2002 , 35, 8621-8630	5.5	41
173	Random packing of model polymers: local structure, topological hindrance and universal scaling. <i>Soft Matter</i> , 2009 , 5, 1762	3.6	40
172	Boundaries steer the contraction of active gels. <i>Nature Communications</i> , 2016 , 7, 13120	17.4	39
172 171	Boundaries steer the contraction of active gels. <i>Nature Communications</i> , 2016 , 7, 13120 Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025	17.4 5.5	39 39
<u> </u>	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics		
171	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025 Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments.	5.5	39
171	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025 Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments. <i>Langmuir</i> , 2015 , 31, 4798-805 Evolution of fivefold local symmetry during crystal nucleation and growth in dense hard-sphere	5.5	39
171 170 169	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025 Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments. <i>Langmuir</i> , 2015 , 31, 4798-805 Evolution of fivefold local symmetry during crystal nucleation and growth in dense hard-sphere packings. <i>Soft Matter</i> , 2012 , 8, 844-858 Magnetization dynamics, rheology, and an effective description of ferromagnetic units in dilute suspension. <i>Physical Review E</i> , 2002 , 66, 021501 Efficient hybrid algorithm for the dynamic creation of wormlike chains in solutions, brushes, melts	5·5 4 3.6	39 38 38
171 170 169	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025 Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments. <i>Langmuir</i> , 2015 , 31, 4798-805 Evolution of fivefold local symmetry during crystal nucleation and growth in dense hard-sphere packings. <i>Soft Matter</i> , 2012 , 8, 844-858 Magnetization dynamics, rheology, and an effective description of ferromagnetic units in dilute suspension. <i>Physical Review E</i> , 2002 , 66, 021501 Efficient hybrid algorithm for the dynamic creation of wormlike chains in solutions, brushes, melts	5·5 4 3.6 2.4	39 38 38 36
171 170 169 168	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025 Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments. <i>Langmuir</i> , 2015 , 31, 4798-805 Evolution of fivefold local symmetry during crystal nucleation and growth in dense hard-sphere packings. <i>Soft Matter</i> , 2012 , 8, 844-858 Magnetization dynamics, rheology, and an effective description of ferromagnetic units in dilute suspension. <i>Physical Review E</i> , 2002 , 66, 021501 Efficient hybrid algorithm for the dynamic creation of wormlike chains in solutions, brushes, melts and glasses. <i>Computer Physics Communications</i> , 1999 , 118, 278-298 Viscosity coefficients for anisotropic, nematic fluids based on structural theories of suspensions.	5.5 4 3.6 2.4 4.2	39 38 38 36 36
171 170 169 168 167	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016 , 49, 9017-9025 Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments. <i>Langmuir</i> , 2015 , 31, 4798-805 Evolution of fivefold local symmetry during crystal nucleation and growth in dense hard-sphere packings. <i>Soft Matter</i> , 2012 , 8, 844-858 Magnetization dynamics, rheology, and an effective description of ferromagnetic units in dilute suspension. <i>Physical Review E</i> , 2002 , 66, 021501 Efficient hybrid algorithm for the dynamic creation of wormlike chains in solutions, brushes, melts and glasses. <i>Computer Physics Communications</i> , 1999 , 118, 278-298 Viscosity coefficients for anisotropic, nematic fluids based on structural theories of suspensions. <i>Journal of Chemical Physics</i> , 1995 , 103, 807-817 Viscoelasticity of carbon nanotube buckypaper: zippinglinzipping mechanism and entanglement	5.5 4 3.6 2.4 4.2 3.9	39 38 38 36 36

163	Primitive-path statistics of entangled polymers: mapping multi-chain simulations onto single-chain mean-field models. <i>New Journal of Physics</i> , 2014 , 16, 015027	2.9	33
162	Projection from an atomistic chain contour to its primitive path. <i>Polymer</i> , 2002 , 43, 477-487	3.9	33
161	Magnetoviscous model fluids. <i>Journal of Physics Condensed Matter</i> , 2003 , 15, S1403-S1423	1.8	31
160	Linear Viscoelastic Behavior of Unentangled Polymer Melts via Non-Equilibrium Molecular Dynamics. <i>Macromolecular Theory and Simulations</i> , 2004 , 13, 748-753	1.5	31
159	Regular and chaotic orientational and rheological behaviour of liquid crystals. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S3835-S3859	1.8	31
158	Phase behavior and structure of Janus fluids. <i>Physical Review E</i> , 2003 , 67, 041209	2.4	31
157	Self-Folding of Charged Single Dendronized Polymers. <i>Advanced Materials</i> , 2008 , 20, 3204-3210	24	30
156	Canonical distribution functions in polymer dynamics. (II). Liquid-crystalline polymers. <i>Physica A:</i> Statistical Mechanics and Its Applications, 2003 , 319, 134-150	3.3	30
155	Thermoresponsive cell culture substrates based on PNIPAM brushes functionalized with adhesion peptides: theoretical considerations of mechanism and design. <i>Langmuir</i> , 2012 , 28, 16623-37	4	29
154	Analytical solution of the SIR-model for the temporal evolution of epidemics. Part A: time-independent reproduction factor. <i>Journal of Physics A: Mathematical and Theoretical</i> , 2020 , 53, 5	05601	29
153	Covid-19 Predictions Using a Gauss Model, Based on Data from April 2 2020 , 2, 197-212	2.1	28
152	Understanding Dynamics in Binary Mixtures of Entangledcis-1,4-Polybutadiene Melts at the Level of Primitive Path Segments by Mapping Atomistic Simulation Data onto the Tube Model. <i>Macromolecules</i> , 2010 , 43, 8239-8250	5.5	28
151	Rheology and Packing of Dendronized Polymers. <i>Macromolecules</i> , 2016 , 49, 7054-7068	5.5	28
150	Microscopic Origin of the Non-Newtonian Viscosity of Semiflexible Polymer Solutions in the Semidilute Regime <i>ACS Macro Letters</i> , 2014 , 3, 136-140	6.6	27
149	Effect of polymer solvent on the mechanical properties of entangled polymer gels: Coarse-grained molecular simulation. <i>Polymer</i> , 2013 , 54, 2555-2564	3.9	27
148	Rheology: From simple and to complex fluids. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1997 , 240, 126-144	3.3	27
147	Magnetoviscosity and orientational order parameters of dilute ferrofluids. Journal of Chemical	2.0	27
	Physics, 2002 , 116, 9078-9088	3.9	<u>-</u> /

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145	Fast equilibration protocol for million atom systems of highly entangled linear polyethylene chains. Journal of Chemical Physics, 2016 , 144, 154901	3.9	27	
144	Ordering and Crystallization of Entangled Polyethylene Melts under Uniaxial Tension: A Molecular Dynamics Study. <i>Macromolecules</i> , 2018 , 51, 9635-9648	5.5	27	
143	Molecularly derived constitutive equation for low-molecular polymer melts from thermodynamically guided simulation. <i>Journal of Rheology</i> , 2011 , 55, 69-93	4.1	26	
142	Formation of double helical and filamentous structures in models of physical and chemical gels. <i>Soft Matter</i> , 2007 , 4, 18-28	3.6	26	
141	Communication: Appearance of undershoots in start-up shear: Experimental findings captured by tumbling-snake dynamics. <i>Journal of Chemical Physics</i> , 2017 , 146, 161101	3.9	25	
140	Structure and rheology of model-ferrofluids under shear flow. <i>Journal of Magnetism and Magnetic Materials</i> , 2005 , 289, 325-327	2.8	25	
139	A simple example for comparing GENERIC with rational non-equilibrium thermodynamics. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2000 , 285, 448-466	3.3	25	
138	Self-assembly of core-polyethylene glycol-lipid shell (CPLS) nanoparticles and their potential as drug delivery vehicles. <i>Nanoscale</i> , 2016 , 8, 14821-35	7.7	25	
137	Computer simulation of dendronized polymers: organization and characterization at the atomistic level. <i>RSC Advances</i> , 2013 , 3, 126-140	3.7	24	
136	Recognition and analysis of local structure in polycrystalline configurations. <i>Computer Physics Communications</i> , 2002 , 145, 371-384	4.2	24	
135	Automated symbolic calculations in nonequilibrium thermodynamics. <i>Computer Physics Communications</i> , 2010 , 181, 2149-2157	4.2	23	
134	Anisotropy of the magnetoviscous effect in ferrofluids. <i>Physical Review E</i> , 2005 , 71, 051201	2.4	23	
133	Modeling of Entangled Polymer Diffusion in Melts and Nanocomposites: A Review. <i>Polymers</i> , 2019 , 11,	4.5	22	
132	Fibers with integrated mechanochemical switches: minimalistic design principles derived from fibronectin. <i>Biophysical Journal</i> , 2012 , 103, 1909-18	2.9	22	
131	Nonaffine deformation of inherent structure as a static signature of cooperativity in supercooled liquids. <i>Physical Review Letters</i> , 2008 , 101, 095501	7.4	22	
130	An extended FENE dumbbell theory for concentration dependent shear-induced anisotropy in dilute polymer solutions. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 1996 , 62, 235-251	2.7	22	
129	Unraveling two-dimensional polymerization in the single crystal. <i>Journal of Applied Crystallography</i> , 2018 , 51, 481-497	3.8	21	
128	Influence of molecular architecture on the entanglement network: topological analysis of linear, long- and short-chain branched polyethylene melts via Monte Carlo simulations. <i>Soft Matter</i> , 2016 , 12, 3770-86	3.6	21	

127	The effect of polymer chain length on the mechanical properties of triblock copolymer gels. <i>Chemical Physics Letters</i> , 2014 , 612, 157-161	2.5	21
126	Computational study on entanglement length and pore size of carbon nanotube buckypaper. <i>Applied Physics Letters</i> , 2012 , 100, 021907	3.4	21
125	From atomistic simulation to the dynamics, structure and helical network formation of dendronized polymers: the Janus chain model. <i>Journal of Chemical Physics</i> , 2007 , 127, 094904	3.9	21
124	Effect of Crosslinking on the Microtribological Behavior of Model Polymer Brushes. <i>Tribology Letters</i> , 2016 , 63, 1	2.8	20
123	Computer simulation of fifth generation dendronized polymers: impact of charge on internal organization. <i>Journal of Physical Chemistry B</i> , 2013 , 117, 6007-17	3.4	20
122	Writhe and mutual entanglement combine to give the entanglement length. <i>Physical Review E</i> , 2013 , 88, 062604	2.4	20
121	The Largest Synthetic Structure with Molecular Precision: Towards a Molecular Object. <i>Angewandte Chemie</i> , 2011 , 123, 763-766	3.6	20
120	Lubricated optical rheometer for the study of two-dimensional complex flows of polymer melts. Journal of Non-Newtonian Fluid Mechanics, 2008, 150, 43-55	2.7	20
119	Self-assembled core-polyethylene glycol-lipid shell nanoparticles demonstrate high stability in shear flow. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 13294-13306	3.6	19
118	Size of graphene sheets determines the structural and mechanical properties of 3D graphene foams. <i>Nanotechnology</i> , 2018 , 29, 104001	3.4	19
117	Synthetic regimes due to packing constraints in dendritic molecules confirmed by labelling experiments. <i>Nature Communications</i> , 2013 , 4, 1993	17.4	19
116	Anisotropic self-diffusion in ferrofluids studied via Brownian dynamics simulations. <i>Physical Review E</i> , 2005 , 72, 031504	2.4	19
115	Influence of Chain Stiffness, Grafting Density and Normal Load on the Tribological and Structural Behavior of Polymer Brushes: A Nonequilibrium-Molecular-Dynamics Study. <i>Polymers</i> , 2016 , 8,	4.5	19
114	Using mesoscopic models to design strong and tough biomimetic polymer networks. <i>Langmuir</i> , 2011 , 27, 13796-805	4	18
113	Derivation of Frank-Ericksen elastic coefficients for polydomain nematics from mean-field molecular theory for anisotropic particles. <i>Journal of Chemical Physics</i> , 2007 , 127, 034903	3.9	18
112	Dynamics of colloidal suspensions of ferromagnetic particles in plane Couette flow: comparison of approximate solutions with Brownian dynamics simulations. <i>Physical Review E</i> , 2003 , 67, 061401	2.4	18
111	NEMD-Computersimulation zur Rheologie von Polymerschmelzen / NEMD Computer Simulation of Polymer Melt Rheology. <i>Applied Rheology</i> , 1995 , 5, 66-71	1.2	18
110	Filamentous networks in phase-separating two-dimensional gels. <i>Europhysics Letters</i> , 2007 , 77, 58007	1.6	17

(2001-2007)

109	From hyperbolic regularization to exact hydrodynamics for linearized Grad's equations. <i>Physical Review E</i> , 2007 , 75, 051204	2.4	17	
108	Analytical solution of the SIR-model for the temporal evolution of epidemics: part B. Semi-time case. <i>Journal of Physics A: Mathematical and Theoretical</i> , 2021 , 54, 175601	2	17	
107	Interactions in dendronized polymers: intramolecular dominates intermolecular. <i>Soft Matter</i> , 2014 , 10, 1032-44	3.6	16	
106	Structure Elucidation of 2D Polymer Monolayers Based on Crystallization Estimates Derived from Tip-Enhanced Raman Spectroscopy (TERS) Polymerization Conversion Data. <i>Journal of the American Chemical Society</i> , 2019 , 141, 9867-9871	16.4	15	
105	A Detailed Examination of the Topological Constraints of Lamellae-Forming Block Copolymers. <i>Macromolecules</i> , 2018 , 51, 2110-2124	5.5	15	
104	Poly(N-isopropylacrylamid)-Phasendiagramme: 50 Jahre Forschung. <i>Angewandte Chemie</i> , 2015 , 127, 15	5 5.& -15	558%	
103	Exact linear hydrodynamics from the Boltzmann equation. <i>Physical Review Letters</i> , 2008 , 100, 214503	7.4	15	
102	Structural changes and viscoplastic behavior of a generic embedded-atom model metal in steady shear flow. <i>Physical Review E</i> , 2004 , 69, 021509	2.4	15	
101	On the Shape and Rheology of Linear Micelles in Dilute Solutions. <i>Journal De Physique II</i> , 1997 , 7, 931-9	46	15	
100	Dynamics of interacting magnetic nanoparticles: effective behavior from competition between Brownian and NBI relaxation. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 22244-22259	3.6	15	
99	Boltzmann equation and hydrodynamic fluctuations. <i>Physical Review E</i> , 2009 , 80, 051202	2.4	14	
98	Crossover between short- and long-time behavior of stress fluctuations and viscoelasticity of liquids. <i>Physical Review E</i> , 2003 , 67, 042201	2.4	14	
97	Thermophysical properties of gases, liquids, and solids composed of particles interacting with a short-range attractive potential. <i>Physical Review E</i> , 2001 , 64, 011201	2.4	14	
96	Combined Experimental and Simulation Studies of Cross-Linked Polymer Brushes under Shear. <i>Macromolecules</i> , 2018 , 51, 10174-10183	5.5	14	
95	Assessing numerical methods for molecular and particle simulation. <i>Soft Matter</i> , 2017 , 13, 8565-8578	3.6	13	
94	Carbon Nanotube Length Governs the Viscoelasticity and Permeability of Buckypaper. <i>Polymers</i> , 2017 , 9,	4.5	13	
93	Beyond-equilibrium molecular dynamics of a rarefied gas subjected to shear flow. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2004 , 120, 175-187	2.7	13	
92	Symbolic test of the Jacobi identity for given generalized P oisson[bracket. <i>Computer Physics Communications</i> , 2001 , 137, 325-340	4.2	13	

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LIST OF PUBLICATIONS

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