

# Martin KrÄƒger

## List of Publications by Year in descending order

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242  
papers

10,461  
citations

36203

51  
h-index

43802

91  
g-index

260  
all docs

260  
docs citations

260  
times ranked

7937  
citing authors

#	ARTICLE	IF	CITATIONS
1	Poly( <i>N</i> -isopropylacrylamide) Phase Diagrams: Fifty Years of Research. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 15342-15367.	7.2	772
2	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.	3.0	387
3	Simple models for complex nonequilibrium fluids. <i>Physics Reports</i> , 2004, 390, 453-551.	10.3	268
4	Shape effect in cellular uptake of PEGylated nanoparticles: comparison between sphere, rod, cube and disk. <i>Nanoscale</i> , 2015, 7, 16631-16646.	2.8	268
5	Topological analysis of polymeric melts: Chain-length effects and fast-converging estimators for entanglement length. <i>Physical Review E</i> , 2009, 80, 031803.	0.8	260
6	Rheology and structural changes of polymer melts via nonequilibrium molecular dynamics. <i>Journal of Rheology</i> , 1993, 37, 1057-1079.	1.3	214
7	Rheological Evidence for a Dynamical Crossover in Polymer Melts via Nonequilibrium Molecular Dynamics. <i>Physical Review Letters</i> , 2000, 85, 1128-1131.	2.9	209
8	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.	2.2	198
9	Primitive Path Networks Generated by Annealing and Geometrical Methods: Insights into Differences. <i>Macromolecules</i> , 2007, 40, 2897-2903.	2.2	188
10	Endocytosis of PEGylated nanoparticles accompanied by structural and free energy changes of the grafted polyethylene glycol. <i>Biomaterials</i> , 2014, 35, 8467-8478.	5.7	176
11	Challenges in Multiscale Modeling of Polymer Dynamics. <i>Polymers</i> , 2013, 5, 751-832.	2.0	173
12	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, 5054-5089.	1.8	160
13	Nanoparticle Effect on the Dynamics of Polymer Chains and Their Entanglement Network. <i>Physical Review Letters</i> , 2012, 109, 118001.	2.9	160
14	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.	2.2	152
15	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.	2.2	146
16	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, 3363-3368.	3.3	139
17	Tuning Polymer Thickness: Synthesis and Scaling Theory of Homologous Series of Dendronized Polymers. <i>Journal of the American Chemical Society</i> , 2009, 131, 11841-11854.	6.6	130
18	Polymer Melts under Uniaxial Elongational Flow: Stress-Optical Behavior from Experiments and Nonequilibrium Molecular Dynamics Computer Simulations. <i>Macromolecules</i> , 1997, 30, 526-539.	2.2	117

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19	Viscoelastic flows studied by smoothed particle dynamics. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2002, 105, 35-51.	1.0	116
20	A predictive multiscale computational framework for viscoelastic properties of linear polymers. <i>Polymer</i> , 2012, 53, 5935-5952.	1.8	115
21	Modeling of Polymer Structure and Conformations in Polymer Nanocomposites from Atomistic to Mesoscale: A Review. <i>Polymer Reviews</i> , 2016, 56, 385-428.	5.3	114
22	The Largest Synthetic Structure with Molecular Precision: Towards a Molecular Object. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 737-740.	7.2	111
23	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.	1.3	108
24	Direct Observation of the Dynamics of Semiflexible Polymers in Shear Flow. <i>Physical Review Letters</i> , 2013, 110, 108302.	2.9	102
25	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.	1.2	101
26	Structure and Dynamics of Dilute Polymer Solutions under Shear Flow via Nonequilibrium Molecular Dynamics. <i>Macromolecules</i> , 1999, 32, 5660-5672.	2.2	99
27	Nanoparticle Geometrical Effect on Structure, Dynamics and Anisotropic Viscosity of Polyethylene Nanocomposites. <i>Macromolecules</i> , 2012, 45, 2099-2112.	2.2	99
28	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.	5.4	97
29	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.	2.2	91
30	Morphology Control of Hairy Nanopores. <i>ACS Nano</i> , 2011, 5, 4737-4747.	7.3	89
31	Molecular simulation guided constitutive modeling on finite strain viscoelasticity of elastomers. <i>Journal of the Mechanics and Physics of Solids</i> , 2016, 88, 204-226.	2.3	87
32	Collapse of Thermoresponsive Brushes and the Tuning of Protein Adsorption. <i>Macromolecules</i> , 2011, 44, 6986-7005.	2.2	85
33	From Dendrimers to Dendronized Polymers and Forests: Scaling Theory and its Limitations. <i>Macromolecules</i> , 2010, 43, 6213-6224.	2.2	80
34	Adsorption of core-shell nanoparticles at liquid-liquid interfaces. <i>Soft Matter</i> , 2011, 7, 7663.	1.2	78
35	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.	1.0	75
36	Molecular dynamics simulations of polymer crystallization under confinement: Entanglement effect. <i>Polymer</i> , 2017, 109, 71-84.	1.8	75

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37	Wormlike micelles under shear flow: A microscopic model studied by nonequilibrium-molecular-dynamics computer simulations. <i>Physical Review E</i> , 1996, 53, 2531-2536.	0.8	74
38	Dynamic structure of unentangled polymer chains in the vicinity of non-attractive nanoparticles. <i>Soft Matter</i> , 2014, 10, 1723.	1.2	73
39	Colloid-Brush Interactions: The Effect of Solvent Quality. <i>Macromolecules</i> , 2011, 44, 3622-3638.	2.2	69
40	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.	1.2	65
41	Dendronized Polymers: Molecular Objects between Conventional Linear Polymers and Colloidal Particles. <i>ACS Macro Letters</i> , 2014, 3, 991-998.	2.3	62
42	Ternary Protein Adsorption onto Brushes: Strong versus Weak. <i>Langmuir</i> , 2009, 25, 11621-11634.	1.6	61
43	A Two-Dimensional Polymer Synthesized at the Air/Water Interface. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 10584-10588.	7.2	61
44	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, 505601.	0.7	61
45	Structure, Dimensions, and Entanglement Statistics of Long Linear Polyethylene Chains. <i>Journal of Physical Chemistry B</i> , 2009, 113, 442-455.	1.2	60
46	Aggregation of polyethylene glycol polymers suppresses receptor-mediated endocytosis of PEGylated liposomes. <i>Nanoscale</i> , 2018, 10, 4545-4560.	2.8	60
47	Primitive chain network study on uncrosslinked and crosslinked cis-polyisoprene polymers. <i>Polymer</i> , 2011, 52, 5867-5878.	1.8	59
48	Chaotic orientational behavior of a nematic liquid crystal subjected to a steady shear flow. <i>Physical Review E</i> , 2002, 66, 040702.	0.8	58
49	Molecular dynamics of model liquid crystals composed of semiflexible molecules. <i>Physical Review E</i> , 1996, 54, 5178-5186.	0.8	56
50	Entanglements and Crystallization of Concentrated Polymer Solutions: Molecular Dynamics Simulations. <i>Macromolecules</i> , 2016, 49, 9017-9025.	2.2	56
51	Chaotic and regular shear-induced orientational dynamics of nematic liquid crystals. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2002, 315, 537-568.	1.2	54
52	Theoretical considerations on mechanisms of harvesting cells cultured on thermoresponsive polymer brushes. <i>Biomaterials</i> , 2012, 33, 4975-4987.	5.7	53
53	Polymer Brushes under Shear: Molecular Dynamics Simulations Compared to Experiments. <i>Langmuir</i> , 2015, 31, 4798-4805.	1.6	53
54	Consistent closure schemes for statistical models of anisotropic fluids. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008, 149, 40-55.	1.0	52

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55	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.	1.0	51
56	Kinetics of gene derepression by ERK signaling. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013, 110, 10330-10335.	3.3	51
57	Boundaries steer the contraction of active gels. <i>Nature Communications</i> , 2016, 7, 13120.	5.8	50
58	Thermomechanical properties of the WCA Lennard-Jones model system in its fluid and solid states. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1998, 250, 58-82.	1.2	49
59	Universal Scaling, Entanglements, and Knots of Model Chain Molecules. <i>Physical Review Letters</i> , 2008, 101, 265702.	2.9	49
60	Ordering and Crystallization of Entangled Polyethylene Melts under Uniaxial Tension: A Molecular Dynamics Study. <i>Macromolecules</i> , 2018, 51, 9635-9648.	2.2	49
61	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.	0.8	48
62	Systematic time-scale-bridging molecular dynamics applied to flowing polymer melts. <i>Physical Review E</i> , 2009, 79, 011802.	0.8	48
63	Entanglements and Dynamics of Polymer Melts near a SWCNT. <i>Macromolecules</i> , 2012, 45, 7274-7281.	2.2	48
64	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.	1.2	47
65	Modeling of Entangled Polymer Diffusion in Melts and Nanocomposites: A Review. <i>Polymers</i> , 2019, 11, 876.	2.0	47
66	Height and Width of Adsorbed Dendronized Polymers: Electron and Atomic Force Microscopy of Homologous Series. <i>Macromolecules</i> , 2011, 44, 6785-6792.	2.2	46
67	Evolution of fivefold local symmetry during crystal nucleation and growth in dense hard-sphere packings. <i>Soft Matter</i> , 2012, 8, 844-858.	1.2	46
68	Covid-19 Predictions Using a Gauss Model, Based on Data from April 2. <i>Physics</i> , 2020, 2, 197-212.	0.5	46
69	Viscosity coefficients for anisotropic, nematic fluids based on structural theories of suspensions. <i>Journal of Chemical Physics</i> , 1995, 103, 807-817.	1.2	45
70	Rotation and Deformation of a Finitely Extendable Flexible Polymer Molecule in a Steady Shear Flow. <i>Macromolecules</i> , 2002, 35, 8621-8630.	2.2	45
71	Random packing of model polymers: local structure, topological hindrance and universal scaling. <i>Soft Matter</i> , 2009, 5, 1762.	1.2	45
72	Viscoelasticity of carbon nanotube buckypaper: zipping/unzipping mechanism and entanglement effects. <i>Soft Matter</i> , 2012, 8, 7822.	1.2	44

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73	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.	3.0	42
74	Magnetization dynamics, rheology, and an effective description of ferromagnetic units in dilute suspension. <i>Physical Review E</i> , 2002, 66, 021501.	0.8	42
75	Dynamics of interacting magnetic nanoparticles: effective behavior from competition between Brownian and Néel relaxation. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 22244-22259.	1.3	41
76	Projection from an atomistic chain contour to its primitive path. <i>Polymer</i> , 2002, 43, 477-487.	1.8	38
77	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.	0.7	38
78	Shear modulus of fluids and solids. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1997, 239, 449-466.	1.2	37
79	Linear Viscoelastic Behavior of Unentangled Polymer Melts via Non-Equilibrium Molecular Dynamics. <i>Macromolecular Theory and Simulations</i> , 2004, 13, 748-753.	0.6	37
80	Effect of polymer solvent on the mechanical properties of entangled polymer gels: Coarse-grained molecular simulation. <i>Polymer</i> , 2013, 54, 2555-2564.	1.8	37
81	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.	1.2	37
82	Regular and chaotic orientational and rheological behaviour of liquid crystals. <i>Journal of Physics Condensed Matter</i> , 2004, 16, S3835-S3859.	0.7	35
83	Atomistic Modeling of Plastic Deformation in Semicrystalline Polyethylene: Role of Interphase Topology, Entanglements, and Chain Dynamics. <i>Macromolecules</i> , 2020, 53, 4605-4617.	2.2	35
84	Magnetoviscosity and orientational order parameters of dilute ferrofluids. <i>Journal of Chemical Physics</i> , 2002, 116, 9078-9088.	1.2	34
85	Canonical distribution functions in polymer dynamics. (II). Liquid-crystalline polymers. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2003, 319, 134-150.	1.2	34
86	Magnetoviscous model fluids. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S1403-S1423.	0.7	34
87	Rheology and Packing of Dendronized Polymers. <i>Macromolecules</i> , 2016, 49, 7054-7068.	2.2	34
88	Reconfigurable artificial microswimmers with internal feedback. <i>Nature Communications</i> , 2021, 12, 4762.	5.8	34
89	On a quantity describing the degree of chain entanglement in linear polymer systems. <i>Macromolecular Theory and Simulations</i> , 1994, 3, 639-647.	0.6	33
90	Rheology: From simple and to complex fluids. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1997, 240, 126-144.	1.2	33

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91	Thermoresponsive Cell Culture Substrates Based on PNIPAM Brushes Functionalized with Adhesion Peptides: Theoretical Considerations of Mechanism and Design. <i>Langmuir</i> , 2012, 28, 16623-16637.	1.6	32
92	Phase behavior and structure of Janus fluids. <i>Physical Review E</i> , 2003, 67, 041209.	0.8	31
93	Self-Folding of Charged Single Dendronized Polymers. <i>Advanced Materials</i> , 2008, 20, 3204-3210.	11.1	31
94	Automated symbolic calculations in nonequilibrium thermodynamics. <i>Computer Physics Communications</i> , 2010, 181, 2149-2157.	3.0	30
95	Molecularly derived constitutive equation for low-molecular polymer melts from thermodynamically guided simulation. <i>Journal of Rheology</i> , 2011, 55, 69-93.	1.3	30
96	Fast equilibration protocol for million atom systems of highly entangled linear polyethylene chains. <i>Journal of Chemical Physics</i> , 2016, 144, 154901.	1.2	30
97	Communication: Appearance of undershoots in start-up shear: Experimental findings captured by tumbling-snake dynamics. <i>Journal of Chemical Physics</i> , 2017, 146, 161101.	1.2	30
98	A simple example for comparing GENERIC with rational non-equilibrium thermodynamics. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2000, 285, 448-466.	1.2	29
99	Understanding Dynamics in Binary Mixtures of Entangled <i>cis</i> -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.	2.2	29
100	Writhe and mutual entanglement combine to give the entanglement length. <i>Physical Review E</i> , 2013, 88, 062604.	0.8	29
101	Microscopic Origin of the Non-Newtonian Viscosity of Semiflexible Polymer Solutions in the Semidilute Regime. <i>ACS Macro Letters</i> , 2014, 3, 136-140.	2.3	29
102	Self-assembly of core-polyethylene glycol-lipid shell (CPLS) nanoparticles and their potential as drug delivery vehicles. <i>Nanoscale</i> , 2016, 8, 14821-14835.	2.8	29
103	Size of graphene sheets determines the structural and mechanical properties of 3D graphene foams. <i>Nanotechnology</i> , 2018, 29, 104001.	1.3	29
104	Unified Analytic Expressions for the Entanglement Length, Tube Diameter, and Plateau Modulus of Polymer Melts. <i>Physical Review Letters</i> , 2020, 124, 147801.	2.9	28
105	Anisotropy of the magnetoviscous effect in ferrofluids. <i>Physical Review E</i> , 2005, 71, 051201.	0.8	27
106	Fibers with Integrated Mechanochemical Switches: Minimalistic Design Principles Derived from Fibronectin. <i>Biophysical Journal</i> , 2012, 103, 1909-1918.	0.2	27
107	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, 3770-3786.	1.2	27
108	Analytical Modeling of the Temporal Evolution of Epidemics Outbreaks Accounting for Vaccinations. <i>Physics</i> , 2021, 3, 386-426.	0.5	27

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109	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.	1.0	26
110	Recognition and analysis of local structure in polycrystalline configurations. <i>Computer Physics Communications</i> , 2002, 145, 371-384.	3.0	26
111	Formation of double helical and filamentous structures in models of physical and chemical gels. <i>Soft Matter</i> , 2008, 4, 18-28.	1.2	26
112	Computer simulation of dendronized polymers: organization and characterization at the atomistic level. <i>RSC Advances</i> , 2013, 3, 126-140.	1.7	26
113	The effect of polymer chain length on the mechanical properties of triblock copolymer gels. <i>Chemical Physics Letters</i> , 2014, 612, 157-161.	1.2	26
114	Structure and rheology of model-ferrofluids under shear flow. <i>Journal of Magnetism and Magnetic Materials</i> , 2005, 289, 325-327.	1.0	25
115	Nonaffine Deformation of Inherent Structure as a Static Signature of Cooperativity in Supercooled Liquids. <i>Physical Review Letters</i> , 2008, 101, 095501.	2.9	25
116	Unraveling two-dimensional polymerization in the single crystal. <i>Journal of Applied Crystallography</i> , 2018, 51, 481-497.	1.9	25
117	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, 254.	2.0	24
118	Self-assembled core-shell polyethylene glycol-lipid shell nanoparticles demonstrate high stability in shear flow. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 13294-13306.	1.3	23
119	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.	6.6	23
120	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.	0.8	22
121	Anisotropic self-diffusion in ferrofluids studied via Brownian dynamics simulations. <i>Physical Review E</i> , 2005, 72, 031504.	0.8	22
122	From hyperbolic regularization to exact hydrodynamics for linearized Grad's equations. <i>Physical Review E</i> , 2007, 75, 051204.	0.8	22
123	Computational study on entanglement length and pore size of carbon nanotube buckypaper. <i>Applied Physics Letters</i> , 2012, 100, .	1.5	22
124	Effect of Crosslinking on the Microtribological Behavior of Model Polymer Brushes. <i>Tribology Letters</i> , 2016, 63, 1.	1.2	22
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.	1.2	21
126	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.	1.2	21



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127	Lubricated optical rheometer for the study of two-dimensional complex flows of polymer melts. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008, 150, 43-55.	1.0	21
128	Exact Linear Hydrodynamics from the Boltzmann Equation. <i>Physical Review Letters</i> , 2008, 100, 214503.	2.9	21
129	Synthetic regimes due to packing constraints in dendritic molecules confirmed by labelling experiments. <i>Nature Communications</i> , 2013, 4, 1993.	5.8	21
130	Boltzmann equation and hydrodynamic fluctuations. <i>Physical Review E</i> , 2009, 80, 051202.	0.8	20
131	Using Mesoscopic Models to Design Strong and Tough Biomimetic Polymer Networks. <i>Langmuir</i> , 2011, 27, 13796-13805.	1.6	20
132	Computer Simulation of Fifth Generation Dendronized Polymers: Impact of Charge on Internal Organization. <i>Journal of Physical Chemistry B</i> , 2013, 117, 6007-6017.	1.2	20
133	Explicit formulae for the peak time of an epidemic from the SIR model. Which approximant to use?. <i>Physica D: Nonlinear Phenomena</i> , 2021, 425, 132981.	1.3	20
134	A Detailed Examination of the Topological Constraints of Lamellae-Forming Block Copolymers. <i>Macromolecules</i> , 2018, 51, 2110-2124.	2.2	19
135	Combined Experimental and Simulation Studies of Cross-Linked Polymer Brushes under Shear. <i>Macromolecules</i> , 2018, 51, 10174-10183.	2.2	19
136	NEMD-Computersimulation zur Rheologie von Polymerschmelzen / NEMD Computer Simulation of Polymer Melt Rheology. <i>Applied Rheology</i> , 1995, 5, 66-71.	3.5	19
137	Dynamics and Wetting Behavior of Core-Shell Soft Particles at a Fluid-Fluid Interface. <i>Langmuir</i> , 2018, 34, 15370-15382.	1.6	18
138	Insights from modeling into structure, entanglements, and dynamics in attractive polymer nanocomposites. <i>Soft Matter</i> , 2021, 17, 6362-6373.	1.2	18
139	Symbolic test of the Jacobi identity for given generalized $\hat{\epsilon}$ -Poisson bracket. <i>Computer Physics Communications</i> , 2001, 137, 325-340.	3.0	17
140	Filamentous networks in phase-separating two-dimensional gels. <i>Europhysics Letters</i> , 2007, 77, 58007.	0.7	17
141	Hyperbolicity of exact hydrodynamics for three-dimensional linearized Grad's equations. <i>Physical Review E</i> , 2007, 76, 022201.	0.8	17
142	Carbon Nanotube Length Governs the Viscoelasticity and Permeability of Buckypaper. <i>Polymers</i> , 2017, 9, 115.	2.0	17
143	Crossover between short- and long-time behavior of stress fluctuations and viscoelasticity of liquids. <i>Physical Review E</i> , 2003, 67, 042201.	0.8	16
144	Structural changes and viscoplastic behavior of a generic embedded-atom model metal in steady shear flow. <i>Physical Review E</i> , 2004, 69, 021509.	0.8	16

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145	Beyond-equilibrium molecular dynamics of a rarefied gas subjected to shear flow. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2004, 120, 175-187.	1.0	16
146	Interactions in dendronized polymers: intramolecular dominates intermolecular. <i>Soft Matter</i> , 2014, 10, 1032.	1.2	16
147	On the Shape and Rheology of Linear Micelles in Dilute Solutions. <i>Journal De Physique II</i> , 1997, 7, 931-946.	0.9	16
148	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.	0.8	15
149	Lubricated cross-slot flow of a low density polyethylene melt. <i>Journal of Non-Newtonian Fluid Mechanics</i> , 2008, 154, 52-64.	1.0	15
150	Pulling-force-induced elongation and alignment effects on entanglement and knotting characteristics of linear polymers in a melt. <i>Physical Review E</i> , 2014, 90, 042602.	0.8	15
151	Miscibility and Nanoparticle Diffusion in Ionic Nanocomposites. <i>Polymers</i> , 2018, 10, 1010.	2.0	15
152	Drag on a spherical particle at the air-liquid interface: Interplay between compressibility, Marangoni flow, and surface viscosities. <i>Physics of Fluids</i> , 2021, 33, .	1.6	15
153	Assessing numerical methods for molecular and particle simulation. <i>Soft Matter</i> , 2017, 13, 8565-8578.	1.2	14
154	Unifying kinetic approach to phoretic forces and torques onto moving and rotating convex particles. <i>Journal of Chemical Physics</i> , 2006, 125, 044105.	1.2	13
155	Non-constant link tension coefficient in the tumbling-snake model subjected to simple shear. <i>Journal of Chemical Physics</i> , 2017, 147, 174903.	1.2	13
156	Mesoscopic model for the viscosities of nematic liquid crystals. <i>Physical Review E</i> , 1999, 60, 4226-4234.	0.8	12
157	Flow of branched polymer melts in a lubricated cross-slot channel: a combined computational and experimental study. <i>Rheologica Acta</i> , 2009, 48, 97-108.	1.1	12
158	Modelling and confocal microscopy of biopolymer mixtures in confined geometries. <i>Soft Matter</i> , 2010, 6, 2713.	1.2	12
159	Solution of the complete Curtiss-Bird model for polymeric liquids subjected to simple shear flow. <i>Journal of Chemical Physics</i> , 2016, 144, 124905.	1.2	12
160	Gaussian Doubling Times and Reproduction Factors of the COVID-19 Pandemic Disease. <i>Frontiers in Physics</i> , 2020, 8, .	1.0	12
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