

# Arantxa Arbe

## List of Publications by Year in descending order

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190  
papers

6,925  
citations

50276

46  
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79698

73  
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192  
all docs

192  
docs citations

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times ranked

3663  
citing authors

#	ARTICLE	IF	CITATIONS
1	Merging of the $\hat{\Gamma}_\pm$ and $\hat{\Gamma}^2$ relaxations in polybutadiene: A neutron spin echo and dielectric study. Physical Review E, 1996, 54, 3853-3869.	2.1	257
2	Crossover from Debye to non-Debye dynamical behavior of the $\hat{\Gamma}_\pm$ relaxation observed by quasielastic neutron scattering in a glass-forming polymer. Physical Review Letters, 1993, 71, 2603-2606.	7.8	194
3	Neutron scattering study of the picosecond dynamics of polybutadiene and polyisoprene. Physical Review E, 1995, 52, 781-795.	2.1	192
4	Correlation between non-Debye behavior and $Q$ behavior of the $\hat{\Gamma}_\pm$ relaxation in glass-forming polymeric systems. Physical Review Letters, 1992, 69, 478-481.	7.8	169
5	Dynamics of Glass-Forming Polymers: "Homogeneous" versus "Heterogeneous" Scenario. Physical Review Letters, 1998, 81, 590-593.	7.8	160
6	Segmental dynamics in miscible polymer blends: recent results and open questions. Soft Matter, 2007, 3, 1474.	2.7	159
7	Effect of Nanoconfinement on Polymer Dynamics: Surface Layers and Interphases. Physical Review Letters, 2013, 110, 108303.	7.8	154
8	How Far Are Single-Chain Polymer Nanoparticles in Solution from the Globular State?. ACS Macro Letters, 2014, 3, 767-772.	4.8	152
9	Neutron Spin Echo in Polymer Systems. , 2005, , .		142
10	Metallo-Folded Single-Chain Nanoparticles with Catalytic Selectivity. ACS Macro Letters, 2014, 3, 439-443.	4.8	130
11	Endowing Single-Chain Polymer Nanoparticles with Enzyme-Mimetic Activity. ACS Macro Letters, 2013, 2, 775-779.	4.8	129
12	Direct Observation of Confined Single Chain Dynamics by Neutron Scattering. Physical Review Letters, 2010, 104, 197801.	7.8	123
13	Self-motion and the $\hat{\Gamma}_\pm$ relaxation in a simulated glass-forming polymer: Crossover from Gaussian to non-Gaussian dynamic behavior. Physical Review E, 2002, 65, 041804.	2.1	121
14	Molecular Motions in Polyisobutylene: A Neutron Spin-Echo and Dielectric Investigation. Macromolecules, 1998, 31, 1133-1143.	4.8	110
15	"Michael" Nanocarriers Mimicking Transient-Binding Disordered Proteins. ACS Macro Letters, 2013, 2, 491-495.	4.8	106
16	How Composition Determines the Properties of Isodimorphic Poly(butylene) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 147 Td (succinate-<i>Crystalline Random Copolymers. Macromolecules, 2015, 48, 43-57.	4.8	105
17	Segmental Dynamics in Poly(vinylethylene)/Polyisoprene Miscible Blends Revisited. A Neutron Scattering and Broad-Band Dielectric Spectroscopy Investigation. Macromolecules, 1999, 32, 7572-7581.	4.8	104
18	Quasielastic neutron scattering in soft matter. Current Opinion in Colloid and Interface Science, 2009, 14, 381-390.	7.4	97

#	ARTICLE	IF	CITATIONS
19	Non-Gaussian Nature of the $\hat{\tau}$ Relaxation of Glass-Forming Polyisoprene. <i>Physical Review Letters</i> , 2002, 89, 245701.	7.8	92
20	Dynamics of poly(ethylene oxide) in a blend with poly(methyl methacrylate): A quasielastic neutron scattering and molecular dynamics simulations study. <i>Physical Review E</i> , 2005, 72, 031808.	2.1	92
21	Advantages of Orthogonal Folding of Single Polymer Chains to Soft Nanoparticles. <i>Macromolecules</i> , 2013, 46, 9748-9759.	4.8	89
22	Study of the Dynamic Structure Factor in the $\hat{\tau}^2$ Relaxation Regime of Polybutadiene. <i>Physical Review Letters</i> , 1996, 76, 1872-1875.	7.8	88
23	Experimental evidence by neutron scattering of a crossover from Gaussian to non-Gaussian behavior in the $\hat{\tau}$ relaxation of polyisoprene. <i>Physical Review E</i> , 2003, 67, 051802.	2.1	82
24	Design and Preparation of Single-Chain Nanocarriers Mimicking Disordered Proteins for Combined Delivery of Dermal Bioactive Cargos. <i>Macromolecular Rapid Communications</i> , 2013, 34, 1681-1686.	3.9	82
25	Intermediate length scale dynamics of polyisobutylene. <i>Physical Review E</i> , 2002, 65, 051803.	2.1	80
26	From Rouse dynamics to local relaxation: A neutron spin echo study on polyisobutylene melts. <i>Journal of Chemical Physics</i> , 1999, 111, 6107-6120.	3.0	78
27	Merging of the Dielectric $\hat{\tau}$ and $\hat{\tau}^2$ Relaxations in Glass-Forming Polymers. <i>Macromolecules</i> , 2001, 34, 503-513.	4.8	77
28	Efficient Route to Compact Single-Chain Nanoparticles: Photoactivated Synthesis via Thiol- $\alpha$ -Yne Coupling Reaction. <i>Macromolecules</i> , 2014, 47, 8270-8280.	4.8	77
29	Study of the dynamics of poly(ethylene oxide) by combining molecular dynamic simulations and neutron scattering experiments. <i>Journal of Chemical Physics</i> , 2009, 130, 094908.	3.0	73
30	Single-chain nanoparticles: opportunities provided by internal and external confinement. <i>Materials Horizons</i> , 2020, 7, 2292-2313.	12.2	72
31	Crossover from Independent to Cooperative Segmental Dynamics in Polymers: Experimental Realization in Poly(Vinyl Chloride). <i>Physical Review Letters</i> , 1997, 78, 1928-1931.	7.8	69
32	Influence of Chain Topology (Cyclic versus Linear) on the Nucleation and Isothermal Crystallization of Poly( $\alpha$ -lactide) and Poly( $\delta$ -lactide). <i>Macromolecules</i> , 2018, 51, 1718-1732.	4.8	68
33	Anomalous relaxation of self-assembled alkyl nanodomains in high-order poly(n-alkyl methacrylates). <i>Soft Matter</i> , 2008, 4, 1792.	2.7	65
34	Concentrated Solutions of Single-Chain Nanoparticles: A Simple Model for Intrinsically Disordered Proteins under Crowding Conditions. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 838-844.	4.6	64
35	Neutron scattering study of the dynamics of a polymer melt under nanoscopic confinement. <i>Journal of Chemical Physics</i> , 2009, 131, 174901.	3.0	62
36	Origin of Internal Viscosity Effects in Flexible Polymers: A Comparative Neutron Spin-Echo and Light Scattering Study on Poly(dimethylsiloxane) and Polyisobutylene. <i>Macromolecules</i> , 2001, 34, 1281-1290.	4.8	61

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37	Dynamics of Water Absorbed in Polyamides. <i>Macromolecules</i> , 2012, 45, 1676-1687.	4.8	61
38	Origin of Dynamic Heterogeneities in Miscible Polymer Blends: A Quasielastic Neutron Scattering Study. <i>Physical Review Letters</i> , 2000, 85, 772-775.	7.8	59
39	Dynamic Confinement Effects in Polymer Blends. A Quasielastic Neutron Scattering Study of the Dynamics of Poly(ethylene oxide) in a Blend with Poly(vinyl acetate). <i>Macromolecules</i> , 2006, 39, 3007-3018.	4.8	56
40	Recent progress on polymer dynamics by neutron scattering: From simple polymers to complex materials. <i>Journal of Polymer Science, Part B: Polymer Physics</i> , 2013, 51, 87-113.	2.1	56
41	The dynamics of the $\hat{1}\pm$ - and $\hat{1}^2$ -relaxations in glass-forming polymers studied by quasielastic neutron scattering and dielectric spectroscopy. <i>Journal of Non-Crystalline Solids</i> , 1994, 172-174, 126-137.	3.1	54
42	Polymer Chain Dynamics in a Random Environment: Heterogeneous Mobilities. <i>Physical Review Letters</i> , 2007, 98, 168301.	7.8	53
43	Dynamics in Poly( <i>n</i> -alkyl methacrylates): A Neutron Scattering, Calorimetric, and Dielectric Study. <i>Macromolecules</i> , 2010, 43, 3107-3119.	4.8	53
44	Efficient Synthesis of Single-Chain Globules Mimicking the Morphology and Polymerase Activity of Metalloenzymes. <i>Macromolecular Rapid Communications</i> , 2015, 36, 1592-1597.	3.9	52
45	Interpretation of anomalous momentum transfer dependences of local chain motion of polymers observed by quasielastic incoherent neutron scattering experiments. <i>Macromolecules</i> , 1992, 25, 6727-6729.	4.8	51
46	On the origin of the non-exponential behaviour of the $\alpha$ -relaxation in glass-forming polymers: incoherent neutron scattering and dielectric relaxation results. <i>Journal of Physics Condensed Matter</i> , 1999, 11, A363-A370.	1.8	50
47	Folding Single Chains to Single-Chain Nanoparticles via Reversible Interactions: What Size Reduction Can One Expect?. <i>Macromolecules</i> , 2017, 50, 1732-1739.	4.8	49
48	Local Structure of Syndiotactic Poly(methyl methacrylate). A Combined Study by Neutron Diffraction with Polarization Analysis and Atomistic Molecular Dynamics Simulations. <i>Macromolecules</i> , 2006, 39, 3947-3958.	4.8	45
49	Self- and Collective Dynamics of Syndiotactic Poly(methyl methacrylate). A Combined Study by Quasielastic Neutron Scattering and Atomistic Molecular Dynamics Simulations. <i>Macromolecules</i> , 2006, 39, 6260-6272.	4.8	45
50	Investigation of the Dielectric $\hat{1}^2$ -Process in Polyisobutylene by Incoherent Quasielastic Neutron Scattering. <i>Macromolecules</i> , 1998, 31, 4926-4934.	4.8	44
51	Structure and dynamics of single-chain nano-particles in solution. <i>Polymer</i> , 2016, 105, 532-544.	3.8	44
52	Carbon-carbon torsional barriers driving the fast dynamics in glass-forming polymers. <i>Physical Review B</i> , 1998, 57, 13508-13513.	3.2	41
53	Dynamic Confinement Effects in Polymer Blends. A Quasielastic Neutron Scattering Study of the Slow Component in the Blend Poly(vinyl acetate)/Poly(ethylene oxide). <i>Macromolecules</i> , 2007, 40, 4568-4577.	4.8	41
54	Poly(butylene succinate- <i>ran</i> - $\hat{1}\mu$ -caprolactone) copolyesters: Enzymatic synthesis and crystalline isodimorphic character. <i>European Polymer Journal</i> , 2017, 95, 795-808.	5.4	41

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55	Quasielastic Neutron Scattering Study on the Dynamics of Poly(alkylene oxide)s. <i>Macromolecules</i> , 2012, 45, 4394-4405.	4.8	40
56	A Solvent-Based Strategy for Tuning the Internal Structure of Metallo-Folded Single-Chain Nanoparticles. <i>Macromolecular Rapid Communications</i> , 2016, 37, 1060-1065.	3.9	39
57	Quasielastic neutron scattering study of hydrogen motions in an aqueous poly(vinyl methyl ether) solution. <i>Journal of Chemical Physics</i> , 2011, 134, 204906.	3.0	37
58	Component dynamics in polyvinylpyrrolidone concentrated aqueous solutions. <i>Journal of Chemical Physics</i> , 2012, 137, 084902.	3.0	36
59	Single Chain Dynamic Structure Factor of Linear Polymers in an All-Polymer Nano-Composite. <i>Macromolecules</i> , 2016, 49, 2354-2364.	4.8	36
60	Neutron scattering and molecular dynamics simulations: synergetic tools to unravel structure and dynamics in polymers. <i>Soft Matter</i> , 2012, 8, 8257.	2.7	35
61	Quasielastic Neutron Scattering Study on the Effect of Blending on the Dynamics of Head-to-Head Poly(propylene) and Poly(ethylene-propylene). <i>Macromolecules</i> , 2006, 39, 1060-1072.	4.8	34
62	Heterogeneous structure of poly(vinyl chloride) as the origin of anomalous dynamical behavior. <i>Journal of Chemical Physics</i> , 2002, 117, 1336-1350.	3.0	33
63	Neutron Spin Echo in Polymer Systems, Chapter 1. , 2005, , 1-221.		33
64	Reply to "Comment on "Merging of the $\hat{I}_\pm$ and $\hat{I}^2$ relaxations in polybutadiene:" A neutron spin echo and dielectric study". <i>Physical Review E</i> , 1999, 60, 1103-1105.	2.1	31
65	The Role of the Topological Constraints in the Chain Dynamics in All-Polymer Nanocomposites. <i>Macromolecules</i> , 2017, 50, 1719-1731.	4.8	31
66	Crowding the Environment of Single-Chain Nanoparticles: A Combined Study by SANS and Simulations. <i>Macromolecules</i> , 2018, 51, 1573-1585.	4.8	31
67	Short and Intermediate Range Order in Poly(alkylene oxide)s. A Neutron Diffraction and Molecular Dynamics Simulation Study. <i>Macromolecules</i> , 2012, 45, 7293-7303.	4.8	29
68	Structure factors in polystyrene: a neutron scattering and MD-simulation study. <i>Physica B: Condensed Matter</i> , 2004, 350, E881-E884.	2.7	28
69	Microscopic Dynamics in Nanocomposites of Poly(ethylene oxide) and Poly(methyl methacrylate) Soft Nanoparticles: A Quasi-Elastic Neutron Scattering Study. <i>Macromolecules</i> , 2014, 47, 304-315.	4.8	28
70	Q-dependence of the relaxation times of the $\hat{I}_\pm$ -relaxation as observed by quasielastic neutron scattering. <i>Journal of Non-Crystalline Solids</i> , 1994, 172-174, 229-233.	3.1	27
71	Structure and Dynamics of Self-Assembled Comb Copolymers: Comparison between Simulations of a Generic Model and Neutron Scattering Experiments. <i>Macromolecules</i> , 2011, 44, 1695-1706.	4.8	27
72	Hydrogen motions in the $\hat{I}_\pm$ -relaxation regime of poly(vinyl ethylene): A molecular dynamics simulation and neutron scattering study. <i>Journal of Chemical Physics</i> , 2004, 121, 3282-3294.	3.0	26

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73	Sub-Tg dynamics in polycarbonate by neutron scattering and its relation with secondary $\hat{\Gamma}^3$ relaxation. <i>Journal of Chemical Physics</i> , 2005, 123, 014907.	3.0	26
74	Short-range order and collective dynamics of poly(vinyl acetate): A combined study by neutron scattering and molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2008, 129, 224903.	3.0	26
75	Modeling the collective relaxation time of glass-forming polymers at intermediate length scales: Application to polyisobutylene. <i>Journal of Chemical Physics</i> , 2013, 139, 044906.	3.0	26
76	Coherent structural relaxation of water from meso- to intermolecular scales measured using neutron spectroscopy with polarization analysis. <i>Physical Review Research</i> , 2020, 2, .	3.6	26
77	Application of SSA thermal fractionation and X-ray diffraction to elucidate comonomer inclusion or exclusion from the crystalline phases in poly(butylene succinate-ran-butylene azelate) random copolymers. <i>Journal of Polymer Science, Part B: Polymer Physics</i> , 2016, 54, 2346-2358.	2.1	25
78	Atomic motions in the $\hat{\Gamma}^2$ -merging region of 1,4-polybutadiene: A molecular dynamics simulation study. <i>Journal of Chemical Physics</i> , 2008, 128, 224905.	3.0	24
79	Characterization of the "simple-liquid" state in a polymeric system: Coherent and incoherent scattering functions. <i>Physical Review E</i> , 2009, 80, 041805.	2.1	24
80	Quasielastic Neutron Scattering and Molecular Dynamics Simulation Study on the Structure Factor of Poly(ethylene-co-propylene). <i>Macromolecules</i> , 2009, 42, 8271-8285.	4.8	24
81	Plasticization and cocrystallization in LLDPE/wax blends. <i>Journal of Polymer Science, Part B: Polymer Physics</i> , 2016, 54, 1469-1482.	2.1	24
82	Direct observation of the crossover from $\hat{\Gamma}^2$ -relaxation to Rouse dynamics in a polymer melt. <i>Europhysics Letters</i> , 2004, 66, 239-245.	2.0	23
83	Atomic motions in poly(vinyl methyl ether): A combined study by quasielastic neutron scattering and molecular dynamics simulations in the light of the mode coupling theory. <i>Journal of Chemical Physics</i> , 2009, 131, 204901.	3.0	23
84	Relaxations and Relaxor-Ferroelectric-Like Response of Nanotubularly Confined Poly(vinylidene fluoride). <i>Journal of Applied Physics</i> , 2010, 107, 044107.	6.7	23
85	Size of Elastic Single-Chain Nanoparticles in Solution and on Surfaces. <i>Macromolecules</i> , 2017, 50, 6323-6331.	4.8	23
86	Space time observation of the $\beta$ -process in polymers by quasielastic neutron scattering. <i>Journal of Physics Condensed Matter</i> , 1999, 11, A297-A306.	1.8	22
87	Methyl group dynamics above the glass transition temperature: a molecular dynamics simulation in polyisoprene. <i>Chemical Physics</i> , 2000, 261, 47-59.	1.9	22
88	Partial Structure Factors in 1,4-Polybutadiene. A Combined Neutron Scattering and Molecular Dynamics Simulations Study. <i>Macromolecules</i> , 2005, 38, 9847-9853.	4.8	22
89	Synthesis and Characterization of Double Crystalline Cyclic Diblock Copolymers of Poly( $\epsilon$ -caprolactone) and Poly( $\delta$ -valerolactone) ( $\text{PCL-}b\text{-PL(D)LA}$ ). <i>Macromolecular Rapid Communications</i> , 2016, 37, 1676-1681.	3.9	22
90	Intermediate length scale dynamics in glass forming polymers: coherent and incoherent quasielastic neutron scattering results on polyisobutylene. <i>Chemical Physics</i> , 2003, 292, 295-309.	1.9	21

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91	Dynamics of Polyethersulfone Phenylene Rings: A Quasielastic Neutron Scattering Study. <i>Macromolecules</i> , 2005, 38, 3999-4013.	4.8	21
92	Study of the structure and dynamics of poly(vinyl pyrrolidone) by molecular dynamics simulations validated by quasielastic neutron scattering and x-ray diffraction experiments. <i>Journal of Chemical Physics</i> , 2011, 134, 054904.	3.0	21
93	Neutron Scattering and X-ray Investigation of the Structure and Dynamics of Poly(ethyl Tj ETQq1 1 0.784314 rgBT /Overlock, 10 Tf 50	4.8	21
94	Phenylene ring dynamics in bisphenol-A-polysulfone by neutron scattering. <i>Journal of Chemical Physics</i> , 2004, 120, 423-436.	3.0	20
95	Glassy Dynamics of Polystyrene by Quasielastic Neutron Scattering. <i>Macromolecules</i> , 2011, 44, 3161-3168.	4.8	20
96	Nanostructuration by Self-Assembly in <i>N</i> -Alkyl Thiazolium and Triazolium Side-Chain Polymethacrylates. <i>Macromolecules</i> , 2015, 48, 7180-7193.	4.8	20
97	Phase behavior of side-chain liquid-crystalline polymers containing biphenyl mesogens with different spacer lengths synthesized via miniemulsion polymerization. <i>Polymer Chemistry</i> , 2016, 7, 4736-4750.	3.9	20
98	Arbeet al.Reply:. <i>Physical Review Letters</i> , 1999, 82, 1336-1336.	7.8	19
99	Positron annihilation and relaxation dynamics from dielectric spectroscopy and nuclear magnetic resonance: <i>cis</i> - <i>trans</i> -1,4-poly(butadiene). <i>Journal of Chemical Physics</i> , 2011, 134, 164507.	3.0	19
100	Dynamic study of polystyrene-block-poly(4-vinylpyridine) copolymer in bulk and confined in cylindrical nanopores. <i>Polymer</i> , 2014, 55, 4057-4066.	3.8	19
101	The role of PLLA-g-montmorillonite nanohybrids in the acceleration of the crystallization rate of a commercial PLA. <i>CrystEngComm</i> , 2016, 18, 9334-9344.	2.6	19
102	Sequential crystallization and morphology of triple crystalline biodegradable PEO-b-PCL-b-PLLA triblock terpolymers. <i>RSC Advances</i> , 2016, 6, 4739-4750.	3.6	19
103	Self-motion and the $\hat{\alpha}$ -relaxation in glass-forming polymers. Molecular dynamic simulation and quasielastic neutron scattering results in polyisoprene. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S1127-S1138.	1.8	18
104	Investigation of a Nanocomposite of 75 wt % Poly(methyl methacrylate) Nanoparticles with 25 wt % Poly(ethylene oxide) Linear Chains: A Quasielastic Neutron Scattering, Calorimetric, and WAXS Study. <i>Macromolecules</i> , 2014, 47, 3005-3016.	4.8	18
105	How Does Microstructural Design Affect the Dynamics and Rheology of Segmented Polyurethanes?. <i>Macromolecules</i> , 2020, 53, 5381-5398.	4.8	18
106	High magnetization FeCo nanoparticles for magnetorheological fluids with enhanced response. <i>Soft Matter</i> , 2021, 17, 840-852.	2.7	18
107	Unexpected PDMS Behavior in Segregated Cylindrical and Spherical Nanophases of PS- <i>b</i> -PDMS Asymmetric Diblock Copolymers. <i>Macromolecules</i> , 2012, 45, 491-502.	4.8	17
108	Effect of polar solvents on the crystalline phase of polyamides. <i>Polymer</i> , 2014, 55, 2867-2881.	3.8	17

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109	Insights into the Network Structure of Cross-Linked Polymers Synthesized via Miniemulsion Nitroxide-Mediated Radical Polymerization. <i>Macromolecules</i> , 2018, 51, 9740-9748.	4.8	17
110	Local Domain Size in Single-Chain Polymer Nanoparticles. <i>ACS Omega</i> , 2018, 3, 8648-8654.	3.5	17
111	Mesoscale Dynamics in Melts of Single-Chain Polymeric Nanoparticles. <i>Macromolecules</i> , 2019, 52, 6935-6942.	4.8	17
112	Insight into the Structure and Dynamics of Polymers by Neutron Scattering Combined with Atomistic Molecular Dynamics Simulations. <i>Polymers</i> , 2020, 12, 3067.	4.5	17
113	Crowding Effects on the Structure and Dynamics of the Intrinsically Disordered Nuclear Chromatin Protein NUPR1. <i>Frontiers in Molecular Biosciences</i> , 2021, 8, 684622.	3.5	17
114	Chain Dynamics of Unentangled Poly(ethylene-<i>alt</i>-propylene) Melts by Means of Neutron Scattering and Fully Atomistic Molecular Dynamics Simulations. <i>Macromolecules</i> , 2011, 44, 3129-3139.	4.8	16
115	Publisher's Note: Effect of Nanoconfinement on Polymer Dynamics: Surface Layers and Interphases [ <i>Phys. Rev. Lett.</i> <b>110</b>, 108303 (2013)]. <i>Physical Review Letters</i> , 2013, 110, .	7.8	16
116	Influence of Solvent on Poly(2-(Dimethylamino)Ethyl Methacrylate) Dynamics in Polymer-Concentrated Mixtures: A Combined Neutron Scattering, Dielectric Spectroscopy, and Calorimetric Study. <i>Macromolecules</i> , 2015, 48, 6724-6735.	4.8	16
117	Effect of Molecular Crowding on Conformation and Interactions of Single-Chain Nanoparticles. <i>Macromolecules</i> , 2019, 52, 4295-4305.	4.8	16
118	Neutron scattering investigation of a diluted blend of poly(ethylene oxide) in polyethersulfone. <i>Journal of Chemical Physics</i> , 2008, 128, 184901.	3.0	15
119	Collective Features in Polyisobutylene. A Study of the Static and Dynamic Structure Factor by Molecular Dynamics Simulations. <i>Macromolecules</i> , 2014, 47, 447-459.	4.8	15
120	Effect of chain stiffness on the structure of single-chain polymer nanoparticles. <i>Journal of Physics Condensed Matter</i> , 2018, 30, 034001.	1.8	15
121	Facile Access to Completely Deuterated Single-Chain Nanoparticles Enabled by Intramolecular Azide Photodecomposition. <i>Macromolecular Rapid Communications</i> , 2019, 40, 1900046.	3.9	15
122	Neutron scattering and the glass transition in polymers – present status and future opportunities. <i>Journal of Non-Crystalline Solids</i> , 2001, 287, 286-296.	3.1	14
123	Atomic motions in the $\beta$ -region of glass-forming polymers: molecular versus mode coupling theory approach. <i>Journal of Physics Condensed Matter</i> , 2007, 19, 205127.	1.8	14
124	Comparative study of $\alpha$ -relaxations in a glass-forming polymer (PVC) by dielectric spectroscopy and quasielastic neutron scattering. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1993, 201, 447-452.	2.6	13
125	Dynamic structure factors due to relaxation processes in glass-forming polymers. <i>Physica B: Condensed Matter</i> , 1997, 241-243, 1005-1012.	2.7	13
126	Hydrogen motions and the $\beta$ -relaxation in glass-forming polymers: Molecular dynamics simulation and quasi-elastic neutron scattering results. <i>Pramana - Journal of Physics</i> , 2004, 63, 25-32.	1.8	13



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127	The decisive influence of local chain dynamics on the overall dynamic structure factor close to the glass transition. <i>Europhysics Letters</i> , 2005, 71, 262-268.	2.0	13
128	Positron annihilation and relaxation dynamics from dielectric spectroscopy: poly(vinylmethylether). <i>Journal of Physics Condensed Matter</i> , 2012, 24, 155104.	1.8	13
129	Applicability of mode-coupling theory to polyisobutylene: A molecular dynamics simulation study. <i>Physical Review E</i> , 2013, 88, 042302.	2.1	13
130	Component dynamics in nanostructured PI-PDMS diblock copolymers with PI segregated in lamellas, cylinders, and spheres. <i>Colloid and Polymer Science</i> , 2014, 292, 1863-1876.	2.1	13
131	Influence of Chain Primary Structure and Topology (Branching) on Crystallization and Thermal Properties: The Case of Polysulfides. <i>Macromolecules</i> , 2019, 52, 2093-2104.	4.8	13
132	Self-Reporting of Folding and Aggregation by Orthogonal Hantzsch Luminophores Within a Single Polymer Chain. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 3534-3539.	13.8	13
133	An unexpected route to aldehyde-decorated single-chain nanoparticles from azides. <i>Polymer Chemistry</i> , 2016, 7, 6570-6574.	3.9	12
134	Temperature and momentum transfer dependence of the dynamics of the $\hat{\tau}$ -relaxation in polymer melts. <i>Physica B: Condensed Matter</i> , 1992, 182, 369-375.	2.7	11
135	DETERMINATION OF FILLER STRUCTURE IN SILICA-FILLED SBR COMPOUNDS BY MEANS OF SAXS AND AFM. <i>Rubber Chemistry and Technology</i> , 2015, 88, 690-710.	1.2	11
136	Applying Polymer Blend Dynamics Concepts to a Simplified Industrial System. A Combined Effort by Dielectric Spectroscopy and Neutron Scattering. <i>Macromolecules</i> , 2018, 51, 6692-6706.	4.8	11
137	Melts of single-chain nanoparticles: A neutron scattering investigation. <i>Journal of Applied Physics</i> , 2020, 127, .	2.5	11
138	Human importin $\hat{\tau}3$ and its N-terminal truncated form, without the importin- $\hat{\tau}2$ -binding domain, are oligomeric species with a low conformational stability in solution. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2020, 1864, 129609.	2.4	11
139	Unraveling the coherent dynamic structure factor of liquid water at the mesoscale by molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2021, 155, 244509.	3.0	11
140	Dynamics of Poly(butylene oxide) Well above the Glass Transition. A Fully Atomistic Molecular Dynamics Simulation Study. <i>Macromolecules</i> , 2013, 46, 1678-1685.	4.8	10
141	A Useful Methodology for Determining the Compaction Degree of Single-Chain Nanoparticles by Conventional SEC. <i>Particle and Particle Systems Characterization</i> , 2016, 33, 373-381.	2.3	10
142	Investigation of the dynamics of aqueous proline solutions using neutron scattering and molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 27739-27754.	2.8	10
143	Cyclic Polyethylene Glycol as Nanoparticle Surface Ligand. <i>ACS Macro Letters</i> , 2020, 9, 1604-1610.	4.8	10
144	Advances in the Multi-Orthogonal Folding of Single Polymer Chains into Single-Chain Nanoparticles. <i>Polymers</i> , 2021, 13, 293.	4.5	10

#	ARTICLE	IF	CITATIONS
145	Ultrafiltration of single-chain polymer nanoparticles through nanopores and nanoslits. <i>Polymer</i> , 2018, 148, 61-67.	3.8	9
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