

Giorgio Celebre

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6496351/publications.pdf>

Version: 2024-02-01

52
papers

599
citations

567281

15
h-index

677142

22
g-index

55
all docs

55
docs citations

55
times ranked

420
citing authors

#	ARTICLE	IF	CITATIONS
1	Structure of biphenyl in a nematic liquid-crystalline solvent. Journal of the Chemical Society, Faraday Transactions, 1991, 87, 2623.	1.7	49
2	Solute-Solvent Interactions and Chiral Induction in Liquid Crystals. Journal of the American Chemical Society, 2005, 127, 11736-11744.	13.7	47
3	Order and Conformation of Biphenyl in Cyanobiphenyl Liquid Crystals: A Combined Atomistic Molecular Dynamics and ¹ H NMR Study. ChemPhysChem, 2014, 15, 1356-1367.	2.1	33
4	Internal rotation potential function for anisole in solution: a liquid crystal NMR study. The Journal of Physical Chemistry, 1992, 96, 2466-2470.	2.9	26
5	On the anisotropic intermolecular potential of biaxial apolar solutes in nematic solvents: Monte Carlo predictions and experimental data. Journal of Chemical Physics, 2001, 115, 9552-9556.	3.0	25
6	Is styrene planar in liquid phases?. Journal of Chemical Physics, 2004, 120, 7075-7084.	3.0	23
7	Conformational Investigation in Solution of a Fluorinated Anti-inflammatory Drug by NMR Spectroscopy in Weakly Ordering Media. Journal of Physical Chemistry B, 2014, 118, 9007-9016.	2.6	22
8	An investigation of the potential governing rotation about the O-CH ₂ bond in 4-chloroethoxybenzene by comparing observed and calculated dipolar couplings obtained for a sample dissolved in a nematic solvent. Journal of the Chemical Society, Faraday Transactions 2, 1988, 84, 1041-1052.	1.1	21
9	The orientational ordering of a biaxial particle in a uniaxial environment by Monte Carlo sampling: a new approach to the problem. Chemical Physics Letters, 2001, 342, 375-381.	2.6	21
10	The orientational behaviour of 1,4-difluorobenzene and p-benzoquinone in ZLI1132 and EBBA nematic solvents. Chemical Physics Letters, 2003, 368, 359-364.	2.6	20
11	The conformational behaviour of naproxen and flurbiprofen in solution by NMR spectroscopy. New Journal of Chemistry, 2015, 39, 9086-9097.	2.8	20
12	The Structure of Acrolein in a Liquid Crystal Phase. Chemistry - A European Journal, 2005, 11, 3599-3608.	3.3	19
13	Short-Range Molecular Interactions Governing the Orientational Ordering of Apolar Molecules Dissolved in Nematic Solvents. Journal of Physical Chemistry B, 2003, 107, 3243-3250.	2.6	18
14	Conformational Distribution of <i>trans</i> -Stilbene in Solution Investigated by Liquid Crystal NMR Spectroscopy and Compared with <i>in Vacuo</i> Theoretical Predictions. Journal of Physical Chemistry B, 2012, 116, 2876-2885.	2.6	17
15	Intrinsic Information Content of NMR Dipolar Couplings: A Conformational Investigation of 1,3-Butadiene in a Nematic Phase. ChemPhysChem, 2006, 7, 1930-1943.	2.1	16
16	A combined LX-NMR and molecular dynamics investigation of the bulk and local structure of ionic liquid crystals. Soft Matter, 2019, 15, 4486-4497.	2.7	16
17	Assessing the stable conformations of ibuprofen in solution by means of Residual Dipolar Couplings. European Journal of Pharmaceutical Sciences, 2017, 106, 113-121.	4.0	15
18	The temperature dependence of biaxiality of solutes dissolved in nematic solvents. Physical Chemistry Chemical Physics, 2000, 2, 1883-1888.	2.8	14

#	ARTICLE	IF	CITATIONS
19	The structure of ethylbenzene as a solute in liquid crystalline solvents via analysis of proton NMR spectra. <i>Physical Chemistry Chemical Physics</i> , 2000, 2, 3405-3413.	2.8	14
20	Smectic order parameters via liquid crystal NMR spectroscopy: Application to a partial bilayer smectic A phase. <i>European Physical Journal E</i> , 2012, 35, 112.	1.6	14
21	On the analysis of Liquid Crystal NMR dipolar coupling data via mixed a priori-maximum entropy methods. <i>Journal of Chemical Physics</i> , 2006, 124, 176101.	3.0	12
22	An Explicit Relationship between the Dielectric Anisotropy and the Average Electric Field Gradient in Nematic Solvents. <i>Journal of Physical Chemistry B</i> , 2007, 111, 2565-2572.	2.6	12
23	Orientalional Mechanisms in Liquid Crystalline Systems. 2. The Contribution to Solute Ordering from the Reaction Field Interaction between the Solute Electric Quadrupole Moment and the Solvent Electric Field Gradient. <i>Journal of Physical Chemistry B</i> , 2010, 114, 235-241.	2.6	12
24	Solvent smectic order parameters from solute nematic order parameters. <i>Journal of Chemical Physics</i> , 2008, 129, 094509.	3.0	9
25	Orientalional Mechanisms in Liquid Crystalline Systems. 1. A Reaction Field Analytical Description of the Interaction between the Electric Quadrupole Moment of a Probe Solute and the Electric Field Gradient of a Nematic Solvent. <i>Journal of Physical Chemistry B</i> , 2010, 114, 228-234.	2.6	9
26	NMR Studies of Solutes in Liquid Crystals: Small Flexible Molecules. , 2003, , 305-324.		8
27	Multitechnique Investigation of Conformational Features of Small Molecules: The Case of Methyl Phenyl Sulfoxide. <i>Journal of Physical Chemistry B</i> , 2008, 112, 2095-2101.	2.6	8
28	Calculated Versus Experimental Force Fields: The Influence in the Structure Determination of Benzene by NMR Spectroscopy in Liquid Crystal Solvents. <i>Molecular Crystals and Liquid Crystals</i> , 2007, 465, 289-299.	0.9	7
29	Doped ionic liquid crystals as effective weakly alignment media for polar solutes. <i>Journal of Magnetic Resonance</i> , 2016, 267, 63-67.	2.1	7
30	Rigid probe solutes in a smectic-liquid crystal: An unconventional route to the latter's positional order parameters. <i>Physical Review E</i> , 2011, 84, 061703.	2.1	6
31	Conformational and Orientalional Determination from Dipolar Couplings in a Weakly Ordering Organic-Based Lyotropic and a Strongly Ordering Thermotropic Liquid Crystal. <i>Molecular Crystals and Liquid Crystals</i> , 2015, 614, 39-53.	0.9	6
32	The combined use of deuterium NMR and computer simulations for conformational investigation of flexible molecules in nematic solutions. <i>Computational and Theoretical Chemistry</i> , 2005, 728, 209-214.	1.5	5
33	The use of residual dipolar couplings for conformational analysis of non-steroidal anti-inflammatory drugs dissolved in weakly ordering media. <i>Liquid Crystals</i> , 2018, 45, 2033-2047.	2.2	5
34	The use of heteronuclear multiple quantum spectra in the automatic analysis of NMR spectra of samples dissolved in liquid crystalline phases. <i>Liquid Crystals</i> , 2001, 28, 1403-1413.	2.2	4
35	Helical solutes orientationally ordered in anisotropic media composed of helical particles: Formulation of a mean torque potential sensitive to P and M chirality as a tool for the assignment of the absolute configuration of enantiomers. <i>Journal of Molecular Liquids</i> , 2019, 288, 111044.	4.9	4
36	Molecular Characterization of the Organic Fraction of Municipal Solid Waste and Compositional Evolution during Oxidative Processes Assessed by HR-MAS ¹³ C NMR Spectroscopy. <i>Applied Sciences (Switzerland)</i> , 2021, 11, 2267.	2.5	4

#	ARTICLE	IF	CITATIONS
37	Metabolic composition and authenticity evaluation of bergamot essential oil assessed by nuclear magnetic resonance spectroscopy. <i>Analytical and Bioanalytical Chemistry</i> , 2022, 414, 2297-2313.	3.7	4
38	Effects of Electrostatic Interactions on Orientational Order of Solutes in Nematic Solvents: NMR Experiments and Theoretical Predictions. <i>Molecular Crystals and Liquid Crystals</i> , 2002, 372, 107-120.	0.9	3
39	The prediction of ordering of parallelepipedal solute particles in nematic solvents: a comparison among different methodologies. <i>Chemical Physics Letters</i> , 2004, 384, 344-349.	2.6	3
40	Orientational ordering of solutes in confined nematic solvents: A possible way to probe director distributions. <i>Physical Review E</i> , 2006, 73, 020702.	2.1	3
41	The stable conformations of 4,4-dichloro-trans-stilbene in solution by liquid crystal NMR spectroscopy. <i>Journal of Molecular Structure</i> , 2013, 1034, 283-288.	3.6	3
42	Detection of Significant Aprotic Solvent Effects on the Conformational Distribution of Methyl 4-Nitrophenyl Sulfoxide: From Gas-Phase Rotational to Liquid-Crystal NMR Spectroscopy. <i>ChemPhysChem</i> , 2015, 16, 2327-2337.	2.1	3
43	Experimental Assessment of the Vibration-Reorientation Contribution to Liquid Crystal NMR Dipolar Couplings: The Case of Tetramethylallene Dissolved in a Nematic Mesophase. <i>Journal of Physical Chemistry B</i> , 2011, 115, 11119-11126.	2.6	2
44	Statistical thermodynamics of thermotropic biaxial nematic liquid crystals: An effective, molecular-field based theoretical description by means of a closed approximate form of the orientational partition function. <i>Journal of Molecular Liquids</i> , 2015, 209, 104-114.	4.9	2
45	Probing the sensitivity of orientational ordering as a way towards absolute enantioselectivity: Helical-particle solutes in helical-particle nematic solvents. <i>Physical Review E</i> , 2018, 98, .	2.1	2
46	Assessing the chirality-dependent conformational distribution of small flexible opposite enantiomers dissolved in weakly ordering enantiopure media by means of liquid crystal NMR techniques. <i>Journal of Molecular Liquids</i> , 2022, 347, 117994.	4.9	2
47	Ordering of Polar Solutes in Nematic Magic Mixtures : Evidences of Dipolar Effects?. <i>Molecular Crystals and Liquid Crystals</i> , 2003, 394, 119-125.	0.9	1
48	Effects of Electrostatic Interactions on Orientational Order of Solutes in Nematic Solvents: NMR Experiments and Theoretical Predictions. <i>Molecular Crystals and Liquid Crystals</i> , 2001, 372, 107-120.	0.3	1
49	Publisher's Note: Rigid probe solutes in a smectic-A liquid crystal: An unconventional route to the latter's positional order parameters [Phys. Rev. E 84, 061703 (2011)]. <i>Physical Review E</i> , 2011, 84, .	2.1	0
50	An analytical altitude- and wavelength-dependent expression of the j_3 rate coefficient governing the photolysis of the atmospheric Ozone. <i>Journal of Mathematical Chemistry</i> , 2018, 56, 1515-1519.	1.5	0
51	A heuristic approach to the assessment of the root mean square quantum noise. <i>Quantum Information Processing</i> , 2022, 21, 1.	2.2	0
52	Chiral symmetry breaking: Frank model for the evolution of homochirality described by population dynamics. <i>Journal of Mathematical Chemistry</i> , 2022, 60, 681-694.	1.5	0