

Francesco Zerbetto

List of Publications by Year in descending order

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399
papers

19,508
citations

16411

64
h-index

17055

122
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422
all docs

422
docs citations

422
times ranked

16033
citing authors

#	ARTICLE	IF	CITATIONS
1	Synthetic Molecular Motors and Mechanical Machines. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 72-191.	7.2	2,428
2	Unidirectional rotation in a mechanically interlocked molecular rotor. <i>Nature</i> , 2003, 424, 174-179.	13.7	862
3	Macroscopic transport by synthetic molecular machines. <i>Nature Materials</i> , 2005, 4, 704-710.	13.3	685
4	Carbon Nanotubes in Electron Donor-Acceptor Nanocomposites. <i>Accounts of Chemical Research</i> , 2005, 38, 871-878.	7.6	453
5	Quantum-chemical investigation of Franck-Condon and Jahn-Teller activity in the electronic spectra of Buckminsterfullerene. <i>Chemical Physics Letters</i> , 1988, 144, 31-37.	1.2	333
6	Interactions in Single Wall Carbon Nanotubes/Pyrene/Porphyrin Nanohybrids. <i>Journal of the American Chemical Society</i> , 2006, 128, 11222-11231.	6.6	320
7	Theoretical analysis of spectra of short polyenes. <i>Chemical Reviews</i> , 1991, 91, 867-891.	23.0	309
8	Influencing intramolecular motion with an alternating electric field. <i>Nature</i> , 2000, 406, 608-611.	13.7	223
9	Patterning through Controlled Submolecular Motion: Rotaxane-Based Switches and Logic Gates that Function in Solution and Polymer Films. <i>Angewandte Chemie - International Edition</i> , 2005, 44, 3062-3067.	7.2	210
10	A Generic Basis for Some Simple Light-Operated Mechanical Molecular Machines. <i>Journal of the American Chemical Society</i> , 2004, 126, 12210-12211.	6.6	199
11	Information Storage Using Supramolecular Surface Patterns. <i>Science</i> , 2003, 299, 531-531.	6.0	193
12	Supramolecular self-assembled fullerene nanostructures. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 5075-5080.	3.3	191
13	Remarkable Positional Discrimination in Bistable Light- and Heat-Switchable Hydrogen-Bonded Molecular Shuttles. <i>Angewandte Chemie - International Edition</i> , 2003, 42, 2296-2300.	7.2	187
14	Photoisomerization of a rotaxane hydrogen bonding template: Light-induced acceleration of a large amplitude rotational motion. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2003, 100, 10-14.	3.3	185
15	Pentagon adjacency as a determinant of fullerene stability. <i>Physical Chemistry Chemical Physics</i> , 1999, 1, 2913-2918.	1.3	178
16	Ordering Fullerene Materials at Nanometer Dimensions. <i>Accounts of Chemical Research</i> , 2005, 38, 38-43.	7.6	177
17	Interpretation of the vibrational structure of the emission and absorption spectra of C ₆₀ . <i>Journal of Chemical Physics</i> , 1992, 97, 6496-6503.	1.2	165
18	Parallel (Face-to-Face) Versus Perpendicular (Edge-to-Face) Alignment of Electron Donors and Acceptors in Fullerene Porphyrin Dyads: The Importance of Orientation in Electron Transfer. <i>Journal of the American Chemical Society</i> , 2001, 123, 9166-9167.	6.6	157

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19	Graphene Can Wreak Havoc with Cell Membranes. ACS Applied Materials & Interfaces, 2015, 7, 4406-4414.	4.0	142
20	Dual-Gate Organic Field-Effect Transistors as Potentiometric Sensors in Aqueous Solution. Advanced Functional Materials, 2010, 20, 898-905.	7.8	136
21	The Devil and Holy Water: Protein and Carbon Nanotube Hybrids. Accounts of Chemical Research, 2013, 46, 2454-2463.	7.6	136
22	Increasing cost of pentagon adjacency for larger fullerenes. Chemical Physics Letters, 1996, 250, 544-548.	1.2	130
23	Kinetics of Place-Exchange Reactions of Thiols on Gold Nanoparticles. Langmuir, 2003, 19, 5172-5174.	1.6	119
24	Switching "On" and "Off" the Expression of Chirality in Peptide Rotaxanes. Journal of the American Chemical Society, 2002, 124, 2939-2950.	6.6	118
25	Directly Observing Micelle Fusion and Growth in Solution by Liquid-Cell Transmission Electron Microscopy. Journal of the American Chemical Society, 2017, 139, 17140-17151.	6.6	118
26	Modulating Charge-Transfer Interactions in Topologically Different Porphyrin-C60 Dyads. Chemistry - A European Journal, 2003, 9, 4968-4979.	1.7	110
27	Baiting Proteins with C ₆₀ . ACS Nano, 2010, 4, 2283-2299.	7.3	104
28	Entropy-Driven Translational Isomerism: A Tristable Molecular Shuttle. Angewandte Chemie - International Edition, 2003, 42, 5886-5889.	7.2	103
29	Extremely Strong and Readily Accessible AAA~DDD Triple Hydrogen Bond Complexes. Journal of the American Chemical Society, 2007, 129, 476-477.	6.6	103
30	Cadiot-Chodkiewicz Active Template Synthesis of Rotaxanes and Switchable Molecular Shuttles with Weak Intercomponent Interactions. Angewandte Chemie - International Edition, 2008, 47, 4392-4396.	7.2	101
31	Singling out the Electrochemistry of Individual Single-Walled Carbon Nanotubes in Solution. Journal of the American Chemical Society, 2008, 130, 7393-7399.	6.6	99
32	A molecular dynamics investigation of structure and dynamics of SDS and SDBS micelles. Soft Matter, 2011, 7, 9148.	1.2	99
33	C62: Theoretical Evidence for a Nonclassical Fullerene with a Heptagonal Ring. The Journal of Physical Chemistry, 1996, 100, 15634-15636.	2.9	97
34	Mechanochemistry: targeted delivery of single molecules. Nature Nanotechnology, 2006, 1, 122-125.	15.6	95
35	Theoretical study of the force fields of the three lowest singlet electronic states of linear polyenes. Journal of Chemical Physics, 1988, 89, 3681-3688.	1.2	93
36	C36, a hexavalent building block for fullerene compounds and solids. Chemical Physics Letters, 1999, 300, 369-378.	1.2	93

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37	Controlling the Frequency of Macrocyclic Ring Rotation in Benzylic Amide [2]Catenanes. <i>Journal of the American Chemical Society</i> , 1998, 120, 6458-6467.	6.6	92
38	QCFF/PI vibrational frequencies of some spherical carbon clusters. <i>Journal of the American Chemical Society</i> , 1991, 113, 6037-6040.	6.6	87
39	Temperature-Dependent Fluorescence of Cu ₅ Metal Clusters: A Molecular Thermometer. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 9662-9665.	7.2	87
40	Molecular Mechanism of Water Bridge Buildup: Field-Induced Formation of Nanoscale Menisci. <i>Langmuir</i> , 2008, 24, 6116-6120.	1.6	86
41	Charge-density oscillation on graphite induced by the interference of electron waves. <i>Physical Review B</i> , 2005, 71, .	1.1	85
42	Energetics of Fullerenes with Four-Membered Rings. <i>The Journal of Physical Chemistry</i> , 1996, 100, 6984-6991.	2.9	84
43	Probing the Structure of Lysozyme-Carbon Nanotube Hybrids with Molecular Dynamics. <i>Chemistry - A European Journal</i> , 2012, 18, 4308-4313.	1.7	84
44	The Effect of Mechanical Interlocking on Crystal Packing: Predictions and Testing. <i>Journal of the American Chemical Society</i> , 2002, 124, 225-233.	6.6	83
45	Effects of Electric Field Stress on a β -Amyloid Peptide. <i>Journal of Physical Chemistry B</i> , 2009, 113, 369-376.	1.2	83
46	Fast photodynamics of azobenzene probed by scanning excited-state potential energy surfaces using slow spectroscopy. <i>Nature Communications</i> , 2015, 6, 5860.	5.8	82
47	Cyclic Voltammetry and Bulk Electronic Properties of Soluble Carbon Nanotubes. <i>Journal of the American Chemical Society</i> , 2004, 126, 1646-1647.	6.6	80
48	Stability, Dynamics, and Lubrication of MoS ₂ Platelets and Nanotubes. <i>Langmuir</i> , 2012, 28, 7393-7400.	1.6	80
49	Tackling the Challenges of Dynamic Experiments Using Liquid-Cell Transmission Electron Microscopy. <i>Accounts of Chemical Research</i> , 2018, 51, 3-11.	7.6	78
50	Wrapping Nanotubes with Micelles, Hemimicelles, and Cylindrical Micelles. <i>Small</i> , 2009, 5, 2191-2198.	5.2	77
51	Dynamics of molecular self-ordering in tetraphenyl porphyrin monolayers on metallic substrates. <i>Nanotechnology</i> , 2009, 20, 275602.	1.3	75
52	Vibrational spectrum and harmonic force field of trimethylamine. <i>The Journal of Physical Chemistry</i> , 1993, 97, 581-595.	2.9	74
53	Electronic Absorption Spectra of Some Alkoxy Radicals. An Experimental and Theoretical Study. <i>Journal of the American Chemical Society</i> , 1995, 117, 2711-2718.	6.6	74
54	Reducing Molecular Shuttling to a Single Dimension. <i>Angewandte Chemie - International Edition</i> , 2000, 39, 350-353.	7.2	74

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55	Role of Substrate in Directing the Self-Assembly of Multicomponent Supramolecular Networks at the Liquid-Solid Interface. ACS Nano, 2012, 6, 8381-8389.	7.3	74
56	C ₆₀ @Lysozyme: Direct Observation by Nuclear Magnetic Resonance of a 1:1 Fullerene Protein Adduct. ACS Nano, 2014, 8, 1871-1877.	7.3	70
57	How Do Benzylic Amide [2]Catenane Rings Rotate?. Journal of the American Chemical Society, 1999, 121, 2364-2379.	6.6	69
58	Light-induced oxygen incision of C ₆₀ . Journal of the Chemical Society Chemical Communications, 1993, 220.	2.0	67
59	An Experimentally Observed Trimetallofullerene Sm ₃ @I _h -C ₈₀ : Encapsulation of Three Metal Atoms in a Cage without a Nonmetallic Mediator. Journal of the American Chemical Society, 2013, 135, 4187-4190.	6.6	67
60	Low-lying electronic excited states of Buckminsterfullerene anions. Journal of the American Chemical Society, 1992, 114, 2909-2913.	6.6	66
61	Modulation of the Reduction Potentials of Fullerene Derivatives. Journal of the American Chemical Society, 2003, 125, 7139-7144.	6.6	66
62	Charging and equilibration of fullerene isomers. Chemical Physics Letters, 1995, 243, 36-41.	1.2	65
63	Energetics of fullerenes with heptagonal rings. Journal of the Chemical Society, Faraday Transactions, 1996, 92, 2203.	1.7	65
64	The missing fluorescence of s ^{trans} butadiene. Journal of Chemical Physics, 1990, 93, 1235-1245.	1.2	64
65	Electrochemical Monitoring of Valence Bond Isomers Interconversion in Bipyridyl-C ₆₁ Anions. Journal of the American Chemical Society, 1995, 117, 6572-6580.	6.6	64
66	The S ₀ (1Ag) ← S ₁ (1B _{2u}) vibronic transition in benzene: Anabinitio study. Journal of Chemical Physics, 1994, 100, 2458-2464.	1.2	60
67	Title is missing!. Angewandte Chemie, 2003, 115, 2398-2402.	1.6	60
68	The Erratic Emission of Pyrene on Gold Nanoparticles. ACS Nano, 2008, 2, 77-84.	7.3	60
69	A Carbon Nano-Onion Ferrocene Donor-Acceptor System: Synthesis, Characterization and Properties. Chemistry - A European Journal, 2009, 15, 4419-4427.	1.7	58
70	A Simple Road for the Transformation of Few-Layer Graphene into MWNTs. Journal of the American Chemical Society, 2012, 134, 13310-13315.	6.6	58
71	Biorecognition in Organic Field Effect Transistors Biosensors: The Role of the Density of States of the Organic Semiconductor. Analytical Chemistry, 2016, 88, 12330-12338.	3.2	58
72	Conformational Self-Recognition as the Origin of Dewetting in Bistable Molecular Surfaces. Journal of Physical Chemistry B, 2001, 105, 10826-10830.	1.2	57

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73	Modeling the Adsorption of Alkanes on an Au(111) Surface. <i>Langmuir</i> , 2003, 19, 7335-7340.	1.6	57
74	Electronic Structure of Carbon Nanotubes with Ultrahigh Curvature. <i>ACS Nano</i> , 2010, 4, 4515-4522.	7.3	57
75	Conformational Selection and Folding-upon-binding of Intrinsically Disordered Protein CP12 Regulate Photosynthetic Enzymes Assembly. <i>Journal of Biological Chemistry</i> , 2012, 287, 21372-21383.	1.6	57
76	Circumstellar carbon chain molecules: A density function theory study of C _n O, n=3-9. <i>Journal of Chemical Physics</i> , 1995, 103, 6343-6349.	1.2	56
77	Amyloid- β fibril disruption by C60 molecular guidance for rational drug design. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 8599.	1.3	56
78	Frequency-Dependent Second-Order Hyperpolarizability of Carbon Clusters: A Semiempirical Investigation. <i>Journal of the American Chemical Society</i> , 1995, 117, 6101-6108.	6.6	55
79	A Tight-Binding Treatment for ¹³ C NMR Spectra of Fullerenes. <i>Journal of Physical Chemistry A</i> , 1999, 103, 8738-8746.	1.1	55
80	From reactants to products via simple hydrogen-bonding networks: Information transmission in chemical reactions. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 4967-4971.	3.3	53
81	Molecular Dynamics of a Dendrimer-Dye Guest-Host System. <i>Journal of the American Chemical Society</i> , 2003, 125, 7388-7393.	6.6	53
82	A computational analysis of the insertion of carbon nanotubes into cellular membranes. <i>Biomaterials</i> , 2011, 32, 7079-7085.	5.7	53
83	Electronic and mechanical coupling between guest and host in carbon peapods. <i>Physical Review B</i> , 2004, 69, .	1.1	52
84	Modeling the Stability and the Motion of DNA Nucleobases on the Gold Surface. <i>Langmuir</i> , 2005, 21, 2512-2518.	1.6	52
85	The vibrational frequencies of C60. <i>Chemical Physics Letters</i> , 1992, 190, 174-178.	1.2	51
86	Theoretical study of the force field of the lowest singlet electronic states of long polyenes. <i>Journal of Chemical Physics</i> , 1989, 91, 6215-6224.	1.2	50
87	The vibrational spectroscopy of C60H36: An experimental and theoretical study. <i>Chemical Physics</i> , 1998, 232, 75-94.	0.9	50
88	The Mechanism of Formation of Amide-Based Interlocked Compounds: Prediction of a New Rotaxane-Forming Motif. <i>Chemistry - A European Journal</i> , 2004, 10, 4960-4969.	1.7	50
89	The Stone-Wales map for C60. <i>Chemical Physics Letters</i> , 1995, 235, 146-151.	1.2	49
90	Rolling up a Graphene Sheet. <i>ChemPhysChem</i> , 2013, 14, 3447-3453.	1.0	49

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91	Vibronic coupling in polyenes: The frequency increase of the active C \rightarrow C ag stretching mode in the absorption spectra. <i>Chemical Physics</i> , 1986, 108, 187-195.	0.9	48
92	Theoretical analysis of the force field of the lowest excited singlet state of trans-stilbene. <i>The Journal of Physical Chemistry</i> , 1989, 93, 5124-5128.	2.9	48
93	Incorporation of Fullerene Derivatives into Smectite Clays: A New Family of Organic-Inorganic Nanocomposites. <i>Journal of the American Chemical Society</i> , 2004, 126, 8561-8568.	6.6	47
94	Shape Governs the Motion of Chemically Propelled Janus Swimmers. <i>Journal of Physical Chemistry C</i> , 2012, 116, 592-598.	1.5	47
95	Structural Motifs and the Stability of Fullerenes. <i>The Journal of Physical Chemistry</i> , 1995, 99, 8076-8081.	2.9	46
96	Clay-Fulleropyrrolidine Nanocomposites. <i>Journal of the American Chemical Society</i> , 2006, 128, 6154-6163.	6.6	46
97	Experimental and Theoretical Study of the Infrared, Raman, and Electronic Spectra of Two Isomers of C ₇₈ of C _{2v} Symmetry. <i>The Journal of Physical Chemistry</i> , 1996, 100, 13399-13407.	2.9	45
98	The hyperpolarisability of an endohedral fullerene: Li@C ₆₀ . <i>Chemical Physics Letters</i> , 1998, 288, 131-137.	1.2	45
99	Assignment and vibrational analysis of the 600 nm absorption band in the phenoxyl radical and some of its derivatives. <i>Canadian Journal of Chemistry</i> , 1993, 71, 1655-1662.	0.6	44
100	The intramolecular vibrations of prototypical polythiophenes. <i>Journal of Chemical Physics</i> , 1996, 104, 9704-9718.	1.2	44
101	C ₃₆ : The Best Fullerene for Covalent Bonding. <i>Journal of the American Chemical Society</i> , 1999, 121, 3218-3219.	6.6	44
102	Water-induced polaron formation at the pentacene surface: Quantum mechanical molecular mechanics simulations. <i>Physical Review B</i> , 2009, 79, .	1.1	44
103	Graphite Oxide and Aromatic Amines: Size Matters. <i>Advanced Functional Materials</i> , 2015, 25, 263-269.	7.8	44
104	Vibronic coupling in the benzyl radical. <i>Chemical Physics Letters</i> , 1985, 115, 253-258.	1.2	43
105	Nanopatterning of carbonaceous structures by field-induced carbon dioxide splitting with a force microscope. <i>Applied Physics Letters</i> , 2010, 96, .	1.5	43
106	In Silico Carborane Docking to Proteins and Potential Drug Targets. <i>Journal of Chemical Information and Modeling</i> , 2011, 51, 1882-1896.	2.5	43
107	The electronic structure and vibrational frequencies of the stable C ₇₆ isomer of D ₂ symmetry. <i>Chemical Physics Letters</i> , 1993, 208, 441-445.	1.2	42
108	Dynamics of Thiolate Chains on a Gold Nanoparticle. <i>Small</i> , 2007, 3, 386-388.	5.2	42

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109	An introduction to bubble dynamics. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 2447.	1.3	42
110	Vibronic coupling in polyenes and their derivatives. Interpretation of the absorption and emission spectra of a derivative of dodecahexaene. <i>Journal of Chemical Physics</i> , 1987, 87, 2505-2512.	1.2	41
111	Quantum chemical and vibronic analysis of the $1^1A_g \rightarrow 1^1A_g$ transition in benzyl and benzyl radicals. <i>Journal of Chemical Physics</i> , 1990, 93, 600-608.	1.2	41
112	Infrared fingerprints of nine fullerene C ₈₂ isomers: a semiempirical prediction. <i>The Journal of Physical Chemistry</i> , 1993, 97, 13575-13579.	2.9	41
113	Fullerene sorting proteins. <i>Nanoscale</i> , 2011, 3, 2873.	2.8	41
114	Atomistic molecular dynamics simulations reveal insights into adsorption, packing, and fluxes of molecules with carbon nanotubes. <i>Journal of Materials Chemistry A</i> , 2014, 2, 12123-12135.	5.2	41
115	Blocking the Passage: C ₆₀ Geometrically Clogs K ⁺ Channels. <i>ACS Nano</i> , 2015, 9, 4827-4834.	7.3	41
116	Absolute intensities of CH-stretching overtones in chloroform and deuteriochloroform. <i>Chemical Physics Letters</i> , 1989, 154, 273-279.	1.2	40
117	Dynamics of molecular inversion: An instanton approach. <i>Journal of Chemical Physics</i> , 1995, 102, 7024-7034.	1.2	40
118	Energetics of C ₂₀ and C ₂₂ Fullerene and Near-Fullerene Carbon Cages. <i>Journal of Physical Chemistry A</i> , 1997, 101, 8339-8344.	1.1	40
119	Adsorption of a Benzylic Amide Macrocycle on a Solid Substrate: XPS and HREELS Characterization of Thin Films Grown on Au(111). <i>Journal of Physical Chemistry B</i> , 2002, 106, 8739-8746.	1.2	40
120	Selective Enhancement of Photoluminescence in Filled Single-Walled Carbon Nanotubes. <i>Advanced Functional Materials</i> , 2012, 22, 3202-3208.	7.8	40
121	Thermodynamics of Binding Between Proteins and Carbon Nanoparticles: The Case of C ₆₀ @Lysozyme. <i>Journal of Physical Chemistry C</i> , 2015, 119, 28077-28082.	1.5	40
122	New Assignments in the 600 nm Region of C ₆₀ : The Origins of the T _{1g} and G _g Transitions. <i>The Journal of Physical Chemistry</i> , 1996, 100, 10849-10853.	2.9	39
123	Large Enhancement of the Nonlinear Optical Response of Reduced Fullerene Derivatives. <i>Chemistry - A European Journal</i> , 2003, 9, 1529-1534.	1.7	39
124	Simple models for hydrophobic hydration. <i>Chemical Society Reviews</i> , 2005, 34, 1012.	18.7	39
125	Application of the Poisson-Nernst-Planck Theory with Space-Dependent Diffusion Coefficients to KcsA. <i>Biophysical Journal</i> , 2006, 91, 3162-3169.	0.2	39
126	Electric Field Effects on Short Fibrils of A β Amyloid Peptides. <i>Journal of Chemical Theory and Computation</i> , 2010, 6, 3516-3526.	2.3	39

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127	Inelastic neutron-scattering study of the intramolecular vibrations of the C70 fullerene. <i>The Journal of Physical Chemistry</i> , 1993, 97, 3641-3643.	2.9	38
128	What Is Adenine Doing in Photolyase?. <i>Journal of Physical Chemistry B</i> , 2010, 114, 4101-4106.	1.2	38
129	Mechanical Interactions in All-Carbon Peapods. <i>Journal of Physical Chemistry B</i> , 2003, 107, 6986-6990.	1.2	37
130	Molecular dynamics study of onset of water gelation around the collagen triple helix. <i>Proteins: Structure, Function and Bioinformatics</i> , 2006, 64, 711-718.	1.5	37
131	Pyridinedicarboxamide Strands Form Double Helices via an Activated Slippage Mechanism. <i>Journal of the American Chemical Society</i> , 2004, 126, 2362-2367.	6.6	36
132	On-the-Fly, Electric-Field-Driven, Coupled Electron~Nuclear Dynamics. <i>Journal of Physical Chemistry A</i> , 2008, 112, 9650-9656.	1.1	36
133	Comparison of synchronous and asynchronous hydrogen transfer mechanisms in free-base porphyrins. <i>Chemical Physics</i> , 1989, 136, 285-295.	0.9	35
134	Theoretical study of the vibronic structure of the $1A_1^+$ and $1B_2$ and $1A_1^+$ and $3B_2$ electronic transitions in cyclopentadiene. <i>Journal of Chemical Physics</i> , 1993, 99, 3721-3729.	1.2	35
135	Growth of <i>p</i> - and <i>n</i> -Dopable Films from Electrochemically Generated C ₆₀ Cations. <i>Journal of the American Chemical Society</i> , 2008, 130, 3788-3796.	6.6	35
136	Splitting CO ₂ with Electric Fields: A Computational Investigation. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 3256-3260.	2.1	34
137	Interaction Model for the Adsorption of Organic Molecules on the Silver Surface. <i>Journal of Physical Chemistry B</i> , 2006, 110, 5595-5601.	1.2	34
138	Quantum-chemical analysis of vibrational intensity distributions in the S ₀ -S ₁ absorption spectra and raman excitation profiles of azulene. <i>Chemical Physics</i> , 1986, 110, 421-430.	0.9	33
139	Theoretical analysis of Franck-Condon and vibronic activity of the a_g and b_{3g} modes in the S ₀ ~S ₁ transitions in anthracene. <i>Chemical Physics</i> , 1988, 127, 17-29.	0.9	33
140	Relative stabilities of C76 isomers. A numerical test of the fullerene isolated-pentagon rule. <i>Chemical Physics Letters</i> , 1994, 226, 219-225.	1.2	33
141	A Density Functional Study of the Vibrations of Three Oligomers of Thiophene. <i>Journal of Physical Chemistry A</i> , 1997, 101, 7283-7291.	1.1	33
142	The Effect of Guest Inclusion on the Crystal Packing of <i>p</i> -tert-Butylcalix[4]arenes. <i>Chemistry - A European Journal</i> , 2002, 8, 4854-4866.	1.7	33
143	Modelling of the Adsorption of C60 on the Au(110) Surface. <i>ChemPhysChem</i> , 2004, 5, 245-248.	1.0	33
144	Proteins as supramolecular hosts for C ₆₀ : a true solution of C ₆₀ in water. <i>Nanoscale</i> , 2018, 10, 9908-9916.	2.8	33

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145	Intermolecular Repulsion through Interfacial Attraction: Toward Engineering of Polymorphs. <i>Journal of the American Chemical Society</i> , 2009, 131, 15655-15659.	6.6	32
146	Controlled Hydrogen-Bond Breaking in a Rotaxane by Discrete Solvation. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 3896-3900.	7.2	32
147	Calcite Single Crystals as Hosts for Atomic-Scale Entrapment and Slow Release of Drugs. <i>Advanced Healthcare Materials</i> , 2015, 4, 1510-1516.	3.9	32
148	The Collapse of Nanobubbles in Water. <i>Journal of the American Chemical Society</i> , 2005, 127, 8020-8021.	6.6	31
149	Unexpected Photophysical Properties of Symmetric Indolylmaleimide Derivatives. <i>Journal of Physical Chemistry A</i> , 2005, 109, 6440-6449.	1.1	31
150	Customizing Properties of β -Chitin in Squid Pen (<i>Gladius</i>) by Chemical Treatments. <i>Marine Drugs</i> , 2014, 12, 5979-5992.	2.2	31
151	Observation and modeling of the recombination kinetics of diphenylmethyl radicals in the cavities of Na-X zeolite. <i>The Journal of Physical Chemistry</i> , 1991, 95, 10018-10024.	2.9	30
152	The infrared and Raman active vibrational frequencies of C ₆₀ hexaanion. <i>Chemical Physics Letters</i> , 1992, 196, 303-310.	1.2	30
153	The Franck-Condon structure of the $1A \rightarrow 1B$ transition of cis- and trans-hexatriene: An ab initio modeling. <i>Journal of Chemical Physics</i> , 1993, 98, 4822-4829.	1.2	30
154	Molecular Structure of Stilbene in the T ₁ State. Transient Resonance Raman Spectra of Stilbene Isotopomers and Quantum Chemical Calculations. <i>The Journal of Physical Chemistry</i> , 1994, 98, 2254-2265.	2.9	30
155	Energetics and isomerisation pathways of a lower fullerene. The Stone-Wales map for C ₄₀ . <i>Journal of the Chemical Society, Faraday Transactions</i> , 1995, 91, 1421-1423.	1.7	30
156	Playing peekaboo with graphene oxide: a scanning electrochemical microscopy investigation. <i>Chemical Communications</i> , 2014, 50, 13117-13120.	2.2	30
157	Graphene Oxide Promotes Site-Selective Allylic Alkylation of Thiophenes with Alcohols. <i>Organic Letters</i> , 2018, 20, 3705-3709.	2.4	30
158	Identification and preparation of stable water dispersions of protein - Carbon nanotube hybrids and efficient design of new functional materials. <i>Carbon</i> , 2019, 147, 70-82.	5.4	30
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