

Tamás Keszthelyi

List of Publications by Year in descending order

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25
papers

483
citations

706676

14
h-index

759306

22
g-index

25
all docs

25
docs citations

25
times ranked

644
citing authors

#	ARTICLE	IF	CITATIONS
1	Quantum-chemistry-aided ligand engineering for potential molecular switches: changing barriers to tune excited state lifetimes. <i>Chemical Communications</i> , 2020, 56, 11831-11834.	2.2	4
2	Bilayer Charge Reversal and Modification of Lipid Organization by Dendrimers as Observed by Sum-Frequency Vibrational Spectroscopy. <i>Langmuir</i> , 2015, 31, 7815-7825.	1.6	11
3	Sodium selective ion channel formation in living cell membranes by polyamidoamine dendrimer. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2013, 1828, 1873-1880.	1.4	20
4	Interaction of Phospholipid Langmuir Monolayers with an Antibiotic Peptide Conjugate. <i>Journal of Physical Chemistry B</i> , 2013, 117, 6969-6979.	1.2	13
5	Organization of the enantiomeric and racemic forms of an amphiphilic resorcinol derivative at the air/water and graphite/phenyloctane interfaces. <i>Chirality</i> , 2012, 24, 155-166.	1.3	11
6	Interaction of Carbon Monoxide with Au(111) Modified by Ion Bombardment: A Surface Spectroscopy Study under Elevated Pressure. <i>Langmuir</i> , 2010, 26, 16312-16324.	1.6	34
7	Characterisation of the membrane affinity of an isoniazide peptide conjugate by tensiometry, atomic force microscopy and sum-frequency vibrational spectroscopy, using a phospholipid Langmuir monolayer model. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 11498.	1.3	14
8	Study of FeOx/Au inverse model catalysts by in situ sum frequency generation vibrational spectroscopy. <i>Reaction Kinetics and Catalysis Letters</i> , 2009, 96, 345-356.	0.6	4
9	Investigation of Solid Surfaces Modified by Langmuir-Blodgett Monolayers Using Sum-Frequency Vibrational Spectroscopy and X-ray Photoelectron Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2006, 110, 8701-8714.	1.2	56
10	Adsorbed and Spread Layers of Poly(ethylene oxide)-Poly(propylene oxide)-Poly(ethylene oxide) Block Copolymers at the Air/Water Interface Studied by Sum-Frequency Vibrational Spectroscopy and Tensiometry. <i>Macromolecules</i> , 2006, 39, 9375-9384.	2.2	32
11	Observation of a Liquid-Gas Phase Transition in Monolayers of Alkyltrimethylammonium Alkyl Sulfates Adsorbed at the Air/Water Interface. <i>Journal of Physical Chemistry B</i> , 2005, 109, 872-878.	1.2	35
12	Electron Delocalization in the Radical Cation of 1,3,6,8-Tetraazatricyclo[4.4.1.1.3,8]dodecane, a 4-Nitrogen-7-Electron System. <i>Journal of the American Chemical Society</i> , 2002, 124, 159-167.	6.6	20
13	Effect of Solvent on the O ₂ (¹ g) → O ₂ (¹ g ₊) Absorption Spectrum: Demonstrating the Importance of Equilibrium vs Nonequilibrium Solvation. <i>Journal of Physical Chemistry A</i> , 2002, 106, 5263-5270.	1.1	18
14	Evidence of Dynamical Jahn-Teller Effect on Triphenylene Radical Cation: Resonance Raman Spectrum and ab Initio Quantum-Chemical Calculations. <i>Journal of Physical Chemistry A</i> , 2000, 104, 9121-9129.	1.1	19
15	O ₂ (¹ g) Absorption and O ₂ (¹ g ₊) Emission in Solution: Quantifying the Stokes Shift. <i>Journal of Physical Chemistry A</i> , 2000, 104, 10550-10555.	1.1	24
16	The Radical Cation and Lowest Rydberg States of 1,4-Diaza[2.2.2]bicyclooctane (DABCO). <i>Journal of Physical Chemistry A</i> , 2000, 104, 1834-1841.	1.1	25
17	2,2'-Bithiophene Radical Cation: An Experimental and Computational Study. <i>Journal of Physical Chemistry A</i> , 2000, 104, 2808-2823.	1.1	26
18	Radiative Transitions of Singlet Oxygen: New Tools, New Techniques and New Interpretations. <i>Photochemistry and Photobiology</i> , 1999, 70, 531-539.	1.3	45

#	ARTICLE	IF	CITATIONS
19	Bithiophene radical cation: resonance Raman spectroscopy and molecular orbital calculations. Chemical Physics Letters, 1998, 282, 171-175.	1.2	7
20	Resonance Raman and quantum chemical studies of short polyene radical cations. Journal of Molecular Structure, 1997, 410-411, 339-342.	1.8	3
21	The radical cation of trans-1,3-pentadiene: resonance Raman spectrum and ab initio calculations. Journal of Molecular Structure, 1996, 379, 211-218.	1.8	7
22	Radical Cation of 1,3-Butadiene: Resonance Raman Spectrum of Deuterated Derivatives and Improved Force Field. The Journal of Physical Chemistry, 1996, 100, 16850-16855.	2.9	17
23	Radical Cations of Monomethyl-Substituted 1,3-Butadienes: Resonance Raman Spectra and Molecular Orbital Calculations. The Journal of Physical Chemistry, 1996, 100, 15785-15793.	2.9	7
24	Radical Cation of 1,3-Butadiene: Resonance Raman Spectrum of the δ Isotopomer. The Journal of Physical Chemistry, 1996, 100, 16843-16849.	2.9	18
25	Resonance Raman Investigation of the Radical Cation of 1,3,5-Hexatriene. The Journal of Physical Chemistry, 1994, 98, 5632-5638.	2.9	13