

Friedrich Temps

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6412993/publications.pdf>

Version: 2024-02-01

43
papers

1,785
citations

279798

23
h-index

265206

42
g-index

44
all docs

44
docs citations

44
times ranked

1871
citing authors

#	ARTICLE	IF	CITATIONS
1	Highly Efficient Reversible $Z \rightarrow E$ Photoisomerization of a Bridged Azobenzene with Visible Light through Resolved S_1 ($n\pi^*$) Absorption Bands. <i>Journal of the American Chemical Society</i> , 2009, 131, 15594-15595.	13.7	445
2	Photofragment velocity map imaging of H atom elimination in the first excited state of pyrrole. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 315-320.	2.8	128
3	Ultrafast Electronic Relaxation in Guanosine is Promoted by Hydrogen Bonding with Cytidine. <i>Journal of the American Chemical Society</i> , 2007, 129, 9272-9273.	13.7	113
4	Photodissociation dynamics of pyrrole: Evidence for mode specific dynamics from conical intersections. <i>Faraday Discussions</i> , 2004, 127, 267.	3.2	105
5	Femtosecond fluorescence up-conversion spectroscopy of adenine and adenosine: experimental evidence for the $\pi\pi^*$ state?. <i>Chemical Physics</i> , 2005, 313, 199-212.	1.9	93
6	Superior $Z \rightarrow E$ and $E \rightarrow Z$ photoswitching dynamics of dihydrodibenzodiazocine, a bridged azobenzene, by S_1 ($n\pi^*$) excitation at $\lambda = 387$ and 490 nm. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 1054-1063.	2.8	86
7	Femtosecond fluorescence up-conversion spectroscopy of a rotation-restricted azobenzene after excitation to the S_1 state. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 1985-1989.	2.8	58
8	Unimolecular dissociation dynamics of highly vibrationally excited DCO ($\chi_f^{\infty} 2A$). II. Calculation of resonance energies and widths and comparison with high-resolution spectroscopic data. <i>Journal of Chemical Physics</i> , 1997, 106, 5359-5378.	3.0	56
9	Unimolecular dissociation dynamics of highly vibrationally excited DCO ($\chi_f^{\infty} 2A$). I. Investigation of dissociative resonance states by stimulated emission pumping spectroscopy. <i>Journal of Chemical Physics</i> , 1997, 106, 5333-5358.	3.0	54
10	Ultraviolet Absorption Induces Hydrogen-Atom Transfer in G-C Watson-Crick DNA Base Pairs in Solution. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 14719-14722.	13.8	54
11	Ultrashort Fluorescence Lifetimes of Hydrogen-Bonded Base Pairs of Guanosine and Cytidine in Solution. <i>Journal of Physical Chemistry B</i> , 2009, 113, 16365-16376.	2.6	46
12	Ultrafast electronic deactivation dynamics of the rare natural nucleobase hypoxanthine. <i>Chemical Physics Letters</i> , 2012, 536, 140-146.	2.6	39
13	Parallel ultrafast $E \rightarrow C$ ring closure and $E \rightarrow Z$ isomerisation in a photochromic furylfulgide studied by femtosecond time-resolved spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 5952.	2.8	35
14	Probing the excited state relaxation dynamics of pyrimidine nucleosides in chloroform solution. <i>Faraday Discussions</i> , 2016, 194, 683-708.	3.2	31
15	Enhanced photoswitching and ultrafast dynamics in structurally modified photochromic fulgides. <i>International Reviews in Physical Chemistry</i> , 2013, 32, 1-38.	2.3	30
16	Ultrafast dynamics of the ESIPt photoswitch N -(3-pyridinyl)-2-pyridinecarboxamide. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 2646-2655.	2.8	29
17	Ultrafast transient absorption spectroscopy of the photo-induced $Z \rightarrow E$ isomerization of a photochromic furylfulgide. <i>Chemical Physics Letters</i> , 2006, 428, 62-67.	2.6	28
18	Sequential photoisomerisation dynamics of the push-pull azobenzene Disperse Red 1. <i>Photochemical and Photobiological Sciences</i> , 2012, 11, 1210-1219.	2.9	28

#	ARTICLE	IF	CITATIONS
19	Ultrafast nonradiative dynamics in electronically excited hexafluorobenzene by femtosecond time-resolved mass spectrometry. <i>Journal of Chemical Physics</i> , 2008, 128, 164314.	3.0	27

20

#	ARTICLE	IF	CITATIONS
37	Note: Energy calibration of a femtosecond photoelectron imaging detector with correction for the ponderomotive shift of atomic ionization energies. <i>Review of Scientific Instruments</i> , 2017, 88, 046101.	1.3	6
38	Real-time observation of multi-mode vibronic coherence in pentafluoropyridine. <i>Journal of Chemical Physics</i> , 2017, 147, 013938.	3.0	5
39	Resonance dynamics of DCO (χ_1^f) simulated with the dynamically pruned discrete variable representation (DP-DVR). <i>Journal of Chemical Physics</i> , 2018, 148, 204309.	3.0	4
40	Rotational state-dependent mixings between resonance states of vibrationally highly excited DCO (χ_1^f). <i>Journal of Chemical Physics</i> , 2004, 120, 10530-10542.	3.0	3
41	Efficient intersystem crossing in 2-aminopurine riboside probed by femtosecond time-resolved transient vibrational absorption spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 20033-20042.	2.8	3
42	Ultrafast Electronic Deactivation Dynamics of Xanthosine Monophosphate. <i>Molecules</i> , 2017, 22, 160.	3.8	2
43	Long-lived Excited States in 7- and 9-Methylpurine Probed by fs Time-Resolved Vibrational Absorption Spectroscopy. , 2020, , .		0