

Wybren Jan Buma

List of Publications by Year in descending order

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185
papers

4,803
citations

101384

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all docs

187
docs citations

187
times ranked

5367
citing authors

#	ARTICLE	IF	CITATIONS
1	Elucidating the photoprotective properties of natural UV screening agents: ZEPHYRUS PFI spectroscopy of methyl sinapate. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 3984-3993.	1.3	5
2	Controlling forward and backward rotary molecular motion on demand. <i>Nature Communications</i> , 2022, 13, 2124.	5.8	15
3	Polycyclic aromatic hydrocarbon growth in a benzene discharge explored by IR-UV action spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 14816-14824.	1.3	6
4	Excited-state dynamics of isolated and (micro)solvated methyl sinapate: the bright and shady sides of a natural sunscreen. <i>Molecular Physics</i> , 2021, 119, e1825850.	0.8	13
5	High-resolution infrared spectroscopy of naphthalene and acenaphthene dimers. <i>Molecular Physics</i> , 2021, 119, e1811908.	0.8	7
6	Vibrational circular dichroism studies of exceptionally strong chirality inducers in liquid crystals. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 10021-10028.	1.3	6
7	Temperature-dependent modulation by biaryl-based monomers of the chain length and morphology of biphenyl-based supramolecular polymers. <i>Chemical Science</i> , 2021, 12, 13001-13012.	3.7	6
8	Allosteric Guest Binding in Chiral Zirconium(IV) Double Decker Porphyrin Cages. <i>European Journal of Organic Chemistry</i> , 2021, 2021, 607-617.	1.2	2
9	Tailoring the optical and dynamic properties of iminothioindoxyl photoswitches through acidochromism. <i>Chemical Science</i> , 2021, 12, 4588-4598.	3.7	13
10	Towards developing novel and sustainable molecular light-to-heat converters. <i>Chemical Science</i> , 2021, 12, 15239-15252.	3.7	18
11	Phenylimino Indolinone: A Green-Light-Responsive π -Type Photoswitch Exhibiting Negative Photochromism. <i>Angewandte Chemie</i> , 2021, 133, 25494.	1.6	2
12	Phenylimino Indolinone: A Green-Light-Responsive π -Type Photoswitch Exhibiting Negative Photochromism. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 25290-25295.	7.2	21
13	Infrared Spectroscopy of Jet-cooled α -GrandPAHs in the 3×10^4 μ m Region. <i>Astrophysical Journal</i> , 2021, 923, 238.	1.6	4
14	Self-Assembly of Supramolecular Polymers of N-Centered Triarylamine Trisamides in the Light of Circular Dichroism: Reaching Consensus between Electrons and Nuclei. <i>Journal of the American Chemical Society</i> , 2020, 142, 1020-1028.	6.6	17
15	Far-IR Absorption of Neutral Polycyclic Aromatic Hydrocarbons (PAHs): Light on the Mechanism of IR-UV Ion Dip Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 8997-9002.	2.1	4
16	Vibrational circular dichroism spectroscopy for probing the expression of chirality in mechanically planar chiral rotaxanes. <i>Chemical Science</i> , 2020, 11, 8469-8475.	3.7	19
17	Photoinduced Forward and Backward Pedalo-Type Motion of a Molecular Switch. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 4741-4746.	2.1	3
18	Photo-activated CO-release in the amino tungsten Fischer carbene complex, $[(CO)_5WC(NC_4H_8)Me]$, picosecond time resolved infrared spectroscopy, time-dependent density functional theory, and an antimicrobial study. <i>Journal of Inorganic Biochemistry</i> , 2020, 208, 111071.	1.5	6

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19	Analysis of Vibrational Circular Dichroism Spectra of Peptides: A Generalized Coupled Oscillator Approach of a Small Peptide Model Using VCDtools. <i>Journal of Physical Chemistry B</i> , 2020, 124, 1665-1677.	1.2	5
20	GUI Implementation of VCDtools, A Program to Analyze Computed Vibrational Circular Dichroism Spectra. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 259-267.	2.5	11
21	Circular Spectropolarimetric Sensing of Vegetation in the Field: Possibilities for the Remote Detection of Extraterrestrial Life. <i>Astrobiology</i> , 2019, 19, 1221-1229.	1.5	22
22	Temperature Control of Sequential Nucleation-Growth Mechanisms in Hierarchical Supramolecular Polymers. <i>Chemistry - A European Journal</i> , 2019, 25, 13008-13016.	1.7	28
23	Taming conformational heterogeneity in and with vibrational circular dichroism spectroscopy. <i>Chemical Science</i> , 2019, 10, 7680-7689.	3.7	40
24	A Tunable, Fullerene-Based Molecular Amplifier for Vibrational Circular Dichroism. <i>Chemistry - A European Journal</i> , 2019, 25, 12560-12566.	1.7	5
25	Vibrational Circular Dichroism of Thiolate-Protected Au ₂₅ Clusters: Accurate Prediction of Spectra and Chirality Transfer within the Mixed Ligand Shell. <i>Journal of Physical Chemistry C</i> , 2019, 123, 22586-22594.	1.5	9
26	Anharmonicity in the mid-infrared spectra of polycyclic aromatic hydrocarbons: molecular beam spectroscopy and calculations. <i>Astronomy and Astrophysics</i> , 2019, 628, A130.	2.1	21
27	Analytical chemistry on many-center chiral compounds based on vibrational circular dichroism: Absolute configuration assignments and determination of contaminant levels. <i>Analytica Chimica Acta</i> , 2019, 1090, 100-105.	2.6	11
28	Opening 2,2-diphenyl-2H-chromene to infrared light. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 11689-11696.	1.3	7
29	Iminothioindoxyl as a molecular photoswitch with 100 nm band separation in the visible range. <i>Nature Communications</i> , 2019, 10, 2390.	5.8	63
30	Frontispiece: Light on the Structural Evolution of Photoresponsive Molecular Switches in Electronically Excited States. <i>Chemistry - A European Journal</i> , 2019, 25, .	1.7	0
31	Taming the Complexity of Donor-Acceptor Stenhouse Adducts: Infrared Motion Pictures of the Complete Switching Pathway. <i>Journal of the American Chemical Society</i> , 2019, 141, 7376-7384.	6.6	66
32	Circular spectropolarimetric sensing of higher plant and algal chloroplast structural variations. <i>Photosynthesis Research</i> , 2019, 140, 129-139.	1.6	12
33	Light on the Structural Evolution of Photoresponsive Molecular Switches in Electronically Excited States. <i>Chemistry - A European Journal</i> , 2019, 25, 6252-6258.	1.7	2
34	Photo- and Electrochemical Properties of a CO ₂ Reducing Ruthenium-Rhenium Quaterpyridine-Based Catalyst. <i>ChemPhotoChem</i> , 2018, 2, 323-331.	1.5	18
35	Enantiospecific Brook Rearrangement of Tertiary Benzylic β -Hydroxysilanes. <i>European Journal of Organic Chemistry</i> , 2018, 2018, 3900-3903.	1.2	11
36	Tailoring Photoisomerization Pathways in Donor-Acceptor Stenhouse Adducts: The Role of the Hydroxy Group. <i>Journal of Physical Chemistry A</i> , 2018, 122, 955-964.	1.1	54

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37	Regional Susceptibility in VCD Spectra to Dynamic Molecular Motions: The Case of a Benzyl Hydroxysilane. <i>ChemPhysChem</i> , 2018, 19, 561-565.	1.0	9
38	Titelbild: Photoinduced Pedal-Type Motion in an Azodicarboxamide-Based Molecular Switch (<i>Angew. Chem.</i>)	1.6	0
39	Fluorescence Correlation Spectroscopy of Labeled Azurin Reveals Photoinduced Electron Transfer between Label and Cu Center. <i>Chemistry - A European Journal</i> , 2018, 24, 646-654.	1.7	3
40	The anharmonic quartic force field infrared spectra of hydrogenated and methylated PAHs. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 1189-1197.	1.3	46
41	Photoinduced Pedal-Type Motion in an Azodicarboxamide-Based Molecular Switch. <i>Angewandte Chemie</i> , 2018, 130, 1810-1814.	1.6	7
42	Photoinduced Pedal-Type Motion in an Azodicarboxamide-Based Molecular Switch. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 1792-1796.	7.2	21
43	Frequency Range Selection Method for Vibrational Spectra. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 6878-6882.	2.1	7
44	Use of Density Functional Based Tight Binding Methods in Vibrational Circular Dichroism. <i>Journal of Physical Chemistry A</i> , 2018, 122, 9435-9445.	1.1	6
45	Solvent Effects on the Actinic Step of Donor-Acceptor Stenhouse Adduct Photoswitching. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 8063-8068.	7.2	70
46	Solvent Effects on the Actinic Step of Donor-Acceptor Stenhouse Adduct Photoswitching. <i>Angewandte Chemie</i> , 2018, 130, 8195-8200.	1.6	21
47	Excited-State Electronic Asymmetry Prevents Photoswitching in Terthiophene Compounds. <i>Inorganic Chemistry</i> , 2018, 57, 9039-9047.	1.9	1
48	Vibrationally-resolved spectroscopic studies of electronically excited states of 1,8-naphthalic anhydride and 1,8-naphthalimide: a delicate interplay between one $\pi\pi^*$ and two $n\pi^*$ states. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 5861-5869.	1.3	4
49	Circular spectropolarimetric sensing of chiral photosystems in decaying leaves. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2017, 189, 303-311.	1.1	24
50	Shedding Light on the Photoisomerization Pathway of Donor-Acceptor Stenhouse Adducts. <i>Journal of the American Chemical Society</i> , 2017, 139, 15596-15599.	6.6	88
51	High-resolution gas-phase spectroscopy of a single-bond axle rotary motor. <i>Tetrahedron</i> , 2017, 73, 4887-4890.	1.0	1
52	Transient two-dimensional vibrational spectroscopy of an operating molecular machine. <i>Nature Communications</i> , 2017, 8, 2206.	5.8	13
53	Interplay of Exciton Coupling and Large-Amplitude Motions in the Vibrational Circular Dichroism Spectrum of Dehydroquinidine. <i>Chemistry - A European Journal</i> , 2016, 22, 704-715.	1.7	20
54	The anharmonic quartic force field infrared spectra of five non-linear polycyclic aromatic hydrocarbons: Benz[a]anthracene, chrysene, phenanthrene, pyrene, and triphenylene. <i>Journal of Chemical Physics</i> , 2016, 145, 084313.	1.2	40

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55	Electron-flux infrared response to varying C–C bond topology in charged aromatic monomers. <i>Nature Communications</i> , 2016, 7, 12633.	5.8	7
56	Energy Transfer between Inorganic Perovskite Nanocrystals. <i>Journal of Physical Chemistry C</i> , 2016, 120, 13310-13315.	1.5	106
57	Direct Observation of a Dark State in the Photocycle of a Light-Driven Molecular Motor. <i>Journal of Physical Chemistry A</i> , 2016, 120, 8606-8612.	1.1	36
58	HIGH-RESOLUTION IR ABSORPTION SPECTROSCOPY OF POLYCYCLIC AROMATIC HYDROCARBONS IN THE 3 $\hat{1}$ / ₄ m REGION: ROLE OF PERIPHERY. <i>Astrophysical Journal</i> , 2016, 831, 58.	1.6	30
59	Water Dissociation upon Adsorption onto Free Iron Clusters Is Size Dependent. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2381-2387.	2.1	15
60	The anharmonic quartic force field infrared spectra of three polycyclic aromatic hydrocarbons: Naphthalene, anthracene, and tetracene. <i>Journal of Chemical Physics</i> , 2015, 143, 224314.	1.2	71
61	HIGH-RESOLUTION IR ABSORPTION SPECTROSCOPY OF POLYCYCLIC AROMATIC HYDROCARBONS: THE REALM OF ANHARMONICITY. <i>Astrophysical Journal</i> , 2015, 814, 23.	1.6	51
62	Interplay between Static and Dynamic Energy Transfer in Biofunctional Upconversion Nanoplatfoms. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 2518-2523.	2.1	39
63	Characterization of Porphyrin-Co(III)- \hat{a} Nitrene Radical \hat{a} ™ Species Relevant in Catalytic Nitrene Transfer Reactions. <i>Journal of the American Chemical Society</i> , 2015, 137, 5468-5479.	6.6	185
64	Ultrafast Excited-State Dynamics of a Cyano-Substituted \hat{a} Proton Sponge \hat{a} . <i>Journal of Physical Chemistry A</i> , 2015, 119, 11233-11240.	1.1	0
65	Water Adsorption on Free Cobalt Cluster Cations. <i>Journal of Physical Chemistry A</i> , 2015, 119, 10828-10837.	1.1	24
66	Elucidating the Structure of Chiral Molecules by using Amplified Vibrational Circular Dichroism: From Theory to Experimental Realization. <i>ChemPhysChem</i> , 2015, 16, 3363-3373.	1.0	17
67	Targeted labeling of an early-stage tumor spheroid in a chorioallantoic membrane model with upconversion nanoparticles. <i>Nanoscale</i> , 2015, 7, 1596-1600.	2.8	11
68	Unraveling the Mechanism of a Reversible Photoactivated Molecular Proton Crane. <i>Journal of Physical Chemistry B</i> , 2014, 118, 12965-12971.	1.2	20
69	Molecular Beam and <i>ab Initio</i> Studies of Photoactive Yellow Protein Chromophores: Influence of the Thioester Functionality and Single Bond Rotation. <i>Journal of Physical Chemistry B</i> , 2014, 118, 12395-12403.	1.2	3
70	Amplified Vibrational Circular Dichroism as a Probe of Local Biomolecular Structure. <i>Journal of the American Chemical Society</i> , 2014, 136, 3530-3535.	6.6	53
71	Excited-State Dynamics of Isolated and Microsolvated Cinnamate-Based UV-B Sunscreens. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 2464-2468.	2.1	91
72	Multispectral upconversion luminescence intensity ratios for ascertaining the tissue imaging depth. <i>Nanoscale</i> , 2014, 6, 9257-9263.	2.8	13

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73	Ultrafast Excited State Dynamics in 9,9- C_2 -Bifluorenylidene. <i>Journal of Physical Chemistry A</i> , 2014, 118, 5961-5968.	1.1	15
74	Water lubricates hydrogen-bonded molecular machines. <i>Nature Chemistry</i> , 2013, 5, 929-934.	6.6	100
75	Critical Landau Velocity in Helium Nanodroplets. <i>Physical Review Letters</i> , 2013, 111, 153002.	2.9	66
76	Excited state dynamics of Photoactive Yellow Protein chromophores elucidated by high-resolution spectroscopy and ab initio calculations. <i>Faraday Discussions</i> , 2013, 163, 321.	1.6	32
77	Conformational Heterogeneity of Methyl 4-Hydroxycinnamate: A Gas-Phase UV-IR Spectroscopic Study. <i>Journal of Physical Chemistry B</i> , 2013, 117, 4798-4805.	1.2	18
78	Non-Equilibrium Isomer Distribution of the Gas-Phase Photoactive Yellow Protein Chromophore. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2259-2263.	2.1	63
79	Time-resolved vibrational spectroscopy of a molecular shuttle. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 1865-1875.	1.3	31
80	Amplification of the linear and nonlinear optical response of a chiral molecular crystal. <i>Journal of Chemical Physics</i> , 2012, 136, 134501.	1.2	18
81	IR Spectroscopy on Jet-Cooled Isolated Two-Station Rotaxanes. <i>Journal of Physical Chemistry A</i> , 2011, 115, 9669-9675.	1.1	32
82	Vibrational and Electronic Spectroscopy of the 4-Hydroxystyrene- CO_2 Cluster and Its Hydrate: A <i>para</i> -Coumaric Acid Impostor. <i>Journal of Physical Chemistry B</i> , 2011, 115, 1275-1281.	1.2	5
83	High-Resolution Spectroscopy of Jet-Cooled 1,1- D_2 -Diphenylethylene: Electronically Excited and Ionic States of a Prototypical Cross-Conjugated System. <i>Journal of Physical Chemistry A</i> , 2011, 115, 9399-9410.	1.1	19
84	Electronic Spectroscopy of Aniline Ions Embedded in Helium Nanodroplets. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 1563-1566.	2.1	28
85	Critical Shell Thickness of Core/Shell Upconversion Luminescence Nanoplatform for FRET Application. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2083-2088.	2.1	124
86	Biphasic Oxidation of Oxy-Hemoglobin in Bloodstains. <i>PLoS ONE</i> , 2011, 6, e21845.	1.1	59
87	Spectroscopy and dynamics of methyl-4-hydroxycinnamate: the influence of isotopic substitution and water complexation. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 4393.	1.3	19
88	Multi-targeting single fiber-optic biosensor based on evanescent wave and quantum dots. <i>Biosensors and Bioelectronics</i> , 2010, 26, 149-154.	5.3	21
89	Red spectral shift and enhanced quantum efficiency in phonon-free photoluminescence from silicon nanocrystals. <i>Nature Nanotechnology</i> , 2010, 5, 878-884.	15.6	294
90	Operation Mechanism of a Molecular Machine Revealed Using Time-Resolved Vibrational Spectroscopy. <i>Science</i> , 2010, 328, 1255-1258.	6.0	95

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91	IR Spectroscopy of Molecular Ions by Nonthermal Ion Ejection from Helium Nanodroplets. <i>Journal of the American Chemical Society</i> , 2010, 132, 14086-14091.	6.6	60
92	Excited State Processes of 2-Butylamino-6-methyl-4-nitropyridine <i>N</i> -oxide in Nonpolar Solvents. A Transient Absorption Spectroscopy Study. <i>Journal of Physical Chemistry A</i> , 2010, 114, 4045-4050.	1.1	7
93	Direct Access to Polyisocyanide Screw Sense Using Vibrational Circular Dichroism. <i>Macromolecules</i> , 2010, 43, 7931-7935.	2.2	37
94	Effect of Surface Related Organic Vibrational Modes in Luminescent Upconversion Dynamics of Rare Earth Ions Doped Nanoparticles. <i>Journal of Nanoscience and Nanotechnology</i> , 2010, 10, 7149-7153.	0.9	6
95	Spectroscopic study of the authentic emitter of AMPPD chemiluminescence in alkaline aqueous solution. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 6789.	1.3	8
96	Absorption spectroscopy of adenine, 9-methyladenine, and 2-aminopurine in helium nanodroplets. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 15600.	1.3	12
97	High-Resolution Excitation and Absorption Spectroscopy of Gas-Phase <i>p</i> -Coumaric Acid: Unveiling an Elusive Chromophore. <i>Journal of the American Chemical Society</i> , 2010, 132, 6315-6317.	6.6	36
98	Interference in acetylene intersystem crossing acts as the molecular analog of Young's double-slit experiment. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009, 106, 2510-2514.	3.3	8
99	Conformational Flexibility of a Rotaxane Thread Probed by Electronic Spectroscopy in Helium Nanodroplets. <i>Journal of the American Chemical Society</i> , 2009, 131, 12902-12903.	6.6	11
100	Multistate Photo-Induced Relaxation and Photoisomerization Ability of Fumaramide Threads: A Computational and Experimental Study. <i>Journal of the American Chemical Society</i> , 2009, 131, 104-117.	6.6	27
101	Stiff, and Sticky in the Right Places: Binding Interactions in Isolated Mechanically Interlocked Molecules Probed by Mid-Infrared Spectroscopy. <i>Journal of the American Chemical Society</i> , 2009, 131, 2428-2429.	6.6	33
102	Two-Dimensional Vibrational Spectroscopy of Rotaxane-Based Molecular Machines. <i>Accounts of Chemical Research</i> , 2009, 42, 1462-1469.	7.6	39
103	Infrared Study of Intercomponent Interactions in a Switchable Hydrogen-Bonded Rotaxane. <i>Chemistry - A European Journal</i> , 2008, 14, 1935-1946.	1.7	45
104	Remedial Mathematics for Quantum Chemistry. <i>Journal of Chemical Education</i> , 2008, 85, 1233.	1.1	5
105	High-Resolution Spectroscopy of Methyl 4-Hydroxycinnamate and Its Hydrogen-Bonded Water Complex. <i>Journal of Physical Chemistry B</i> , 2008, 112, 4427-4434.	1.2	37
106	Tagging multiphoton ionization events by two-dimensional photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , 2007, 126, 204312.	1.2	4
107	Photoelectron studies on vibronic coupling in pyrazine. <i>Journal of Chemical Physics</i> , 2007, 127, 104301.	1.2	6
108	Structure and Photophysics of 2-(2-Pyridyl)benzindoles: The Role of Intermolecular Hydrogen Bonds. <i>Journal of Physical Chemistry A</i> , 2007, 111, 11400-11409.	1.1	22

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109	Heterovibrational Interactions, Cooperative Hydrogen Bonding, and Vibrational Energy Relaxation Pathways in a Rotaxane. <i>Journal of Physical Chemistry C</i> , 2007, 111, 6798-6804.	1.5	15
110	Fluorescent, molecularly imprinted thin-layer films based on a common polymer. <i>Journal of Applied Polymer Science</i> , 2007, 105, 229-235.	1.3	20
111	Vibronic spectra of the lower excited singlet states of styrene: A Time Dependent Density Functional Theory study. <i>Chemical Physics Letters</i> , 2007, 435, 224-229.	1.2	18
112	Femtosecond Studies of Charge-Transfer Mediated Proton Transfer in 2-Butylamino-6-methyl-4-nitropyridineN-Oxide. <i>Journal of Physical Chemistry A</i> , 2006, 110, 7086-7091.	1.1	19
113	Femtosecond Spectroscopic Studies of the One- and Two-Photon Excited-State Dynamics of 2,2,17,17-Tetramethyloctadeca-5,9,13-trien-3,7,11,15-tetrayne: A Trimeric Oligodiacetylene. <i>Journal of Physical Chemistry A</i> , 2006, 110, 11435-11439.	1.1	6
114	A time-dependent density functional study of vibrationally resolved excitation, emission, and ionization spectra of the S1 state of phenol. <i>Chemical Physics Letters</i> , 2006, 420, 459-464.	1.2	13
115	Combined experimental-theoretical study of the lower excited singlet states of paravinyl phenol, an analog of the paracoumaric acid chromophore. <i>Journal of Chemical Physics</i> , 2006, 125, 204303.	1.2	14
116	The asymmetric nature of charge transfer states of the cyano-substituted proton sponge. <i>Chemical Physics Letters</i> , 2005, 401, 189-195.	1.2	8
117	From The Cover: Probing the structure of a rotaxane with two-dimensional infrared spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 13378-13382.	3.3	49
118	Femtosecond Excited State Studies of the Two-Center Three-Electron Bond Driven Twisted Internal Charge Transfer Dynamics in 1,8-Bis(dimethylamino)naphthalene. <i>Journal of Physical Chemistry A</i> , 2005, 109, 3535-3541.	1.1	17
119	Comment on "Gas-Phase Photochemistry of the Photoactive Yellow Protein Chromophoretrans-p-Coumaric Acid". <i>Journal of Physical Chemistry A</i> , 2005, 109, 6135-6136.	1.1	21
120	Ionization Potentials of Fluoroindoles and the Origin of Nonexponential Tryptophan Fluorescence Decay in Proteins. <i>Journal of the American Chemical Society</i> , 2005, 127, 4104-4113.	6.6	85
121	Photophysics of 1,8-Bis(dimethylamino)naphthalene in Solution: Internal Charge Transfer with a Twist. <i>Journal of Physical Chemistry A</i> , 2004, 108, 10623-10631.	1.1	16
122	Rydberg-Valence Interactions in Monoolefins: Dispersing Electronic Properties in 1,1-Bicyclohexylidene. <i>ChemPhysChem</i> , 2003, 4, 97-101.	1.0	2
123	Spectroscopy and dynamics of excited states in maleimide and N-methyl maleimide: Ionic projection and ab initio calculations. <i>Journal of Chemical Physics</i> , 2003, 118, 10944-10955.	1.2	17
124	Vibronic coupling in excited states of acetone. <i>Journal of Chemical Physics</i> , 2002, 116, 547-560.	1.2	25
125	High-Resolution Excited-State Photoelectron Spectroscopy of the Lower Rydberg States of Jet-Cooled C2H4 and C2D4. <i>Journal of Physical Chemistry A</i> , 2002, 106, 3727-3737.	1.1	15
126	Isolated Building Blocks of Photonic Materials: High-Resolution Excited-State Photoelectron Spectroscopy of Jet-Cooled Tetramethylethylene and 1,1-Bicyclohexylidene. <i>Journal of Physical Chemistry A</i> , 2002, 106, 5249-5262.	1.1	13

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127	Isolated Building Blocks of Photonic Materials: High-Resolution Spectroscopy of Excited States of Jet-Cooled Push-Pull Stilbenes. <i>Journal of Physical Chemistry A</i> , 2002, 106, 2446-2456.	1.1	31
128	Excited and ionic states of formamide: An excited-state photoelectron spectroscopy and ab initio study. <i>Journal of Chemical Physics</i> , 2002, 117, 8270-8280.	1.2	20
129	Structure and Photophysics of an Old, New Molecule: 1,3,6,8-Tetraazatricyclo[4.4.1.1.3,8]dodecane. <i>Journal of the American Chemical Society</i> , 2002, 124, 149-158.	6.6	10
130	Fluorescence Excitation Spectroscopy of the 3p Rydberg States of 1-Azabicyclo[2.2.2]octane and 1-Azaadamantane. <i>Journal of Physical Chemistry A</i> , 2000, 104, 729-734.	1.1	4
131	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
132	A Perspective on Symmetry Breaking and Vibronic Coupling in cis-1,3,5-Hexatriene. <i>Journal of the American Chemical Society</i> , 2000, 122, 7418-7419.	6.6	12
133	Scaling of the second hyperpolarisabilities of conjugated carbon systems: polyynes versus polyenes and fullerenes. <i>Chemical Physics Letters</i> , 1999, 313, 426-430.	1.2	16
134	Modeling the Spectroscopy of the Lowest Excited Singlet State of cis,trans-1,3,5,7-Octatetraene: The Role of Symmetry Breaking and Vibronic Interactions. <i>Journal of Physical Chemistry A</i> , 1999, 103, 2220-2226.	1.1	15
135	Vibronic interactions in s-trans-butadiene. <i>Chemical Physics Letters</i> , 1998, 287, 275-281.	1.2	8
136	The large $1\pi^*1\sigma_g^* \rightarrow 2\pi^*1\sigma_g^* \dots C$ and $C\pi^*C$ stretch vibronic interaction in all-trans polyenes. <i>Chemical Physics Letters</i> , 1998, 289, 118-124.	1.2	7
137	Experimental and theoretical studies of the low-lying electronic states of the simplest benzylic amide [2]catenane. <i>Chemical Physics</i> , 1998, 238, 421-428.	0.9	7
138	Resonance enhanced multiphoton ionization photoelectron spectroscopy of gerade excited Rydberg states of the xenon dimer. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , 1998, 97, 147-158.	0.8	12
139	Photoionization and photodissociation dynamics of H ₂ after (3+1) resonance-enhanced multiphoton ionization via the $B^1\Sigma_u^+$ state. <i>Journal of Chemical Physics</i> , 1998, 109, 8319-8329.	1.2	17
140	The spectroscopy of high Rydberg states of ammonia. <i>Journal of Chemical Physics</i> , 1998, 108, 6667-6680.	1.2	38
141	Resonance-enhanced multiphoton ionization photoelectron spectroscopy of Rydberg states of N ₂ O below the $X^2\Sigma^+$ ionization limit. <i>Journal of Chemical Physics</i> , 1998, 109, 7844-7850.	1.2	18
142	Two Ground State Conformers of the Proton Sponge 1,8-Bis(dimethylamino)naphthalene Revealed by Fluorescence Spectroscopy and ab Initio Calculations. <i>Journal of the American Chemical Society</i> , 1998, 120, 4840-4844.	6.6	49
143	Resonance-enhanced multiphoton-ionization photoelectron spectroscopy of even-parity autoionizing Rydberg states of atomic sulphur. <i>Journal of Chemical Physics</i> , 1997, 106, 6831-6838.	1.2	8
144	Rotationally resolved multiphoton ionization photoelectron spectroscopy of the $[a^1\pi^*]3d^1\epsilon^*2\pi^*$ and $[a^1\pi^*]5p^1\epsilon^*2\pi^*$ Rydberg states of the SH radical. <i>Journal of Chemical Physics</i> , 1997, 107, 2782-2792.	1.2	9

#	ARTICLE	IF	CITATIONS
145	The Lowest Excited Singlet States of 1-Azaadamantane and 1-Azabicyclo[2.2.2]octane: Fluorescence Excitation Spectroscopy and Density Functional Calculations. <i>Journal of the American Chemical Society</i> , 1997, 119, 11523-11533.	6.6	21
146	Evidence of Stringlike Behavior in all-trans-Octatetraene. <i>Journal of the American Chemical Society</i> , 1996, 118, 9178-9179.	6.6	7
147	Resonance-enhanced multiphoton-ionization photoelectron spectroscopy of even-parity Rydberg states of atomic sulfur. <i>Physical Review A</i> , 1996, 54, 5126-5132.	1.0	8
148	Dynamics of high- n Rydberg states employed in zero kinetic energy-pulsed field ionization spectroscopy via the $F\epsilon\%1\%2$, $D\epsilon\%1\%$, and $f\epsilon\%3\%2$ Rydberg states of HCl. <i>Journal of Chemical Physics</i> , 1996, 105, 5702-5710.	1.2	12
149	(3 + 1) resonance enhanced multiphoton ionization-photoelectron spectroscopy on the E, F, and G Rydberg states of ClO. <i>Chemical Physics Letters</i> , 1996, 259, 213-218.	1.2	15
150	Resonance enhanced multiphoton ionisation (REMPI) and REMPI-photoelectron spectroscopy of carbonyl sulphide and carbon disulphide. <i>International Journal of Mass Spectrometry and Ion Processes</i> , 1996, 159, 1-11.	1.9	14
151	Resonance enhanced multiphoton ionization spectroscopy of carbonyl sulphide. <i>Journal of Chemical Physics</i> , 1996, 105, 2141-2152.	1.2	34
152	Zero-kinetic-energy pulsed-field ionization spectroscopy of the $a\epsilon\%1\%$ state of SH ⁺ (SD ⁺). <i>Journal of Chemical Physics</i> , 1996, 104, 521-527.	1.2	23
153	Zero kinetic energy-pulsed field ionization and resonance enhanced multiphoton ionization photoelectron spectroscopy: Ionization dynamics of Rydberg states in HBr. <i>Journal of Chemical Physics</i> , 1996, 104, 4911-4919.	1.2	35
154	Characterization of spin-orbit autoionizing Rydberg states excited via one-photon absorption from the $F\epsilon\%1\%2$ Rydberg state of HBr. <i>Journal of Chemical Physics</i> , 1996, 105, 2978-2991.	1.2	13
155	Two-photon resonance enhanced multiphoton ionization photoelectron spectroscopy of the SH (SD) radical below and above the lowest ionization threshold. <i>Journal of Chemical Physics</i> , 1996, 105, 6688-6712.	1.2	16
156	Resonance enhanced multiphoton ionization spectroscopy of carbon disulphide. <i>Journal of Chemical Physics</i> , 1996, 104, 6117-6129.	1.2	48
157	Two-photon resonance enhanced MPI-PES above the lowest ionization threshold. Observation of the $[a\epsilon\%1\%2]$ state of the SH (SD) radical. <i>Chemical Physics Letters</i> , 1995, 239, 326-331.	1.2	16
158	Rotationally resolved photoelectron spectroscopy of the $[a\epsilon\%1\%3]$ Rydberg state of the SH radical. <i>Journal of Chemical Physics</i> , 1995, 103, 3262-3264.	1.2	14
159	Four-photon excitation of autoionizing states of Ar, Kr, and Xe between the $2P_{3/2}$ and $2P_{1/2}$ ionic limits. <i>Physical Review A</i> , 1995, 51, 1097-1109.	1.0	43
160	A resonance enhanced multiphoton ionization study of the gerade excited states of Xe ₂ with a Xe $\epsilon\%1\%0$ +Xe* $\epsilon\%6s[3/2]$ 1 dissociation limit. <i>Journal of Chemical Physics</i> , 1995, 102, 4020-4026.	1.2	16
161	Vibronic activity in trans,trans-1,3,5,7 octatetraene: The S $\epsilon\%1\%$ S $\epsilon\%1\%$ spectrum. <i>Journal of Chemical Physics</i> , 1995, 103, 10492-10501.	1.2	29
162	Resonance-enhanced multiphoton ionisation spectroscopy of methanethiol. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1995, 91, 2715.	1.7	13

#	ARTICLE	IF	CITATIONS
163	Resonance enhanced multiphoton ionisation spectroscopy of dimethyl sulfide. Journal of the Chemical Society, Faraday Transactions, 1995, 91, 3339.	1.7	20
164	Three-photon excitation of autoionizing states of Ar, Kr, and Xe between the $2P_{3/2}$ and $2P_{1/2}$ ionic limits. Physical Review A, 1994, 49, 3322-3332.	1.0	42
165	(3+1) resonance enhanced multiphoton ionization photoelectron spectroscopy on nf Rydberg states of carbon dioxide. Journal of Chemical Physics, 1994, 101, 9303-9325.	1.2	20
166	One- and two-color two-photon resonance enhanced multiphoton ionization spectroscopy of the $d^{\infty}1^{\infty}+$ state of NH. Journal of Chemical Physics, 1994, 100, 7984-7994.	1.2	13
167	Resonance-enhanced multiphoton ionisation spectroscopy of thiirane. Journal of the Chemical Society, Faraday Transactions, 1994, 90, 3591.	1.7	11
168	A (3+1) resonance enhanced multiphoton ionization study of the $C^{\infty}1^{\infty}+$ and $E^{\infty}1^{\infty}$ states of CO: Polarization dependence used to probe electronic excitation routes and electronic character. Journal of Chemical Physics, 1993, 99, 5061-5070.	1.2	11
169	Resonance enhanced multiphoton ionization photoelectron spectroscopy on nanosecond and picosecond time scales of Rydberg states of methyl iodide. Journal of Chemical Physics, 1993, 99, 836-853.	1.2	60
170	Resonance enhanced multiphoton ionization photoelectron spectroscopy and pulsed field ionization via the $F^{\infty}1^{\infty}2(v^{\infty}=0)$ and $f^{\infty}3^{\infty}2(v^{\infty}=0)$ Rydberg states of HCl. Journal of Chemical Physics, 1993, 99, 3252-3261.	1.2	46
171	The lowest excited singlet state of isolated 1-phenyl-1,3-butadiene and 1-phenyl-1,3,5-hexatriene. Journal of Chemical Physics, 1992, 96, 4860-4868.	1.2	13
172	The $2^{\infty}1A_g$ state of isolated cis,trans-1,3,5,7-octatetraene: Two-color resonance enhanced two-photon ionization studies. Journal of Chemical Physics, 1992, 96, 399-407.	1.2	24
173	Ab initio calculations on vibronic coupling in the lower triplet states of pyrimidine. Journal of the American Chemical Society, 1992, 114, 9544-9551.	6.6	13
174	Lowest energy excited singlet state of isolated cis-hexatriene. Journal of Chemical Physics, 1991, 94, 6367-6376.	1.2	76
175	Lowest energy excited singlet states of isomers of alkyl substituted hexatrienes. Journal of Chemical Physics, 1991, 94, 4691-4698.	1.2	32
176	Conformational instability of the lowest triplet state of the benzene nucleus. I. The unsubstituted molecule. Journal of Chemical Physics, 1990, 93, 3733-3745.	1.2	27
177	Conformational instability of the lowest triplet state of the benzene nucleus. II. p-Xylene, the influence of substituents. Journal of Chemical Physics, 1990, 93, 3746-3751.	1.2	9
178	Ab initio calculations on the structure of pyridine in its lowest triplet state. Journal of the American Chemical Society, 1990, 112, 5447-5451.	6.6	19
179	Location of the $2^{\infty}1A_g$ state in hexatriene. Journal of Chemical Physics, 1990, 92, 4622-4623.	1.2	63
180	An electron spin-echo envelope modulation study of the lowest triplet state of pyridine-d ₅ : Spin-density distribution and structure. Journal of Chemical Physics, 1989, 91, 6549-6565.	1.2	35

#	ARTICLE	IF	CITATIONS
181	Conformational instability of the lowest triplet state of benzene: the result of ab initio calculations. Journal of the American Chemical Society, 1989, 111, 86-87.	6.6	13
182	Electron spin echo spectroscopy of the photo-excited triplet state of pyridine. Journal of Molecular Structure, 1988, 173, 249-252.	1.8	1
183	The reaction center triplet state of the photosynthetic bacterium Rhodobacter sphaeroides R26: Electron spin echo spectroscopy of a single crystal. Chemical Physics Letters, 1987, 142, 231-236.	1.2	7
184	Pseudo-jahn-teller distortion of pyridine in its lowest triplet state. Chemical Physics Letters, 1986, 127, 189-192.	1.2	18
185	The triplet state of pyridine: A magnetic resonance study using electron-spin-echo spectroscopy. Chemical Physics Letters, 1985, 117, 203-207.	1.2	26