

Jacek Waluk

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/356804/publications.pdf>

Version: 2024-02-01

285
papers

7,367
citations

46918

47
h-index

98622

67
g-index

303
all docs

303
docs citations

303
times ranked

5480
citing authors

#	ARTICLE	IF	CITATIONS
1	Influence of bulky substituents on single-molecule SERS sensitivity. <i>Journal of Chemical Physics</i> , 2022, 156, 014201.	1.2	4
2	Solving the Puzzle of Unusual Excited-State Proton Transfer in 2,5-Bis(6-methyl-2-benzoxazolyl)phenol. <i>Journal of Physical Chemistry A</i> , 2022, 126, 1823-1836.	1.1	1
3	Phosphorescence and Photophysical Parameters of Porphycene in Cryogenic Matrices. <i>Photochem</i> , 2022, 2, 217-224.	1.3	0
4	Energy Relaxation of Porphycene in Atomic and Molecular Cryogenic Matrices. <i>Photochem</i> , 2022, 2, 299-307.	1.3	0
5	Scouting for strong light-matter coupling signatures in Raman spectra. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 16837-16846.	1.3	14
6	Synthesis and Photostability of Cyclooctatetraene-Substituted Free Base Porphyrins. <i>Chemistry</i> , 2021, 3, 104-115.	0.9	2
7	Controlling Emissive Properties by Intramolecular Hydrogen Bonds: Alkyl and Aryl meso-substituted Porphycenes. <i>Chemistry - A European Journal</i> , 2021, 27, 6324-6333.	1.7	8
8	Substituent screening effect on single-molecule photostability: comparison of three differently substituted porphycenes. <i>Methods and Applications in Fluorescence</i> , 2021, 9, 035004.	1.1	2
9	Tailoring Tautomerization of Single Phthalocyanine Molecules through Modification of Chromophore Photophysics by the Purcell Effect of an Optical Microcavity. <i>Journal of Physical Chemistry C</i> , 2021, 125, 14932-14939.	1.5	3
10	Spectral and photophysical modifications of porphyrins attached to core-shell nanoparticles. Theory and experiment. <i>Methods and Applications in Fluorescence</i> , 2021, 9, 045003.	1.1	3
11	SERS-based sensor for the detection of sexually transmitted pathogens in the male swab specimens: A new approach for clinical diagnosis. <i>Biosensors and Bioelectronics</i> , 2021, 189, 113358.	5.3	17
12	Photoinduced and ground state conversions in a cyclic $\hat{2}$ -thioxoketone. <i>RSC Advances</i> , 2021, 12, 681-689.	1.7	2
13	Influence of local microenvironment on the double hydrogen transfer in porphycene. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 17117-17128.	1.3	4
14	Magnetic Circular Dichroism of <i>meso</i> -Phenyl-Substituted Pd-Octaethylporphyrins. <i>Journal of Physical Chemistry A</i> , 2020, 124, 8144-8158.	1.1	6
15	Fluorinated Porphycenes: Synthesis, Spectroscopy, Photophysics, and Tautomerism. <i>ChemPlusChem</i> , 2020, 85, 2197-2206.	1.3	5
16	2 + 2 Can Make Nearly a Thousand! Comparison of Di- and Tetra- <i>Meso</i> -Alkyl-Substituted Porphycenes. <i>Journal of Physical Chemistry A</i> , 2020, 124, 4594-4604.	1.1	11
17	Matrix isolation studies of vibrational structure of hemiporphycene. <i>Journal of Molecular Structure</i> , 2020, 1218, 128497.	1.8	0
18	Photoeradication of bacteria with porphycenes: Substituent effects on the efficiency. <i>European Journal of Medicinal Chemistry</i> , 2020, 200, 112472.	2.6	6

#	ARTICLE	IF	CITATIONS
19	Multimode Vibrational Strong Coupling of Methyl Salicylate to a Fabry-Pérot Microcavity. <i>Journal of Physical Chemistry B</i> , 2020, 124, 5709-5716.	1.2	19
20	Towards More Photostable, Brighter, and Less Phototoxic Chromophores: Synthesis and Properties of Porphyrins Functionalized with Cyclooctatetraene. <i>Chemistry - A European Journal</i> , 2020, 26, 16666-16675.	1.7	9
21	Porphycene Protonation: A Fast and Reversible Reaction Enabling Optical Transduction for Acid Sensing. <i>ChemPhotoChem</i> , 2020, 4, 5264-5270.	1.5	4
22	Fluorescence quantum yield determination using simultaneous double-beam absorption measurement. <i>Measurement: Journal of the International Measurement Confederation</i> , 2020, 165, 108159.	2.5	4
23	Triarylisocyanurate-Based Fluorescent Two-Photon Absorbers. <i>ChemPlusChem</i> , 2020, 85, 411-425.	1.3	5
24	Nanoplasmonic sensor for foodborne pathogens detection. Towards development of ISO-SERS methodology for taxonomic affiliation of <i>Campylobacter</i> spp.. <i>Journal of Biophotonics</i> , 2020, 13, e201960227.	1.1	12
25	Porphycene Films Grown on Highly Oriented Pyrolytic Graphite: Unveiling Structure-Property Relationship through Combined Reflectance Anisotropy Spectroscopy and Atomic Force Microscopy Investigations. <i>Proceedings (mdpi)</i> , 2020, 56, 44.	0.2	1
26	Spectrofluorometer: the excitation beam intensity calibration using a single standard solution. <i>Methods and Applications in Fluorescence</i> , 2020, 8, 047001.	1.1	0
27	Frontispiece: Towards More Photostable, Brighter, and Less Phototoxic Chromophores: Synthesis and Properties of Porphyrins Functionalized with Cyclooctatetraene. <i>Chemistry - A European Journal</i> , 2020, 26, .	1.7	0
28	Fluorescence studies of porphycene in various cryogenic environments. <i>Low Temperature Physics</i> , 2019, 45, 656-662.	0.2	2
29	Goodbye to Quinine in Sulfuric Acid Solutions as a Fluorescence Quantum Yield Standard. <i>Analytical Chemistry</i> , 2019, 91, 5389-5394.	3.2	19
30	Photoinduced oxidation of an indole derivative: 2-(1H-indol-2-yl)-[1,5]naphthyridine. <i>Photochemical and Photobiological Sciences</i> , 2019, 18, 2225-2231.	1.6	4
31	Near-Field Spectral Response of Optically Excited Scanning Tunneling Microscope Junctions Probed by Single-Molecule Action Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 2068-2074.	2.1	11
32	Antiaromatic or Nonaromatic? ¹ H-, ⁶ H-, ^{2,6} H-, ^{2,5} -Dipyrrolo-1,5(2,6)-dipyridinacyclooctaphane-3,7-diene: a Porphycene Derivative with 4 N Electrons. <i>Journal of Physical Chemistry A</i> , 2019, 123, 2727-2733.	1.1	5
33	Nature of Large Temporal Fluctuations of Hydrogen Transfer Rates in Single Molecules. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1211-1215.	2.1	20
34	Quantum tunneling in real space: Tautomerization of single porphycene molecules on the (111) surface of Cu, Ag, and Au. <i>Journal of Chemical Physics</i> , 2018, 148, 102330.	1.2	29
35	Combined effect of hydrogen bonding interactions and freezing of rotameric equilibrium on the enhancement of photostability. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 13306-13315.	1.3	12
36	Two Macrocycles in One Shot: Synthesis, Spectroscopy, Photophysics, and Tautomerism of 23-Oxahemiporphycene and 21-Oxacorrole-5-carbaldehyde. <i>Chemistry - A European Journal</i> , 2018, 24, 9884-9891.	1.7	3

#	ARTICLE	IF	CITATIONS
37	Near-Field Enhanced Photochemistry of Single Molecules in a Scanning Tunneling Microscope Junction. <i>Nano Letters</i> , 2018, 18, 152-157.	4.5	32
38	Anharmonicity in a double hydrogen transfer reaction studied in a single porphycene molecule on a Cu(110) surface. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 12112-12119.	1.3	3
39	Unusual effects in single molecule tautomerization: hemiporphycene. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 26591-26596.	1.3	5
40	Supersonic jet spectroscopy of parent hemiporphycene: Structural assignment and vibrational analysis for S0 and S1 electronic states. <i>Journal of Chemical Physics</i> , 2018, 149, 134307.	1.2	3
41	Versatile Approach for Reliable Determination of Both High and Low Values of Luminescence Quantum Yields. <i>Analytical Chemistry</i> , 2018, 90, 10139-10143.	3.2	11
42	Spectroscopic and microscopic investigations of tautomerization in porphycenes: condensed phases, supersonic jets, and single molecule studies. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 4921-4937.	1.3	24
43	Non-typical fluorescence studies of excited and ground state proton and hydrogen transfer. <i>Methods and Applications in Fluorescence</i> , 2017, 5, 014007.	1.1	3
44	New class of easily-synthesizable and modifiable organic materials for applications in luminescent devices. <i>Dyes and Pigments</i> , 2017, 138, 267-277.	2.0	13
45	Plasmon-Mediated Surface Engineering of Silver Nanowires for Surface-Enhanced Raman Scattering. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 2774-2779.	2.1	38
46	Surface-enhanced Raman spectroscopy introduced into the International Standard Organization (ISO) regulations as an alternative method for detection and identification of pathogens in the food industry. <i>Analytical and Bioanalytical Chemistry</i> , 2017, 409, 1555-1567.	1.9	49
47	Direct Observation of Double Hydrogen Transfer via Quantum Tunneling in a Single Porphycene Molecule on a Ag(110) Surface. <i>Journal of the American Chemical Society</i> , 2017, 139, 12681-12687.	6.6	49
48	SERS-based Immunoassay in a Microfluidic System for the Multiplexed Recognition of Interleukins from Blood Plasma: Towards Picogram Detection. <i>Scientific Reports</i> , 2017, 7, 10656.	1.6	75
49	Antimicrobial photodynamic therapy by means of porphycene photosensitizers. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2017, 174, 84-89.	1.7	29
50	Improved Method of Fluorescence Quantum Yield Determination. <i>Analytical Chemistry</i> , 2017, 89, 8650-8655.	3.2	38
51	Detection of a weak ring current in a nonaromatic porphyrin nanoring using magnetic circular dichroism. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 32556-32565.	1.3	8
52	Polymer mat prepared via Forcespinning [®] as a SERS platform for immobilization and detection of bacteria from blood plasma. <i>Materials Science and Engineering C</i> , 2017, 71, 345-350.	3.8	28
53	Spectroscopy and Tautomerization Studies of Porphycenes. <i>Chemical Reviews</i> , 2017, 117, 2447-2480.	23.0	130
54	Resonance Raman spectroscopy study of protonated porphyrin. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2017, 173, 350-355.	2.0	14

#	ARTICLE	IF	CITATIONS
55	Linear Dichroism, Instrumentation. , 2017, , 601-603.		0
56	Linear Dichroism, Applications. , 2017, , 595-600.		3
57	Force-induced tautomerization in a single molecule. Nature Chemistry, 2016, 8, 935-940.	6.6	111
58	Detection and identification of human fungal pathogens using surface-enhanced Raman spectroscopy and principal component analysis. Analytical Methods, 2016, 8, 8427-8434.	1.3	47
59	Highly efficient SERS-based detection of cerebrospinal fluid neopterin as a diagnostic marker of bacterial infection. Analytical and Bioanalytical Chemistry, 2016, 408, 4319-4327.	1.9	28
60	Non-classical effects in proton or hydrogen transfer. Pure and Applied Chemistry, 2016, 88, 1063-1071.	0.9	1
61	Parent, Unsubstituted Hemiporphycene: Synthesis and Properties. Chemistry - A European Journal, 2016, 22, 17311-17320.	1.7	20
62	Rapid detection and identification of bacterial meningitis pathogens in ex vivo clinical samples by SERS method and principal component analysis. Analytical Methods, 2016, 8, 4521-4529.	1.3	38
63	Near Infrared Phosphorescent, Non-oxidizable Palladium and Platinum Perfluoro-phthalocyanines. ChemPhysChem, 2016, 17, 1123-1135.	1.0	12
64	ABO blood groups' antigen-antibody interactions studied using SERS spectroscopy: towards blood typing. Analytical Methods, 2016, 8, 1463-1472.	1.3	13
65	Structure, NMR and Electronic Spectra of [<i>m.n</i>]Paracyclophanes with Varying Bridges Lengths (<i>m, n</i> = 2-4). Journal of Physical Chemistry A, 2016, 120, 724-736.	1.1	10
66	Direct Observation of Photoinduced Tautomerization in Single Molecules at a Metal Surface. Nano Letters, 2016, 16, 1034-1041.	4.5	67
67	Influence of alkyl substituents in corphycene on geometry, electronic structure, hydrogen bonding, and tautomerization. Journal of Porphyrins and Phthalocyanines, 2016, 20, 367-377.	0.4	5
68	Simulations of fluorescence enhancement and emission profile changes in porphyrin attached to gold-silica core-shell nanoparticles. Methods and Applications in Fluorescence, 2016, 4, 014002.	1.1	6
69	Single molecule Raman spectra of porphycene isotopologues. Nanoscale, 2016, 8, 3337-3349.	2.8	25
70	Evidence for Dominant Role of Tunneling in Condensed Phases and at High Temperatures: Double Hydrogen Transfer in Porphycenes. Journal of Physical Chemistry Letters, 2016, 7, 283-288.	2.1	43
71	ZnO oxide films for ultrasensitive, rapid, and label-free detection of neopterin by surface-enhanced Raman spectroscopy. Analyst, The, 2015, 140, 5090-5098.	1.7	12
72	Tailored gold nanostructure arrays as catalysts for oxygen reduction in alkaline media and a single molecule SERS platform. Nanoscale, 2015, 7, 10767-10774.	2.8	15

#	ARTICLE	IF	CITATIONS
73	Hot Carrier-Induced Tautomerization within a Single Porphycene Molecule on Cu(111). ACS Nano, 2015, 9, 7287-7295.	7.3	72
74	Solvent-Controlled Excited State Relaxation Path of 4-Acetyl-4-(dimethylamino)biphenyl. Journal of Physical Chemistry B, 2015, 119, 7294-7307.	1.2	5
75	Solvent-Induced Changes in Photophysics and Photostability of Indole-Naphthyridines. Journal of Physical Chemistry B, 2015, 119, 7283-7293.	1.2	13
76	Tautomerism in Porphycenes: Analysis of Rate-Affecting Factors. Journal of Physical Chemistry B, 2015, 119, 2292-2301.	1.2	40
77	In Search for the Best Environment for Single Molecule Studies: Photostability of Single Terrylendiimide Molecules in Various Polymer Matrices. Journal of Physical Chemistry Letters, 2015, 6, 2477-2482.	2.1	14
78	Detection of Hepatitis B virus antigen from human blood: SERS immunoassay in a microfluidic system. Biosensors and Bioelectronics, 2015, 66, 461-467.	5.3	132
79	Substituent and Solvent Effects on the Excited State Deactivation Channels in Anils and Boranils. Chemistry - A European Journal, 2015, 21, 1312-1327.	1.7	45
80	Towards improved precision in the quantification of surface-enhanced Raman scattering (SERS) enhancement factors: a renewed approach. Analyst, The, 2015, 140, 489-496.	1.7	13
81	Spectroscopic Study of Jet-Cooled Deuterated Porphycenes: Unusual Isotopic Effects on Proton Tunneling. Journal of Physical Chemistry B, 2015, 119, 2193-2203.	1.2	28
82	Electrochemical pathway for the quantification of SERS enhancement factor. Electrochemistry Communications, 2014, 49, 103-106.	2.3	5
83	Enhancing fluorescence by using pluronic block copolymers as carriers of monomeric porphycenes. Methods and Applications in Fluorescence, 2014, 2, 024003.	1.1	6
84	Controlling intramolecular hydrogen transfer in a porphycene molecule with single atoms or molecules located nearby. Nature Chemistry, 2014, 6, 41-46.	6.6	204
85	Resonance Raman and FTIR spectra of Mg-porphyrazines. Journal of Molecular Structure, 2014, 1058, 197-204.	1.8	4
86	Michael Kasha: From Photochemistry and Flowers to Spectroscopy and Music. Angewandte Chemie - International Edition, 2014, 53, 14316-14324.	7.2	30
87	Arresting consecutive steps of a photochromic reaction: studies of β^2 -thioxoketones combining laser photolysis with NMR detection. Physical Chemistry Chemical Physics, 2014, 16, 9128-9137.	1.3	11
88	Electrospun polymer mat as a SERS platform for the immobilization and detection of bacteria from fluids. Analyst, The, 2014, 139, 5061-5064.	1.7	41
89	A new algorithm for identification of components in a mixture: application to Raman spectra of solid amino acids. Analyst, The, 2014, 139, 5755-5764.	1.7	3
90	Synthesis, spectroscopy, and photophysics of porphyrins attached to gold nanoparticles via one or two linkers. Journal of Porphyrins and Phthalocyanines, 2014, 18, 686-697.	0.4	3

#	ARTICLE	IF	CITATIONS
91	Structure, Electronic States, and Anion-Binding Properties of Cyclo[4]naphthobipyrroles. <i>Journal of Physical Chemistry A</i> , 2014, 118, 1038-1046.	1.1	14
92	Nanostructured silver-gold bimetallic SERS substrates for selective identification of bacteria in human blood. <i>Analyst</i> , 2014, 139, 1037.	1.7	110
93	Raman Spectra of Solid Amino Acids: Spectral Correlation Analysis as the First Step Towards Identification by Raman Spectroscopy. <i>Challenges and Advances in Computational Chemistry and Physics</i> , 2014, , 329-354.	0.6	5
94	7-Hydroxyquinoline-8-carbaldehydes. 1. Ground- and Excited-State Long-Range Prototropic Tautomerization. <i>Journal of Physical Chemistry A</i> , 2013, 117, 9127-9146.	1.1	31
95	7-Hydroxyquinoline-8-carbaldehydes. 2. Prototropic Equilibria. <i>Journal of Physical Chemistry A</i> , 2013, 117, 9147-9155.	1.1	11
96	Arresting Tautomerization in a Single Molecule by the Surrounding Polymer: 2,7,12,17-Tetraphenyl Porphycene. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 3967-3971.	2.1	25
97	Thermally and Vibrationally Induced Tautomerization of Single Porphycene Molecules on a Cu(110) Surface. <i>Physical Review Letters</i> , 2013, 111, 246101.	2.9	93
98	Double Hydrogen Transfer in Low Symmetry Porphycenes. <i>Zeitschrift Fur Physikalische Chemie</i> , 2013, 227, 1009-1020.	1.4	7
99	Vibrations of porphycene in the S and S1 electronic states: Single vibronic level dispersed fluorescence study in a supersonic jet. <i>Journal of Chemical Physics</i> , 2013, 138, 174201.	1.2	25
100	Spectroscopy and Photophysics of Bifunctional Proton Donor-Acceptor Indole Derivatives. <i>Journal of Physical Chemistry A</i> , 2013, 117, 4898-4906.	1.1	3
101	Polymorphism, Hydrogen Bond Properties, and Vibrational Structure of 1H-Pyrrolo[3,2-h]Quinoline Dimers. <i>Journal of Atomic, Molecular, and Optical Physics</i> , 2012, 2012, 1-11.	0.5	3
102	Porphycenes. , 2012, , 809-829.		1
103	Vibrations and hydrogen bonding in porphycene. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 5489.	1.3	41
104	1 <i>H</i> -Pyrrolo[3,2- <i>h</i>]quinoline: A Benchmark Molecule for Reliable Calculations of Vibrational Frequencies, IR Intensities, and Raman Activities. <i>Journal of Physical Chemistry A</i> , 2012, 116, 11973-11986.	1.1	13
105	Tautomerization in 2,7,12,17-Tetraphenylporphycene and 9-Amino-2,7,12,17-tetraphenylporphycene: Influence of Asymmetry on the Direction of the Transition Moment. <i>Chemistry - A European Journal</i> , 2012, 18, 13160-13167.	1.7	23
106	Three Modes of Proton Transfer in One Chromophore: Photoinduced Tautomerization in 1 <i>H</i> -Pyrrolo[5 <i>a</i>]pyridines, Their Dimers and Alcohol Complexes. <i>ChemPhysChem</i> , 2012, 13, 3661-3671.	1.0	25
107	Electronic states of cyclophanes with small bridges. <i>Journal of Chemical Physics</i> , 2012, 136, 074201.	1.2	8
108	The long and winding road to new porphycenes. <i>Journal of Porphyrins and Phthalocyanines</i> , 2012, 16, 589-602.	0.4	28

#	ARTICLE	IF	CITATIONS
109	The dynamics and origin of the unrelaxed fluorescence of free-base tetraphenylporphyrin. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2012, 234, 100-106.	2.0	23
110	Highly reproducible, stable and multiply regenerated surface-enhanced Raman scattering substrate for biomedical applications. <i>Journal of Materials Chemistry</i> , 2011, 21, 8662.	6.7	65
111	Excited-State Proton Transfer in <i>syn</i> -2-(2-Pyridyl)pyrrole Occurs on the Nanosecond Time Scale in the Gas Phase. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2114-2117.	2.1	10
112	1,4-Bis(1,3-dioxo-2-indenylidene)cyclohexane: polymorphism, gas phase oxidation and enol form mediated radical formation in the solid state. <i>CrystEngComm</i> , 2011, 13, 3170-3174.	1.3	4
113	Photoluminescence Properties of Porous Silicon with CdSe/ZnS Quantum Dots. <i>Medziagotyra</i> , 2011, 17, .	0.1	0
114	Ground and Excited State Double Hydrogen Transfer in Symmetric and Asymmetric Potentials: Comparison of 2,7,12,17-Tetra β -n-propylporphycene with 9 β -Acetoxy-2,7,12,17-tetra β -n-propylporphycene. <i>Chemistry - A European Journal</i> , 2011, 17, 3672-3678.	1.7	32
115	Bridging the Gap between Porphyrins and Porphycenes: Substituent-Position-Sensitive Tautomerism and Photophysics in <i>meso</i> -Diphenyloctaethylporphyrins. <i>Chemistry - A European Journal</i> , 2011, 17, 10039-10049.	1.7	18
116	Structure, vibrations, and hydrogen bond parameters of dibenzotetraaza[14]annulene. <i>Journal of Molecular Structure</i> , 2010, 976, 215-225.	1.8	10
117	SERS Active Surface Based on Au-Coated Porous GaN. , 2010, , .		1
118	On the origin of fluorescence quenching of pyridylindoles by hydroxylic solvents. <i>Photochemical and Photobiological Sciences</i> , 2010, 9, 923-930.	1.6	20
119	Unusual, Solvent Viscosity-Controlled Tautomerism and Photophysics: <i>Meso</i> -Alkylated Porphycenes. <i>Journal of the American Chemical Society</i> , 2010, 132, 13472-13485.	6.6	63
120	Car π -Parrinello Molecular Dynamics Study of the Intramolecular Vibrational Mode-Sensitive Double Proton-Transfer Mechanisms in Porphycene. <i>Journal of Physical Chemistry A</i> , 2010, 114, 2313-2318.	1.1	28
121	Structure and Hydrogen-Bond Vibrations of Water Complexes of Azaaromatic Compounds: 7-(3-Pyridyl)indole. <i>Journal of Physical Chemistry A</i> , 2010, 114, 3270-3279.	1.1	16
122	Partitioning and Localization of Environment-Sensitive 2-(2-Pyridyl)- and 2-(2-Pyrimidyl)-Indoles in Lipid Membranes: A Joint Refinement Using Fluorescence Measurements and Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2010, 114, 13574-13584.	1.2	36
123	From the Gas Phase to a Lipid Membrane Environment: DFT and MD Simulations of Structure and Dynamics of Hydrogen-Bonded Solvates of Bifunctional Heteroazaaromatic Compounds. <i>Challenges and Advances in Computational Chemistry and Physics</i> , 2010, , 35-75.	0.6	2
124	Ground- and Excited-State Tautomerization Rates in Porphycenes. <i>Chemistry - A European Journal</i> , 2009, 15, 4851-4856.	1.7	60
125	Mode-Selective Promotion and Isotope Effects of Concerted Double-Hydrogen Tunneling in Porphycene Embedded in Superfluid Helium Nanodroplets. <i>ChemPhysChem</i> , 2009, 10, 761-765.	1.0	50
126	On the Origin of Radiationless Transitions in Porphycenes. <i>Journal of Physical Chemistry A</i> , 2009, 113, 7714-7716.	1.1	40

#	ARTICLE	IF	CITATIONS
127	Tautomerization in Condensed Phases and in Isolated Molecules. <i>Israel Journal of Chemistry</i> , 2009, 49, 175-185.	1.0	6
128	Polarized Spectroscopy Studies of Single Molecules of Porphycenes: Tautomerism and Orientation. <i>Journal of Physical Chemistry C</i> , 2009, 113, 11514-11519.	1.5	45
129	Proton transfer with a twist? Femtosecond Dynamics of 2-(2-pyridyl)indole in Condensed Phase and in Supersonic Jets. <i>Angewandte Chemie - International Edition</i> , 2008, 47, 6037-6040.	7.2	54
130	Distribution and favorable binding sites of pyrroloquinoline and its analogues in a lipid bilayer studied by molecular dynamics simulations. <i>Biophysical Chemistry</i> , 2008, 136, 128-135.	1.5	20
131	Electronic and Vibrational Relaxation of Porphycene in Solution. <i>Journal of Physical Chemistry A</i> , 2008, 112, 10753-10757.	1.1	19
132	Separation of Different Hydrogen-Bonded Clusters by Femtosecond UV-Ionization-Detected Infrared Spectroscopy: 1H-Pyrrolo[3,2-h]quinoline·(H ₂ O) _{n=1,2} Complexes. <i>Journal of Physical Chemistry A</i> , 2008, 112, 1150-1156.	1.1	35
133	Unusually Slow Intermolecular Proton-Deuteron Exchange in Porphycene. <i>Zeitschrift Fur Physikalische Chemie</i> , 2008, 222, 1165-1173.	1.4	26
134	Efficient synthesis of porphycene. <i>Journal of Porphyrins and Phthalocyanines</i> , 2007, 11, 596-600.	0.4	31
135	Matrix isolation spectroscopy and molecular dynamics simulations for 2,7,12,17-tetra-tert-butylporphycene in argon and xenon. <i>Journal of Chemical Physics</i> , 2007, 127, 134501.	1.2	4
136	Mode-Selective Excited-State Proton Transfer in 2-(2-pyridyl)pyrrole Isolated in a Supersonic Jet. <i>Journal of the American Chemical Society</i> , 2007, 129, 2738-2739.	6.6	61
137	Fluorescence quenching in cyclic hydrogen-bonded complexes of 1H-pyrrolo[3,2-h]quinoline with methanol: cluster size effect. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 3276.	1.3	34
138	Vibrational Gating of Double Hydrogen Tunneling in Porphycene. <i>Journal of the American Chemical Society</i> , 2007, 129, 1335-1341.	6.6	82
139	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
140	Monothiodibenzoylmethane: Structural and vibrational assignments. <i>Vibrational Spectroscopy</i> , 2007, 43, 53-63.	1.2	12
141	Ground and excited state vibrations of 2-(2-pyridyl)pyrrole. <i>Journal of Molecular Structure</i> , 2007, 844-845, 286-299.	1.8	9
142	Photochromism in p-methylbenzoylthioacetone and related 1 ² -thioxoketones. <i>Chemical Physics</i> , 2007, 338, 11-22.	0.9	10
143	Conformation-Dependent Photophysics of Bifunctional Hydrogen Bond Donor/Acceptor Molecules. <i>Acta Physica Polonica A</i> , 2007, 112, S-105-S-120.	0.2	11
144	Evidence for Two Forms, Double Hydrogen Tunneling, and Proximity of Excited States in Bridge-Substituted Porphycenes: A Supersonic Jet Studies. <i>Journal of the American Chemical Society</i> , 2006, 128, 2577-2586.	6.6	61

#	ARTICLE	IF	CITATIONS
145	Excited-State Proton Transfer through Water Bridges and Structure of Hydrogen-Bonded Complexes in 1H-Pyrrolo[3,2-h]quinoline: A $\epsilon\%$ Adiabatic Time-Dependent Density Functional Theory Study. <i>Journal of Physical Chemistry A</i> , 2006, 110, 11958-11967.	1.1	74
146	Detection and Structural Characterization of Clusters with Ultrashort-Lived Electronically Excited States: IR Absorption Detected by Femtosecond Multiphoton Ionization. <i>Journal of the American Chemical Society</i> , 2006, 128, 10000-10001.	6.6	42
147	Ground- and Excited-State Tautomerism in Porphycenes. <i>Accounts of Chemical Research</i> , 2006, 39, 945-952.	7.6	103
148	Inverted Sapphyrin: A New Family of Doubly N-Confused Expanded Porphyrins. <i>Journal of the American Chemical Society</i> , 2006, 128, 12640-12641.	6.6	73
149	Ultrafast dynamics of alkyl-substituted porphycenes in solution. <i>Chemical Physics Letters</i> , 2006, 422, 142-146.	1.2	15
150	Excited state proton transfer in 2,9-(di-2-pyridyl)-4,7-di(t-butyl)carbazole. <i>Chemical Physics Letters</i> , 2006, 423, 288-292.	1.2	13
151	From Bifunctional Nucleophilic Behavior of DBU to a New Heterocyclic Fluorescent Platform. <i>Organic Letters</i> , 2006, 8, 4747-4750.	2.4	26
152	Photochromism and polarization spectroscopy of p-methyl(thiobenzoyl)acetone. <i>Chemical Physics</i> , 2006, 328, 205-215.	0.9	9
153	Energy relaxation paths in matrix-isolated excited molecules: Comparison of porphycene with dibenzoporphycenes. <i>Chemical Physics Letters</i> , 2005, 416, 128-132.	1.2	17
154	Electronic Structure, Spectra, and Magnetic Circular Dichroism of Cyclohexa-, Cyclohepta-, and Cyclooctapyrrole. <i>Chemistry - A European Journal</i> , 2005, 11, 4179-4184.	1.7	35
155	ELECTRONIC STATES OF ANTHANTHRENE. LINEAR AND MAGNETIC CIRCULAR DICHROISM, FLUORESCENCE ANISOTROPY, AND QUANTUM CHEMICAL CALCULATIONS. <i>Polycyclic Aromatic Compounds</i> , 2005, 25, 23-45.	1.4	7
156	Molecular dynamics simulations of matrix deposition. III. Site structure analysis for porphycene in argon and xenon. <i>Journal of Chemical Physics</i> , 2005, 123, 064706.	1.2	11
157	Vibrational spectroscopy of hydroxy-heterobiaryls. IR-active modes of [2,2-bipyridyl]-3,3-diol. <i>Photochemical and Photobiological Sciences</i> , 2005, 4, 143-148.	1.6	5
158	Imaging of Tautomerism in a Single Molecule. <i>Journal of the American Chemical Society</i> , 2005, 127, 5302-5303.	6.6	74
159	Molecular dynamics and density functional theory simulations of matrix deposition. II. Absolute site structure assignment for porphyrin in xenon. <i>Journal of Chemical Physics</i> , 2004, 121, 12017-12025.	1.2	8
160	Thioacetylacetone: Structural and Vibrational Assignments. <i>ChemPhysChem</i> , 2004, 5, 495-502.	1.0	21
161	Hydrogen-Bonding-Induced Phenomena in Bifunctional Heterozaaromatics. <i>ChemInform</i> , 2004, 35, no.	0.1	0
162	Time-resolved fluorescence studies of porphycene isolated in low-temperature gas matrices. <i>Chemical Physics Letters</i> , 2004, 394, 410-414.	1.2	11

#	ARTICLE	IF	CITATIONS
163	Intra- and intermolecular fluorescence quenching in 7-(pyridyl)indoles. <i>Chemical Physics Letters</i> , 2004, 400, 379-383.	1.2	42
164	Conformational equilibria and photoinduced tautomerization in 2-(2- π -pyridyl)pyrrole. <i>Chemical Physics Letters</i> , 2004, 400, 279-285.	1.2	33
165	From purely organic to metallo-organic chiral magnetic materials. <i>Polyhedron</i> , 2003, 22, 2349-2354.	1.0	23
166	Molecular dynamics simulations of matrix deposition. I. Site structure analysis for porphyrin in argon and xenon. <i>Journal of Chemical Physics</i> , 2003, 119, 7318-7327.	1.2	15
167	Hydrogen-Bonding-Induced Phenomena in Bifunctional Heteroazaaromatics. <i>Accounts of Chemical Research</i> , 2003, 36, 832-838.	7.6	185
168	Electronic spectroscopy and photophysics of 2-(N-methyl-N-isopropylamino)-5-cyanopyridine and related compounds Dedicated to Professor Dr Z. R. Grabowski and Professor Dr J. Wirz on the occasions of their 75th and 60th birthdays.. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 1027-1031.	1.3	12
169	Excited states of 4-dimethylaminopyridines: Magnetic circular dichroism and computational studies. <i>Photochemical and Photobiological Sciences</i> , 2003, 2, 187-194.	1.6	16
170	Determination of triplet formation efficiency from kinetic profiles of the ground state recovery Dedicated to Professor Jean Kossanyi on the occasion of his 70th birthday.. <i>Photochemical and Photobiological Sciences</i> , 2003, 2, 267.	1.6	20
171	Kinetic description of dioxygen binding to human hemoglobin on the 1-100 ns time scale. , 2002, 4749, 355.		0
172	Magnetic Circular Dichroism of Octaethylcorrphycene and Its Doubly Protonated and Deprotonated Forms. <i>Journal of Physical Chemistry A</i> , 2002, 106, 8139-8145.	1.1	24
173	An Experimental Test of C-N Bond Twisting in the TICT State: Syn ^o Anti Photoisomerization in 2-(N-Methyl-N-isopropylamino)-5-cyanopyridine. <i>Journal of the American Chemical Society</i> , 2002, 124, 2406-2407.	6.6	76
174	High-Resolution Spectroscopic Study of Matrix-Isolated Reactive Intermediates: A Vibrational Assignments for 3-Fluoro-o-Benzyne and Perfluoro-o-Benzyne ⁺ . <i>Journal of Physical Chemistry A</i> , 2002, 106, 6730-6737.	1.1	6
175	Fluorescence Quenching by Pyridine and Derivatives Induced by Intermolecular Hydrogen Bonding to Pyrrole-Containing Heteroaromatics. <i>Journal of Physical Chemistry A</i> , 2002, 106, 2158-2163.	1.1	58
176	Intramolecular charge-transfer properties of a molecule with a large donor group: the case of 4- π -(pyren-1-yl)benzonitrile. <i>Physical Chemistry Chemical Physics</i> , 2002, 4, 4334-4339.	1.3	22
177	Metal Complexes of Porphycene, Corrphycene, and Hemiporphycene: Stability and Coordination Chemistry. <i>Chemistry - A European Journal</i> , 2002, 8, 3485.	1.7	69
178	Magnetic Circular Dichroism of 5,10,15,20-Tetraphenylsapphyrin. <i>ChemPhysChem</i> , 2002, 3, 849-855.	1.0	8
179	In search for phototautomerization in solid dipyrido[2,3-a:3 π ,2 π -i]carbazole. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2002, 154, 61-68.	2.0	12
180	Magnetic circular dichroism of neutral and ionic forms of octaethylhemiporphycene. <i>Chemical Physics</i> , 2002, 282, 37-49.	0.9	18

#	ARTICLE	IF	CITATIONS
181	Photoinduced double proton transfer in water complexes of 1H-pyrrolo[3,2-h]quinoline and dipyrido[2,3-a:3â€²,2â€²-i]carbazole. <i>Chemical Physics Letters</i> , 2002, 366, 329-335.	1.2	22
182	Electronic states of the phenoxy radical. <i>Journal of Chemical Physics</i> , 2001, 115, 9733-9738.	1.2	82
183	Vibrations of the Phenoxy Radical. <i>Journal of the American Chemical Society</i> , 2001, 123, 11253-11261.	6.6	75
184	Raman Spectrum of the Phenyl Radical. <i>Journal of Physical Chemistry A</i> , 2001, 105, 10520-10524.	1.1	28
185	Electronic States of a Novel Smaragdyrin Isomer: Polarized Spectroscopy and Theoretical Studies. <i>Journal of Physical Chemistry A</i> , 2001, 105, 4992-4999.	1.1	13
186	The structure of the phototransformation product of monothiodibenzoylmethane. <i>Chemical Physics Letters</i> , 2001, 350, 502-508.	1.2	24
187	Vibrations of nitrous oxide: Matrix isolation Fourier transform infrared spectroscopy of twelve N2O isotopomers. <i>Journal of Chemical Physics</i> , 2001, 115, 1757-1764.	1.2	57
188	Tetrazete (N4). Can it be prepared and observed?. <i>Chemical Physics Letters</i> , 2000, 328, 227-233.	1.2	59
189	Electronic states of diphenyl- and dipyridyl-s-tetrazines: linear and magnetic circular dichroism, and quantum chemical calculations. <i>Chemical Physics</i> , 2000, 254, 135-149.	0.9	21
190	Ground and excited state tautomerization in 9-acetoxy-2,7,12,17-tetra-n-propylporphycene. <i>Chemical Physics Letters</i> , 2000, 323, 534-541.	1.2	31
191	Relaxation in excited states of porphycene in low-temperature argon and nitrogen matrices. <i>Chemical Physics Letters</i> , 2000, 318, 79-84.	1.2	12
192	Title is missing!. <i>Journal of Fluorescence</i> , 2000, 10, 41-48.	1.3	13
193	Molecular Dynamics and DFT Studies of Intermolecular Hydrogen Bonds between Bifunctional Heteroazaaromatic Molecules and Hydroxylic Solvents. <i>Journal of Physical Chemistry A</i> , 2000, 104, 9542-9555.	1.1	55
194	Solvent-Induced syn ⁺ anti Rotamerization of 2-(2â€²-Pyridyl)indole and the Structure of its Alcohol Complexes. <i>Journal of the American Chemical Society</i> , 2000, 122, 2818-2827.	6.6	64
195	(Sub)picosecond Fluorescence Upconversion Studies of Intermolecular Proton Transfer of Dipyrido[2,3-a:3â€²,2â€²-i]carbazole and Related Compounds. <i>Journal of Physical Chemistry A</i> , 2000, 104, 7167-7175.	1.1	47
196	Electronic transition moment directions in indoloindoles: the use of orientation amplifiers. <i>Journal of Molecular Structure</i> , 1999, 475, 141-151.	1.8	6
197	The Electronic Structure of Carcinogenic Dibenzopyrenes: Linear Dichroism, Fluorescence Polarization Spectroscopy and Quantum Mechanical Calculations. <i>Photochemistry and Photobiology</i> , 1999, 69, 158-166.	1.3	4
198	Linear Dichroism, Applications*. , 1999, , 1340-1345.		0

#	ARTICLE	IF	CITATIONS
199	Spectroscopy and Photophysics of a Highly Nonplanar Expanded Porphyrin: 4,9,13,18,22,27-Hexaethyl-5,8,14,17,23,26-hexamethyl-2,11,20-triphenylrosarin. <i>Chemistry - A European Journal</i> , 1999, 5, 3039-3045.	1.7	27
200	Photoinduced Double Proton Transfer: Inter- and Intramolecular Cases. <i>Israel Journal of Chemistry</i> , 1999, 39, 309-318.	1.0	40
201	Role of Ground State Structure in Photoinduced Tautomerization in Bifunctional Proton Donor-Acceptor Molecules: 1H-Pyrrolo[3,2-h]quinoline and Related Compounds. <i>Journal of the American Chemical Society</i> , 1999, 121, 11179-11188.	6.6	60
202	Light-Induced Tautomerization in Porphyrin Isomers. <i>Acta Physica Polonica A</i> , 1999, 95, 49-62.	0.2	5
203	Linear Dichroism, Instrumentation*. , 1999, , 1346-1348.		0
204	Excited-state intramolecular proton transfer in anthralin.. <i>Chemical Physics Letters</i> , 1998, 291, 51-56.	1.2	23
205	Proton tunnelling in porphycene seeded in a supersonic jet. <i>Chemical Physics Letters</i> , 1998, 296, 549-556.	1.2	65
206	Diversity of excited state deactivation paths in heteroazaaromatics with multiple intermolecular hydrogen bonds. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1998, 102, 469-475.	0.9	19
207	Vibrational spectroscopy of hydroxy-heterobiaryls. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1998, 54, 1291-1305.	2.0	9
208	Spectroscopy and Photophysics of Tetraalkyldibenzoporphycenes. <i>Journal of Physical Chemistry A</i> , 1998, 102, 4966-4971.	1.1	37
209	Electronic Spectra of Porphycenes in Rare Gas and Nitrogen Matrices. <i>Journal of Physical Chemistry A</i> , 1998, 102, 9999-10006.	1.1	37
210	Compact multifunctional spectrofluorimeter with a novel design for anisotropy measurements. <i>Review of Scientific Instruments</i> , 1998, 69, 2242-2245.	0.6	10
211	Intermolecular Excited State Double Proton Transfer in Dipyridocarbazole:Alcohol Complexes. <i>Journal of Physical Chemistry A</i> , 1997, 101, 5839-5845.	1.1	50
212	Photophysics of trans-stilbene analogues: Indolo[3,2-b]indole and its heterosubstituted sulfur and selenium derivatives. <i>Chemical Physics</i> , 1997, 216, 179-192.	0.9	18
213	Stilbene-like molecules: sulfur- and selenium-heterosubstituted indolo[3,2-b]indoles. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 1997, 105, 149-152.	2.0	5
214	Luminescence anisotropy and linear dichroism studies of large polycyclic aromatic hydrocarbons based on the perylene chromophore. <i>Journal of Luminescence</i> , 1997, 72-74, 517-519.	1.5	3
215	Vibrational spectroscopy of hydroxy-heterobiaryls. Low frequency modes of 2-(2-hydroxyphenyl)-3-pyridinol. <i>Journal of Molecular Structure</i> , 1997, 408-409, 363-366.	1.8	0
216	Matrix-isolated products of cyanoacetylene dissociation. <i>Journal of Molecular Structure</i> , 1997, 408-409, 473-476.	1.8	30

#	ARTICLE	IF	CITATIONS
217	Electronic transition moment directions and tautomerization of 9,10,19,20-tetra-n-propylporphycene. <i>Chemical Physics Letters</i> , 1997, 271, 341-348.	1.2	13
218	Electronic spectra and symmetry of metalloporphyrins in low-temperature rare gas and nitrogen matrices. <i>Chemical Physics Letters</i> , 1997, 272, 405-411.	1.2	25
219	Solvent-Controlled Excited State Behavior: 2-(2-Pyridyl)indoles in Alcohols. <i>Journal of the American Chemical Society</i> , 1996, 118, 3508-3518.	6.6	116
220	Exciton coupling in various substituted aryl-phthalimide bichromophoric systems. <i>Tetrahedron</i> , 1996, 52, 13201-13214.	1.0	17
221	Determination of Structure of Nonplanar Polycyclic Aromatic Hydrocarbons by Polarized Spectroscopy. <i>Polycyclic Aromatic Compounds</i> , 1996, 9, 283-290.	1.4	4
222	Proton and electron transfer in hydrogen-bonded systems. <i>Journal of Molecular Liquids</i> , 1995, 64, 49-56.	2.3	8
223	Electronic states of symmetrically disubstituted s-tetrazines. <i>Chemical Physics</i> , 1995, 200, 201-213.	0.9	31
224	Electronic and vibrational spectroscopy of phototautomerizing heteroazaaromatics with intramolecular hydrogen bonds. <i>Journal of Molecular Structure</i> , 1995, 349, 277-280.	1.8	4
225	Polarized Infrared Spectra of Photooriented Matrix-Isolated Free-Base Porphyrin Isotopomers. <i>The Journal of Physical Chemistry</i> , 1995, 99, 14254-14260.	2.9	41
226	Electronic States and Structure of Dibenzo[j,lm]-phenanthro[5,4,3-abcd]perylene. <i>Acta Physica Polonica A</i> , 1995, 88, 295-304.	0.2	6
227	Spectroscopy and Photophysics of 3'-[p-(dimethylamino)phenyl]spiro-[fluorene-9,4'-oxazolidine]-2',5'-dione. <i>Acta Physica Polonica A</i> , 1995, 88, 445-454.	0.2	7
228	Distance Dependence of Excited-State Double Proton Transfer in Porphycenes Studied by Fluorescence Polarization. <i>The Journal of Physical Chemistry</i> , 1994, 98, 4530-4535.	2.9	50
229	Excited charge transfer states in 4-aminopyrimidines, 4-(dimethylanilino)pyrimidine and 4-(dimethylamino)pyridine. <i>Chemical Physics</i> , 1994, 188, 247-265.	0.9	73
230	The molecular and electronic structure of dibenzo[g,p]chrysene: A twisted case. <i>Theoretica Chimica Acta</i> , 1994, 89, 301-309.	0.9	10
231	Excited state tautomerization in porphycenes studied by polarized spectroscopy. <i>Journal of Luminescence</i> , 1994, 60-61, 867-869.	1.5	4
232	Mechanisms of fluorescence quenching by hydrogen bonding in various aza aromatics. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 1994, 80, 157-160.	2.0	50
233	Spectral studies of organic chromophores in polymeric matrices. , 1993, , .		0
234	<title>Site structure in the electronic and IR spectra of porphyrin</title>. , 1992, , .		0

#	ARTICLE	IF	CITATIONS
235	Electronic states of benzo[a]pyrene. Linear and magnetic circular dichroism, polarized fluorescence, and quantum chemical calculations. <i>Journal of the American Chemical Society</i> , 1992, 114, 1942-1949.	6.6	21
236	Ground state structures of molecules prepared for phototautomerization: 2,2'-bipyridyl-3,3'-diol and 2,2'-bipyridyl-3-ol. <i>Journal of Crystallographic and Spectroscopic Research</i> , 1992, 22, 563-572.	0.3	30
237	Secondary fluorescences of 4'-(1-pyrenyl)acetophenone and p-dimethylaminoacetophenone. <i>International Journal of Radiation Applications and Instrumentation Nuclear Tracks and Radiation Measurements</i> , 1992, 39, 149-153.	0.0	0
238	Quenching of fluorescence of 2-(2-pyridyl) indoles upon complexation with alcohols. <i>Chemical Physics Letters</i> , 1992, 195, 556-562.	1.2	41
239	ELECTRONIC ABSORPTION PROPERTIES OF SYMMETRICAL DIALKOXYANTHRACENES. LINEAR DICHROISM AND MAGNETIC CIRCULAR DICHROISM*. <i>Photochemistry and Photobiology</i> , 1992, 55, 335-347.	1.3	13
240	Excited state relaxation processes in the case of some acetophenone derivatives. <i>Journal of Chemical Sciences</i> , 1992, 104, 143-152.	0.7	8
241	Magnetic circular dichroism of metallotetrapyrins. <i>Journal of Organic Chemistry</i> , 1991, 56, 2735-2742.	1.7	11
242	Electronic states of porphycenes. <i>Journal of the American Chemical Society</i> , 1991, 113, 5511-5527.	6.6	142
243	Perimeter model and magnetic circular dichroism of porphyrin analogs. <i>Journal of Organic Chemistry</i> , 1991, 56, 2729-2735.	1.7	92
244	Determination of Orientation Averages for Partially Aligned low Symmetry Molecules. <i>Spectroscopy Letters</i> , 1991, 24, 951-957.	0.5	1
245	Evidence for conformational equilibrium in an unsubstituted benzenoid hydrocarbon: tetrabenz[a,c,d,f,l,m]perylene. <i>The Journal of Physical Chemistry</i> , 1991, 95, 8660-8663.	2.9	11
246	Fourier transform fluorescence and phosphorescence of porphine in rare gas matrixes. <i>The Journal of Physical Chemistry</i> , 1991, 95, 1963-1969.	2.9	67
247	Magnetic circular dichroism of cyclic π -electron systems. 29. Bicyclic phospholium and arsolium cations with phosphorus-sulfur, arsenic-sulfur, phosphorus-nitrogen, and arsenic-nitrogen p - π bonding. <i>Organometallics</i> , 1990, 9, 1085-1088.	1.1	11
248	Electronic states of chrysene: linear and magnetic circular dichroism and quantum chemical calculations. <i>The Journal of Physical Chemistry</i> , 1990, 94, 1800-1806.	2.9	25
249	FT visible absorption spectroscopy of porphine in noble gas matrixes. <i>Journal of Molecular Spectroscopy</i> , 1990, 140, 373-389.	0.4	43
250	Site-population conserving and site-population altering photo-orientation of matrix-isolated free-base porphine by double proton transfer: IR dichroism and vibrational symmetry assignments. <i>Chemical Physics</i> , 1989, 136, 165-180.	0.9	50
251	The nature of the excited states of p-nitro-N,N-dimethylaniline. <i>Journal of Luminescence</i> , 1989, 44, 149-160.	1.5	16
252	Transition moment directions and molecular structure of some p-cyano-N,N-dimethylaniline derivatives. <i>Chemical Physics</i> , 1989, 138, 105-113.	0.9	26

#	ARTICLE	IF	CITATIONS
253	Effective electronegativities of phosphorus, arsenic, and antimony in a .pi. system. Evidence from magnetic circular dichroism. <i>Organometallics</i> , 1989, 8, 2804-2808.	1.1	70
254	Electronic states and magnetic circular dichroism of cyclic π systems with antiaromatic perimeters. <i>Pure and Applied Chemistry</i> , 1989, 61, 2117-2128.	0.9	9
255	The electronic spectra of some aza-derivatives of 1,2-benzanthracene. <i>Chemical Physics</i> , 1988, 124, 103-112.	0.9	4
256	Luminescence of neutral and protonated aminoquinoxalines. <i>Journal of Luminescence</i> , 1988, 40-41, 211-212.	1.5	0
257	Luminescence of hydrogen-bonded 2-aminopyridine dimers. <i>Journal of Molecular Structure</i> , 1988, 177, 415-419.	1.8	6
258	Excited states of indoloquinoxalines. <i>Spectrochimica Acta Part A: Molecular Spectroscopy</i> , 1988, 44, 1335-1340.	0.1	5
259	Ground- and excited-state protonation of aminoquinoxalines. <i>The Journal of Physical Chemistry</i> , 1988, 92, 6930-6935.	2.9	20
260	The electronic spectrum of benz[a]anthracene. Linear and magnetic circular dichroism and fluorescence polarization studies. <i>Chemical Physics</i> , 1987, 116, 411-420.	0.9	18
261	A complete determination of transition moment directions from fluorescence spectroscopy and IR and UV linear dichroism: 1,2-benzanthracene. <i>Chemical Physics Letters</i> , 1987, 135, 515-520.	1.2	21
262	Modification of photophysical behaviour by hydrogen bonding: Indoloquinoxaline and its methylated derivatives. <i>Chemical Physics Letters</i> , 1987, 133, 368-372.	1.2	24
263	Photophysics of pseudoazulenes: 7-azaindole derivatives. <i>Journal of Photochemistry and Photobiology</i> , 1987, 39, 49-58.	0.6	6
264	Excited-state double proton transfer in 1-azacarbazole-alcohol complexes. <i>The Journal of Physical Chemistry</i> , 1986, 90, 3868-3871.	2.9	46
265	A probe for substituent hyperconjugative power: MCD (magnetic circular dichroism) of the benzene Lb band. <i>Pure and Applied Chemistry</i> , 1986, 58, 39-53.	0.9	14
266	Excited-state double proton transfer in the solid state: the dimers of 1-azacarbazole. <i>The Journal of Physical Chemistry</i> , 1986, 90, 3866-3868.	2.9	44
267	Linear and magnetic circular dichroism spectra of 1,2:7,8-dibenzchrysene: out-of-plane polarized intensity. <i>Chemical Physics Letters</i> , 1986, 123, 102-107.	1.2	8
268	Solvent-Dependent Photophysics of Indoloquinoxaline. <i>Journal of Molecular Structure</i> , 1986, 142, 159-162.	1.8	14
269	Determination of the energy barrier origin of the excited state double proton transfer in 7-azaindole: Alcohol complexes. <i>Journal of Molecular Structure</i> , 1984, 114, 329-332.	1.8	26
270	Viscosity and temperature effects in excited state double proton transfer: Luminescence of 1-azacarbazole dimers in solid state and solution. <i>Journal of Molecular Structure</i> , 1984, 114, 359-362.	1.8	8

#	ARTICLE	IF	CITATIONS
271	Spectroscopy of doubly hydrogen-bonded 7-azaindole. Reinvestigation of the excited state reaction. <i>Journal of Luminescence</i> , 1984, 29, 65-81.	1.5	71
272	Spectroscopy of doubly hydrogen-bonded 7-azaindole. Reinvestigation of the excited state reaction. <i>Journal of Luminescence</i> , 1984, 29, 65-81.	1.5	20
273	Viscosity vs. temperature effects in excited-state double proton transfer. Comparison of 1-azacarbazole with 7-azaindole. <i>The Journal of Physical Chemistry</i> , 1984, 88, 1160-1162.	2.9	46
274	Excited-state dipole moment changes in some pseudoazulenes. <i>Chemical Physics Letters</i> , 1983, 94, 58-62.	1.2	12
275	Applications of magnetic circular dichroism (MCD) spectroscopy: electronic structure of the thiotriazolium cation, S4N3 ⁺ . <i>Inorganic Chemistry</i> , 1982, 21, 556-558.	1.9	8
276	Magnetic circular dichroism of cyclic π -electron systems. 22. Derivatives of the trisulfur trinitride anion. <i>Inorganic Chemistry</i> , 1982, 21, 832-834.	1.9	8
277	Magnetic circular dichroism of cyclic π -electron systems. 21. The trisulfur trinitride anion. <i>Inorganic Chemistry</i> , 1981, 20, 963-965.	1.9	15
278	Magnetic circular dichroism of cyclic π -electron systems. 23. Monocyclic π -excessive nitrogen heterocycles. 1,4-Dihydro-1,4-diazocines. <i>Journal of Organic Chemistry</i> , 1981, 46, 3306-3311.	1.7	11
279	Determination of positional isomers of methylpyrenes and other polycyclic aromatic hydrocarbons by magnetic circular dichroism. <i>Analytical Chemistry</i> , 1981, 53, 236-239.	3.2	6
280	Red edge excitation study of cooperative double proton transfer in 7-azaindole. <i>Journal of Luminescence</i> , 1981, 24-25, 519-522.	1.5	26
281	Charge density flow as a driving force of distortion in excited protonated azaaromatics. <i>Chemical Physics Letters</i> , 1980, 70, 175-179.	1.2	17
282	Proton addition to excited diazaphenanthrenes topology determined molecular properties. <i>Journal of Luminescence</i> , 1980, 21, 277-291.	1.5	18
283	Influence of protonation upon excited states of diazaphthalenes. Semi-empirical study of phthalazine and quinoxaline. <i>Chemical Physics Letters</i> , 1979, 63, 579-583.	1.2	7
284	Are protonated ortho-diazines planar in excited states?. <i>Journal of Luminescence</i> , 1979, 18-19, 201-204.	1.5	11
285	Tautomerization in Porphycenes. , 0, , 245-271.		2