Javier CatalÃ¡n

List of Publications by Year
in descending order

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152	5,663	71102	85541
papers	citations	41 h-index	71 g-index

155	155	155	4446
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Carbon disulfide solvent, helping to clarify the Stokes shift in diphenylpolyenes. Journal of Physical Organic Chemistry, 2021, 34, e4147.	1.9	2
2	Solvatochromic correlation analysis of monomolecular SN1/E1 heterolysis reactions of tertiary haloalkanes. Journal of Molecular Liquids, 2021, 324, 114699.	4.9	4
3	On the empirical scales of organic solvents established using probe/homomorph pairs. Journal of Physical Organic Chemistry, 2021, 34, e4206.	1.9	8
4	Photophysics of the electronic states S 0 and S 1 for the coplanar molecular structures of the Î±,Ï‰â€diphenylpolyenes DPH and DPO. Journal of Physical Organic Chemistry, 2021, 34, e4256.	1.9	0
5	On the mirror symmetry between the absorption and emission of complex molecules in solution. Journal of Physical Organic Chemistry, 2020, 33, e4034.	1.9	3
6	Comments on â€œQuantifying solvent effects through QSPR: A new look over different model equationsâ€• Journal of Molecular Liquids, 2020, 298, 111922.	4.9	11
7	Solvatochromism in urea/water and urea-derivative/water solutions. Physical Chemistry Chemical Physics, 2020, 22, 25165-25176.	2.8	6
8	On the photophysics of a polyene dissolved in n â€alkanes when the temperature drops from 293 to 77 K. Journal of Physical Organic Chemistry, 2020, 33, e4097.	1.9	2
9	On the first electronic transitions in molecular spectra of conjugated diphenylpolyenes: A reappraisal. Chemical Physics, 2019, 525, 110422.	1.9	9
10	On the hydrophobic effect in waterâ€“alcohol mixtures. Chemical Physics, 2019, 527, 110467.	1.9	11
11	On the photoâ€activation of the S 0 Âât'ÂS 1 transition in polyenes. Journal of Physical Organic Chemistry, 2019, 32, e3933.	1.9	4
12	Kasha's rule: a reappraisal. Physical Chemistry Chemical Physics, 2019, 21, 10061-10069.	2.8	97
13	Thermochromism of pure alkanols and water versus its polarizability. Chemical Physics, 2019, 522, 99-103.	1.9	1
14	On the dual fluorescence of Î±,Ï‰â€diphenylpolyenes from five to seven polyene double bonds. Journal of Physical Organic Chemistry, 2019, 32, e3906.	1.9	3
15	On the photophysical model for polyenes: Experimental evidence of the 1,8â€diphenyloctaâ€1,3,5,7â€tetraene molecule dissolved in nâ€octane, cyclooctane, and 2,2,4â€trimethylpentane. Journal of Physical Organic Chemistry, 2018, 31, e3794.	1.9	7
16	Molecule 1-Methyl-5-nitroindoline Probes the Structural Change of Liquid Water with Temperature. ACS Omega, 2018, 3, 18930-18934.	3.5	6
17	Existence of Two Fluorescence Bands in all-trans-Polyenes with Six and Seven Double Bonds. Journal of Physical Chemistry A, 2018, 122, 6391-6395.	2.5	3
18	Spectroscopy of 1,6-diphenyl-1,3,5-hexatriene (DPH) dissolved in three hexane structural isomers, and its consequences on the photophysical model of polyenes. Physical Chemistry Chemical Physics, 2017, 19, 10657-10662.	2.8	8

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19	Liquid water changes its structure at 43 Â°C. Chemical Physics Letters, 2017, 679, 86-89.	2.6	12
20	Solvation by Glycerol at Temperatures from 353 to 77 K: Its Solvatochromic Characterization and Use to Block the Molecular Structure of Conformationally Flexible Structures. Journal of Physical Chemistry A, 2017, 121, 7114-7120.	2.5	9
21	Inversion of the 1¹B_uand 2¹A_gelectronic states of all-trans-1,6-diphenyl-1,3,5-hexatriene in carbon disulfide. Physical Chemistry Chemical Physics, 2017, 19, 27099-27104.	2.8	11
22	Influence of solvent basicity on DMABN photophysics. Journal of Physical Organic Chemistry, 2017, 30, e3613.	1.9	2
23	On the dimerization of unsubstituted Î±,Ï‰-diphenylpolyenes at low concentrations in inert solvents. Journal of Physical Organic Chemistry, 2016, 29, 414-420.	1.9	2
24	The first UV absorption band of<scp>l</scp>-tryptophan is not due to two simultaneous orthogonal electronic transitions differing in the dipole moment. Physical Chemistry Chemical Physics, 2016, 18, 15170-15176.	2.8	12
25	Compounds with Ï€(loc)â€‰â†’â€‰Ï€*(deloc) electronic transitions and their solvatochromism. Journal of Physical Organic Chemistry, 2015, 28, 497-503.	1.9	10
26	Fluorosolvatochromism of monomethyl indoles: further evidence in support of a new photophysical model for the indole chromophore. Journal of Physical Organic Chemistry, 2015, 28, 329-336.	1.9	6
27	On Saltiel's isopolarizability approach and its applicability to diphenylpolyenes. Chemical Physics Letters, 2015, 635, 56-59.	2.6	6
28	The first UV absorption band for indole is not due to two simultaneous orthogonal electronic transitions differing in dipole moment. Physical Chemistry Chemical Physics, 2015, 17, 12515-12520.	2.8	7
29	Reply to the comment on â€œOn Saltiel's isopolarizability approach and its applicability to diphenylpolyenesâ€•by J. CatalÃ¡n, Chem. Phys. Lett. 635 (2015) 56. Chemical Physics Letters, 2015, 641, 104-105.	2.6	1
30	SOLVENT EFFECTS BASED ON PURE SOLVENT SCALES. , 2014, , 581-622.		7
31	On dimers and complexes of tetracene and the Kasha's molecular excitonic model. Journal of Physical Organic Chemistry, 2014, 27, 456-462.	1.9	1
32	Can the dipolarity of the medium induce the formation of charge transfer structures? An unexpected finding in the photophysics of DMABN. Physical Chemistry Chemical Physics, 2014, 16, 7734.	2.8	13
33	A Spectroscopic Rule from the Solvatochromism of Aromatic Solutes in Nonpolar Solvents. Journal of Physical Chemistry B, 2014, 118, 5168-5176.	2.6	16
34	On the solvatochromism, dimerization and tautomerism of indazole. Arkivoc, 2014, 2014, 57-70.	0.5	1
35	Reply to the comment on â€œOn the dual emission of p-dimethylaminobenzonitrile and its photophysical implicationsâ€•by J. CatalÃ¡n, Phys. Chem. Chem. Phys., 2013, 15, 8811â€“8820. Physical Chemistry Chemical Physics, 2013, 15, 16978.	2.8	5
36	On the dual emission of p-dimethylaminobenzonitrile and its photophysical implications. Physical Chemistry Chemical Physics, 2013, 15, 8811.	2.8	34

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37	On the use of Î²â€¢carotene as a probe for solvent polarizability. Journal of Physical Organic Chemistry, 2013, 26, 948-952.	1.9	6
38	Is the LEâ†'TICT process in the S1 excited state of 9,9â€²-bisanthracenyl influenced by the viscosity or the dipolarity of the solvent?. Journal of Luminescence, 2013, 143, 635-639.	3.1	4
39	Analysis of the Solvatochromism of 9,9â€²-Biaryl Compounds Using a Pure Solvent Dipolarity Scale. Journal of Physical Chemistry A, 2012, 116, 4726-4734.	2.5	30
40	On the fluorescence of methyl salicylate: the significance of its IMHB. Physical Chemistry Chemical Physics, 2012, 14, 8903.	2.8	15
41	Do stilbazolium betaine dyes exhibit inverted solvatochromism by changes in solvent dipolarity?. Dyes and Pigments, 2012, 95, 180-187.	3.7	23
42	Questioning the photophysical model for the indole chromophore in the light of evidence obtained by controlling the non-specific effect of the medium with 1-chlorobutane as solvent. Physical Chemistry Chemical Physics, 2011, 13, 15022.	2.8	8
43	Photophysics of the 6H-Indolo[2,3-b]quinoline Molecule: The Excited-State Double-Proton-Transfer Process in Perspective. Journal of Physical Chemistry A, 2011, 115, 1900-1907.	2.5	3
44	On the solvatochromism of the n â†"Ï€* electronic transitions in ketones. Physical Chemistry Chemical Physics, 2011, 13, 4072.	2.8	31
45	Generalized solvent scales as a tool for investigating solvent dependence of spectroscopic and kinetic parameters. Application to fluorescent BODIPY dyes. Photochemical and Photobiological Sciences, 2010, 9, 996-1008.	2.9	100
46	On the Molecular Structure and UV/vis Spectroscopic Properties of the Solvatochromic and Thermochromic Pyridinium-N-Phenolate Betaine Dye B30. Journal of Physical Chemistry A, 2010, 114, 6226-6234.	2.5	64
47	Activation Energy of the Two-Proton Phototautomerism in 7-Azaindole Dimer and Its Medium-Dependence. Journal of Physical Chemistry A, 2010, 114, 5666-5673.	2.5	14
48	Proton Phototransfers in Doubly Hydrogen Bonded Dimers: The Photophysics of 6,7,8,9-Tetrahydro-5H-pyrido[2,3-b]indole Dimers. Journal of Physical Chemistry A, 2010, 114, 811-816.	2.5	6
49	On the 7-azaindole in acetonitrile anhydrous solutions as an inappropriate photophysical model for DNA base pairs. Nature Precedings, 2009, , .	0.1	0
50	Toward a Generalized Treatment of the Solvent Effect Based on Four Empirical Scales: Dipolarity (SdP,) Tj ETQq0 0 0 rgBT /Overlock 10 T Chemistry B, 2009, 113, 5951-5960.	2.6	581
51	On the chromism of polyenes. Chemical Physics Letters, 2008, 457, 87-90.	2.6	8
52	Fluorescence Spectroscopy and Amplified Spontaneous Emission (ASE) of Phenylimidazoles: Predicted Vibronic Coupling Along the Excited-State Intramolecular Proton Transfer in 2-(2â€²-Hydroxyphenyl)imidazoles. Journal of Physical Chemistry A, 2008, 112, 5555-5565.	2.5	14
53	On the Inoperativeness of the ESIPT Process in the Emission of 1-Hydroxy-2-acetonaphthone:â€‰ A Reappraisal. Journal of Physical Chemistry A, 2008, 112, 904-914.	2.5	24
54	On the Photophysics of Polyenes. 1. Bathochromic Shifts in Their 1Ag â†' 1Bu Electronic Transitions Caused by the Polarizability of the Medium. Journal of Physical Chemistry A, 2008, 112, 5653-5657.	2.5	15

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55	On the concerted mechanism of photo-induced biprotonic transfer in C _{2h} 7-azaindole dimer. Proceedings of the National Academy of Sciences of the United States of America, 2008, 105, E78; author reply E79.	7.1	12
56	New contributions to the photophysical model for all-trans-polyenesfrom ttbP4, a nonphotolabile octatetraene. Journal of Chemical Physics, 2008, 128, 104504.	3.0	10
57	The photophysics of all-trans polyenes from ttbP5, a nonphotolabile pentaene. Journal of Chemical Physics, 2008, 129, 014505.	3.0	6
58	On the Doubly hydrogen bonded dimer of 7-azaindole (0.1 M) as a model for DNA base pairs in acetonitrile solutions at rt. Nature Precedings, 2008, , .	0.1	0
59	Photophysics of 1-Azacarbazole Dimers:â€‰ A Reappraisal. Journal of Physical Chemistry A, 2007, 111, 8774-8779.	2.5	14
60	On the molecular conformation of bisaromatic systems the case of 2-phenyl-2H-benzotriazoles. Chemical Physics, 2007, 340, 32-42.	1.9	5
61	The emission of Î±,Ï‰-diphenylpolyenes: A model involving several molecular structures. Chemical Physics, 2007, 335, 69-78.	1.9	9
62	On the photophysics of all-trans polyenes: Hexatriene versus octatetraene. Journal of Chemical Physics, 2006, 124, 034306.	3.0	28
63	Excited State Proton Transfer in 3-Methyl-7-Azaindole Dimer. Symmetry Control. Journal of Physical Chemistry A, 2006, 110, 9116-9122.	2.5	21
64	Excited-state proton phototransfer in the (3-methyl-7-azaindole)-(7-azaindole) heterodimer. Chemical Physics Letters, 2006, 419, 164-167.	2.6	14
65	Polarization of the T1â†’S0 phosphorescence and S0â†’Sn phosphorescence excitation of aromatic hydrocarbons prototype for Ï€,Ï€* states. A reappraisal. Chemical Physics, 2005, 316, 253-259.	1.9	5
66	Molecular structure distortions and the Mullikenâ€“Rieke rule: The case of t-stilbene. Chemical Physics Letters, 2005, 416, 165-170.	2.6	17
67	Electronic Energy Levels in all-trans Long Linear Polyenes: The Case of the 3,20-Di(tert-butyl)-2,2,21,21-tetramethyl-all-trans-3,5,7,9,11,13,15,17,19-docosanonaen (ttbp9) Conforming to Kasha's Rule. Chemistry - A European Journal, 2005, 11, 3915-3920.	3.3	15
68	Study of 7-azaindole in its first four singlet states. Journal of Chemical Physics, 2005, 122, 244320.	3.0	13
69	The molecular symmetry and electronic spectroscopy of 7-azaindole dimer: Its proton-transfer channels. Journal of Chemical Physics, 2005, 123, 114302.	3.0	31
70	Empirical Treatment of the Inductive and Dispersive Components of Soluteâ^'Solvent Interactions: The Solvent Polarizability (SP) Scale. European Journal of Organic Chemistry, 2004, 2004, 4694-4702.	2.4	140
71	Analysis of mixed solvent effects on the properties of singlet oxygen (). Chemical Physics, 2004, 300, 33-39.	1.9	32
72	On the absorption and emission spectra for the purine chromophore in weakly perturbative environments. Chemical Physics, 2004, 303, 205-218.	1.9	9

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73	First reported evidence that solvent polarity induces an â†" inversion in the indole chromophore. Chemical Physics Letters, 2003, 368, 717-723.	2.6	23
74	DFT study of ionization potentials for aza-substituted aromatic rings. International Journal of Quantum Chemistry, 2003, 91, 432-437.	2.0	22
75	Characterization of binary solvent mixtures: the waterâ€"acetonitrile mixture. Organic and Biomolecular Chemistry, 2003, 1, 575-580.	2.8	37
76	ON THE ABSORPTION SPECTRUM OF C60IN THE VISIBLE SPECTRAL REGION. Fullerenes Nanotubes and Carbon Nanostructures, 2002, 10, 171-180.	2.1	3
77	The concerted mechanism of photo-induced biprotonic transfer in 7-azaindole dimers: A model for the secondary evolution of the classic C2h dimer and comparison of four mechanisms. Proceedings of the National Academy of Sciences of the United States of America, 2002, 99, 5799-5803.	7.1	56
78	The concerted mechanism of photo-induced biprotonic transfer in 7-azaindole dimers: Structure, quantum-theoretical analysis, and simultaneity principles. Proceedings of the National Academy of Sciences of the United States of America, 2002, 99, 5793-5798.	7.1	54
79	Characterization of Binary Solvent Mixtures of DMSO with Water and Other Cosolvents. Journal of Organic Chemistry, 2001, 66, 5846-5852.	3.2	183
80	On the Origin of Nonvertical Triplet Excitation Transfer:Â The Relative Role of Double-Bond and Phenyl-Vinyl Torsions in the Stilbenes. Journal of Physical Chemistry A, 2001, 105, 6273-6276.	2.5	27
81	Understanding the solvatochromism of 10-hydroxybenzo[h]quinoline. An appraisal of a polarity calibrator. Chemical Physics, 2001, 270, 1-12.	1.9	14
82	The singular coincidence of fluorescence spectra of the anionic and cationic species formed by the respective deprotonated and protonated pyrido-pyrrolo bases. International Journal of Quantum Chemistry, 2000, 77, 118-127.	2.0	41
83	Conformity of the 7-azaindole dimer cationic potential with photoionization/Coulomb-explosion MS observations and the concerted biprotonic transfer mechanism. Chemical Physics Letters, 2000, 318, 629-636.	2.6	51
84	Photophysics of 7-Azaindole, Its Doubly-H-Bonded Base-Pair, and Corresponding Proton-Transfer-Tautomer Dimeric Species, via Defining Experimental and Theoretical Results. Journal of Physical Chemistry A, 2000, 104, 10812-10820.	2.5	86
85	Medium-Controlled Aggregation of trans-Stilbene. Journal of the American Chemical Society, 2000, 122, 2377-2378.	13.7	43
86	Solvatochromism of fluorophores with an intramolecular hydrogen bond and their use as probes in biomolecular cavity sites. International Journal of Quantum Chemistry, 1999, 72, 421-438.	2.0	51
87	Extending the Solvent Acidiy Scale to Highly Acidic Organic Solvents: The Unique Photophysical Behaviour of 3,6-Diethyltetrazine. European Journal of Organic Chemistry, 1999, 1999, 885-891.	2.4	74
88	The Six-Membered Intramolecular Hydrogen Bond Position as a Switch for Inducing an Excited State Intramolecular Proton Transfer (ESIPT) in Esters ofo-Hydroxynaphthoic Acids. Journal of Physical Chemistry A, 1999, 103, 10921-10934.	2.5	68
89	Correlation of Solvolysis Rates 50 Years Later. Journal of Organic Chemistry, 1999, 64, 6512-6514.	3.2	23
90	Gas-phase protolysis between a neutral BrÃ¸nsted acid and a neutral BrÃ¸nsted base?. Chemical Physics Letters, 1998, 293, 511-514.	2.6	15

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91	On the first triplet state of benzotriazole-like ultraviolet stabilizers. Chemical Physics Letters, 1998, 297, 549-552.	2.6	3
92	On the TICT Mechanism of 9,9â€²-Biaryl Compounds. European Journal of Organic Chemistry, 1998, 1998, 1697-1704.	2.4	9
93	Inter-ring Torsional Modulation in Molecular Lasers. Ultraviolet Lasing via Amplified Spontaneous Emission Spectroscopy of Phenylimidazoles. Journal of Physical Chemistry A, 1997, 101, 5284-5291.	2.5	33
94	Photophysical Study of Pyridoxal 5'-Phosphate and Its Schiff Base with n-Hexylamine. Photochemistry and Photobiology, 1997, 66, 810-816.	2.5	6
95	Spectroscopy of Amplified Spontaneous Emission Laser Spikes in Phenyloxazoles. Prototype Classes. Journal of Physical Chemistry A, 1997, 101, 3260-3272.	2.5	45
96	Photophysics of the 2-(2â€²-hydroxyphenyl)perimidines: Photostability studies. Journal of Luminescence, 1997, 75, 17-26.	3.1	10
97	A Generalized Solvent Acidity Scale: The Solvatochromism of oâ€tertâ€Butylstilbazolium Betaine Dye and Its Homomorph o,oâ€²â€Diâ€tertâ€butylstilbazolium Betaine Dye. Liebigs Annalen, 1997, 1997, 1941-1949.	0.8	144
98	Importance of aromaticity on the relative stabilities of indazole annular tautomers: an ab initio study. Journal of the Chemical Society Perkin Transactions II, 1996, , 57-60.	0.9	55
99	A Generalized Solvent Basicity Scale: The Solvatochromism of 5â€Nitroindoline and Its Homomorph 1â€Methylâ€5â€nitroindoline. Liebigs Annalen, 1996, 1996, 1785-1794.	0.8	174
100	Chemical physics of excitation dynamics via amplified spontaneous emission (ASE) laser spike spectroscopy in substituted phenyloxazoles. Chemical Physics Letters, 1996, 263, 154-160.	2.6	11
101	THE INFLUENCE OF MOLECULAR CONFORMATION ON THE STABILITY OF ULTRAVIOLET STABILIZERS TOWARD DIRECT AND DYE-SENSITIZED PHOTOIRRADIATION: THE CASE OF 2-(2'-HYDROXY-5'-METHYPHENYL)BENZOTRIAZOLE (TIN P). Photochemistry and Photobiology, 1995, 61, 118-123.	2.5	48
102	Progress towards a generalized solvent polarity scale: The solvatochromism of 2â€(dimethylamino)â€7â€nitrofluorene and its homomorph 2â€fluoroâ€7â€nitrofluorene. Liebigs Annalen, 1995, 1995, 241-252.	0.8	187
103	Solvent dipolarity/polarizability (SPP) of alcoholic solvents. Liebigs Annalen, 1995, 1995, 793-795.	0.8	62
104	The Colors of C60 Solutions. Angewandte Chemie International Edition in English, 1995, 34, 105-107.	4.4	47
105	Toward the Photostability Mechanism of Intramolecular Hydrogen Bond Systems. 4. 3(5)-(1'-Hydroxy-2'-naphthyl)pyrazoles and 3(5)-(2'-Hydroxy-1'-naphthyl)pyrazoles. Journal of Organic Chemistry, 1995, 60, 3427-3439.	3.2	20
106	2-Arylperimidine derivatives. Part 1. Synthesis, NMR spectroscopy, X-ray crystal and molecular structures. Journal of the Chemical Society Perkin Transactions II, 1995, , 1389-1398.	0.9	25
107	Calorimetric quantification of the hydrogen-bond acidity of solvents and its relationship with solvent polarity. Journal of the Chemical Society Perkin Transactions II, 1995, , 2301-2305.	0.9	15
108	Acidity and Basicity of Indazole and its N-Methyl Derivatives in the Ground and in the Excited State. The Journal of Physical Chemistry, 1994, 98, 10606-10612.	2.9	68

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109	Influence of Lone Pair Repulsion vs Resonance Energy on the Relative Stabilities of Molecular Structures: A Theoretical Approach to the Equilibrium between 1H- and 2H-Benzotriazole Tautomers. Journal of Organic Chemistry, 1994, 59, 2799-2802.	3.2	69
110	Solvatochromism of Sterically Hindered Stilbazolium Betaines and Its Relationship to Reichardt's E_T(30) Scale: The Problem of the Measurement of the Polarity vs the Acidity of Alcohols. Chemische Berichte, 1993, 126, 2445-2448.	0.2	18
111	Experimental (13C and15N NMR spectroscopy) and theoretical (6-31G) study of the protonation ofN-methylazoles andN-methylbenzazoles. Magnetic Resonance in Chemistry, 1993, 31, 791-800.	1.9	45
112	N-aminoazoles. Part 2. Basicity and protonation site of N-aminoazoles: an experimental (pKa,13C and15N) Tj ETQq0 0 0 rgBT /Overlock 1 Transactions II, 1993, , 1687-1699.	0.9	46
113	Fluorescence of fullerenes (C60 and C70). Journal of the American Chemical Society, 1993, 115, 9249-9252.	13.7	105
114	Toward the photostability mechanism of intramolecular hydrogen bond systems. The photophysics of 1'-hydroxy-2'-acetonaphthone. Journal of the American Chemical Society, 1993, 115, 4321-4325.	13.7	101
115	Correction. Toward the Photostability Mechanism of Intramolecular Hydrogen Bond Systems. The Photophysics of 1'-Hydroxy-2'-acetonaphthone. Journal of the American Chemical Society, 1993, 115, 8519-8519.	13.7	1
116	Structure of benzotriazole in the gas phase: a UV experimental study. Journal of Organic Chemistry, 1993, 58, 5276-5277.	3.2	44
117	The role of the torsion of the phenyl moiety in the mechanism of stimulated ultraviolet light generation in 2â€phenylbenzazoles. Journal of Chemical Physics, 1992, 96, 2005-2016.	3.0	58
118	Photophysical properties of some 2-(2'-hydroxyaryl)benzotriazoles: dramatic effect of an ortho-located bulky tert-butyl group. Journal of the American Chemical Society, 1992, 114, 964-966.	13.7	36
119	Towards a solvent acidity scale: the calorimetry of the N-methyl imidazole probe. Journal of the Chemical Society Perkin Transactions II, 1992, , 1181-1185.	0.9	23
120	New ultraviolet stabilizers: 3- and 5-(2'-hydroxyphenyl)pyrazoles. Journal of the American Chemical Society, 1992, 114, 5039-5048.	13.7	139
121	Interaction of formamide with stilbazolium betaines: Steric effects in amides. Journal of Physical Organic Chemistry, 1992, 5, 609-613.	1.9	2
122	Effect ofN-methylation on the conformation of 2-phenylbenzimidazole, 2-phenylindole and three related rotationally constrained indoles. Magnetic Resonance in Chemistry, 1992, 30, 800-802.	1.9	5
123	Effect of the replacement of a methyl by a trifluoromethyl group on the acid-base properties of pyrazoles. Journal of Organic Chemistry, 1991, 56, 3942-3947.	3.2	35
124	Thermodynamic basicity vs. kinetic basicity of diazoles (imidazoles and pyrazoles). Journal of Organic Chemistry, 1991, 56, 179-183.	3.2	31
125	Analysis of the solvent effect on the photophysics properties of 6-propionyl-2-(dimethylamino)naphthalene (PRODAN). Journal of Fluorescence, 1991, 1, 215-223.	2.5	107
126	Lone-pair charges and structural effects. Journal of Physical Organic Chemistry, 1990, 3, 255-259.	1.9	1

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127	Photoinduced intramolecular proton transfer as the mechanism of ultraviolet stabilizers: a reappraisal. Journal of the American Chemical Society, 1990, 112, 747-759.	13.7	198
128	The structure of N-aminoazoles: an experimental (X-ray and15N NMR) and theoretical study. Journal of the Chemical Society Perkin Transactions II, 1990, , 237-244.	0.9	40
129	Toward a solvent basicity scale: the calorimetry of the pyrrole probe. Journal of the American Chemical Society, 1990, 112, 1678-1681.	13.7	72
130	The tautomerism of 1,2,3-triazole, 3(5)-methylpyrazole and their cations. Journal of Computational Chemistry, 1989, 10, 426-433.	3.3	51
131	Calorimetric study of the effect ofN-methylation in azoles: Loss of an active centre of solvation. Journal of Physical Organic Chemistry, 1989, 2, 646-652.	1.9	11
132	An AB initio comparative study of the electronic properties of sulfonamides and amides. Computational and Theoretical Chemistry, 1989, 184, 115-129.	1.5	34
133	Experimental and theoretical study of the R3P+-X- bond. Case of betaines derived from N-iminophosphoranes and alkyl isocyanates. Journal of the American Chemical Society, 1989, 111, 355-363.	13.7	58
134	Tautomerism and aromaticity in 1,2,3-triazoles: the case of benzotriazole. Journal of the American Chemical Society, 1989, 111, 7348-7353.	13.7	119
135	Polarizability effects on the aqueous solution basicity of substituted pyridines. Journal of Organic Chemistry, 1988, 53, 1137-1140.	3.2	48
136	Basicity and acidity of azoles: the annelation effect in azoles. Journal of the American Chemical Society, 1988, 110, 4105-4111.	13.7	127
137	Basicity and Acidity of Azoles. Advances in Heterocyclic Chemistry, 1987, , 187-274.	1.7	187
138	Electrostatic proximity effects in the relative basicities and acidities of pyrazole, imidazole, pyridazine, and pyrimidine. Journal of the American Chemical Society, 1986, 108, 3237-3239.	13.7	120
139	Basicity of azoles. VII. Basicity of Câ€aminopyrazoles in relation to tautomeric and protonation studies. Journal of Heterocyclic Chemistry, 1985, 22, 997-1000.	2.6	24
140	Basicity of azoles. IV. Empirical relationships between basicity and ionization potential for aromatic five membered rings containing nitrogen or oxygen. Journal of Heterocyclic Chemistry, 1984, 21, 269-270.	2.6	24
141	The Relative Basicities of Imidazole and Benzimidazole. Angewandte Chemie International Edition in English, 1983, 22, 323-324.	4.4	16
142	On the relative basicities of imidazole and benzimidazole. Angewandte Chemie International Edition in English, 1983, 22, 411-418.	4.4	1
143	Basicity of azoles. Part 2. Theoretical study of the basicity of methylpyrazoles and methylimidazoles. Journal of the Chemical Society Perkin Transactions II, 1983, , 1869-1874.	0.9	35
144	Ãœber die relative BasizitÃt von Imidazol und Benzimidazol. Angewandte Chemie, 1983, 95, 323-324.	2.0	2

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145	Prediction of proton affinities and protonation sites using a multivariate linear correlation. Journal of the Chemical Society Perkin Transactions II, 1982, , 1409-1418.	0.9	18
146	Intrinsic acidities of meta- and para-substituted phenols from calculated molecular properties. Journal of the Chemical Society Perkin Transactions II, 1979, , 1632-1636.	0.9	12
147	Prediction of proton affinities and preferred protonation sites in benzene derivatives, from 1s orbital energies. Journal of the Chemical Society Perkin Transactions II, 1979, , 1627-1631.	0.9	15
148	Theoretical study of the intramolecular hydrogen bonding in benzene derivatives. Advances in Molecular Relaxation and Interaction Processes, 1978, 12, 265-287.	0.5	25
149	Proton transfer in salicylic acid excited states. Human Development, 1976, 8, 87-94.	0.8	14
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