

Roger A Klein

List of Publications by Year in descending order

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Version: 2024-02-01

48
papers

3,980
citations

279798

23
h-index

214800

47
g-index

50
all docs

50
docs citations

50
times ranked

4350
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Definition of the hydrogen bond (IUPAC Recommendations 2011). <i>Pure and Applied Chemistry</i> , 2011, 83, 1637-1641. | 1.9 | 1,449 |
| 2 | Defining the hydrogen bond: An account (IUPAC Technical Report). <i>Pure and Applied Chemistry</i> , 2011, 83, 1619-1636. | 1.9 | 856 |
| 3 | What is a hydrogen bond? Mutually consistent theoretical and experimental criteria for characterizing H-bonding interactions. <i>Molecular Physics</i> , 2012, 110, 565-579. | 1.7 | 195 |
| 4 | Anti- π -Electrostatic Hydrogen Bonds. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 11214-11217. | 13.8 | 171 |
| 5 | Electron Density Topological Analysis of Hydrogen Bonding in Glucopyranose and Hydrated Glucopyranose. <i>Journal of the American Chemical Society</i> , 2002, 124, 13931-13937. | 13.7 | 110 |
| 6 | Ab initio conformational studies on diols and binary diol-water systems using DFT methods. Intramolecular hydrogen bonding and 1:1 complex formation with water. <i>Journal of Computational Chemistry</i> , 2002, 23, 585-599. | 3.3 | 104 |
| 7 | What is a hydrogen bond? Resonance covalency in the supramolecular domain. <i>Chemistry Education Research and Practice</i> , 2014, 15, 276-285. | 2.5 | 97 |
| 8 | Hydrogen bonding in diols and binary diol-water systems investigated using DFT methods. II. Calculated infrared OH-stretch frequencies, force constants, and NMR chemical shifts correlate with hydrogen bond geometry and electron density topology. A reevaluation of geometrical criteria for hydrogen bonding. <i>Journal of Computational Chemistry</i> , 2003, 24, 1120-1131. | 3.3 | 79 |
| 9 | Modified van der Waals atomic radii for hydrogen bonding based on electron density topology. <i>Chemical Physics Letters</i> , 2006, 425, 128-133. | 2.6 | 71 |
| 10 | Ab Initio Calculations of ^{17}O NMR-Chemical Shifts for Water. The Limits of PCM Theory and the Role of Hydrogen-Bond Geometry and Cooperativity. <i>Journal of Physical Chemistry A</i> , 2004, 108, 5851-5863. | 2.5 | 70 |
| 11 | Crystalline ethane-1,2-diol does not have intra-molecular hydrogen bonding: Experimental and theoretical charge density studies. <i>Journal of Molecular Structure</i> , 2010, 964, 126-133. | 3.6 | 66 |
| 12 | Mass spectrometric localization of carbon-carbon double bonds: A critical review of recent methods. <i>Chemistry and Physics of Lipids</i> , 1986, 39, 285-311. | 3.2 | 60 |
| 13 | Hydrogen bonding in strained cyclic vicinal diols: The birth of the hydrogen bond. <i>Chemical Physics Letters</i> , 2006, 429, 633-637. | 2.6 | 48 |
| 14 | Isolation and Structural Characterization of Sialic Acid-Containing Glycopeptides of the O-Glycosidic Type from the Urine of Two Patients with an Hereditary Deficiency in α -N-Acetylgalactosaminidase Activity. <i>Biological Chemistry Hoppe-Seyler</i> , 1989, 370, 661-672. | 1.4 | 46 |
| 15 | Improved General Understanding of the Hydrogen-Bonding Phenomena: A Reply. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 2600-2602. | 13.8 | 42 |
| 16 | MS and NMR analysis of the cross-reacting determinant glycan from <i>Trypanosoma brucei brucei</i> MITat 1.6 variant specific glycoprotein. <i>Biochemical and Biophysical Research Communications</i> , 1987, 146, 1055-1063. | 2.1 | 38 |
| 17 | UNIQUAC interaction parameters for molecules with -OH group on adjacent carbon atoms in aqueous solution determined by molecular mechanics α -glycols, glycerol and glucose. <i>Fluid Phase Equilibria</i> , 1997, 132, 117-137. | 2.5 | 35 |
| 18 | Thermodynamics and membrane processes. <i>Quarterly Reviews of Biophysics</i> , 1982, 15, 667-759. | 5.7 | 28 |

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|----|--|-----|-----------|
| 19 | Threonine as a Preferred Source of 2-Carbon Units for Lipid Synthesis in <i>Trypanosoma brucei</i> . <i>Biochemical Society Transactions</i> , 1976, 4, 48-50. | 3.4 | 27 |
| 20 | The aqueous solution structure of a lipoteichoic acid from <i>Streptococcus pneumoniae</i> strain R6 containing 2,4-diamino-2,4,6-trideoxy-galactose: evidence for conformational mobility of the galactopyranose ring. <i>Carbohydrate Research</i> , 1996, 281, 79-98. | 2.3 | 26 |
| 21 | The use of 13-methyltetradecanoic acid as an indicator of adipose tissue turnover. <i>Lipids</i> , 1980, 15, 572-579. | 1.7 | 24 |
| 22 | Pople versus Dunning basis-sets for group IA metal hydrides and some other second row hydrides: The case against a De Facto standard. <i>Chemical Physics Letters</i> , 2006, 419, 254-258. | 2.6 | 24 |
| 23 | Lack of intramolecular hydrogen bonding in glucopyranose: Vicinal hydroxyl groups exhibit negative cooperativity. <i>Chemical Physics Letters</i> , 2006, 433, 165-169. | 2.6 | 24 |
| 24 | The critical role of force-fields in property prediction. <i>New Journal of Chemistry</i> , 1999, 23, 153-164. | 2.8 | 22 |
| 25 | UNIQUAC interaction parameters for alkane/amine systems determined by Molecular Mechanics. <i>Fluid Phase Equilibria</i> , 1996, 115, 59-72. | 2.5 | 21 |
| 26 | Reporter resonances in the NMR spectra of oligosaccharides containing sialic acid linked to galactopyranose rings. <i>Carbohydrate Research</i> , 1994, 254, 289-294. | 2.3 | 19 |
| 27 | The aqueous solution structure of the tetrasaccharide-ribitol repeat-unit from the lipoteichoic acid of <i>Streptococcus pneumoniae</i> strain R6 determined using a combination of NMR spectroscopy and computer calculations. <i>Carbohydrate Research</i> , 1994, 256, 189-222. | 2.3 | 18 |
| 28 | Improved General Understanding of the Hydrogen Bonding Phenomena: A Reply. <i>Angewandte Chemie</i> , 2015, 127, 2636-2638. | 2.0 | 17 |
| 29 | Barriers to internal rotation around the C-N bond in 3-(o-aryl)-5-methyl-rhodanines using NMR spectroscopy and computational studies. Electron density topological analysis of the transition states. <i>Organic and Biomolecular Chemistry</i> , 2004, 2, 2426-2436. | 2.8 | 14 |
| 30 | The Enzymic Hydrolysis of Acetyl-Coenzyme A by Trypanosomatid Flagellates. <i>Biochemical Society Transactions</i> , 1976, 4, 285-287. | 3.4 | 13 |
| 31 | A study of the membrane attachment site of the membrane-form variant surface glycoprotein from <i>Trypanosoma brucei brucei</i> using lipid vesicles as a model of the plasma membrane. <i>Molecular and Biochemical Parasitology</i> , 1986, 20, 191-197. | 1.1 | 13 |
| 32 | The role of threonine in the metabolism of acetyl coenzyme A by <i>Trypanosoma brucei brucei</i> . <i>Comparative Biochemistry and Physiology Part B: Comparative Biochemistry</i> , 1983, 74, 277-281. | 0.2 | 12 |
| 33 | Bromoacetyl-L-carnitine: Biochemical and antitrypanosomal actions against <i>Trypanosoma brucei brucei</i> . <i>Biochemical Pharmacology</i> , 1983, 32, 3447-3451. | 4.4 | 11 |
| 34 | Binary Diol-Water Systems Studied by ¹⁷ O Nuclear Magnetic Resonance Spectroscopy. Interpretation of the Effect of Diol Structure on ¹⁷ O-Water Chemical Shift. Formation of Networks of Water Molecules Stabilized by Weak C-H...O Interactions. <i>Journal of Physical Chemistry A</i> , 2001, 105, 9298-9304. | 2.5 | 10 |
| 35 | Carnitine stimulates ATP synthesis in <i>Trypanosoma brucei brucei</i> . <i>FEBS Letters</i> , 1982, 141, 271-274. | 2.8 | 8 |
| 36 | Pyruvate kinase: A carnitine regulated site of ATP production in <i>Trypanosoma brucei brucei</i> . <i>Comparative Biochemistry and Physiology Part B: Comparative Biochemistry</i> , 1984, 78, 595-599. | 0.2 | 8 |

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|----|---|-----|-----------|
| 37 | Characterization of DNA and protein synthesis in <i>Trypanosoma brucei</i> by using inhibitors. <i>Biochemical Society Transactions</i> , 1983, 11, 366-367. | 3.4 | 6 |
| 38 | Double bond location in fatty acids. A critical analysis of the feasibility of using specifically deuterated pyrrolidides for mass spectral analysis. <i>Biological Mass Spectrometry</i> , 1986, 13, 429-437. | 0.5 | 6 |
| 39 | Lectin interactions with the variant surface glycoprotein from <i>Trypanosoma brucei brucei</i> incorporated into liposomes. <i>Biochemical and Biophysical Research Communications</i> , 1986, 141, 1274-1278. | 2.1 | 4 |
| 40 | Threonine uptake in <i>Trypanosoma brucei</i> . <i>Molecular and Biochemical Parasitology</i> , 1984, 11, 215-223. | 1.1 | 3 |
| 41 | Single point calculations using geometries derived at a lower level of theory: Caveats to be observed in using compound levels of theory or "double-barrelling". <i>Chemical Physics Letters</i> , 2006, 423, 413-416. | 2.6 | 3 |
| 42 | Re-examination of the Solvolysis of Acyl Chlorides. Rethinking the Role of Both Carbonyl and Chloride. <i>Journal of Physical Chemistry A</i> , 2004, 108, 9328-9330. | 2.5 | 2 |
| 43 | Stereospecificity of the threonine dehydrogenase from bloodstream <i>Trypanosoma brucei</i> . <i>Comparative Biochemistry and Physiology Part B: Comparative Biochemistry</i> , 1980, 66, 143-146. | 0.2 | 1 |
| 44 | Aspects of conformational mobility in charged oligosaccharides. <i>Computational and Theoretical Chemistry</i> , 1997, 395-396, 313-332. | 1.5 | 1 |
| 45 | Reply to the comment on "Pople versus Dunning basis-sets for group IA metal hydrides and some other second row hydrides: The case against a De Facto standard" by R.A. Klein and M.A. Zottola [<i>Chem. Phys. Lett.</i> 419 (2006) 254-258]. <i>Chemical Physics Letters</i> , 2006, 430, 464-465. | 2.6 | 1 |
| 46 | The Movement of Molecules Across Membranes: The Thermodynamic Analysis of the Dependence on Structure, Pressure, and Temperature. , 1983, , 301-317. | | 1 |
| 47 | The measurement of motility in micro-organisms and its biochemical significance. <i>Biochemical Society Transactions</i> , 1984, 12, 627-630. | 3.4 | 0 |
| 48 | The trypanocidal activity of bromoacetyl-L-carnitine and \pm -difluoromethylornithine in <i>Trypanosoma brucei brucei</i> infected rats. <i>Biochemical Society Transactions</i> , 1986, 14, 1093-1094. | 3.4 | 0 |