

Richard F W Bader

List of Publications by Year in descending order

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69
papers

17,411
citations

43973

48
h-index

98622

67
g-index

71
all docs

71
docs citations

71
times ranked

9702
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | A quantum theory of molecular structure and its applications. <i>Chemical Reviews</i> , 1991, 91, 893-928. | 23.0 | 5,550 |
| 2 | Calculation of the average properties of atoms in molecules. II. <i>Journal of Computational Chemistry</i> , 1982, 3, 317-328. | 1.5 | 1,317 |
| 3 | Properties of atoms in molecules: atomic volumes. <i>Journal of the American Chemical Society</i> , 1987, 109, 7968-7979. | 6.6 | 1,120 |
| 4 | The Lewis Model and Beyond. <i>Journal of Physical Chemistry A</i> , 1999, 103, 304-314. | 1.1 | 944 |
| 5 | Hydrogenâ€“Hydrogen Bonding: A Stabilizing Interaction in Molecules and Crystals. <i>Chemistry - A European Journal</i> , 2003, 9, 1940-1951. | 1.7 | 688 |
| 6 | Bond Paths Are Not Chemical Bonds. <i>Journal of Physical Chemistry A</i> , 2009, 113, 10391-10396. | 1.1 | 544 |
| 7 | Identifying and Analyzing Intermolecular Bonding Interactions in van der Waals Moleculesâ€“. <i>The Journal of Physical Chemistry</i> , 1996, 100, 10892-10911. | 2.9 | 411 |
| 8 | VIBRATIONALLY INDUCED PERTURBATIONS IN MOLECULAR ELECTRON DISTRIBUTIONS. <i>Canadian Journal of Chemistry</i> , 1962, 40, 1164-1175. | 0.6 | 343 |
| 9 | Theoretical analysis of hydrocarbon properties. 1. Bonds, structures, charge concentrations, and charge relaxations. <i>Journal of the American Chemical Society</i> , 1987, 109, 985-1001. | 6.6 | 298 |
| 10 | Electron Delocalization and the Fermi Hole. <i>Journal of the American Chemical Society</i> , 1996, 118, 4959-4965. | 6.6 | 271 |
| 11 | Topological analysis of magnetically induced molecular current distributions. <i>Journal of Chemical Physics</i> , 1993, 99, 3669-3682. | 1.2 | 264 |
| 12 | Theoretical analysis of hydrocarbon properties. 2. Additivity of group properties and the origin of strain energy. <i>Journal of the American Chemical Society</i> , 1987, 109, 1001-1012. | 6.6 | 248 |
| 13 | Pauli Repulsions Exist Only in the Eye of the Beholder. <i>Chemistry - A European Journal</i> , 2006, 12, 2896-2901. | 1.7 | 240 |
| 14 | A physical basis for the VSEPR model of molecular geometry. <i>Journal of the American Chemical Society</i> , 1988, 110, 7329-7336. | 6.6 | 228 |
| 15 | Toward a theory of chemical reactivity based on the charge density. <i>Journal of the American Chemical Society</i> , 1985, 107, 6788-6795. | 6.6 | 210 |
| 16 | The Quantum Mechanical Basis of Conceptual Chemistry. <i>Monatshefte FÃ¼r Chemie</i> , 2005, 136, 819-854. | 0.9 | 210 |
| 17 | Atomic Charges Are Measurable Quantum Expectation Values: A Rebuttal of Criticisms of QTAIM Charges. <i>Journal of Physical Chemistry A</i> , 2004, 108, 8385-8394. | 1.1 | 209 |
| 18 | The Electron Pair. <i>The Journal of Physical Chemistry</i> , 1996, 100, 15398-15415. | 2.9 | 206 |

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|----|--|-----|-----------|
| 19 | Molecular fragments or chemical bonds. <i>Accounts of Chemical Research</i> , 1975, 8, 34-40. | 7.6 | 202 |
| 20 | Quantum Topology: Theory of Molecular Structure and its Change. <i>Israel Journal of Chemistry</i> , 1980, 19, 8-29. | 1.0 | 184 |
| 21 | Quantum topology of molecular charge distributions. II. Molecular structure and its change. <i>Journal of Chemical Physics</i> , 1979, 70, 4316-4329. | 1.2 | 183 |
| 22 | Definition of Molecular Structure: By Choice or by Appeal to Observation?. <i>Journal of Physical Chemistry A</i> , 2010, 114, 7431-7444. | 1.1 | 176 |
| 23 | Bonding to Titanium. <i>Inorganic Chemistry</i> , 2001, 40, 5603-5611. | 1.9 | 156 |
| 24 | Properties of atoms in molecules: Atomic polarizabilities. <i>Journal of Chemical Physics</i> , 1990, 93, 7213-7224. | 1.2 | 155 |
| 25 | Properties of atoms in molecules: Magnetic susceptibilities. <i>Journal of Chemical Physics</i> , 1993, 99, 3683-3693. | 1.2 | 149 |
| 26 | Origin of rotation and inversion barriers. <i>Journal of the American Chemical Society</i> , 1990, 112, 6530-6536. | 6.6 | 148 |
| 27 | Where To Draw the Line in Defining a Molecular Structure. <i>Organometallics</i> , 2004, 23, 6253-6263. | 1.1 | 145 |
| 28 | Properties of Atoms in Molecules: \hat{A} Group Additivity. <i>Journal of Physical Chemistry A</i> , 2000, 104, 5579-5589. | 1.1 | 125 |
| 29 | Properties of atoms in molecules: electrophilic aromatic substitution. <i>The Journal of Physical Chemistry</i> , 1989, 93, 2946-2956. | 2.9 | 122 |
| 30 | Properties of Atoms in Molecules: \hat{A} Caged Atoms and the Ehrenfest Force. <i>Journal of Chemical Theory and Computation</i> , 2005, 1, 403-414. | 2.3 | 119 |
| 31 | An Experimentalist's Reply to "What Is an Atom in a Molecule?". <i>Journal of Physical Chemistry A</i> , 2006, 110, 6365-6371. | 1.1 | 119 |
| 32 | Towards the development of the quantum mechanics of a subspace. <i>Journal of Chemical Physics</i> , 1975, 63, 3945. | 1.2 | 117 |
| 33 | Core Distortions and Geometries of the Difluorides and Dihydrides of Ca, Sr, and Ba. <i>Inorganic Chemistry</i> , 1995, 34, 2407-2414. | 1.9 | 114 |
| 34 | The nature of the structure difference between light and heavy water and the origin of the solvent isotope effect. <i>Tetrahedron</i> , 1960, 10, 182-199. | 1.0 | 112 |
| 35 | Properties of Atoms in Molecules: \hat{A} Atoms Forming Molecules. <i>Journal of Physical Chemistry A</i> , 2000, 104, 1779-1794. | 1.1 | 112 |
| 36 | The zero-flux surface and the topological and quantum definitions of an atom in a molecule. <i>Theoretical Chemistry Accounts</i> , 2001, 105, 276-283. | 0.5 | 102 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 37 | Geometry of the Fluorides, Oxofluorides, Hydrides, and Methanides of Vanadium(V), Chromium(VI), and Molybdenum(VI): A Understanding the Geometry of Non-VSEPR Molecules in Terms of Core Distortion. <i>Inorganic Chemistry</i> , 1996, 35, 3954-3963. | 1.9 | 98 |
| 38 | Atoms in molecules as non-overlapping, bounded, space-filling open quantum systems. <i>Foundations of Chemistry</i> , 2013, 15, 253-276. | 0.4 | 94 |
| 39 | Effect of electron correlation on the topological properties of molecular charge distributions. <i>Journal of Chemical Physics</i> , 1988, 88, 3792-3804. | 1.2 | 93 |
| 40 | Subspace quantum mechanics and the variational principle. <i>Journal of Chemical Physics</i> , 1978, 68, 3667-3679. | 1.2 | 92 |
| 41 | Quantum topology. IV. Relation between the topological and energetic stabilities of molecular structures. <i>Journal of Chemical Physics</i> , 1981, 74, 5162-5167. | 1.2 | 89 |
| 42 | Everyman's Derivation of the Theory of Atoms in Molecules. <i>Journal of Physical Chemistry A</i> , 2007, 111, 7966-7972. | 1.1 | 76 |
| 43 | Subspace quantum dynamics and the quantum action principle. <i>Journal of Chemical Physics</i> , 1978, 68, 3680-3691. | 1.2 | 75 |
| 44 | Atoms-in-molecules study of the genetically encoded amino acids. III. Bond and atomic properties and their correlations with experiment including mutation-induced changes in protein stability and genetic coding. <i>Proteins: Structure, Function and Bioinformatics</i> , 2003, 52, 360-399. | 1.5 | 75 |
| 45 | Structural homeomorphism between the electronic charge density and the nuclear potential of a molecular system. <i>Physical Review A</i> , 1980, 21, 1-11. | 1.0 | 69 |
| 46 | Proton Spin-Spin Coupling and Electron Delocalization. <i>Journal of Physical Chemistry A</i> , 2002, 106, 7369-7375. | 1.1 | 59 |
| 47 | Forces in molecules. <i>Faraday Discussions</i> , 2007, 135, 79-95. | 1.6 | 57 |
| 48 | Atoms in molecules in external fields. <i>Journal of Chemical Physics</i> , 1989, 91, 6989-7001. | 1.2 | 54 |
| 49 | On the non-existence of parallel universes in chemistry. <i>Foundations of Chemistry</i> , 2011, 13, 11-37. | 0.4 | 49 |
| 50 | Trigonal Bipyramidal and Related Molecules of the Main Group Elements: Investigation of Apparent Exceptions to the VSEPR Model through the Analysis of the Laplacian of the Electron Density. <i>Inorganic Chemistry</i> , 1994, 33, 2115-2121. | 1.9 | 46 |
| 51 | Properties of atoms in crystals: Dielectric polarization. <i>International Journal of Quantum Chemistry</i> , 2001, 85, 592-607. | 1.0 | 46 |
| 52 | Recognizing a triple bond between main group atoms. <i>Theoretical Chemistry Accounts</i> , 2001, 105, 365-373. | 0.5 | 43 |
| 53 | Atoms-in-molecules study of the genetically encoded amino acids. II. Computational study of molecular geometries. <i>Proteins: Structure, Function and Bioinformatics</i> , 2002, 48, 519-538. | 1.5 | 42 |
| 54 | Atomic properties and the reactivity of carbenes. <i>Canadian Journal of Chemistry</i> , 1986, 64, 1496-1508. | 0.6 | 41 |

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|----|---|-----|-----------|
| 55 | Comment on the Comparative Use of the Electron Density and Its Laplacian. Chemistry - A European Journal, 2006, 12, 7769-7772. | 1.7 | 39 |
| 56 | Nearsightedness of Electronic Matter As Seen by a Physicist and a Chemist. Journal of Physical Chemistry A, 2008, 112, 13717-13728. | 1.1 | 39 |
| 57 | Properties of atoms in molecules: dipole moments and substituent effects in ethyl and carbonyl compounds. The Journal of Physical Chemistry, 1988, 92, 6219-6227. | 2.9 | 38 |
| 58 | Extending the VSEPR model through the properties of the Laplacian of the charge density. Canadian Journal of Chemistry, 1989, 67, 1842-1846. | 0.6 | 36 |
| 59 | Chemistry and the near-sighted nature of the one-electron density matrix. International Journal of Quantum Chemistry, 1995, 56, 409-419. | 1.0 | 30 |
| 60 | QTAIM Study on the Degenerate Cope Rearrangements of 1,5-Hexadiene and Semibullvalene. Journal of Physical Chemistry A, 2009, 113, 3254-3265. | 1.1 | 28 |
| 61 | Worlds Apart in Chemistry: A Personal Tribute to J. C. Slater. Journal of Physical Chemistry A, 2011, 115, 12667-12676. | 1.1 | 28 |
| 62 | Why define atoms in real space?. International Journal of Quantum Chemistry, 1994, 49, 299-308. | 1.0 | 26 |
| 63 | The Mechanism of Ethylene Oxide Formation from 2-Chloroethanol ^{1,2} . Journal of the American Chemical Society, 1959, 81, 2353-2359. | 6.6 | 24 |
| 64 | THE USE OF THE HELLMANN-“FEYNMAN THEOREM TO CALCULATE MOLECULAR ENERGIES. Canadian Journal of Chemistry, 1960, 38, 2117-2127. | 0.6 | 21 |
| 65 | Energy additivity in branched and cyclic hydrocarbons. Canadian Journal of Chemistry, 2009, 87, 1583-1592. | 0.6 | 7 |
| 66 | The Lagrangian Approach to Chemistry. , 0, , 35-59. | | 5 |
| 67 | Atomic and Group Properties in the Alkanes. , 0, , 1-77. | | 3 |
| 68 | Atoms in Medicinal Chemistry. Methods and Principles in Medicinal Chemistry, 2005, , 199-231. | 0.3 | 2 |
| 69 | Topology of Electron Density and Open Quantum Systems. NATO ASI Series Series B: Physics, 1995, , 237-272. | 0.2 | 1 |