

Richard F W Bader

List of Publications by Year in descending order

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69
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43973

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71
times ranked

9702
citing authors

#	ARTICLE	IF	CITATIONS
1	Atoms in molecules as non-overlapping, bounded, space-filling open quantum systems. <i>Foundations of Chemistry</i> , 2013, 15, 253-276.	0.4	94
2	Worlds Apart in Chemistry: A Personal Tribute to J. C. Slater. <i>Journal of Physical Chemistry A</i> , 2011, 115, 12667-12676.	1.1	28
3	On the non-existence of parallel universes in chemistry. <i>Foundations of Chemistry</i> , 2011, 13, 11-37.	0.4	49
4	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
5	Energy additivity in branched and cyclic hydrocarbons. <i>Canadian Journal of Chemistry</i> , 2009, 87, 1583-1592.	0.6	7
6	QTAIM Study on the Degenerate Cope Rearrangements of 1,5-Hexadiene and Semibullvalene. <i>Journal of Physical Chemistry A</i> , 2009, 113, 3254-3265.	1.1	28
7	Bond Paths Are Not Chemical Bonds. <i>Journal of Physical Chemistry A</i> , 2009, 113, 10391-10396.	1.1	544
8	Nearsightedness of Electronic Matter As Seen by a Physicist and a Chemist. <i>Journal of Physical Chemistry A</i> , 2008, 112, 13717-13728.	1.1	39
9	Forces in molecules. <i>Faraday Discussions</i> , 2007, 135, 79-95.	1.6	57
10	Everyman's Derivation of the Theory of Atoms in Molecules. <i>Journal of Physical Chemistry A</i> , 2007, 111, 7966-7972.	1.1	76
11	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
12	Pauli Repulsions Exist Only in the Eye of the Beholder. <i>Chemistry - A European Journal</i> , 2006, 12, 2896-2901.	1.7	240
13	Comment on the Comparative Use of the Electron Density and Its Laplacian. <i>Chemistry - A European Journal</i> , 2006, 12, 7769-7772.	1.7	39
14	Atoms in Medicinal Chemistry. <i>Methods and Principles in Medicinal Chemistry</i> , 2005, , 199-231.	0.3	2
15	The Quantum Mechanical Basis of Conceptual Chemistry. <i>Monatshefte für Chemie</i> , 2005, 136, 819-854.	0.9	210
16	Properties of Atoms in Molecules: Caged Atoms and the Ehrenfest Force. <i>Journal of Chemical Theory and Computation</i> , 2005, 1, 403-414.	2.3	119
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	Where To Draw the Line in Defining a Molecular Structure. <i>Organometallics</i> , 2004, 23, 6253-6263.	1.1	145

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19	Hydrogenâ€“Hydrogen Bonding: A Stabilizing Interaction in Molecules and Crystals. Chemistry - A European Journal, 2003, 9, 1940-1951.	1.7	688
20	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. Proteins: Structure, Function and Bioinformatics, 2003, 52, 360-399.	1.5	75
21	Proton Spinâˆ“Spin Coupling and Electron Delocalization. Journal of Physical Chemistry A, 2002, 106, 7369-7375.	1.1	59
22	Atoms-in-molecules study of the genetically encoded amino acids. II. Computational study of molecular geometries. Proteins: Structure, Function and Bioinformatics, 2002, 48, 519-538.	1.5	42
23	Bonding to Titanium. Inorganic Chemistry, 2001, 40, 5603-5611.	1.9	156
24	Recognizing a triple bond between main group atoms. Theoretical Chemistry Accounts, 2001, 105, 365-373.	0.5	43
25	The zero-flux surface and the topological and quantum definitions of an atom in a molecule. Theoretical Chemistry Accounts, 2001, 105, 276-283.	0.5	102
26	Properties of atoms in crystals: Dielectric polarization. International Journal of Quantum Chemistry, 2001, 85, 592-607.	1.0	46
27	Properties of Atoms in Molecules:Â Atoms Forming Molecules. Journal of Physical Chemistry A, 2000, 104, 1779-1794.	1.1	112
28	Properties of Atoms in Molecules:Â Group Additivity. Journal of Physical Chemistry A, 2000, 104, 5579-5589.	1.1	125
29	The Lewis Model and Beyond. Journal of Physical Chemistry A, 1999, 103, 304-314.	1.1	944
30	Electron Delocalization and the Fermi Hole. Journal of the American Chemical Society, 1996, 118, 4959-4965.	6.6	271
31	Identifying and Analyzing Intermolecular Bonding Interactions in van der Waals Moleculesâ€. The Journal of Physical Chemistry, 1996, 100, 10892-10911.	2.9	411
32	Geometry of the Fluorides, Oxofluorides, Hydrides, and Methanides of Vanadium(V), Chromium(VI), and Molybdenum(VI):Â Understanding the Geometry of Non-VSEPR Molecules in Terms of Core Distortion. Inorganic Chemistry, 1996, 35, 3954-3963.	1.9	98
33	The Electron Pair. The Journal of Physical Chemistry, 1996, 100, 15398-15415.	2.9	206
34	Chemistry and the near-sighted nature of the one-electron density matrix. International Journal of Quantum Chemistry, 1995, 56, 409-419.	1.0	30
35	Core Distortions and Geometries of the Difluorides and Dihydrides of Ca, Sr, and Ba. Inorganic Chemistry, 1995, 34, 2407-2414.	1.9	114
36	Topology of Electron Density and Open Quantum Systems. NATO ASI Series Series B: Physics, 1995, , 237-272.	0.2	1

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37	Why define atoms in real space?. International Journal of Quantum Chemistry, 1994, 49, 299-308.	1.0	26
38	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. Inorganic Chemistry, 1994, 33, 2115-2121.	1.9	46
39	Properties of atoms in molecules: Magnetic susceptibilities. Journal of Chemical Physics, 1993, 99, 3683-3693.	1.2	149
40	Topological analysis of magnetically induced molecular current distributions. Journal of Chemical Physics, 1993, 99, 3669-3682.	1.2	264
41	A quantum theory of molecular structure and its applications. Chemical Reviews, 1991, 91, 893-928.	23.0	5,550
42	Origin of rotation and inversion barriers. Journal of the American Chemical Society, 1990, 112, 6530-6536.	6.6	148
43	Properties of atoms in molecules: Atomic polarizabilities. Journal of Chemical Physics, 1990, 93, 7213-7224.	1.2	155
44	Atoms in molecules in external fields. Journal of Chemical Physics, 1989, 91, 6989-7001.	1.2	54
45	Properties of atoms in molecules: electrophilic aromatic substitution. The Journal of Physical Chemistry, 1989, 93, 2946-2956.	2.9	122
46	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
47	A physical basis for the VSEPR model of molecular geometry. Journal of the American Chemical Society, 1988, 110, 7329-7336.	6.6	228
48	Effect of electron correlation on the topological properties of molecular charge distributions. Journal of Chemical Physics, 1988, 88, 3792-3804.	1.2	93
49	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
50	Properties of atoms in molecules: atomic volumes. Journal of the American Chemical Society, 1987, 109, 7968-7979.	6.6	1,120
51	Theoretical analysis of hydrocarbon properties. 1. Bonds, structures, charge concentrations, and charge relaxations. Journal of the American Chemical Society, 1987, 109, 985-1001.	6.6	298
52	Theoretical analysis of hydrocarbon properties. 2. Additivity of group properties and the origin of strain energy. Journal of the American Chemical Society, 1987, 109, 1001-1012.	6.6	248
53	Atomic properties and the reactivity of carbenes. Canadian Journal of Chemistry, 1986, 64, 1496-1508.	0.6	41
54	Toward a theory of chemical reactivity based on the charge density. Journal of the American Chemical Society, 1985, 107, 6788-6795.	6.6	210

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55	Calculation of the average properties of atoms in molecules. II. Journal of Computational Chemistry, 1982, 3, 317-328.	1.5	1,317
56	Quantum topology. IV. Relation between the topological and energetic stabilities of molecular structures. Journal of Chemical Physics, 1981, 74, 5162-5167.	1.2	89
57	Structural homeomorphism between the electronic charge density and the nuclear potential of a molecular system. Physical Review A, 1980, 21, 1-11.	1.0	69
58	Quantum Topology: Theory of Molecular Structure and its Change. Israel Journal of Chemistry, 1980, 19, 8-29.	1.0	184
59	Quantum topology of molecular charge distributions. II. Molecular structure and its change. Journal of Chemical Physics, 1979, 70, 4316-4329.	1.2	183
60	Subspace quantum dynamics and the quantum action principle. Journal of Chemical Physics, 1978, 68, 3680-3691.	1.2	75
61	Subspace quantum mechanics and the variational principle. Journal of Chemical Physics, 1978, 68, 3667-3679.	1.2	92
62	Towards the development of the quantum mechanics of a subspace. Journal of Chemical Physics, 1975, 63, 3945.	1.2	117
63	Molecular fragments or chemical bonds. Accounts of Chemical Research, 1975, 8, 34-40.	7.6	202
64	VIBRATIONALLY INDUCED PERTURBATIONS IN MOLECULAR ELECTRON DISTRIBUTIONS. Canadian Journal of Chemistry, 1962, 40, 1164-1175.	0.6	343
65	The nature of the structure difference between light and heavy water and the origin of the solvent isotope effect. Tetrahedron, 1960, 10, 182-199.	1.0	112
66	THE USE OF THE HELLMANN-FEYNMAN THEOREM TO CALCULATE MOLECULAR ENERGIES. Canadian Journal of Chemistry, 1960, 38, 2117-2127.	0.6	21
67	The Mechanism of Ethylene Oxide Formation from 2-Chloroethanol. Journal of the American Chemical Society, 1959, 81, 2353-2359.	6.6	24
68	Atomic and Group Properties in the Alkanes. , 0, , 1-77.		3
69	The Lagrangian Approach to Chemistry. , 0, , 35-59.		5