

Amnon Stanger

List of Publications by Year in descending order

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70
papers

3,535
citations

186209

28
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58
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73
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docs citations

73
times ranked

2275
citing authors

#	ARTICLE	IF	CITATIONS
1	Nucleus-Independent Chemical Shifts (NICS): Distance Dependence and Revised Criteria for Aromaticity and Antiaromaticity. <i>Journal of Organic Chemistry</i> , 2006, 71, 883-893.	1.7	636
2	Magnetic criteria of aromaticity. <i>Chemical Society Reviews</i> , 2015, 44, 6597-6615.	18.7	354
3	The NICS-XY-Scan: Identification of Local and Global Ring Currents in Multi-Ring Systems. <i>Chemistry - A European Journal</i> , 2014, 20, 5673-5688.	1.7	252
4	What is aromaticity: a critique of the concept of aromaticity "can it really be defined?". <i>Chemical Communications</i> , 2009, , 1939.	2.2	186
5	Obtaining Relative Induced Ring Currents Quantitatively from NICS. <i>Journal of Organic Chemistry</i> , 2010, 75, 2281-2288.	1.7	146
6	Is the Mills Nixon effect real?. <i>Journal of the American Chemical Society</i> , 1991, 113, 8277-8280.	6.6	131
7	The Impact of Antiaromatic Subunits in [4n+2] π -Systems: Bispentalenes with [4n+2] π -Electron Perimeters and Antiaromatic Character. <i>Journal of the American Chemical Society</i> , 2015, 137, 7178-7188.	6.6	115
8	Novel route to carbon-silicon double bonds via a Peterson-type reaction. <i>Organometallics</i> , 1992, 11, 2326-2328.	1.1	102
9	Can Substituted Cyclopentadiene Become Aromatic or Antiaromatic?. <i>Chemistry - A European Journal</i> , 2006, 12, 2745-2751.	1.7	100
10	NICS " Past and Present. <i>European Journal of Organic Chemistry</i> , 2020, 2020, 3120-3127.	1.2	85
11	Cyclopropenylcarbinol Derivatives as New Versatile Intermediates in Organic Synthesis: Application to the Formation of Enantiomerically Pure Alkylidenecyclopropane Derivatives. <i>Chemistry - A European Journal</i> , 2009, 15, 8449-8464.	1.7	67
12	Strain-Induced Bond Localization. The Heteroatom Case. <i>Journal of the American Chemical Society</i> , 1998, 120, 12034-12040.	6.6	64
13	Are carbenium ions stabilized or destabilized by .alpha.-silyl substitution? The solvolysis of 2-(trimethylsilyl)-2-adamantyl p-nitrobenzoate. <i>Journal of the American Chemical Society</i> , 1985, 107, 2806-2807.	6.6	61
14	En Route to Archimedene: Total Synthesis of C _{3h} -Symmetric [7]Phenylene. <i>Organic Letters</i> , 2003, 5, 549-552.	2.4	59
15	Enantio- and Diastereoselective Tandem Zn-Promoted Brook Rearrangement/Ene Allene Carbocyclization Reaction. <i>Organic Letters</i> , 2009, 11, 1853-1856.	2.4	57
16	The first demonstration of solvolytic generation of a simple silicenium ion (R ₃ Si ⁺). Access via 1,2-methyl migration. <i>Journal of the American Chemical Society</i> , 1987, 109, 272-273.	6.6	54
17	The origin of the symmetrical structure of benzene. Is the .sigma. or the .pi. frame responsible? An ab-initio study of the effect of HCC bond angle distortion. <i>Journal of Organic Chemistry</i> , 1988, 53, 4889-4890.	1.7	49
18	An Excursion from Normal to Inverted C-C Bonds Shows a Clear Demarcation between Covalent and Charge-Shift C-C Bonds. <i>ChemPhysChem</i> , 2009, 10, 2658-2669.	1.0	46

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19	A Nickel-Anthracene Complex Having η^3 and η^4 Coordination in One Crystal. <i>Angewandte Chemie International Edition in English</i> , 1993, 32, 1475-1477.	4.4	42
20	The anomeric effect at silicon. <i>Journal of Organometallic Chemistry</i> , 1988, 346, 305-313.	0.8	40
21	Is the haptotropic rearrangement in bis(tributylphosphine)(anthracene)nickel inter- or intramolecular? Determination of the molecularity by a spin saturation transfer approach. <i>Organometallics</i> , 1991, 10, 2979-2982.	1.1	40
22	Synthesis and fluxional behavior of [bis(trialkylphosphine)nickel]anthracene (alkyl = Et, Bu). <i>Organometallics</i> , 1992, 11, 317-320.	1.1	40
23	Reexamination of NICS _{zz} : Height Dependence, Off-Center Values, and Integration. <i>Journal of Physical Chemistry A</i> , 2019, 123, 3922-3927.	1.1	40
24	The predictive power of aromaticity: quantitative correlation between aromaticity and ionization potentials and HOMO-LUMO gaps in oligomers of benzene, pyrrole, furan, and thiophene. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 14808-14817.	1.3	38
25	The crystal structures of (R ₃ P) ₂ Ni-anthracene (R = Et, Bu). <i>Journal of Organometallic Chemistry</i> , 1992, 430, 235-243.	0.8	35
26	Strain induced bond localization in strained aromatic compounds with extended π systems. <i>Journal of Computational Chemistry</i> , 2001, 22, 1377-1386.	1.5	34
27	α -Silicon-substituted vinyl cations. A theoretical ab initio investigation. <i>Journal of Organic Chemistry</i> , 1982, 47, 1462-1468.	1.7	32
28	Hexabromotricyclobutabenzene and Hexabromohexaradialene: Their Nickel-Mediated One-Pot Syntheses and Crystal Structure. <i>Chemistry - A European Journal</i> , 1997, 3, 208-211.	1.7	30
29	Structure, Deformation Electron Densities, Photoelectron Spectra, and Reactivity of 3,4-Dihydro-1H-cyclobuta[a]cyclopropa[d]benzene. <i>Angewandte Chemie International Edition in English</i> , 1989, 28, 206-208.	4.4	29
30	Double-Stranded Cycles: Toward C ₈₄ 's Belt Region. <i>Journal of Organic Chemistry</i> , 2007, 72, 424-430.	1.7	29
31	Bimolecular substitution at carbon in neopentyl-like silylcarbinyl sulfonates. <i>Journal of the American Chemical Society</i> , 1982, 104, 6852-6854.	6.6	28
32	Syntheses of Syn and Anti Doublebonded [5]Phenylene. <i>Organic Letters</i> , 2004, 6, 2249-2252.	2.4	26
33	Towards a Fully Conjugated, Double-Stranded Cycle: A Mass Spectrometric and Theoretical Study. <i>Chemistry - A European Journal</i> , 2008, 14, 1628-1637.	1.7	26
34	Concurrence between Current Density, Nucleus-Independent Chemical Shifts, and Aromatic Stabilization Energy: The Case of Isomeric [4]- and [5]Phenylenes. <i>Journal of Organic Chemistry</i> , 2013, 78, 7544-7553.	1.7	26
35	NICS _{xy} Scan Predictions of Local, Semi-Global, and Global Ring Currents in Annulated Pentalene and β -Indacene Cores Compared to First-Principles Current Density Maps. <i>ChemPhysChem</i> , 2020, 21, 65-82.	1.0	26
36	NICS ² Nucleus-independent Chemical Shift. , 2021, , 99-154.		25

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37	The Different Aromatic Characters of Some Localized Benzene Derivatives. <i>Journal of Physical Chemistry A</i> , 2008, 112, 12849-12854.	1.1	24
38	Inter- vs. intramolecular rearrangement of a (Bu ₃) ₂ Ni moiety in its 9-alkyl and 9,10-dialkyl anthracene complexes. Limiting conditions and isomer stabilities. <i>Journal of Organometallic Chemistry</i> , 1996, 515, 183-191.	0.8	21
39	Evidence for metal induced bond localization in cyclobutabenzene: The crystal and molecular structures of λ -6-Cr(CO) ₃ and λ -4-Fe(CO) ₃ complexes of cyclobutabenzene. <i>Journal of Organometallic Chemistry</i> , 1997, 542, 19-24.	0.8	20
40	Photo- and Thermal Haptotropism in Cyclopentadienylcobalt Complexes of Linear Phenylenes: Intercyclobutadiene Metal Migration. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 9853-9857.	7.2	20
41	Nickel-Mediated Cyclobutabenzene Syntheses. <i>trans</i> -7,8-Dibromo- cyclobutabenzene: 1 Their One-Pot Preparation, X-ray Structure, and Diels-Alder Reactions. <i>Journal of Organic Chemistry</i> , 1996, 61, 2549-2552.	1.7	18
42	The Competition for Electrons: Aromatic Stabilization in a Six-Membered Ring vs Cyclobutadiene-Iron Complex. The Molecular Structure of Tris(tricarbonylironcyclobutadiene)benzene. <i>Journal of Organic Chemistry</i> , 1998, 63, 247-253.	1.7	18
43	Evidence for Fully Conjugated Double-Stranded Cycles. <i>Chemistry - A European Journal</i> , 2011, 17, 12163-12174.	1.7	17
44	Interplay of Aromaticity and Antiaromaticity in N-Doped Nanographenes. <i>Journal of Physical Chemistry A</i> , 2020, 124, 695-703.	1.1	17
45	An MO-Based Identification of Charge-Shift Bonds. <i>ChemPhysChem</i> , 2012, 13, 2377-2381.	1.0	16
46	A one-pot method for the preparation of (R ₃ P) ₂ Ni ⁰ L complexes. <i>Journal of Organometallic Chemistry</i> , 1993, 458, 233-236.	0.8	14
47	Stereoselective Synthesis of Metalated Cyclobutyl Derivatives. <i>Advanced Synthesis and Catalysis</i> , 2009, 351, 1005-1008.	2.1	13
48	Tetraazaacenes Containing Four-Membered Rings in Different Oxidation States. Are They Aromatic? A Computational Study. <i>Journal of Organic Chemistry</i> , 2014, 79, 11644-11650.	1.7	13
49	The Planar Cyclooctatetraene Bridge in Bis-Metallic Macrocycles: Isolating or Conjugating?. <i>Inorganic Chemistry</i> , 2017, 56, 2287-2296.	1.9	13
50	Flat corannulene: when a transition state becomes a stable molecule. <i>Chemical Science</i> , 2020, 11, 13015-13025.	3.7	13
51	Crystal Structure of 1,2-Diphenyl-5,7-di-tert-butylspiro[2.5]octa-1,4,7-trien-6-one, a Possible Model for Diphenylvinylidenephonium Ions. <i>Journal of Organic Chemistry</i> , 1999, 64, 4370-4375.	1.7	12
52	Substituent effect and multisite protonation in the fragmentation of alkyl benzoates. <i>Journal of Mass Spectrometry</i> , 2002, 37, 336-342.	0.7	12
53	Aromatic Stabilization Energy and Magnetic Properties in Fulvalenes: Is There a Connection Between These Two Aromaticity Indices?. <i>Journal of Organic Chemistry</i> , 2013, 78, 12374-12380.	1.7	12
54	Comment on "Origin of the Nonplanarity of Tetrafluoro Cyclobutadiene, C ₄ F ₄ ". <i>Journal of Physical Chemistry A</i> , 2007, 111, 5116-5118.	1.1	11

