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List of Publications by Year
in descending order

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58 papers	844 citations	471509 17 h-index	580821 25 g-index
65 all docs	65 docs citations	65 times ranked	782 citing authors

#	Article	IF	Citations
1	Triplet state (anti)aromaticity of some monoheterocyclic analogues of benzene, naphthalene and anthracene. New Journal of Chemistry, 2021, 45, 5060-5074.	2.8	6
2	Singlet/Triplet State Anti/Aromaticity of CyclopentadienylCation: Sensitivity to Substituent Effect. Chemistry, 2021, 3, 765-782.	2.2	6
3	Revival of HÃ¼ckel Aromatic (Poly)benzenoid Subunits in Triplet State Polycyclic Aromatic Hydrocarbons by Silicon Substitution. Chemistry - an Asian Journal, 2021, , .	3.3	3
4	Synthesis of 2,3-Dihydro-4-pyridones and 4-Pyridones by the Cyclization Reaction of Ester-Tethered Enaminones. Journal of Organic Chemistry, 2020, 85, 13495-13507.	3.2	8
5	Substituent Effect on Triplet State Aromaticity of Benzene. Journal of Organic Chemistry, 2020, 85, 4289-4297.	3.2	35
6	Triplet-State Structures, Energies, and Antiaromaticity of BN Analogues of Benzene and Their Benzo-Fused Derivatives. Journal of Organic Chemistry, 2019, 84, 13582-13594.	3.2	16
7	The effect of two types of dibenzoannulation of pentalene on molecular energies and magnetically induced currents. Physical Chemistry Chemical Physics, 2019, 21, 3250-3263.	2.8	9
8	A DFT Study of the Modulation of the Antiaromatic and Openâ€Shell Character of Dibenzo[a,f]pentalene by Employing Three Strategies: Additional Benzoannulation, BN/CC Isosterism, and Substitution. Chemistry - A European Journal, 2019, 25, 9747-9757.	3.3	19
9	Does aromaticity account for an enhanced thermodynamic stability? The case of monosubstituted azaborines and the stereoelectronic chameleonism of the NH₂ group. Physical Chemistry Chemical Physics, 2019, 21, 9465-9476.	2.8	5
10	Can Variations of ¹Hâ€…NMR Chemical Shifts in Benzene Substituted with an Electronâ€Accepting (NO₂)/Donating (NH₂) Group be Explained in Terms of Resonance Effects of Substituents?. Chemistry - an Asian Journal, 2018, 13, 877-881.	3.3	7
11	Analysis of Stability and (Anti)aromaticity of BNâ€Dibenzo[a,e]pentalenes. European Journal of Organic Chemistry, 2018, 2018, 6230-6240.	2.4	4
12	Mono BN-substituted analogues of naphthalene: a theoretical analysis of the effect of BN position on stability, aromaticity and frontier orbital energies. New Journal of Chemistry, 2018, 42, 12968-12976.	2.8	15
13	Silica Gel as a Promoter of Sequential Azaâ€Michael/Michael Reactions of Amines and Propiolic Esters: Solventâ€and Metalâ€Free Synthesis of Polyfunctionalized Conjugated Dienes. Chemistry - an Asian Journal, 2018, 13, 1811-1835.	3.3	8
14	Electron Delocalization in Electronâ€Deficient Alkenes and Pushâ€Pull Alkenes. ChemistrySelect, 2017, 2, 42-50.	1.5	4
15	Theoretical study of azido gauche effect and its origin. New Journal of Chemistry, 2017, 41, 4644-4661.	2.8	2
16	4Ïâ€Electron Bâ€“N Monocycles: Stability and (Anti)aromaticity. European Journal of Organic Chemistry, 2017, 2017, 5163-5169.	2.4	10
17	Quantification of the pushâ€“pull effect in 2-alkylidene-4-oxothiazolidines by using NMR spectral data and barriers to rotation around the Cî€€ bond. New Journal of Chemistry, 2016, 40, 6364-6373.	2.8	6
18	Aromaticity of Diazaborines and Their Protonated Forms. Journal of Organic Chemistry, 2016, 81, 197-205.	3.2	16

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19	Theoretical analysis of the rotational barrier in ethane: cause and consequences. Structural Chemistry, 2015, 26, 989-996.	2.0	16
20	The effect of steric repulsion on the torsional potential of n-butane: a theoretical study. Tetrahedron, 2015, 71, 5119-5123.	1.9	9
21	Energy decomposition analysis of gauche preference in 2-haloethanol, 2-haloethylamine (halogen = F,) Tj ETQq1 1 0.784314 rgBT /Ov 22980-22995.	3.6	18
22	Origin of Fluorine/Sulfur Gauche Effect of Î²-Fluorinated Thiol, Sulfoxide, Sulfone, and Thionium Ion. Journal of Organic Chemistry, 2015, 80, 10197-10207.	3.2	20
23	A theoretical study on borenium ion affinities toward ammonia, formaldehyde and chloride anions. RSC Advances, 2015, 5, 75895-75910.	3.6	2
24	New insight into the anisotropic effects in solution-state NMR spectroscopy. RSC Advances, 2014, 4, 308-321.	3.6	63
25	Aromaticity and Stability of Azaborines. Chemistry - A European Journal, 2014, 20, 16558-16565.	3.3	66
26	Gauche preference in 1,2-difluoroethane originates from both orbital and electrostatic stabilization interactions. RSC Advances, 2014, 4, 43834-43838.	3.6	20
27	4-Oxothiazolidines with Exocyclic C=C Double Bond(s): Synthesis, Structure, Reactions and Biological Activity. Current Organic Chemistry, 2014, 18, 1108-1148.	1.6	14
28	Unusual mode of reactivity of 2-alkylidene-4-oxothiazolidine S-oxides under the Pummerer reaction conditions. Tetrahedron, 2013, 69, 9819-9825.	1.9	5
29	Cyclic Ï€ Electron Delocalization in Fluoroborazines. Journal of Physical Chemistry A, 2013, 117, 11540-11547.	2.5	9
30	Substituent effects on cyclic electron delocalization in symmetric B- and N-trisubstituted borazine derivatives. RSC Advances, 2013, 3, 24108.	3.6	11
31	¹H NMR Chemical Shifts of Cyclopropane and Cyclobutane: A Theoretical Study. Journal of Organic Chemistry, 2013, 78, 1504-1507.	3.2	20
32	2-Alkylidene-4-oxothiazolidine S-oxides: synthesis andÂstereochemistry. Tetrahedron, 2013, 69, 6436-6447.	1.9	11
33	Magnetic Anisotropy of the Cï£¿C Single Bond. Chemistry - A European Journal, 2013, 19, 4249-4254.	3.3	8
34	Synthesis of thiazolidine-fused heterocycles via exo-mode cyclizations of vinylogous N-acyliminium ions. Organic and Biomolecular Chemistry, 2012, 10, 575-589.	2.8	14
35	Density Functional Calculations of the Anisotropic Effects of Borazine and 1,3,2,4â€Ðiazadiboretidine. ChemPhysChem, 2012, 13, 3803-3811.	2.1	9
36	Is the Conventional Interpretation of the Anisotropic Effects of Cï£¾C Double Bonds and Aromatic Rings in NMR Spectra in Terms of the Ï€â€Electron Shielding/Deshielding Contributions Correct?. Chemistry - A European Journal, 2012, 18, 370-376.	3.3	29

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37	Quantification of the Aromaticity of 2-Alkylidenethiazolines Subjected to Pushâ€“Pull Activity. Journal of Organic Chemistry, 2011, 76, 3861-3871.	3.2	15
38	Catalytic oxidations of enolizable ketones using 2-alkylidene-4-oxothiazolidine vinyl bromide. Tetrahedron, 2011, 67, 8000-8008.	1.9	17
39	endo-Mode cyclizations of vinylogous N-acyliminium ions as a route to the synthesis of condensed thiazolidines. Tetrahedron, 2011, 67, 9541-9554.	1.9	13
40	Thermal solid-state Z/E isomerization of 2-alkylidene-4-oxothiazolidines: Effects of non-covalent interactions. Journal of the Serbian Chemical Society, 2011, 76, 317-328.	0.8	6
41	Transformations of 2-alkylidene-4-oxothiazolidine vinyl bromides initiated by bromophilic attack of neutral and anionic nucleophiles. Tetrahedron, 2010, 66, 6873-6884.	1.9	12
42	Structure, configuration, conformation and quantification of the pushâ€“pull effect of 2-alkylidene-4-thiazolidinones and 2-alkylidene-4,5-fused bicyclic thiazolidine derivatives. Tetrahedron, 2010, 66, 8958-8967.	1.9	17
43	Bromophilic substitution/carbophilic substitution cascade reactions of Î±,Î±-dibromo-2-methoxyacetophenone with C-, N- and O-nucleophiles. Tetrahedron Letters, 2010, 51, 4851-4855.	1.4	11
44	Reactions of ortho-substituted Î±,Î±-dibromoacetophenones with nucleophiles: first examples of combined carbophilic and bromophilic attack on Câ€“Br bonds. Tetrahedron Letters, 2009, 50, 700-703.	1.4	17
45	Cyclic voltammetry study of (5-ethoxycarbonylmethylidene-4-oxothiazolidin-2-ylidene)-N-phenylethanamide. Russian Journal of Physical Chemistry A, 2009, 83, 1571-1576.	0.6	4
46	High-Yield Synthesis of Substituted and Unsubstituted Pyridinium Salts Containing a 4-Oxothiazolidine Moiety. Synthesis, 2008, 2008, 2117-2121.	2.3	5
47	Carbonâ€“bromine cleavage by dimethyl sulfoxide: the key step of C(5) functionalization of pushâ€“pull 2-alkylidene-4-oxothiazolidine vinyl bromides. Tetrahedron Letters, 2007, 48, 1695-1698.	1.4	5
48	1H NMR Dynamic study of thermal Z/E isomerization of 5-substituted 2-alkylidene-4-oxothiazolidine derivatives: Barriers to rotation about CC bond. Journal of Molecular Structure, 2006, 800, 85-92.	3.6	17
49	2-Alkylidene-4-oxothiazolidine Vinyl Bromides: Versatile Precursors for C(5) Functionalization via Pyridine-Assisted Bromine Transfer. Synlett, 2006, 2006, 0729-0732.	1.8	9
50	Nucleophilic Functionalization of 2-Alkylidene-4-Oxothiazolidines at C(5)-Position Induced by Formation of Novel Pyridinium Salt. Phosphorus, Sulfur and Silicon and the Related Elements, 2005, 180, 1411-1415.	1.6	7
51	Stereocontrolled Synthesis of New Tetrahydrofuro[2,3-d]thiazole Derivatives via Activated Vinylogous Iminium Ions. Heterocycles, 2005, 65, 2635.	0.7	23
52	Configurational isomerization of push-pull thiazolidinone derivatives controlled by intermolecular and intramolecular RAHB: 1 H NMR dynamic investigation of concentration and temperature effects. Journal of Physical Organic Chemistry, 2004, 17, 118-123.	1.9	30
53	Ethyl (E)-2-(4-oxo-1,3-thiazinan-2-ylidene)ethanoate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o484-o485.	0.2	0
54	Thionation of N-methyl- and N-unsubstituted thiazolidine enaminones. Journal of the Serbian Chemical Society, 2004, 69, 909-918.	0.8	11

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55	A novel and efficient 4-oxothiazolidine-1,2-dithiole rearrangement induced by Lawesson's reagent. Tetrahedron Letters, 2003, 44, 7087-7090.	1.4	25
56	High regioselectivity in the heterocyclization of Î²-oxonitriles to 4-oxothiazolidines: X-ray structure proof. Tetrahedron, 2003, 59, 7803-7810.	1.9	36
57	Hydrogen bonding in push-pull 5-substituted-2-alkylidene-4-oxothiazolidines: 1H-MNR spectroscopic study. Journal of the Serbian Chemical Society, 2003, 68, 1-7.	0.8	10
58	Stereo- and regiocontrol of electrophile-initiated rearrangement of pushâ€“pull 5-substituted 4-oxothiazolidine derivatives. Tetrahedron, 2001, 57, 5833-5841.	1.9	31