Jorge Echeverra

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

42
papers

3,400
h-index

45
g-index

45
ext. papers

5.8
avg, IF

L-index

#	Paper	IF	Citations
42	From Simple Alkenes and CO2 to Fluorinated Carboxylic Acids: Computational Studies and Predictions. <i>European Journal of Organic Chemistry</i> , 2022 , 2022, e202101243	3.2	
41	Cooperative Effects between Hydrogen Bonds and C?OIIIS Interactions in the Crystal Structures of Sulfoxides. <i>Crystal Growth and Design</i> , 2021 , 21, 2481-2487	3.5	3
40	A Multifunctional Dysprosium-Carboxylato 2D Metall-Organic Framework. <i>Angewandte Chemie - International Edition</i> , 2021 , 60, 12001-12006	16.4	10
39	Delocalized Bonding in LiX Rings: Probing the Limits of the Covalent and Ionic Bonding Models. <i>Inorganic Chemistry</i> , 2021 , 60, 345-356	5.1	2
38	Interplay of halogen and hydrogen bonding in a series of heteroleptic iron(III) complexes. <i>CrystEngComm</i> , 2021 , 23, 4069-4076	3.3	1
37	Intermolecular Interactions between Thiocyanato Ligands in Metal Complexes. <i>Crystal Growth and Design</i> , 2021 , 21, 1636-1644	3.5	1
36	Methyl groups as widespread Lewis bases in noncovalent interactions. <i>Nature Communications</i> , 2021 , 12, 5030	17.4	3
35	Non-covalent interactions induced supramolecular architecture of Hg(NCS)2 with 3-pyridinecarbaldehyde nicotinoylhydrazone. <i>Inorganica Chimica Acta</i> , 2020 , 509, 119700	2.7	5
34	Understanding the Interplay of Dispersion, Charge Transfer, and Electrostatics in Noncovalent Interactions: The Case of Bromine Carbonyl Short Contacts. <i>Crystal Growth and Design</i> , 2020 , 20, 7180-7	71 8 7	9
33	Experimental and theoretical study of Pb?S and Pb?O Ehole interactions in the crystal structures of Pb(II) complexes. <i>CrystEngComm</i> , 2019 , 21, 6018-6025	3.3	9
32	Effect of the Substituents on the Nature and Strength of Lone-Pair@arbonyl Interactions in Acyl Halides. <i>Crystal Growth and Design</i> , 2019 , 19, 6511-6518	3.5	6
31	The n -E interaction in metal complexes. <i>Chemical Communications</i> , 2018 , 54, 3061-3064	5.8	42
30	Intermolecular Carbonyl Carbonyl Interactions in Transition-Metal Complexes. <i>Inorganic Chemistry</i> , 2018 , 57, 5429-5437	5.1	26
29	Taming a monomeric [Cu(IICH)] complex with silylene. <i>Chemical Science</i> , 2018 , 9, 4333-4337	9.4	17
28	Noncovalent Interactions in Succinic and Maleic Anhydride Derivatives. <i>Crystal Growth and Design</i> , 2018 , 18, 506-512	3.5	13
27	Understanding the Molecule-Electrode Interface for Molecular Spintronic Devices: A Computational and Experimental Study. <i>Molecules</i> , 2018 , 23,	4.8	3
26	The interplay of non-covalent interactions determining the antiparallel conformation of (isocyanide)gold(I) dimers. <i>CrystEngComm</i> , 2018 , 20, 3987-3993	3.3	8

(2011-2018)

25	In(iii)In(iii) short contacts: an unnoticed metallophilic interaction?. <i>Chemical Communications</i> , 2018 , 54, 6312-6315	5.8	7
24	Abundance and Strength of MHIIIHI (M = Al, Ga, In) Dihydrogen Bonds. <i>Crystal Growth and Design</i> , 2017 , 17, 2097-2103	3.5	18
23	Dihydrogen intermolecular contacts in group 13 compounds: HH or EH (E = B, Al, Ga) interactions?. <i>Dalton Transactions</i> , 2017 , 46, 2844-2854	4.3	21
22	Mercurophilic interactions: a theoretical study on the importance of ligands. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 11645-11654	3.6	25
21	Zinc-Zinc Double Bonds: A Theoretical Study. Angewandte Chemie - International Edition, 2017, 56, 1015	1±16041 5	558
20	Attractive PHHP interactions revealed by state-of-the-art ab initio calculations. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 28044-28055	3.6	10
19	Alkyl groups as electron density donors in Ehole bonding. CrystEngComm, 2017, 19, 6289-6296	3.3	15
18	Frustrated Lewis Trios and Long-Range Hole Interactions: A Combined Structural and Theoretical Study of LB-AX ???LB and LB???AX ???LB (A=B, Al, Ga, In) Systems. <i>ChemPhysChem</i> , 2017 , 18, 2864-2872	3.2	10
17	Intermolecular interactions in group 14 hydrides: Beyond C?HIIIH?C contacts. <i>International Journal of Quantum Chemistry</i> , 2017 , 117, e25432	2.1	23
16	The silane-methane dimer revisited: more than a dispersion-bound system?. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 32663-32669	3.6	9
15	ZincZinc Double Bonds: A Theoretical Study. <i>Angewandte Chemie</i> , 2017 , 129, 10285-10289	3.6	0
14	Supramolecular Rotor and Translator at Work: On-Surface Movement of Single Atoms. <i>ACS Nano</i> , 2015 , 9, 8394-400	16.7	25
13	Current-Driven Supramolecular Motor with In Situ Surface Chiral Directionality Switching. <i>Nano Letters</i> , 2015 , 15, 4793-8	11.5	49
12	The Design of a Single-Molecule Motor. Advances in Atom and Single Molecule Machines, 2015, 81-94	О	1
11	One-way rotation of a molecule-rotor driven by a shot noise. <i>Nanoscale</i> , 2014 , 6, 2793-9	7.7	21
10	Seeding molecular rotators on a passivated silicon surface. <i>ChemPhysChem</i> , 2014 , 15, 271-5	3.2	11
9	Understanding the Nature of the CHIIIHC Interactions in Alkanes. <i>Journal of Chemical Theory and Computation</i> , 2013 , 9, 1977-91	6.4	90
8	Dihydrogen contacts in alkanes are subtle but not faint. <i>Nature Chemistry</i> , 2011 , 3, 323-30	17.6	199

7	Concurrent symmetries: the interplay between local and global molecular symmetries. <i>Chemistry - A European Journal</i> , 2011 , 17, 359-67	4.8	14
6	The trigonal prism in coordination chemistry. <i>Chemistry - A European Journal</i> , 2010 , 16, 10380-96	4.8	52
5	New perspectives on polyhedral molecules and their crystal structures. <i>Journal of Physical Organic Chemistry</i> , 2010 , 23, 1080-1087	2.1	9
4	Jahn-Teller distortions of six-coordinate Cull compounds: cis or trans?. <i>Chemical Communications</i> , 2009 , 4242-4	5.8	25
3	Application of symmetry operation measures in structural inorganic chemistry. <i>Inorganic Chemistry</i> , 2008 , 47, 10965-70	5.1	20
2	Molecules and crystals with both icosahedral and cubic symmetry. <i>Chemical Communications</i> , 2008 , 271	7 ₅ 285	39
1	Covalent radii revisited. <i>Dalton Transactions</i> , 2008 , 2832-8	4.3	2540