

Jorge Echeverra

List of Publications by Citations

Source: <https://exaly.com/author-pdf/6497016/jorge-echeverria-publications-by-citations.pdf>

Version: 2024-04-29

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

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
citations

17
h-index

45
g-index

45
ext. papers

3,829
ext. citations

5.8
avg. IF

5.4
L-index

#	Paper	IF	Citations
42	Covalent radii revisited. <i>Dalton Transactions</i> , 2008 , 2832-8	4.3	2540
41	Dihydrogen contacts in alkanes are subtle but not faint. <i>Nature Chemistry</i> , 2011 , 3, 323-30	17.6	199
40	Understanding the Nature of the CH \cdots HC Interactions in Alkanes. <i>Journal of Chemical Theory and Computation</i> , 2013 , 9, 1977-91	6.4	90
39	The trigonal prism in coordination chemistry. <i>Chemistry - A European Journal</i> , 2010 , 16, 10380-96	4.8	52
38	Current-Driven Supramolecular Motor with In Situ Surface Chiral Directionality Switching. <i>Nano Letters</i> , 2015 , 15, 4793-8	11.5	49
37	The n- π interaction in metal complexes. <i>Chemical Communications</i> , 2018 , 54, 3061-3064	5.8	42
36	Molecules and crystals with both icosahedral and cubic symmetry. <i>Chemical Communications</i> , 2008 , 2717-25	5.25	39
35	Intermolecular Carbonyl \cdots Carbonyl Interactions in Transition-Metal Complexes. <i>Inorganic Chemistry</i> , 2018 , 57, 5429-5437	5.1	26
34	Mercuriphilic interactions: a theoretical study on the importance of ligands. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 11645-11654	3.6	25
33	Supramolecular Rotor and Translator at Work: On-Surface Movement of Single Atoms. <i>ACS Nano</i> , 2015 , 9, 8394-400	16.7	25
32	Jahn-Teller distortions of six-coordinate CuII compounds: cis or trans?. <i>Chemical Communications</i> , 2009 , 4242-4	5.8	25
31	Intermolecular interactions in group 14 hydrides: Beyond C \cdots H \cdots H \cdots C contacts. <i>International Journal of Quantum Chemistry</i> , 2017 , 117, e25432	2.1	23
30	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
29	One-way rotation of a molecule-rotor driven by a shot noise. <i>Nanoscale</i> , 2014 , 6, 2793-9	7.7	21
28	Application of symmetry operation measures in structural inorganic chemistry. <i>Inorganic Chemistry</i> , 2008 , 47, 10965-70	5.1	20
27	Abundance and Strength of M \cdots H \cdots H \cdots C (M = Al, Ga, In) Dihydrogen Bonds. <i>Crystal Growth and Design</i> , 2017 , 17, 2097-2103	3.5	18
26	Taming a monomeric [Cu(η -CH)] complex with silylene. <i>Chemical Science</i> , 2018 , 9, 4333-4337	9.4	17

25	Alkyl groups as electron density donors in H-bonding. <i>CrystEngComm</i> , 2017 , 19, 6289-6296	3.3	15
24	Concurrent symmetries: the interplay between local and global molecular symmetries. <i>Chemistry - A European Journal</i> , 2011 , 17, 359-67	4.8	14
23	Noncovalent Interactions in Succinic and Maleic Anhydride Derivatives. <i>Crystal Growth and Design</i> , 2018 , 18, 506-512	3.5	13
22	Seeding molecular rotators on a passivated silicon surface. <i>ChemPhysChem</i> , 2014 , 15, 271-5	3.2	11
21	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
20	Frustrated Lewis Trios and Long-Range H-bond Interactions: A Combined Structural and Theoretical Study of LB-AX \cdots LB and LB \cdots AX \cdots LB (A=B, Al, Ga, In) Systems. <i>ChemPhysChem</i> , 2017 , 18, 2864-2872	3.2	10
19	A Multifunctional Dysprosium-Carboxylato 2D Metall-Organic Framework. <i>Angewandte Chemie - International Edition</i> , 2021 , 60, 12001-12006	16.4	10
18	Experimental and theoretical study of Pb \cdots S and Pb \cdots O H-bond interactions in the crystal structures of Pb(II) complexes. <i>CrystEngComm</i> , 2019 , 21, 6018-6025	3.3	9
17	The silane-methane dimer revisited: more than a dispersion-bound system?. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 32663-32669	3.6	9
16	New perspectives on polyhedral molecules and their crystal structures. <i>Journal of Physical Organic Chemistry</i> , 2010 , 23, 1080-1087	2.1	9
15	Understanding the Interplay of Dispersion, Charge Transfer, and Electrostatics in Noncovalent Interactions: The Case of Bromine \cdots Carbonyl Short Contacts. <i>Crystal Growth and Design</i> , 2020 , 20, 7180-7187	3.5	9
14	Zinc-Zinc Double Bonds: A Theoretical Study. <i>Angewandte Chemie - International Edition</i> , 2017 , 56, 10151-10155	16.1	8
13	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
12	In(iii) \cdots In(iii) short contacts: an unnoticed metallophilic interaction?. <i>Chemical Communications</i> , 2018 , 54, 6312-6315	5.8	7
11	Effect of the Substituents on the Nature and Strength of Lone-Pair \cdots Carbonyl Interactions in Acyl Halides. <i>Crystal Growth and Design</i> , 2019 , 19, 6511-6518	3.5	6
10	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
9	Understanding the Molecule-Electrode Interface for Molecular Spintronic Devices: A Computational and Experimental Study. <i>Molecules</i> , 2018 , 23,	4.8	3
8	Cooperative Effects between Hydrogen Bonds and C=O \cdots S Interactions in the Crystal Structures of Sulfoxides. <i>Crystal Growth and Design</i> , 2021 , 21, 2481-2487	3.5	3

7	Methyl groups as widespread Lewis bases in noncovalent interactions. <i>Nature Communications</i> , 2021 , 12, 5030	17.4	3
6	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
5	The Design of a Single-Molecule Motor. <i>Advances in Atom and Single Molecule Machines</i> , 2015 , 81-94	0	1
4	Interplay of halogen and hydrogen bonding in a series of heteroleptic iron(III) complexes. <i>CrystEngComm</i> , 2021 , 23, 4069-4076	3.3	1
3	Intermolecular Interactions between Thiocyanato Ligands in Metal Complexes. <i>Crystal Growth and Design</i> , 2021 , 21, 1636-1644	3.5	1
2	Zinc-Zinc Double Bonds: A Theoretical Study. <i>Angewandte Chemie</i> , 2017 , 129, 10285-10289	3.6	0
1	From Simple Alkenes and CO ₂ to Fluorinated Carboxylic Acids: Computational Studies and Predictions. <i>European Journal of Organic Chemistry</i> , 2022 , 2022, e202101243	3.2	