

# Zeferino GÃ³mez-Sandoval

## List of Publications by Year in descending order

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36  
papers

349  
citations

759190

12  
h-index

888047

17  
g-index

37  
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37  
docs citations

37  
times ranked

543  
citing authors

#	ARTICLE	IF	CITATIONS
1	Synthesis, X-ray and complete assignments of 1H and 13C nuclear magnetic resonance data for novel dichloro-1,4-dihydro-1,4-epoxynaphthalene derivatives. <i>Journal of Molecular Structure</i> , 2021, 1224, 129287.	3.6	2
2	Tamoxifen Sensitizes Acute Lymphoblastic Leukemia Cells to Cannabidiol by Targeting Cyclophilin-D and Altering Mitochondrial Ca <sup>2+</sup> Homeostasis. <i>International Journal of Molecular Sciences</i> , 2021, 22, 8688.	4.1	5
3	Ground state geometries and stability of impurity doped clusters: LinY (n=13). The role of yttrium atom in electronic and magnetic properties. <i>Chemical Physics Letters</i> , 2021, 779, 138884.	2.6	3
4	Synthesis, crystal structure, antioxidant activity and dft study of 2-aryl-2,3-dihydro-4H-[1,3]thiazino[3,2-a]benzimidazol-4-One. <i>Journal of Molecular Structure</i> , 2020, 1199, 127036.	3.6	17
5	In silico structure-based design of GABA B receptor agonists using a combination of docking and QSAR. <i>Chemical Biology and Drug Design</i> , 2019, 94, 1782-1798.	3.2	7
6	Synthesis, structural investigation, antibacterial and DFT studies of complexes derived from a cholesteryl dithiophosphonate ligand with some thio-metallolane and thio-metalloocene heterocycles of As(III) and Sb(III). <i>Inorganica Chimica Acta</i> , 2019, 495, 118943.	2.4	4
7	Cytotoxic Acetogenins from the Roots of <i>Annona purpurea</i> . <i>International Journal of Molecular Sciences</i> , 2019, 20, 1870.	4.1	14
8	Magnesium oxide clusters as promising candidates for hydrogen storage. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 23102-23110.	2.8	19
9	Organotin(IV) compounds derived from ibuprofen and cinnamic acids, an alternative into design of anti-inflammatory by the cyclooxygenases (COX-1 and COX-2) pathway. <i>Journal of Organometallic Chemistry</i> , 2018, 862, 58-70.	1.8	25
10	h function: A protonic take on the numerical Fukui function as a graphical descriptor for deprotonation. <i>International Journal of Quantum Chemistry</i> , 2018, 118, e25532.	2.0	2
11	Exploring the Structure, Energetic, and Magnetic Properties of Neutral Small Lithium Clusters Doped with Yttrium: Supermagnetic Atom Research. <i>ACS Omega</i> , 2018, 3, 11252-11261.	3.5	5
12	In silico receptor-based drug design of X,Y-benzenesulfonamide derivatives as selective COX-2 inhibitors. <i>Comptes Rendus Chimie</i> , 2017, 20, 169-180.	0.5	12
13	Conformational analysis of N <sup>+</sup> BH <sub>3</sub> <sup>+</sup> , N <sup>+</sup> BF <sub>3</sub> <sup>+</sup> , and N-CH <sub>3</sub> <sup>+</sup> complexes with ibuprofen-derivative amides. <i>Heteroatom Chemistry</i> , 2017, 28, e21368.	0.7	1
14	Silicon containing ibuprofen derivatives with antioxidant and anti-inflammatory activities: An in vivo and in silico study. <i>European Journal of Pharmacology</i> , 2017, 814, 18-27.	3.5	10
15	<i>Lactobacillus plantarum</i> WCFS1 Î <sup>2</sup> -Fructosidase: Evidence for an Open Funnel-Like Channel Through the Catalytic Domain with Importance for the Substrate Selectivity. <i>Applied Biochemistry and Biotechnology</i> , 2016, 180, 1056-1075.	2.9	3
16	Synthesis and Biological Screening of Silicon-Containing Ibuprofen Derivatives: A Study of Their NF-Î <sup>2</sup> Inhibitory Activity, Cytotoxicity, and Their Ability to Bind IKKÎ <sup>2</sup> . <i>Australian Journal of Chemistry</i> , 2016, 69, 662.	0.9	7
17	Computational study of the structure, bonding and reactivity of selected helical metallocenes. <i>Inorganica Chimica Acta</i> , 2015, 438, 203-207.	2.4	6
18	Antiradical capacity of a series of organotin(IV) compounds: A chemical reactivity study in the Density Functional Theory framework. <i>Inorganica Chimica Acta</i> , 2014, 413, 143-148.	2.4	5

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19	QSAR study of the DPPH• radical scavenging activity of coumarin derivatives and xanthine oxidase inhibition by molecular docking. <i>Open Chemistry</i> , 2014, 12, 1067-1080.	1.9	13
20	Evaluation of the antiradical activity of hyperjovinol-A utilizing donor-acceptor maps. <i>Journal of Molecular Modeling</i> , 2014, 20, 2337.	1.8	26
21	Shape entropy's response to molecular ionization. <i>Journal of Molecular Modeling</i> , 2013, 19, 1677-1683.	1.8	11
22	High magnetic moments on binary yttrium-alkali superatoms. <i>Chemical Physics Letters</i> , 2013, 583, 97-102.	2.6	18
23	Synthesis and in Vitro Antioxidant Activity Evaluation of 3-Carboxycoumarin Derivatives and QSAR Study of Their DPPH• Radical Scavenging Activity. <i>Molecules</i> , 2012, 17, 14882-14898.	3.8	26
24	Antitumor structure-activity relationship in bis-stannoxane derivatives from pyridine dicarboxylic and benzoic acids. <i>Inorganica Chimica Acta</i> , 2012, 392, 229-235.	2.4	7
25	Estudio de usabilidad de visualizaci3n molecular educativa en un tel3fono inteligente. <i>Quimica Nova</i> , 2012, 35, 648-653.	0.3	3
26	Can an eight $\pi$ -electron bare ring be planar?. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 20615.	2.8	14
27	X-ray, DFT, FTIR and NMR structural study of 2,3-dihydro-2-(R-phenylacylidene)-1,3,3-trimethyl-1H-indole. <i>Journal of Molecular Structure</i> , 2011, 987, 106-118.	3.6	21
28	Synthesis of Novel Pyridinium Betaine Precursors from exo-Norbornene Dicarboximides. <i>Letters in Organic Chemistry</i> , 2011, 8, 249-257.	0.5	1
29	Renin gene haplotype diversity and linkage disequilibrium in two Mexican and one German population samples. <i>JRAAS - Journal of the Renin-Angiotensin-Aldosterone System</i> , 2011, 12, 231-237.	1.7	6
30	A Helicoid Ferrocene. <i>Inorganic Chemistry</i> , 2009, 48, 2714-2716.	4.0	22
31	Integraci3n de visualizaci3n cient3fica molecular en el sal3n de clases. <i>Quimica Nova</i> , 2008, 31, 2184-2189.	0.3	0
32	Density Functional Study of 2-[(R-Phenyl)amine]-1,4-naphthalenediones. <i>Journal of Chemical Theory and Computation</i> , 2007, 3, 894-904.	5.3	4
33	Density Functional Study of the Structure and Properties of Cu9 and Cu9-. <i>Journal of Chemical Theory and Computation</i> , 2007, 3, 905-913.	5.3	16
34	2-[(R-phenyl)amine]-1,4-naphthalenediones as photosystem I electron acceptors. Structure-activity relationship of m- and p-PAN compounds with QSAR analysis. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2006, 83, 105-113.	3.8	7
35	First principle $f$ - $f$ energy separation. <i>Theoretical Chemistry Accounts</i> , 2005, 114, 137-144.	1.4	1
36	Separation of $f$ and $f$ Energies. <i>Journal of Physical Chemistry A</i> , 2005, 109, 1257-1259.	2.5	6