

# Ana Cristina Mora Tello

## List of Publications by Year in descending order

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Version: 2024-02-01

11  
papers

68  
citations

1937685

4  
h-index

1720034

7  
g-index

11  
all docs

11  
docs citations

11  
times ranked

94  
citing authors

#	ARTICLE	IF	CITATIONS
1	Microwave-Driven Hexagonal-to-Monoclinic Transition in BiPO <sub>4</sub> : An In-Depth Experimental Investigation and First-Principles Study. <i>Inorganic Chemistry</i> , 2020, 59, 7453-7468.	4.0	24
2	Zinc-substituted Ag <sub>2</sub> CrO <sub>4</sub> : A material with enhanced photocatalytic and biological activity. <i>Journal of Alloys and Compounds</i> , 2020, 835, 155315.	5.5	16
3	Structure, Photoluminescence Emissions, and Photocatalytic Activity of Ag <sub>2</sub> SeO <sub>3</sub> : A Joint Experimental and Theoretical Investigation. <i>Inorganic Chemistry</i> , 2021, 60, 5937-5954.	4.0	10
4	Multi-dimensional architecture of Ag <sub>1-x</sub> Ag <sub>2</sub> WO <sub>4</sub> crystals: insights into microstructural, morphological, and photoluminescence properties. <i>CrystEngComm</i> , 2020, 22, 7903-7917.	2.6	9
5	The 1,2-hydrogen shift reaction for monohalogenophosphanes PH <sub>2</sub> X and HPX (X = F, Cl). <i>Molecular Physics</i> , 2016, 114, 2999-3014.	1.7	3
6	On polarization functions for Gaussian basis sets. <i>Journal of Molecular Modeling</i> , 2020, 26, 293.	1.8	3
7	Ag <sub>2</sub> WO <sub>4</sub> under microwave, electron beam and femtosecond laser irradiations: Unveiling the relationship between morphology and photoluminescence emissions. <i>Journal of Alloys and Compounds</i> , 2022, 903, 163840.	5.5	3
8	Design approach of a molecular nanotransmitter using quantum dynamics: A case study on ethane. , 2010, , .		0
9	On the Use of an Interpolation Approach for the Choice of Gaussian Polarization Functions. <i>Journal of the Brazilian Chemical Society</i> , 2018, , .	0.6	0
10	Accurate atomic electron affinities calculated by using anionic Gaussian basis sets. <i>Theoretical Chemistry Accounts</i> , 2020, 139, 1.	1.4	0
11	Generation, contraction, and polarisation of Gaussian basis sets for atomic and molecular calculations using the generator coordinate method with polynomial discretisation: atoms from Na through Cl. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 16989-16997.	2.8	0