

# Mariela M Nolasco

## List of Publications by Year in descending order

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Version: 2024-02-01

47  
papers

1,970  
citations

304743

22  
h-index

243625

44  
g-index

48  
all docs

48  
docs citations

48  
times ranked

2749  
citing authors

#	ARTICLE	IF	CITATIONS
1	Identification of microplastics using Raman spectroscopy: Latest developments and future prospects. <i>Water Research</i> , 2018, 142, 426-440.	11.3	512
2	Inelastic neutron scattering study of relin: shedding light on the hydrogen bonding network of deep eutectic solvents. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 17998-18009.	2.8	132
3	Inside PEF: Chain Conformation and Dynamics in Crystalline and Amorphous Domains. <i>Macromolecules</i> , 2018, 51, 3515-3526.	4.8	110
4	Photo-Click Chemistry to Design Highly Efficient Lanthanide $\beta^2$ -Diketonate Complexes Stable under UV Irradiation. <i>Chemistry of Materials</i> , 2013, 25, 586-598.	6.7	96
5	Engineering highly efficient Eu(III)-based tri-ureasil hybrids toward luminescent solar concentrators. <i>Journal of Materials Chemistry A</i> , 2013, 1, 7339.	10.3	95
6	Characterization of Microplastics by Raman Spectroscopy. <i>Comprehensive Analytical Chemistry</i> , 2017, , 119-151.	1.3	84
7	Computationally-Assisted Approach to the Vibrational Spectra of Molecular Crystals: Study of Hydrogen-Bonding and Pseudo-Polymorphism. <i>ChemPhysChem</i> , 2006, 7, 2150-2161.	2.1	71
8	Excimer Formation in a Terbium Metal-Organic Framework Assists Luminescence Thermometry. <i>Chemistry of Materials</i> , 2017, 29, 9547-9554.	6.7	65
9	Probing Pseudopolymorphic Transitions in Pharmaceutical Solids using Raman Spectroscopy: Hydration and Dehydration of Theophylline. <i>Journal of Pharmaceutical Sciences</i> , 2007, 96, 1366-1379.	3.3	56
10	Dependence of the Lifetime upon the Excitation Energy and Intramolecular Energy Transfer Rates: The $^5D_0 \rightarrow ^3F_4$ Emission Case. <i>Chemistry - A European Journal</i> , 2012, 18, 12130-12139.	3.3	54
11	New chloro and triphenylsiloxy derivatives of dioxomolybdenum(VI) chelated with pyrazolylpyridine ligands: Catalytic applications in olefin epoxidation. <i>Journal of Molecular Catalysis A</i> , 2007, 261, 79-87.	4.8	52
12	C-H...O Hydrogen Bonds in Cyclohexenone Reveal the Spectroscopic Behavior of Csp <sup>3</sup> H and Csp <sup>2</sup> H Donors. <i>ChemPhysChem</i> , 2005, 6, 496-502.	2.1	50
13	Chemistry and Catalytic Activity of Molybdenum(VI)-Pyrazolylpyridine Complexes in Olefin Epoxidation. Crystal Structures of Monomeric Dioxo, Dioxo-1/4-oxo, and Oxodiperoxo Derivatives. <i>Inorganic Chemistry</i> , 2011, 50, 525-538.	4.0	50
14	Hydrogen-Bond Dynamics of C-H...O Interactions: The Chloroform...Acetone Case. <i>Chemistry - A European Journal</i> , 2010, 16, 9010-9017.	3.3	38
15	Synthesis, Structural Elucidation, and Catalytic Properties in Olefin Epoxidation of the Polymeric Hybrid Material [Mo <sub>3</sub> O <sub>9</sub> (2-[3(5)-Pyrazolyl]pyridine)] <sub>n</sub> . <i>Inorganic Chemistry</i> , 2014, 53, 2652-2665.	4.0	38
16	Intriguing light-emission features of ketoprofen-based Eu(III) adduct due to a strong electron-phonon coupling. <i>Journal of Luminescence</i> , 2016, 170, 357-363.	3.1	34
17	A Combined Theoretical-Experimental Study of the Inclusion of Niobocene Dichloride in Native and Permethylated $\beta^2$ -Cyclodextrins. <i>Organometallics</i> , 2007, 26, 4220-4228.	2.3	32
18	Effect of hydrogen bonding in the vibrational spectra of <i>trans</i> -cinnamic acid. <i>Journal of Raman Spectroscopy</i> , 2009, 40, 394-400.	2.5	32

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19	Hydrogen Bond Dynamics of Cellulose through Inelastic Neutron Scattering Spectroscopy. <i>Biomacromolecules</i> , 2018, 19, 1305-1313.	5.4	28
20	Light emission of a polyfluorene derivative containing complexed europium ions. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 26238-26248.	2.8	26
21	Efficient Visible-Light-Excitable $\text{Eu}^{3+}$ Complexes for Red Organic Light-Emitting Diodes. <i>European Journal of Inorganic Chemistry</i> , 2020, 2020, 1260-1270.	2.0	25
22	What a difference a methyl group makes – probing choline-urea molecular interactions through urea structure modification. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 18278-18289.	2.8	24
23	Asymmetric Monomer, Amorphous Polymer? Structure-Property Relationships in 2,4-FDCA and 2,4-PEF. <i>Macromolecules</i> , 2020, 53, 1380-1387.	4.8	24
24	The role of 4,7-disubstituted phenanthroline ligands in energy transfer of europium(iii) complexes: a DFT study. <i>New Journal of Chemistry</i> , 2011, 35, 2435.	2.8	21
25	Modelling the luminescence of extended solids: an example of a highly luminescent MCM-41 impregnated with a $\text{Eu}^{3+}$ -diketonate complex. <i>Journal of Materials Chemistry C</i> , 2014, 2, 9701-9711.	5.5	20
26	Poly(4-styrene sulfonic acid)/bacterial cellulose membranes: Electrochemical performance in a single-chamber microbial fuel cell. <i>Bioresource Technology Reports</i> , 2020, 9, 100376.	2.7	20
27	C-H...O Hydrogen bonding in 4-phenyl-benzaldehyde: A comprehensive crystallographic, spectroscopic and computational study. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 3027.	2.8	19
28	Vibrational Study on the Local Structure of Post-Synthesis and Hybrid Mesoporous Materials: Are There Fundamental Distinctions?. <i>Chemistry - A European Journal</i> , 2007, 13, 7874-7882.	3.3	19
29	A green-emitting $\beta$ -substituted $\beta$ -diketonate $\text{Tb}^{3+}$ phosphor for ultraviolet LED-based solid-state lighting. <i>Journal of Coordination Chemistry</i> , 2014, 67, 4076-4089.	2.2	19
30	Water in Deep Eutectic Solvents: New Insights From Inelastic Neutron Scattering Spectroscopy. <i>Frontiers in Physics</i> , 2022, 10, .	2.1	17
31	Understanding the vibrational spectra of crystalline isoniazid: Raman, IR and INS spectroscopy and solid-state DFT study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2018, 204, 452-459.	3.9	13
32	Crystal structure landscapes from combined vibrational spectroscopy and ab initio calculations: 4-(Dimethylamino)benzaldehyde. <i>Computational and Theoretical Chemistry</i> , 2010, 946, 65-69.	1.5	12
33	Understanding the Structure and Dynamics of Nanocellulose-Based Composites with Neutral and Ionic Poly(methacrylate) Derivatives Using Inelastic Neutron Scattering and DFT Calculations. <i>Molecules</i> , 2020, 25, 1689.	3.8	12
34	Influence of the Crystal Structure on the Luminescence Properties of Mixed Eu,La-(1,10-Phenanthroline) Complexes. <i>European Journal of Inorganic Chemistry</i> , 2015, 2015, 4861-4868.	2.0	10
35	Spectroscopic and thermal studies on the inclusion of <i>trans</i> -cinnamic acid and a number of its hydroxyl-derivatives with $\beta$ , $\gamma$ and $\delta$ -cyclodextrins molecules. <i>Journal of Raman Spectroscopy</i> , 2009, 40, 687-695.	2.5	7
36	Modelling the Luminescence of Phosphonate Lanthanide-Organic Frameworks. <i>European Journal of Inorganic Chemistry</i> , 2015, 2015, 1254-1260.	2.0	7

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37	Vibrational Dynamics of Crystalline 4-Phenylbenzaldehyde from INS Spectra and Periodic DFT Calculations. <i>Molecules</i> , 2020, 25, 1374.	3.8	6
38	Intermolecular C-H...O interactions in cyclopentanone: An inelastic neutron scattering study. <i>Chemical Physics</i> , 2013, 427, 117-123.	1.9	5
39	Coordination polymers based on a glycine-derivative ligand. <i>CrystEngComm</i> , 2014, 16, 8119-8137.	2.6	5
40	High Emission Quantum Yield Tb <sup>3+</sup> -Activated Organic-Inorganic Hybrids for UV-Down-Shifting Green Light-Emitting Diodes. <i>European Journal of Inorganic Chemistry</i> , 2020, 2020, 1736-1742.	2.0	5
41	Insights into phase stability of anhydrous/hydrate systems: a Raman-based methodology. <i>Journal of Raman Spectroscopy</i> , 2010, 41, 340-349.	2.5	4
42	Exploring C-H...O hydrogen bonds in dihydrocoumarin from combined vibrational spectroscopy and DFT calculations. <i>Chemical Physics Letters</i> , 2012, 551, 86-91.	2.6	4
43	Effects of hydrogen-bond and cooperativity in the vibrational spectra of Luminol. <i>Vibrational Spectroscopy</i> , 2013, 64, 119-125.	2.2	4
44	Preformulation Studies of the $\beta$ -Cyclodextrin and Montelukast Inclusion Compound Prepared by Comilling. <i>Journal of Pharmaceutical Sciences</i> , 2019, 108, 1837-1847.	3.3	4
45	New Insights on the Vibrational Dynamics of 2-Methoxy-, 4-Methoxy- and 4-Ethoxy-Benzaldehyde from INS Spectra and Periodic DFT Calculations. <i>Materials</i> , 2021, 14, 4561.	2.9	4
46	Vibrational dynamics of 4-fluorobenzaldehyde from periodic DFT calculations. <i>Chemical Physics Letters: X</i> , 2019, 2, 100006.	2.1	3
47	Vibrational Dynamics in crystalline 4-(dimethylamino) benzaldehyde: Inelastic Neutron Scattering and Periodic DFT Study. <i>Materials</i> , 2022, 15, 475.	2.9	2