

Lourdes Infantes

List of Publications by Year in descending order

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70
papers

2,614
citations

331259

21
h-index

189595

50
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83
all docs

83
docs citations

83
times ranked

2628
citing authors

#	ARTICLE	IF	CITATIONS
1	The structure and flexibility analysis of the <i>Arabidopsis</i> synaptotagmin 1 reveal the basis of its regulation at membrane contact sites. <i>Life Science Alliance</i> , 2021, 4, e202101152.	1.3	9
2	Fluorescence mechanism switching from ICT to PET by substituent chemical manipulation: Macrophage cytoplasm imaging probes. <i>Dyes and Pigments</i> , 2020, 175, 108172.	2.0	11
3	New Multicomponent Forms of the Antiretroviral Nevirapine with Improved Dissolution Performance. <i>Crystal Growth and Design</i> , 2020, 20, 688-698.	1.4	9
4	Optimization and comparison of statistical tools for the prediction of multicomponent forms of a molecule: the antiretroviral nevirapine as a case study. <i>CrystEngComm</i> , 2020, 22, 7460-7474.	1.3	10
5	Environment-Sensitive Probes for Illuminating Amyloid Aggregation <i>In Vitro</i> and in Zebrafish. <i>ACS Sensors</i> , 2020, 5, 2792-2799.	4.0	21
6	Smart lanthanide antennas for sensing water. <i>Chemical Communications</i> , 2020, 56, 5484-5487.	2.2	20
7	A structural study of new tetrakis(1H-pyrazol-1-yl)methanes. <i>Tetrahedron</i> , 2019, 75, 130690.	1.0	2
8	New Quinolylnitrones for Stroke Therapy: Antioxidant and Neuroprotective (<i>Z</i>)- <i>N</i> - <i>tert</i> -Butyl-1-(2-chloro-6-methoxyquinolin-3-yl)methanimine Oxide as a New Lead-Compound for Ischemic Stroke Treatment. <i>Journal of Medicinal Chemistry</i> , 2019, 62, 2184-2201.	2.9	35
9	Chiral Microneedles from an Achiral Bis(boron dipyrromethene): Spontaneous Mirror Symmetry Breaking Leading to a Promising Photoluminescent Organic Material. <i>Langmuir</i> , 2019, 35, 5021-5028.	1.6	6
10	An Example of Polynomial Expansion: The Reaction of 3(5)-Methyl-1H-Pyrazole with Chloroform and Characterization of the Four Isomers. <i>Molecules</i> , 2019, 24, 568.	1.7	6
11	The structure of four thallium tris(1H-pyrazol-1-yl)hydroborates in the solid state by X-ray crystallography and in solution by NMR and DFT-GIAO calculations. <i>Inorganica Chimica Acta</i> , 2018, 483, 402-410.	1.2	6
12	Deciphering the Inhibition of the Neuronal Calcium Sensor 1 and the Guanine Exchange Factor Ric8a with a Small Phenothiazine Molecule for the Rational Generation of Therapeutic Synapse Function Regulators. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 5910-5921.	2.9	10
13	Interference of the complex between NCS-1 and Ric8a with phenothiazines regulates synaptic function and is an approach for fragile X syndrome. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E999-E1008.	3.3	40
14	Structure of Ligand-Bound Intermediates of Crop ABA Receptors Highlights $\hat{A}P2C$ as Necessary ABA Co-receptor. <i>Molecular Plant</i> , 2017, 10, 1250-1253.	3.9	49
15	The structures of two scorpionates: thallium tetrakis(3-phenyl-1H-pyrazol-1-yl)borate and potassium tetrakis(3-cyclopropyl-1H-pyrazol-1-yl)borate. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2016, 72, 819-825.	0.2	4
16	Experimental and Theoretical Studies on the Rearrangement of 2-Oxoazepane \hat{A} -Amino Acids into 2-Oxopiperidine $\hat{A}^{2,3}$ -Amino Acids: An Example of Intramolecular Catalysis. <i>Chemistry - A European Journal</i> , 2015, 21, 2489-2500.	1.7	3
17	Synthesis of Enantiopure 3-Hydroxypiperidines from Sulfinyl Dienyl Amines by Diastereoselective Intramolecular Cyclization and [2,3]-Sigmatropic Rearrangement. <i>Journal of Organic Chemistry</i> , 2015, 80, 7674-7692.	1.7	13
18	The reaction of 2-amino-4-H-pyrans with <i>N</i> -bromosuccinimide. <i>Molecular Diversity</i> , 2015, 19, 103-122.	2.1	2

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19	Prediction of H-Bonding Motifs for Pyrazoles and Oximes Using the Cambridge Structural Database. , 2015, , 269-280.		1
20	Silver Triflate-Catalyzed Cyclization of 2-Amino-6-propargyl-amineazines Leading to Iminoimidazoazines. Advanced Synthesis and Catalysis, 2014, 356, 1235-1241.	2.1	9
21	Efficient Light Harvesters Based on the 10-(1,3-Dithiol-2-ylidene)anthracene Core. Organic Letters, 2013, 15, 4166-4169.	2.4	18
22	Sulfoxide-Directed Enantioselective Synthesis of Functionalized Tetrahydropyridines. Organic Letters, 2013, 15, 4936-4939.	2.4	18
23	Silver-Catalyzed Cyclization of <i>N</i> -(Prop-2-yn-1-yl)pyridin-2-amines. European Journal of Organic Chemistry, 2013, 2013, 35-39.	1.2	44
24	The Structure and Dynamic Properties of 1H-Pyrazole-4-Carboxylic Acids in the Solid State. Zeitschrift Fur Physikalische Chemie, 2013, 227, 841-856.	1.4	3
25	Highly Functionalized 1,2-Diamino Compounds through Reductive Amination of Amino Acid-Derived β -Keto Esters. PLoS ONE, 2013, 8, e53231.	1.1	5
26	Azepane Quaternary Amino Acids As Effective Inducers of 3 ₁₀ Helix Conformations. Journal of Organic Chemistry, 2012, 77, 9833-9839.	1.7	13
27	The Structural Domains of Pseudomonas aeruginosa Phosphorylcholine Phosphatase Cooperate in Substrate Hydrolysis: 3D Structure and Enzymatic Mechanism. Journal of Molecular Biology, 2012, 423, 503-514.	2.0	6
28	A practical two-step synthesis of imidazo[1,2-a]pyridines from <i>N</i> -(prop-2-yn-1-yl)pyridin-2-amines. Chemical Communications, 2011, 47, 5043.	2.2	39
29	Quaternary β -2-Oxoazepane β -Amino Acids: Synthesis from Ornithine-Derived β -Lactams and Incorporation into Model Dipeptides. Journal of Organic Chemistry, 2011, 76, 6592-6603.	1.7	33
30	Controlling Optical Properties and Function of BODIPY by Using Asymmetric Substitution Effects. Chemistry - A European Journal, 2010, 16, 14094-14105.	1.7	38
31	Synthesis of (E)-diethyl 6,6-(diazene-1,2-diyl)bis(5-cyano-2-methyl-4-phenylnicotinates), a new type of 2,2-azopyridine dyes. Tetrahedron Letters, 2010, 51, 6278-6281.	0.7	8
32	Modification of pancreatic lipase properties by directed molecular evolution. Protein Engineering, Design and Selection, 2010, 23, 365-373.	1.0	14
33	Further Evidence for 2-Alkyl-2-carboxyazetidines as β -Turn Inducers. Journal of Organic Chemistry, 2009, 74, 8203-8211.	1.7	19
34	The molecular structure of 1,3-dimethyl-4-phenyl-1H-pyrazole-5-carboxylic acid. Arkivoc, 2008, 2008, 74-84.	0.3	3
35	Synthesis, Photophysical Properties, and Laser Behavior of 3-Amino and 3-Acetamido BODIPY Dyes. Organic Letters, 2007, 9, 4183-4186.	2.4	60
36	Organic crystal hydrates: what are the important factors for formation. CrystEngComm, 2007, 9, 65-71.	1.3	175

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37	Structural Changes in the BODIPY Dye PM567 Enhancing the Laser Action in Liquid and Solid Media. <i>Advanced Functional Materials</i> , 2007, 17, 3088-3098.	7.8	56
38	Knowledge-based approaches to crystal design. <i>CrystEngComm</i> , 2006, , .	1.3	3
39	Substituent Effects on Enthalpies of Formation of Nitrogen Heterocycles: 2-Substituted Benzimidazoles and Related Compounds. <i>Journal of Physical Chemistry A</i> , 2006, 110, 2535-2544.	1.1	17
40	Water Oligomers in Crystal Hydrates "What's News and What Isn't?. <i>Angewandte Chemie - International Edition</i> , 2006, 45, 32-36.	7.2	265
41	Classification of hydrogen-bond motives in crystals of NH-pyrazoles: a mixed empirical and theoretical approach. <i>Arkivoc</i> , 2006, 2006, 15-30.	0.3	4
42	Hydrogen bond competition between chemical groups: new methodology and the Cambridge Structural Database. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2005, 220, .	0.4	39
43	The Unusual Transformation of an Aromatic 1H-Imidazole into a Non-Aromatic 2H-Imidazole. <i>Structural Chemistry</i> , 2005, 16, 485-490.	1.0	12
44	Neural Network Prediction of Secondary Structure in Crystals: Hydrogen-Bond Systems in Pyrazole Derivatives. <i>Crystal Growth and Design</i> , 2005, 5, 191-200.	1.4	13
45	Hydrogen bond capacity of organic functional groups: a CSD derived database. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2005, 61, c482-c482.	0.3	0
46	Prediction of H-Bonding Motifs for Pyrazoles and Oximes Using the Cambridge Structural Database. <i>Structural Chemistry</i> , 2004, 15, 173-184.	1.0	33
47	The probable number of hydrogen-bonded contacts for chemical groups in organic crystal structures Electronic supplementary information (ESI) available: details of the calculated properties of atoms and groups in Tables 1 and 2. See http://www.rsc.org/suppdata/cc/b4/b402939a/ . <i>Chemical Communications</i> , 2004, , 1166.	2.2	38
48	CSDContact: a database of hydrogen-bond contacts for chemical groups. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2004, 60, s95-s95.	0.3	0
49	The Structure of N1-Hydroxylophine N3-Oxide (=1-Hydroxy-2,4,5- triphenyl-1H-imidazole 3-Oxide) in the Solid State. <i>Helvetica Chimica Acta</i> , 2003, 86, 1026-1039.	1.0	10
50	1,2,4,5-Tetrazines vs. Carboxylic Acid Dimers: Molecular Chemistry vs. Supramolecular Chemistry. <i>Helvetica Chimica Acta</i> , 2003, 86, 1205-1221.	1.0	16
51	Structure of a 4-Nitroso-5-aminopyrazole and Its Salts: Tautomerism, Protonation, and E/Z Isomerism. <i>Journal of Organic Chemistry</i> , 2003, 68, 8831-8837.	1.7	21
52	Extended motifs from water and chemical functional groups in organic molecular crystals. <i>CrystEngComm</i> , 2003, 5, 480.	1.3	430
53	The annular tautomerism of 4(5)-phenylimidazole. <i>Perkin Transactions II RSC</i> , 2002, , 564-568.	1.1	20
54	Water clusters in organic molecular crystals. <i>CrystEngComm</i> , 2002, 4, 454.	1.3	593

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55	Water clusters in organic molecular crystals in the CSD. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2002, 58, c330-c330.	0.3	0
56	2,4,6-Tris(azol-1-yl)-1,3,5-triazines: A New Class of Multidentate Ligands. <i>Heterocycles</i> , 2001, 55, 905.	0.4	18
57	3(5),4-Dimethyl- and 3,4,5-trimethylpyrazole at 200 K. X-ray crystallography and quantum-chemical analysis. <i>Acta Crystallographica Section B: Structural Science</i> , 1999, 55, 441-447.	1.8	10
58	Aminopyrazoles and their conjugated acids. An X-ray study of 3,5-dimethyl-4-aminopyrazole and the Picrate of 3(5)-aminopyrazole. <i>Journal of Heterocyclic Chemistry</i> , 1999, 36, 595-600.	1.4	16
59	Packing Modes in Eight 3-Ethoxycarbonylpyrazole Derivatives. Influence of the Substituents on the Crystal Structure and Annular Tautomerism. <i>Heterocycles</i> , 1999, 50, 227.	0.4	21
60	Tautomerism of NH-pyrazolinones in the solid state: the case of 3(5)-ethoxycarbonyl-5(3)-hydroxypyrazole. <i>Journal of Molecular Structure</i> , 1998, 447, 71-79.	1.8	17
61	The Structure of Aminoazoles and Its Relationship with Aromaticity. Crystal and Molecular Structure of Two Polymorphic Forms of 4-Aminopyrazole. <i>Heterocycles</i> , 1998, 49, 157.	0.4	12
62	Structure and tautomerism of 3(5)-amino-5(3)-arylpyrazoles in the solid state and in solution: An X-ray and NMR study. <i>Tetrahedron</i> , 1997, 53, 10783-10802.	1.0	45
63	Solid-state structure of NH-pyrazolium hydrochlorides and hydrobromides by X-ray crystallography and CPMAS NMR. <i>Journal of Molecular Structure</i> , 1997, 415, 81-92.	1.8	10
64	Mixed crystals of pyrazoles and benzoic acids. Part 1. The molecular structure of 3,5-dimethylpyrazole-2,4,6-trimethylbenzoic acid co-crystals. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1996, , 349-353.	0.9	13
65	Gas-phase (ion cyclotron resonance spectrometric) and solid-state (crystallographic) studies of highly substituted pyrazoles. <i>Journal of Physical Organic Chemistry</i> , 1996, 9, 79-86.	0.9	7
66	Host-guest chemistry. The structure and proton disorder of the three-component crystal formed by 3(5)-methyl-4-nitropyrazole, (R, R)-(?)-trans-4,5-bis (hydroxydiphenylmethyl)-2,2-dimethyl-1,3-dioxolane and toluene. , 1996, 9, 611-618.		6
67	An orientation on solving supramolecular compounds by vector search. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 1996, 52, C73-C73.	0.3	0
68	New synthetic approaches to condensed pyridazinones: alkylpyridazinyl carbonitriles as building blocks for the synthesis of condensed pyridazinones. <i>Tetrahedron</i> , 1995, 51, 12745-12762.	1.0	55
69	Optical resolution of 1,3-dimethyl-5-phenyl-1 ^H -2-pyrazoline by diastereoisomeric complex formation with an optically active host compound: X-ray and molecular structure of the complex. <i>Journal of the Chemical Society Chemical Communications</i> , 1995, , 1453-1454.	2.0	21
70	Structure-Based Modulation of the Ligand Sensitivity of a Tomato Dimeric Abscisic Acid Receptor Through a Glu to Asp Mutation in the Latch Loop. <i>Frontiers in Plant Science</i> , 0, 13, .	1.7	2