

Filomena Martins

List of Publications by Year in descending order

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52
papers

1,193
citations

471061

17
h-index

414034

32
g-index

52
all docs

52
docs citations

52
times ranked

1125
citing authors

#	ARTICLE	IF	CITATIONS
1	In vitro Evaluation of Isoniazid Derivatives as Potential Agents Against Drug-Resistant Tuberculosis. <i>Frontiers in Pharmacology</i> , 2022, 13, .	1.6	2
2	The separation between solvent polarizability and solvent dipolarity: Revisiting the Kamlet-Abraham-Taft model equation. <i>Journal of Molecular Liquids</i> , 2022, 362, 119656.	2.3	5
3	Properties of the <i>tert</i> -butyl halide solvolysis transition states. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 3311-3320.	1.3	4
4	Paving the Way to Fight Multi-Drug Resistant Tuberculosis. <i>Biophysical Journal</i> , 2021, 120, 284a.	0.2	0
5	Designing new antitubercular isoniazid derivatives with improved reactivity and membrane trafficking abilities. <i>Biomedicine and Pharmacotherapy</i> , 2021, 144, 112362.	2.5	11
6	Standardization of antimicrobial testing of dental devices. <i>Dental Materials</i> , 2020, 36, e59-e73.	1.6	33
7	Probing Substrate/Catalyst Effects Using QSPR Analysis on Friedel-Crafts Acylation Reactions over Hierarchical BEA Zeolites. <i>Molecules</i> , 2020, 25, 5682.	1.7	4
8	Reply to the short communication "Comments on Quantifying solvent effects through QSPR: A new look over different model equations". <i>Journal of Molecular Liquids</i> , 2020, 310, 113108.	2.3	1
9	Cinnamic Derivatives as Antitubercular Agents: Characterization by Quantitative Structure-Activity Relationship Studies. <i>Molecules</i> , 2020, 25, 456.	1.7	9
10	Lipophilicity assessment of some isoniazid derivatives active against <i>Mycobacterium tuberculosis</i> . <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2020, 599, 124820.	2.3	3
11	Quantifying solvent effects through QSPR: A new look over different model equations. <i>Journal of Molecular Liquids</i> , 2019, 291, 111244.	2.3	15
12	Zooming in with QSPR on Friedel-Crafts acylation reactions over modified BEA zeolites. <i>Molecular Catalysis</i> , 2019, 476, 110495.	1.0	8
13	Using solvatochromic probes to investigate intermolecular interactions in 1,4-dioxane/methanol/acetonitrile solvent mixtures. <i>Journal of Molecular Liquids</i> , 2018, 266, 259-268.	2.3	8
14	The role of ethanol-water solvent mixtures in N719 sensitization of electrodeposited ZnO nanorods. <i>Journal of Solid State Electrochemistry</i> , 2018, 22, 2779-2787.	1.2	1
15	Volumetric and refractive index study of the ternary mixture methanol/formamide/acetonitrile at 298.15 K. <i>Journal of Molecular Liquids</i> , 2017, 234, 463-468.	2.3	6
16	Kinetic study of Friedel-Crafts acylation reactions over hierarchical MCM-22 zeolites. <i>Molecular Catalysis</i> , 2017, 434, 175-183.	1.0	19
17	Revisiting the Reactions of <i>tert</i> -BuX (X = Br, I) with Monoalcohols: A Mechanistic Analysis through Numerical Integration and Nonlinear Regression Methods. <i>International Journal of Chemical Kinetics</i> , 2017, 49, 100-111.	1.0	0
18	Insights on the Mechanism of Action of INH-C ₁₀ as an Antitubercular Prodrug. <i>Molecular Pharmaceutics</i> , 2017, 14, 4597-4605.	2.3	15

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19	Structure-property relationships in protic ionic liquids: a study of solvent-solvent and solvent-solute interactions. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 28133-28138.	1.3	26
20	Structure-property relationships in protic ionic liquids: a thermochemical study. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 19928-19936.	1.3	15
21	Synthesis and Biological Evaluation of Hybrid 1,5- and 2,5-Disubstituted Indoles as Potentially New Antitubercular Agents. <i>Medicinal Chemistry</i> , 2017, 13, 439-447.	0.7	5
22	Use of quantitative structure-property relationships to study the solvation process of 18-crown-6. <i>Thermochimica Acta</i> , 2015, 604, 140-144.	1.2	15
23	Molecular Details of INH-C ₁₀ Binding to <i>wt</i> KatG and Its S315T Mutant. <i>Molecular Pharmaceutics</i> , 2015, 12, 898-909.	2.3	12
24	Design, synthesis and biological evaluation of novel isoniazid derivatives with potent antitubercular activity. <i>European Journal of Medicinal Chemistry</i> , 2014, 81, 119-138.	2.6	97
25	UV-Vis spectroscopic study of preferential solvation and intermolecular interactions in methanol/1-propanol/acetonitrile by means of solvatochromic probes. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2014, 124, 470-479.	2.0	23
26	Comparison of Multiple Linear Regressions and Neural Networks based QSAR models for the design of new antitubercular compounds. <i>European Journal of Medicinal Chemistry</i> , 2013, 70, 831-845.	2.6	43
27	Solution enthalpies of 1,4-dioxane: Study of solvent effects through quantitative structure-property relationships. <i>Thermochimica Acta</i> , 2013, 574, 85-87.	1.2	12
28	A new approach for the extraction of pollutants from wastewaters handled by the graphic industry. <i>Journal of Environmental Management</i> , 2013, 122, 99-104.	3.8	0
29	QSAR Based Design of New Antitubercular Compounds: Improved Isoniazid Derivatives Against Multidrug-Resistant TB. <i>Current Pharmaceutical Design</i> , 2013, 20, 4427-4454.	0.9	18
30	Acidity and Hydrophobicity of Several New Potential Antitubercular Drugs: Isoniazid and Benzimidazole Derivatives. <i>Journal of Chemical & Engineering Data</i> , 2012, 57, 330-338.	1.0	43
31	Solution enthalpies of hydroxylic compounds. <i>Journal of Thermal Analysis and Calorimetry</i> , 2012, 108, 761-767.	2.0	9
32	Densities and refractive indices for the ternary mixture methanol/propan-1-ol/acetonitrile. <i>Journal of Molecular Liquids</i> , 2012, 170, 30-36.	2.3	14
33	Algorithms for Skin Permeability Using Hydrogen Bond Descriptors: the Problem of Steroids. <i>Journal of Pharmacy and Pharmacology</i> , 2011, 49, 858-865.	1.2	104
34	QSAR modeling of antitubercular activity of diverse organic compounds. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2011, 107, 69-74.	1.8	24
35	Solvent effects on solution enthalpies of adamantyl derivatives. <i>Journal of Thermal Analysis and Calorimetry</i> , 2010, 100, 483-491.	2.0	15
36	Enthalpies of Solution of 1-Butyl-3-methylimidazolium Tetrafluoroborate in 15 Solvents at 298.15 K. <i>Journal of Chemical & Engineering Data</i> , 2010, 55, 616-620.	1.0	16

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37	Modeling Preferential Solvation in Ternary Solvent Systems. <i>Journal of Physical Chemistry B</i> , 2009, 113, 3071-3079.	1.2	18
38	Application of Quantitative Structure-Activity Relationships to the Modeling of Antitubercular Compounds. 1. The Hydrazide Family. <i>Journal of Medicinal Chemistry</i> , 2008, 51, 612-624.	2.9	56
39	The Influence of Carbon-Carbon Multiple Bonds on the Solvolyses of Tertiary Alkyl Halides: a Grunwald-Winstein Analysis. <i>International Journal of Molecular Sciences</i> , 2008, 9, 1704-1716.	1.8	6
40	Determination of solvation and specific interaction enthalpies of adamantane derivatives in aprotic solvents. <i>Journal of Chemical Thermodynamics</i> , 2007, 39, 1201-1205.	1.0	19
41	Design of an Excel Spreadsheet To Estimate Rate Constants, Determine Associated Errors, and Choose Curve's Extent. <i>Journal of Chemical Education</i> , 2006, 83, 1879.	1.1	12
42	Solvent and temperature effects on ion association and mobility of 2,6-lutidinium chloride in non-aqueous solvents. <i>Molecular Physics</i> , 2006, 104, 1905-1913.	0.8	8
43	Thermochemistry of 1-bromoadamantane in binary mixtures of water-aprotic solvent. <i>Thermochimica Acta</i> , 2006, 441, 27-29.	1.2	12
44	Solution enthalpies of 1-bromoadamantane in monoalcohols at 298.15K. <i>Thermochimica Acta</i> , 2006, 444, 83-85.	1.2	13
45	Solvation effects in the heterolyses of 3-methylpentanes (X=Cl, Br, I). <i>Journal of Physical Organic Chemistry</i> , 2004, 17, 1061-1066.	0.9	11
46	Human Skin Permeation and Partition: General Linear Free-Energy Relationship Analyses. <i>Journal of Pharmaceutical Sciences</i> , 2004, 93, 1508-1523.	1.6	182
47	Structural characterization of the ternary solvent mixture methanol-acetonitrile-1-propanol. <i>Journal of Physical Organic Chemistry</i> , 2002, 15, 623-630.	0.9	31
48	Hydrogen bonding. 47. Characterization of the ethylene glycol-heptane partition system: Hydrogen bond acidity and basicity of peptides. <i>Journal of Pharmaceutical Sciences</i> , 1999, 88, 241-247.	1.6	58
49	Hydrogen bonding part 46: a review of the correlation and prediction of transport properties by an LFER method: physicochemical properties, brain penetration and skin permeability. <i>Pest Management Science</i> , 1999, 55, 78-88.	0.7	89
50	Hydrogen bonding part 46: a review of the correlation and prediction of transport properties by an LFER method: physicochemical properties, brain penetration and skin permeability. , 1999, 55, 78.		15
51	Hydrogen bonding part 46: A review of the correlation and prediction of transport properties by an LFER method: physicochemical properties, brain penetration and skin permeability??. <i>Pest Management Science</i> , 1999, 55, 78-88.	0.7	12
52	Enthalpies of solution and intermolecular forces.tert-butyl halides in hydroxylic solvents. <i>Journal of Physical Organic Chemistry</i> , 1992, 5, 93-100.	0.9	16