Filomena Martins

List of Publications by Year in descending order

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| 52 | 1,193 | 17 h-index | 32 |
|----------|----------------|--------------|----------------|
| papers | citations | | g-index |
| 52 | 52 | 52 | 1125 |
| all docs | docs citations | times ranked | citing authors |

| # | Article | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | In vitro Evaluation of Isoniazid Derivatives as Potential Agents Against Drug-Resistant Tuberculosis. Frontiers in Pharmacology, 2022, 13 , . | 3.5 | 2 |
| 2 | The separation between solvent polarizability and solvent dipolarity: Revisiting the Kamlet-Abraham-Taft model equation. Journal of Molecular Liquids, 2022, 362, 119656. | 4.9 | 5 |
| 3 | Properties of the <i>tert</i> -butyl halide solvolysis transition states. Physical Chemistry Chemical Physics, 2021, 23, 3311-3320. | 2.8 | 4 |
| 4 | Paving the Way to Fight Multi-Drug Resistant Tuberculosis. Biophysical Journal, 2021, 120, 284a. | 0.5 | 0 |
| 5 | Designing new antitubercular isoniazid derivatives with improved reactivity and membrane trafficking abilities. Biomedicine and Pharmacotherapy, 2021, 144, 112362. | 5.6 | 11 |
| 6 | Standardization of antimicrobial testing of dental devices. Dental Materials, 2020, 36, e59-e73. | 3.5 | 33 |
| 7 | Probing Substrate/Catalyst Effects Using QSPR Analysis on Friedel-Crafts Acylation Reactions over Hierarchical BEA Zeolites. Molecules, 2020, 25, 5682. | 3.8 | 4 |
| 8 | Reply to the short communication "Comments on Quantifying solvent effects through QSPR: A new look over different model equationsâ€. Journal of Molecular Liquids, 2020, 310, 113108. | 4.9 | 1 |
| 9 | Cinnamic Derivatives as Antitubercular Agents: Characterization by Quantitative Structure–Activity Relationship Studies. Molecules, 2020, 25, 456. | 3.8 | 9 |
| 10 | Lipophilicity assessment of some isoniazid derivatives active against Mycobacterium tuberculosis. Colloids and Surfaces A: Physicochemical and Engineering Aspects, 2020, 599, 124820. | 4.7 | 3 |
| 11 | Quantifying solvent effects through QSPR: A new look over different model equations. Journal of Molecular Liquids, 2019, 291, 111244. | 4.9 | 15 |
| 12 | Zooming in with QSPR on Friedel-Crafts acylation reactions over modified BEA zeolites. Molecular Catalysis, 2019, 476, 110495. | 2.0 | 8 |
| 13 | Using solvatochromic probes to investigate intermolecular interactions in 1,4-dioxane/methanol/acetonitrile solvent mixtures. Journal of Molecular Liquids, 2018, 266, 259-268. | 4.9 | 8 |
| 14 | The role of ethanol-water solvent mixtures in N719 sensitization of electrodeposited ZnO nanorods. Journal of Solid State Electrochemistry, 2018, 22, 2779-2787. | 2.5 | 1 |
| 15 | Volumetric and refractive index study of the ternary mixture methanol/formamide/acetonitrile at 298.15 K. Journal of Molecular Liquids, 2017, 234, 463-468. | 4.9 | 6 |
| 16 | Kinetic study of Friedel-Crafts acylation reactions over hierarchical MCM-22 zeolites. Molecular Catalysis, 2017, 434, 175-183. | 2.0 | 19 |
| 17 | Revisiting the Reactions of <i>tâ€</i> BuX (X = Br, I) with Monoalcohols: A Mechanistic Analysis through Numerical Integration and Nonlinear Regression Methods. International Journal of Chemical Kinetics, 2017, 49, 100-111. | 1.6 | 0 |
| 18 | Insights on the Mechanism of Action of INH-C ₁₀ as an Antitubercular Prodrug. Molecular Pharmaceutics, 2017, 14, 4597-4605. | 4.6 | 15 |

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|----|--|-----|-----------|
| 19 | Structure–property relationships in protic ionic liquids: a study of solvent–solvent and solvent–solute interactions. Physical Chemistry Chemical Physics, 2017, 19, 28133-28138. | 2.8 | 26 |
| 20 | Structure–property relationships in protic ionic liquids: a thermochemical study. Physical Chemistry Chemical Physics, 2017, 19, 19928-19936. | 2.8 | 15 |
| 21 | Synthesis and Biological Evaluation of Hybrid 1,5- and 2,5-Disubstituted Indoles as Potentially New Antitubercular Agents. Medicinal Chemistry, 2017, 13, 439-447. | 1.5 | 5 |
| 22 | Use of quantitative structure–property relationships to study the solvation process of 18-crown-6. Thermochimica Acta, 2015, 604, 140-144. | 2.7 | 15 |
| 23 | Molecular Details of INH-C ₁₀ Binding to <i>wt</i> KatG and Its S315T Mutant. Molecular Pharmaceutics, 2015, 12, 898-909. | 4.6 | 12 |
| 24 | Design, synthesis and biological evaluation of novel isoniazid derivatives with potent antitubercular activity. European Journal of Medicinal Chemistry, 2014, 81, 119-138. | 5.5 | 97 |
| 25 | UV–Vis spectroscopic study of preferential solvation and intermolecular interactions in methanol/1-propanol/acetonitrile by means of solvatochromic probes. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2014, 124, 470-479. | 3.9 | 23 |
| 26 | Comparison of Multiple Linear Regressions and Neural Networks based QSAR models for the design of new antitubercular compounds. European Journal of Medicinal Chemistry, 2013, 70, 831-845. | 5.5 | 43 |
| 27 | Solution enthalpies of 1,4-dioxane: Study of solvent effects through quantitative structure–property relationships. Thermochimica Acta, 2013, 574, 85-87. | 2.7 | 12 |
| 28 | A new approach for the extraction of pollutants from wastewaters handled by the graphic industry. Journal of Environmental Management, 2013, 122, 99-104. | 7.8 | 0 |
| 29 | QSAR Based Design of New Antitubercular Compounds: Improved Isoniazid Derivatives Against Multidrug-Resistant TB. Current Pharmaceutical Design, 2013, 20, 4427-4454. | 1.9 | 18 |
| 30 | Acidity and Hydrophobicity of Several New Potential Antitubercular Drugs: Isoniazid and Benzimidazole Derivatives. Journal of Chemical & Engineering Data, 2012, 57, 330-338. | 1.9 | 43 |
| 31 | Solution enthalpies of hydroxylic compounds. Journal of Thermal Analysis and Calorimetry, 2012, 108, 761-767. | 3.6 | 9 |
| 32 | Densities and refractive indices for the ternary mixture methanol/propan-1-ol/acetonitrile. Journal of Molecular Liquids, 2012, 170, 30-36. | 4.9 | 14 |
| 33 | Algorithms for Skin Permeability Using Hydrogen Bond Descriptors: the Problem of Steroids. Journal of Pharmacy and Pharmacology, 2011, 49, 858-865. | 2.4 | 104 |
| 34 | QSAR modeling of antitubercular activity of diverse organic compounds. Chemometrics and Intelligent Laboratory Systems, 2011, 107, 69-74. | 3.5 | 24 |
| 35 | Solvent effects on solution enthalpies of adamantyl derivatives. Journal of Thermal Analysis and Calorimetry, 2010, 100, 483-491. | 3.6 | 15 |
| 36 | Enthalpies of Solution of 1-Butyl-3-methylimidazolium Tetrafluoroborate in 15 Solvents at 298.15 K. Journal of Chemical & Data, 2010, 55, 616-620. | 1.9 | 16 |

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|----|--|-------------|-----------|
| 37 | Modeling Preferential Solvation in Ternary Solvent Systems. Journal of Physical Chemistry B, 2009, 113, 3071-3079. | 2.6 | 18 |
| 38 | Application of Quantitative Structureâ^'Activity Relationships to the Modeling of Antitubercular Compounds. 1. The Hydrazide Family. Journal of Medicinal Chemistry, 2008, 51, 612-624. | 6.4 | 56 |
| 39 | The Influence of Carbon-Carbon Multiple Bonds on the Solvolyses of Tertiary Alkyl Halides: a Grunwald-Winstein Analysis. International Journal of Molecular Sciences, 2008, 9, 1704-1716. | 4.1 | 6 |
| 40 | Determination of solvation and specific interaction enthalpies of adamantane derivatives in aprotic solvents. Journal of Chemical Thermodynamics, 2007, 39, 1201-1205. | 2.0 | 19 |
| 41 | Design of an Excel Spreadsheet To Estimate Rate Constants, Determine Associated Errors, and Choose Curve's Extent. Journal of Chemical Education, 2006, 83, 1879. | 2.3 | 12 |
| 42 | Solvent and temperature effects on ion association and mobility of 2,6-lutidinium chloride in non-aqueous solvents. Molecular Physics, 2006, 104, 1905-1913. | 1.7 | 8 |
| 43 | Thermochemistry of 1-bromoadamantane in binary mixtures of water–aprotic solvent. Thermochimica Acta, 2006, 441, 27-29. | 2.7 | 12 |
| 44 | Solution enthalpies of 1-bromoadamantane in monoalcohols at 298.15K. Thermochimica Acta, 2006, 444, 83-85. | 2.7 | 13 |
| 45 | Solvation effects in the heterolyses of 3â€Xâ€3â€methylpentanes (X = Cl, Br, I). Journal of Physical Orga Chemistry, 2004, 17, 1061-1066. | inic 1.9 | 11 |
| 46 | Human Skin Permeation and Partition: General Linear Freeâ€Energy Relationship Analyses. Journal of Pharmaceutical Sciences, 2004, 93, 1508-1523. | 3.3 | 182 |
| 47 | Structural characterization of the ternary solvent mixture methanol-acetonitrile-1-propanol. Journal of Physical Organic Chemistry, 2002, 15, 623-630. | 1.9 | 31 |
| 48 | Hydrogen bonding. 47. Characterization of the ethylene glycol–heptane partition system: Hydrogen bond acidity and basicity of peptides. Journal of Pharmaceutical Sciences, 1999, 88, 241-247. | 3.3 | 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. Pest Management Science, 1999, 55, 78-88. | 0.4 | 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. Pest Management Science, 1999, 55, 78-88. | 0.4 | 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??. Pest Management Science, 1999, 55, 78-88. | 0.4 | 12 |
| 52 | Enthalpies of solution and intermolecular forces.tert-butyl halides in hydroxylic solvents. Journal of Physical Organic Chemistry, 1992, 5, 93-100. | 1.9 | 16 |