## German L Perlovich

## List of Publications by Citations

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226
papers
citations

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h-index
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ext. papers

3,702
h-index

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avg, IF

5.95
L-index

| #   | Paper  | IF  | Citations |
|-----|--|-----|-----------|
| 226 | Thermodynamics of solutions III: comparison of the solvation of (+)-naproxen with other NSAIDs. <i>European Journal of Pharmaceutics and Biopharmaceutics</i> , <b>2004</b> , 57, 411-20   | 5.7 | 200       |
| 225 | The Polymorphism of Glycine. Thermochemical and structural aspects. <i>Magyar Apr ad Kalem byek</i> , <b>2001</b> , 66, 699-715  | О   | 171       |
| 224 | Thermodynamics of solutions. II. Flurbiprofen and diflunisal as models for studying solvation of drug substances. <i>European Journal of Pharmaceutical Sciences</i> , <b>2003</b> , 19, 423-32  | 5.1 | 128       |
| 223 | The Vapour Pressure and the Enthalpy of Sublimation: Determination by inert gas flow method. <i>Magyar Apr</i> Dad KDlemDyek, <b>1999</b> , 57, 225-234  | О   | 104       |
| 222 | Thermodynamics of sublimation, crystal lattice energies, and crystal structures of racemates and enantiomers: (+)- and (+/-)-ibuprofen. <i>Journal of Pharmaceutical Sciences</i> , <b>2004</b> , 93, 654-66   | 3.9 | 74        |
| 221 | Thermodynamic characteristics of cocrystal formation and melting points for rational design of pharmaceutical two-component systems. <i>CrystEngComm</i> , <b>2015</b> , 17, 7019-7028   | 3.3 | 68        |
| 220 | Towards an understanding of the molecular mechanism of solvation of drug molecules: a thermodynamic approach by crystal lattice energy, sublimation, and solubility exemplified by paracetamol, acetanilide, and phenacetin. <i>Journal of Pharmaceutical Sciences</i> , <b>2006</b> , 95, 2158-69 | 3.9 | 67        |
| 219 | N-(2-Chlorophenyl)benzenesulfonamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , <b>2006</b> , 62, o780-o782  |     | 66        |
| 218 | Pharmaceutical cocrystals of diflunisal and diclofenac with theophylline. <i>Molecular Pharmaceutics</i> , <b>2014</b> , 11, 3707-15   | 5.6 | 62        |
| 217 | Solvation of drugs as a key for understanding partitioning and passive transport exemplified by NSAIDs. <i>Current Drug Delivery</i> , <b>2004</b> , 1, 213-26   | 3.2 | 50        |
| 216 | Salicylamide cocrystals: screening, crystal structure, sublimation thermodynamics, dissolution, and solid-state DFT calculations. <i>Journal of Physical Chemistry B</i> , <b>2014</b> , 118, 6803-14  | 3.4 | 49        |
| 215 | Thermodynamics of solubility, sublimation and solvation processes of parabens. <i>European Journal of Pharmaceutical Sciences</i> , <b>2005</b> , 24, 25-33  | 5.1 | 49        |
| 214 | Thermodynamic and Structural Aspects of Some Fenamate Molecular Crystals. <i>Crystal Growth and Design</i> , <b>2009</b> , 9, 3265-3272  | 3.5 | 46        |
| 213 | Polymorphism of paracetamol. Journal of Thermal Analysis and Calorimetry, 2007, 89, 767-774  | 4.1 | 46        |
| 212 | Pharmaceutical salts of ciprofloxacin with dicarboxylic acids. <i>European Journal of Pharmaceutical Sciences</i> , <b>2015</b> , 77, 112-21   | 5.1 | 44        |
| 211 | Crystallization and Polymorphism of Felodipine. Crystal Growth and Design, 2012, 12, 4022-4030   | 3.5 | 44        |
| 210 | Solubility, lipophilicity and membrane permeability of some fluoroquinolone antimicrobials. <i>European Journal of Pharmaceutical Sciences</i> , <b>2016</b> , 93, 29-37   | 5.1 | 43        |

| 209 | Cocrystal screening of hydroxybenzamides with benzoic acid derivatives: a comparative study of thermal and solution-based methods. <i>European Journal of Pharmaceutical Sciences</i> , <b>2014</b> , 65, 56-64   | 5.1              | 41 |  |
|-----|---|------------------|----|--|
| 208 | Sulfonamide Molecular Crystals: Structure, Sublimation Thermodynamic Characteristics, Molecular Packing, Hydrogen Bonds Networks. <i>Crystal Growth and Design</i> , <b>2013</b> , 13, 4002-4016  | 3.5              | 41 |  |
| 207 | Solvation and hydration characteristics of ibuprofen and acetylsalicylic acid. <i>AAPS PharmSci</i> , <b>2004</b> , 6, E3   |                  | 41 |  |
| 206 | Novel 1,2,4-thiadiazole derivatives as potent neuroprotectors: approach to creation of bioavailable drugs. <i>Molecular Pharmaceutics</i> , <b>2012</b> , 9, 2156-67  | 5.6              | 40 |  |
| 205 | Sulfonamides as a subject to study molecular interactions in crystals and solutions: sublimation, solubility, solvation, distribution and crystal structure. <i>International Journal of Pharmaceutics</i> , <b>2008</b> , 349, 300-13  | 6.5              | 39 |  |
| 204 | Sublimation of Molecular Crystals: Prediction of Sublimation Functions on the Basis of HYBOT Physicochemical Descriptors and Structural Clusterization. <i>Crystal Growth and Design</i> , <b>2010</b> , 10, 2707-27  | 7∮2 <sup>5</sup> | 37 |  |
| 203 | Partial molar volumes of some drug and pro-drug substances in 1-octanol at T=298.15K. <i>Journal of Chemical Thermodynamics</i> , <b>2010</b> , 42, 429-435   | 2.9              | 37 |  |
| 202 | Novel drugdrug cocrystals of carbamazepine with para-aminosalicylic acid: screening, crystal structures and comparative study of carbamazepine cocrystal formation thermodynamics. <i>CrystEngComm</i> , <b>2017</b> , 19, 4273-4286  | 3.3              | 35 |  |
| 201 | Two-component molecular crystals: evaluation of the formation thermodynamics based on melting points and sublimation data. <i>CrystEngComm</i> , <b>2017</b> , 19, 2870-2883  | 3.3              | 34 |  |
| 200 | Drug-drug cocrystals of antituberculous 4-aminosalicylic acid: Screening, crystal structures, thermochemical and solubility studies. <i>European Journal of Pharmaceutical Sciences</i> , <b>2017</b> , 99, 228-239   | 5.1              | 34 |  |
| 199 | Cocrystals of the antiandrogenic drug bicalutamide: screening, crystal structures, formation thermodynamics and lattice energies. <i>CrystEngComm</i> , <b>2016</b> , 18, 4818-4829   | 3.3              | 33 |  |
| 198 | Sulfonamide Molecular Crystals: Thermodynamic and Structural Aspects. <i>Crystal Growth and Design</i> , <b>2011</b> , 11, 1067-1081  | 3.5              | 32 |  |
| 197 | Thermodynamics of solutions I: benzoic acid and acetylsalicylic acid as models for drug substances and the prediction of solubility. <i>Pharmaceutical Research</i> , <b>2003</b> , 20, 471-8   | 4.5              | 32 |  |
| 196 | Thermodynamics of solutions IV: Solvation of ketoprofen in comparison with other NSAIDs. <i>Journal of Pharmaceutical Sciences</i> , <b>2003</b> , 92, 2502-11  | 3.9              | 32 |  |
| 195 | Influence of Secondary Interactions on the Structure, Sublimation Thermodynamics, and Solubility of Salicylate:4-Hydroxybenzamide Cocrystals. Combined Experimental and Theoretical Study. <i>Journal of Physical Chemistry B</i> , <b>2015</b> , 119, 10466-77                 | 3.4              | 31 |  |
| 194 | Thermodynamic properties of flufenamic and niflumic acidsspecific and non-specific interactions in solution and in crystal lattices, mechanism of solvation, partitioning and distribution. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , <b>2007</b> , 45, 679-87 | 3.5              | 31 |  |
| 193 | Energetic aspects of diclofenac acid in crystal modifications and in solutionsmechanism of solvation, partitioning and distribution. <i>Journal of Pharmaceutical Sciences</i> , <b>2007</b> , 96, 1031-42  | 3.9              | 30 |  |
| 192 | Fenamate Cocrystals with 4,4?-Bipyridine: Structural and Thermodynamic Aspects. <i>Crystal Growth and Design</i> , <b>2015</b> , 15, 228-238  | 3.5              | 27 |  |

| 191 | Weak Interactions Cause Packing Polymorphism in Pharmaceutical Two-Component Crystals. The Case Study of the Salicylamide Cocrystal. <i>Crystal Growth and Design</i> , <b>2017</b> , 17, 1425-1437  | 3.5            | 26 |  |
|-----|--|----------------|----|--|
| 190 | Thermodynamic and structural study of tolfenamic acid polymorphs. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , <b>2009</b> , 50, 831-40  | 3.5            | 26 |  |
| 189 | Driving forces and the influence of the buffer composition on the complexation reaction between ibuprofen and HPCD. <i>European Journal of Pharmaceutical Sciences</i> , <b>2003</b> , 20, 197-200   | 5.1            | 26 |  |
| 188 | Thermodynamics of porphyrin sublimation. <i>Journal of Porphyrins and Phthalocyanines</i> , <b>2000</b> , 04, 699-706  | <b>5</b> 1.8   | 26 |  |
| 187 | Cocrystal formation, crystal structure, solubility and permeability studies for novel 1,2,4-thiadiazole derivative as a potent neuroprotector. <i>European Journal of Pharmaceutical Sciences</i> , <b>2017</b> , 109, 31-39   | 5.1            | 25 |  |
| 186 | Thermodynamic and structural aspects of sulfonamide crystals and solutions. <i>Journal of Pharmaceutical Sciences</i> , <b>2009</b> , 98, 4738-55  | 3.9            | 25 |  |
| 185 | Towards an understanding of the molecular mechanism of solvation of drug molecules: a thermodynamic approach by crystal lattice energy, sublimation, and solubility exemplified by hydroxybenzoic acids. <i>Journal of Pharmaceutical Sciences</i> , <b>2006</b> , 95, 1448-58 | 3.9            | 24 |  |
| 184 | New Solid Forms of the Antiviral Drug Arbidol: Crystal Structures, Thermodynamic Stability, and Solubility. <i>Molecular Pharmaceutics</i> , <b>2015</b> , 12, 4154-65   | 5.6            | 23 |  |
| 183 | Saccharin salts of biologically active hydrazone derivatives. New Journal of Chemistry, 2015, 39, 8614-86  | 5 <b>2</b> 326 | 23 |  |
| 182 | Polymorphism and solvatomorphism of bicalutamide. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2013</b> , 111, 655-662  | 4.1            | 23 |  |
| 181 | Towards the rational design of novel drugs based on solubility, partitioning/distribution, biomimetic permeability and biological activity exemplified by 1,2,4-thiadiazole derivatives.<br>MedChemComm, 2017, 8, 162-175  | 5              | 23 |  |
| 180 | Aqueous Solubilities, Infinite Dilution Activity Coefficients and Octanol Water Partition Coefficients of Tricyclic Analogs of Acyclovir. <i>Journal of Solution Chemistry</i> , <b>1999</b> , 28, 731-745   | 1.8            | 23 |  |
| 179 | Volumetric properties of tetraphenylporphine, their metallo-complexes and some substituted tetraphenylporphines in benzene solution. <i>Journal of Solution Chemistry</i> , <b>1996</b> , 25, 135-153  | 1.8            | 23 |  |
| 178 | Polymorphism of felodipine co-crystals with 4,4?-bipyridine. <i>CrystEngComm</i> , <b>2014</b> , 16, 6603-6611   | 3.3            | 22 |  |
| 177 | Sublimation thermodynamics of four fluoroquinolone antimicrobial compounds. <i>Journal of Chemical Thermodynamics</i> , <b>2017</b> , 105, 37-43   | 2.9            | 22 |  |
| 176 | Extent and mechanism of solvation and partitioning of isomers of substituted benzoic acids: a thermodynamic study in the solid state and in solution. <i>Journal of Pharmaceutical Sciences</i> , <b>2008</b> , 97, 3883-96  | 3.9            | 22 |  |
| 175 | Influence of position and size of substituents on the mechanism of partitioning: a thermodynamic study on acetaminophens, hydroxybenzoic acids, and parabens. <i>AAPS PharmSciTech</i> , <b>2008</b> , 9, 205-16   | 3.9            | 22 |  |
| 174 | Inclusion complex formation of <code>\(\text{B}\) and <code>\(\text{E}\)</code>yclodextrins with riboflavin and alloxazine in aqueous solution: thermodynamic study. Journal of Inclusion Phenomena and Macrocyclic Chemistry, <b>2011</b>, 69, 167</code>                     | 7-172          | 21 |  |

## (2020-2009)

| 173 | Novel isothiourea derivatives as potent neuroprotectors and cognition enhancers: synthesis, biological and physicochemical properties. <i>Journal of Medicinal Chemistry</i> , <b>2009</b> , 52, 1845-52            | 8.3                 | 21 |  |
|-----|---|---------------------|----|--|
| 172 | The difference between partitioning and distribution from a thermodynamic point of view: NSAIDs as an example. <i>European Journal of Pharmaceutical Sciences</i> , <b>2006</b> , 27, 150-7                         | 5.1                 | 21 |  |
| 171 | Specific features of supramolecular organisation and hydrogen bonding in proline cocrystals: a case study of fenamates and diclofenac. <i>CrystEngComm</i> , <b>2018</b> , 20, 6970-6981                            | 3.3                 | 21 |  |
| 170 | Formation Thermodynamics of Two-Component Molecular Crystals: Polymorphism, Stoichiometry, and Impact of Enantiomers. <i>Crystal Growth and Design</i> , <b>2020</b> , 20, 5526-5537                                | 3.5                 | 20 |  |
| 169 | Vapor pressure and sublimation thermodynamics of aminobenzoic acid, nicotinic acid, and related amido-derivatives. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2016</b> , 123, 841-849                  | 4.1                 | 20 |  |
| 168 | Thermodynamic aspects of solubility and partitioning processes of some sulfonamides in the solvents modeling biological media. <i>Journal of Chemical Thermodynamics</i> , <b>2014</b> , 69, 56-65                  | 2.9                 | 20 |  |
| 167 | Acetamidobenzoic acid isomers: Studying sublimation and fusion processes and their relation with crystal structures. <i>Thermochimica Acta</i> , <b>2014</b> , 583, 72-77   | 2.9                 | 20 |  |
| 166 | Thermodynamic study of sublimation, solubility, solvation, and distribution processes of atenolol and pindolol. <i>Molecular Pharmaceutics</i> , <b>2007</b> , 4, 929-35  | 5.6                 | 20 |  |
| 165 | Three Polymorphic Forms of Ciprofloxacin Maleate: Formation Pathways, Crystal Structures, Calculations, and Thermodynamic Stability Aspects. <i>Crystal Growth and Design</i> , <b>2016</b> , 16, 6556-6567         | 3.5                 | 20 |  |
| 164 | Synthesis, pharmacology, crystal properties, and quantitative solvation studies from a drug transport perspective for three new 1,2,4-thiadiazoles. <i>Journal of Pharmaceutical Sciences</i> , <b>2010</b> , 99, 3 | 75 <del>4</del> -68 | 19 |  |
| 163 | Studying thermodynamic aspects of sublimation, solubility and solvation processes and crystal structure analysis of some sulfonamides. <i>International Journal of Pharmaceutics</i> , <b>2007</b> , 334, 115-24    | 6.5                 | 19 |  |
| 162 | Interrelation between thermochemical and structural data of polymorphs exemplified by diflunisal. <i>Journal of Pharmaceutical Sciences</i> , <b>2002</b> , 91, 1036-45   | 3.9                 | 19 |  |
| 161 | Two-component molecular crystals: relationship between the entropy term and the molecular volume of co-crystal formation. <i>CrystEngComm</i> , <b>2018</b> , 20, 3634-3637   | 3.3                 | 19 |  |
| 160 | Synthesis, biological activity, distribution and membrane permeability of novel spiro-thiazines as potent neuroprotectors. <i>European Journal of Medicinal Chemistry</i> , <b>2014</b> , 77, 8-17                  | 6.8                 | 18 |  |
| 159 | Design of pharmaceutical cocrystals for drug solubility improvement. <i>Russian Journal of General Chemistry</i> , <b>2014</b> , 84, 407-414  | 0.7                 | 18 |  |
| 158 | Thermodynamic and Structural Aspects of Hydrated and Unhydrated Phases of 4-Hydroxybenzamide. <i>Crystal Growth and Design</i> , <b>2007</b> , 7, 2643-2648   | 3.5                 | 18 |  |
| 157 | The Melting Process of Acetylsalicylic Acid Single Crystals <b>2001</b> , 63, 653-661   |                     | 18 |  |
| 156 | Cocrystals of Fluconazole with Aromatic Carboxylic Acids: Competition between Anhydrous and Hydrated Solid Forms. <i>Crystal Growth and Design</i> , <b>2020</b> , 20, 1218-1228                                    | 3.5                 | 18 |  |

| 155 | Pharmaceutical Salts of Biologically Active Hydrazone Compound Salinazid: Crystallographic, Solubility, and Thermodynamic Aspects. <i>Crystal Growth and Design</i> , <b>2016</b> , 16, 2605-2617                             | 3.5 | 18 |
|-----|---|-----|----|
| 154 | Effects of the crystal structure and thermodynamic stability on solubility of bioactive compounds: DFT study of isoniazid cocrystals. <i>Computational and Theoretical Chemistry</i> , <b>2016</b> , 1092, 1-11               | 2   | 18 |
| 153 | New Pharmaceutical Cocrystal Forms of Flurbiprofen: Structural, Physicochemical, and Thermodynamic Characterization. <i>Crystal Growth and Design</i> , <b>2019</b> , 19, 5751-5761   | 3.5 | 17 |
| 152 | Prediction of Sublimation Functions of Molecular Crystals Based on Melting Points: Cocrystal Formation Thermodynamics Application. <i>Crystal Growth and Design</i> , <b>2017</b> , 17, 4110-4117                             | 3.5 | 17 |
| 151 | Impact of Sulfonamide Structure on Solubility and Transfer Processes in Biologically Relevant Solvents. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2014</b> , 59, 4217-4226                                       | 2.8 | 17 |
| 150 | Inclusion complex of antiasthmatic compound with 2-hydroxypropyl-Etyclodextrin: Preparation and physicochemical properties. <i>Journal of Molecular Liquids</i> , <b>2017</b> , 237, 185-192                                  | 6   | 16 |
| 149 | Solid Forms of Ciprofloxacin Salicylate: Polymorphism, Formation Pathways, and Thermodynamic Stability. <i>Crystal Growth and Design</i> , <b>2019</b> , 19, 2979-2990  | 3.5 | 16 |
| 148 | The impact of structural modification of 1,2,4-thiadiazole derivatives on thermodynamics of solubility and hydration processes. <i>Physical Chemistry Chemical Physics</i> , <b>2015</b> , 17, 20889-96                       | 3.6 | 16 |
| 147 | Diversity of felodipine solvates: structure and physicochemical properties. <i>CrystEngComm</i> , <b>2015</b> , 17, 4089-4097   | 3.3 | 16 |
| 146 | Diversity of crystal structures and physicochemical properties of ciprofloxacin and norfloxacin salts with fumaric acid. <i>CrystEngComm</i> , <b>2018</b> , 20, 755-767  | 3.3 | 16 |
| 145 | A thermodynamic study of sublimation, dissolution and distribution processes of anti-inflammatory drug Clonixin. <i>Journal of Chemical Thermodynamics</i> , <b>2019</b> , 132, 281-288                                       | 2.9 | 16 |
| 144 | The Pore-Lipid Interface: Role of Amino-Acid Determinants of Lipophilic Access by Ivabradine to the hERG1 Pore Domain. <i>Molecular Pharmacology</i> , <b>2019</b> , 96, 259-271  | 4.3 | 15 |
| 143 | Thermodynamic and structural aspects of novel 1,2,4-thiadiazoles in solid and biological mediums. <i>Molecular Pharmaceutics</i> , <b>2011</b> , 8, 1807-20   | 5.6 | 15 |
| 142 | Thermodynamics of potassium diclofenac salt aqueous solutions at various temperatures. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2011</b> , 104, 279-289  | 4.1 | 15 |
| 141 | Redetermination and H-atom refinement of (S)-(+)-ibuprofen. <i>Acta Crystallographica Section E: Structure Reports Online</i> , <b>2003</b> , 59, o1357-o1358   |     | 15 |
| 140 | Structural and energetic aspects of adamantane and memantine derivatives of sulfonamide molecular crystals: experimental and theoretical characterisation. <i>CrystEngComm</i> , <b>2018</b> , 20, 3476-3489                  | 3.3 | 14 |
| 139 | Crystal architecture and physicochemical properties of felodipine solvates. <i>CrystEngComm</i> , <b>2013</b> , 15, 6054  | 3.3 | 14 |
| 138 | Physicochemical properties/descriptors governing the solubility and partitioning of chemicals in water-solvent-gas systems. Part 2. Solubility in 1-octanol. <i>SAR and QSAR in Environmental Research</i> , 2007, 18, 543-78 | 3.5 | 14 |

| 137 | Intermolecular interactions and permeability of 5-fluorouracil cocrystals with a series of isomeric hydroxybenzoic acids: a combined theoretical and experimental study. <i>CrystEngComm</i> , <b>2019</b> , 21, 5095-   | -517035 | 13 |  |
|-----|--|---------|----|--|
| 136 | Inclusion complexes of hydroxypropyl-Ecyclodextrin with novel cytotoxic compounds: Solubility and thermodynamic properties. <i>Fluid Phase Equilibria</i> , <b>2014</b> , 384, 68-72   | 2.5     | 13 |  |
| 135 | Novel 1,2,4-thiadiazole derivatives: crystal structure, conformational analysis, hydrogen bond networks, calculations, and thermodynamic characteristics of crystal lattices. <i>Journal of Physical Chemistry B</i> , <b>2013</b> , 117, 10414-29   | 3.4     | 13 |  |
| 134 | A study of the inclusion complex of bioactive thiadiazole derivative with 2-hydroxypropyl-Eyclodextrin: Preparation, characterization and physicochemical properties. <i>Journal of Molecular Liquids</i> , <b>2019</b> , 273, 653-662   | 6       | 13 |  |
| 133 | Thermodynamic approaches to the challenges of solubility in drug discovery and development. <i>Molecular Pharmaceutics</i> , <b>2014</b> , 11, 1-11  | 5.6     | 12 |  |
| 132 | Vapor pressures and thermodynamic sublimation of antitubercular drugs. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2015</b> , 120, 1053-1060   | 4.1     | 12 |  |
| 131 | Thermochemistry of Drugs. Experimental and First-Principles Study of Fenamates. <i>Journal of Chemical &amp; Chemica</i> | 2.8     | 12 |  |
| 130 | Melting points of one- and two-component molecular crystals as effective characteristics for rational design of pharmaceutical systems. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials,</i> <b>2020</b> , 76, 696-706  | 1.8     | 12 |  |
| 129 | Novel cocrystals of itraconazole: Insights from phase diagrams, formation thermodynamics and solubility. <i>International Journal of Pharmaceutics</i> , <b>2021</b> , 599, 120441   | 6.5     | 12 |  |
| 128 | Adamantane derivatives of sulfonamide molecular crystals: structure, sublimation thermodynamic characteristics, molecular packing, and hydrogen bond networks. <i>CrystEngComm</i> , <b>2015</b> , 17, 753-763   | 3.3     | 11 |  |
| 127 | Adamantane derivatives of sulfonamides: sublimation, solubility, solvation and transfer processes in biologically relevant solvents. <i>Physical Chemistry Chemical Physics</i> , <b>2016</b> , 18, 9281-94  | 3.6     | 11 |  |
| 126 | Hydrogen Bond Donor/Acceptor Ratios of the Coformers: Do They Really Matter for the Prediction of Molecular Packing in Cocrystals? The Case of Benzamide Derivatives with Dicarboxylic Acids. <i>Crystal Growth and Design</i> , <b>2018</b> , 18, 5254-5269   | 3.5     | 11 |  |
| 125 | Cocrystals of a 1,2,4-thiadiazole-based potent neuroprotector with gallic acid: solubility, thermodynamic stability relationships and formation pathways. <i>Physical Chemistry Chemical Physics</i> , <b>2018</b> , 20, 14469-14481   | 3.6     | 11 |  |
| 124 | Thermodynamic studies of Fenbufen, Diflunisal, and Flurbiprofen: sublimation, solution and solvation of biphenyl substituted drugs. <i>International Journal of Pharmaceutics</i> , <b>2008</b> , 357, 100-7   | 6.5     | 11 |  |
| 123 | Solution thermodynamics of pyrazinamide, isoniazid, and p-aminobenzoic acid in buffers and octanol. <i>Journal of Chemical Thermodynamics</i> , <b>2015</b> , 91, 396-403  | 2.9     | 10 |  |
| 122 | Thermodynamic and structural aspects of hydroxybenzamide molecular crystals study. <i>Thermochimica Acta</i> , <b>2013</b> , 551, 57-61  | 2.9     | 10 |  |
| 121 | Enhancement of dissolution behavior of antiarthritic drug leflunomide using solid dispersion methods. <i>Thermochimica Acta</i> , <b>2017</b> , 656, 123-128   | 2.9     | 10 |  |
| 120 | Crystal Structures, Thermal Analysis, and Dissolution Behavior of New Solid Forms of the Antiviral Drug Arbidol with Dicarboxylic Acids. <i>Crystals</i> , <b>2015</b> , 5, 650-669  | 2.3     | 10 |  |

| 119 | Polymorphs and solvates of felodipine: analysis of crystal structures and thermodynamic aspects of sublimation and solubility processes. <i>CrystEngComm</i> , <b>2012</b> , 14, 8577   | 3.3 | 10 |
|-----|---|-----|----|
| 118 | Thermodynamic aspects of solubility, solvation and partitioning processes of some sulfonamides. <i>Journal of Chemical Thermodynamics</i> , <b>2011</b> , 43, 683-689   | 2.9 | 10 |
| 117 | 5-Fluorouracil Cocrystals with Lipophilic Hydroxy-2-Naphthoic Acids: Crystal Structures, Theoretical Computations, and Permeation Studies. <i>Crystal Growth and Design</i> , <b>2020</b> , 20, 923-933   | 3.5 | 10 |
| 116 | A combined experimental and theoretical study of miconazole salts and cocrystals: crystal structures, DFT computations, formation thermodynamics and solubility improvement. <i>Physical Chemistry Chemical Physics</i> , <b>2021</b> , 23, 12456-12470 | 3.6 | 10 |
| 115 | Pharmaceutical Salts of Fenbendazole with Organic Counterions: Structural Analysis and Solubility Performance. <i>Crystal Growth and Design</i> , <b>2021</b> , 21, 4516-4530   | 3.5 | 10 |
| 114 | Thermodynamic properties of Nalidixic and Oxolinic acids: Experimental and computational study. <i>Thermochimica Acta</i> , <b>2019</b> , 682, 178411   | 2.9 | 9  |
| 113 | Design of 4-aminobenzoic acid two-component molecular crystals: prediction and experiments. <i>CrystEngComm</i> , <b>2019</b> , 21, 2119-2129   | 3.3 | 9  |
| 112 | Poorly soluble drugs: disbalance of thermodynamic characteristics of crystal lattice and solvation. <i>RSC Advances</i> , <b>2016</b> , 6, 77870-77886  | 3.7 | 9  |
| 111 | Solubility and Solution Thermodynamics of Novel Bicyclic Derivatives of 1,3-Selenazine in Biological Relevant Solvents. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2014</b> , 59, 2298-2304   | 2.8 | 9  |
| 110 | Partition coefficients and thermodynamics of transfer of novel drug-like spiro-derivatives in model biological solutions. <i>Journal of Chemical Thermodynamics</i> , <b>2013</b> , 61, 11-17   | 2.9 | 9  |
| 109 | Thermodynamics of Solubility Processes of Novel Drug-like Spiro-Derivatives in Model Biological Solutions. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2012</b> , 57, 1996-2003  | 2.8 | 9  |
| 108 | Monte Carlo studies of drug nucleation 1: formation of crystalline clusters of bicalutamide in water.<br>Journal of Physical Chemistry B, <b>2011</b> , 115, 3062-72  | 3.4 | 9  |
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| 69 | Comparative analysis of solubilization and complexation characteristics for new antifungal compound with cyclodextrins. Impact of cyclodextrins on distribution process. <i>European Journal of Pharmaceutical Sciences</i> , <b>2020</b> , 154, 105531 | 5.1             | 5                 |
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| 66 | Sulfasalazine: Dissolution and Distribution in Pharmaceutically Relevant Mediums. <i>Journal of Chemical &amp; Data</i> , <b>2017</b> , 62, 123-128   | 2.8             | 4                 |

## (1998-2020)

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| 64 | New derivatives of hydrogenated pyrido[4,3-b]indoles as potential neuroprotectors: Synthesis, biological testing and solubility in pharmaceutically relevant solvents. <i>Saudi Pharmaceutical Journal</i> , <b>2018</b> , 26, 801-809 | 4.4 | 4 |
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| 55 | New diclofenac choline hydrate salt: Synthesis, characterization and solubility. <i>Journal of Molecular Structure</i> , <b>2019</b> , 1198, 126922  | 3.4 | 3 |
| 54 | Comparative analysis of experimental methods for determining thermodynamic parameters of formation of multi-component molecular crystals: Benefits and limitations. <i>Journal of Molecular Liquids</i> , <b>2019</b> , 295, 111644    | 6   | 3 |
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| 28 | Physico-chemical study of bioactive N-(5-ethyl-1,3,4-thiadiazole-2-yl)-4-nitrobenzamide: Sublimation, solubility and distribution. <i>Thermochimica Acta</i> , <b>2017</b> , 657, 72-78  | 2.9   | 1 |
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| 22 | Redetermination of 3-hydroxybenzamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , <b>2007</b> , 63, o2359-o2360   |       | 1 |
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| 18 | Isavuconazole: Thermodynamic Evaluation of Processes Sublimation, Dissolution and Partition in Pharmaceutically Relevant Media. <i>Molecules</i> , <b>2021</b> , 26,   | 4.8   | 1 |
| 17 | Sublimation thermodynamics of antifungal drugs: Tioconazole, miconazole and climbazole. <i>Fluid Phase Equilibria</i> , <b>2021</b> , 544-545, 113098  | 2.5   | 1 |
| 16 | Solubility and lipophilicity of antiarrhythmic drug Dofetilide in modeling physiological media.<br>Journal of Chemical Thermodynamics, <b>2021</b> , 161, 106512   | 2.9   | 1 |
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