

Maciej Przybyłek

List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

29
papers

359
citations

13
h-index

18
g-index

31
ext. papers

480
ext. citations

3.8
avg, IF

4.35
L-index

#	Paper	IF	Citations
29	Natural Deep Eutectic Solvents as Agents for Improving Solubility, Stability and Delivery of Curcumin. <i>Pharmaceutical Research</i> , 2019 , 36, 116	4.5	65
28	Distinguishing Cocrystals from Simple Eutectic Mixtures: Phenolic Acids as Potential Pharmaceutical Coformers. <i>Crystal Growth and Design</i> , 2018 , 18, 3524-3534	3.5	28
27	Selection of effective cocrystals former for dissolution rate improvement of active pharmaceutical ingredients based on lipoaffinity index. <i>European Journal of Pharmaceutical Sciences</i> , 2017 , 107, 87-96	5.1	21
26	Color prediction from first principle quantum chemistry computations: a case of alizarin dissolved in methanol. <i>New Journal of Chemistry</i> , 2012 , 36, 1836	3.6	20
25	Applicability of Phenolic Acids as Effective Enhancers of Cocrystal Solubility of Methylxanthines. <i>Crystal Growth and Design</i> , 2017 , 17, 2186-2193	3.5	19
24	Propensity of salicylamide and ethenzamide cocrystallization with aromatic carboxylic acids. <i>European Journal of Pharmaceutical Sciences</i> , 2016 , 85, 132-40	5.1	19
23	Formation of chlorinated breakdown products during degradation of sunscreen agent, 2-ethylhexyl-4-methoxycinnamate in the presence of sodium hypochlorite. <i>Environmental Science and Pollution Research</i> , 2016 , 23, 1886-97	5.1	19
22	Experimental and theoretical studies on the photodegradation of 2-ethylhexyl 4-methoxycinnamate in the presence of reactive oxygen and chlorine species. <i>Open Chemistry</i> , 2014 , 12, 612-623	1.6	19
21	On the origin of surface imposed anisotropic growth of salicylic and acetylsalicylic acids crystals during droplet evaporation. <i>Journal of Molecular Modeling</i> , 2015 , 21, 49	2	16
20	Solubility advantage of sulfanilamide and sulfacetamide in natural deep eutectic systems: experimental and theoretical investigations. <i>Drug Development and Industrial Pharmacy</i> , 2019 , 45, 1120-1129	3.6	15
19	Exploring the cocrystallization potential of urea and benzamide. <i>Journal of Molecular Modeling</i> , 2016 , 22, 103	2	15
18	Application of Multivariate Adaptive Regression Splines (MARSplines) Methodology for Screening of Dicarboxylic Acid Cocrystal Using 1D and 2D Molecular Descriptors. <i>Crystal Growth and Design</i> , 2019 , 19, 3876-3887	3.5	13
17	Utilization of oriented crystal growth for screening of aromatic carboxylic acids cocrystallization with urea. <i>Journal of Crystal Growth</i> , 2016 , 433, 128-138	1.6	13
16	On the origin of surfaces-dependent growth of benzoic acid crystal inferred through the droplet evaporation method. <i>Structural Chemistry</i> , 2015 , 26, 705-712	1.8	12
15	Solubility of sulfanilamide in binary solvents containing water: Measurements and prediction using Buchowski-Ksiazczak solubility model. <i>Journal of Molecular Liquids</i> , 2020 , 319, 114342	6	11
14	Reaction of aniline with ammonium persulphate and concentrated hydrochloric acid: Experimental and DFT studies. <i>Chemical Papers</i> , 2012 , 66,	1.9	9
13	Experimental and theoretical solubility advantage screening of bi-component solid curcumin formulations. <i>Journal of Drug Delivery Science and Technology</i> , 2019 , 50, 125-135	4.5	8

12	Studies on the formation of formaldehyde during 2-ethylhexyl 4-(dimethylamino)benzoate demethylation in the presence of reactive oxygen and chlorine species. <i>Environmental Science and Pollution Research</i> , 2017 , 24, 8049-8061	5.1	5
11	Application of Multivariate Adaptive Regression Splines (MARSplines) for Predicting Hansen Solubility Parameters Based on 1D and 2D Molecular Descriptors Computed from SMILES String. <i>Journal of Chemistry</i> , 2019 , 2019, 1-15	2.3	5
10	Solvent Screening for Solubility Enhancement of Theophylline in Neat, Binary and Ternary NADES Solvents: New Measurements and Ensemble Machine Learning. <i>International Journal of Molecular Sciences</i> , 2021 , 22,	6.3	5
9	Studies on the solid-liquid equilibria and intermolecular interactions Urea binary mixtures with Sulfanilamide and Sulfacetamide. <i>Journal of Chemical Thermodynamics</i> , 2021 , 153, 106308	2.9	5
8	Experimental and theoretical studies on the Sulfamethazine-Urea and Sulfamethizole-Urea solid-liquid equilibria. <i>Journal of Drug Delivery Science and Technology</i> , 2021 , 61, 102186	4.5	4
7	Experimental and Theoretical Screening for Green Solvents Improving Sulfamethizole Solubility. <i>Materials</i> , 2021 , 14,	3.5	3
6	Thermodynamic Characteristics of Phenacetin in Solid State and Saturated Solutions in Several Neat and Binary Solvents. <i>Molecules</i> , 2021 , 26,	4.8	3
5	Predicting Value of Binding Constants of Organic Ligands to Beta-Cyclodextrin: Application of MARSplines and Descriptors Encoded in SMILES String. <i>Symmetry</i> , 2019 , 11, 922	2.7	2
4	Thermodynamics and Intermolecular Interactions of Nicotinamide in Neat and Binary Solutions: Experimental Measurements and COSMO-RS Concentration Dependent Reactions Investigations. <i>International Journal of Molecular Sciences</i> , 2021 , 22,	6.3	2
3	The use of fast molecular descriptors and artificial neural networks approach in organochlorine compounds electron ionization mass spectra classification. <i>Environmental Science and Pollution Research</i> , 2019 , 26, 28188-28201	5.1	1
2	Application 2D Descriptors and Artificial Neural Networks for Beta-Glucosidase Inhibitors Screening. <i>Molecules</i> , 2020 , 25,	4.8	1
1	New Screening Protocol for Effective Green Solvents Selection of Benzamide, Salicylamide and Ethenzamide. <i>Molecules</i> , 2022 , 27, 3323	4.8	0