

Wang Mingliang

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Preferential solvation and solute-solvent interactions of posaconazole in mixtures of (ethyl) Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 347 Td (106661.	2.0	0
2	Research on solubility modelling of actarit in different solvents: Solvent effect, Hansen solubility parameter, molecular interactions and solution thermodynamics. Journal of Molecular Liquids, 2022, 347, 117965.	4.9	9
3	Investigation on co-solvency, solvent effect, Hansen solubility parameter and preferential solvation of fenbufen dissolution and models correlation. Journal of Molecular Liquids, 2022, 348, 118415.	4.9	14
4	Research on Solubility Measurement, Solvent Effects, Preferential Solvation, and Model Correlation of Sofosbuvir Form A in Different Pure and Binary Solvents. Journal of Chemical & Engineering Data, 2022, 67, 748-760.	1.9	1
5	Solid-liquid phase equilibrium of SOV-I2 in various pure and binary mixed solvents: Solubility determination, solvent effect, preferential solvation and model correlation. Journal of Chemical Thermodynamics, 2022, 171, 106794.	2.0	1
6	Research on dissolution of actarit in aqueous mixtures: Solubility determination and correlation, preferential solvation, solvent effect and thermodynamics. Journal of Molecular Liquids, 2022, 358, 119141.	4.9	13
7	Dual-Mode Aptasensor Assembled by a WO ₃ /Fe ₂ O ₃ Heterojunction for Paper-Based Colorimetric Prediction/Photoelectrochemical Multicomponent Analysis. ACS Applied Materials & Interfaces, 2021, 13, 3645-3652.	8.0	42
8	Solubility Measurement and the Correlation of Cilostazol in Pure and 1,4-Dioxane + Ethanol Binary Solvents from $T = 273.15$ to 313.15 K. Journal of Chemical & Engineering Data, 2021, 66, 2895-2900.	1.9	2
9	Solubility of 5-Fluorocytosine in Different Pure and Binary Mixed Solvents: Measurement, Model Correlation, Solvent Effect, and Preferential Solvation. Journal of Chemical & Engineering Data, 2021, 66, 3090-3100.	1.9	1
10	Investigation on the Hansen solubility parameter, solvent effect and thermodynamic analysis of indapamide dissolution and molecular dynamics simulation. Journal of Molecular Liquids, 2021, 334, 116489.	4.9	18
11	Dissolution behavior and preferential solvation of diacerein in mixtures of (ethyl) Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 347 Td (106482.	2.0	7
12	Solubility modelling, solvent effect, preferential solvation and solution thermodynamic of thymine form AH A° in ten mono solvents and two solvent mixtures. Journal of Molecular Liquids, 2020, 300, 112257.	4.9	27
13	Determination, Construction, and Evaluation of Ternary and Quaternary Solid-Liquid Phase Equilibrium of Uric Acid, Adenine, and Guanine in Water. Journal of Chemical & Engineering Data, 2020, 65, 2133-2143.	1.9	2
14	Construction and evaluation of ternary solid-liquid phase diagram of pyraclostrobin (form IV) and its intermediate in ethanol and N,N-dimethylformamide. Journal of Chemical Thermodynamics, 2019, 128, 1-9.	2.0	8
15	Binary Solid-Liquid Solubility Determination and Model Correlation of Quizalofop-ethyl in Different Pure Solvents. Journal of Chemical & Engineering Data, 2019, 64, 1611-1621.	1.9	17
16	Solubility Increment and Thermodynamic Analysis of Bioactive Antofloxacin Hydrochloride in Aqueous ChCl/PTS Deep Eutectic Solvent and Cosolvent Mixtures. Journal of Chemical & Engineering Data, 2019, 64, 5748-5754.	1.9	5
17	Bilinear Staphylococcus aureus detection based on suspension immunoassay. Talanta, 2019, 192, 154-159.	5.5	12
18	Investigation of Charge-Transfer Interaction in Mixed Stack Donor-Acceptor Cocrystals Toward Tunable Solid-State Emission Characteristics. Crystal Growth and Design, 2018, 18, 6001-6008.	3.0	51

#	ARTICLE	IF	CITATIONS
19	Research on Dissolution Capability of Several Antofloxacin Salts. Journal of Chemical & Engineering Data, 2018, 63, 3018-3026.	1.9	2
20	A novel one-pot approach to oxidative aromatization and bromination of pyrazolidin-3-one with HBr-H ₂ O ₂ system. Heterocyclic Communications, 2018, 24, 165-169.	1.2	3
21	Molecular Marriage via Charge Transfer Interaction in Organic Charge Transfer Co-Crystals toward Solid-State Fluorescence Modulation. Crystal Growth and Design, 2017, 17, 1251-1257.	3.0	65
22	Understanding charge transfer stacking mode in cocrystals involving Tetrachloro- <i>p</i> -benzoquinone via experimental (SXR, DSC, TGA, DRS) studies and hirshfeld surfaces analysis. Crystal Research and Technology, 2017, 52, 1600329.	1.3	2
23	Study of H-bonded assemblies of the solvates of anthracene derivatives: guest effect on the crystal symmetry and spectroscopic properties. Supramolecular Chemistry, 2017, 29, 497-505.	1.2	8
24	Co-crystals with Delayed Fluorescence Assembled by 1,4-Diiodotetrafluorobenzene and Polycyclic Aromatic Compounds via Halogen Bonds. ChemistrySelect, 2017, 2, 6323-6330.	1.5	10
25	Efficient Luminescent Microtubes of Charge-Transfer Organic Cocrystals Involving 1,2,4,5-Tetracyanobenzene, Carbazole Derivatives, and Pyrene Derivatives. Crystal Growth and Design, 2017, 17, 6684-6691.	3.0	25
26	Understanding Charge-Transfer Interaction Mode in Cocrystals and Solvates of 1-Phenyl-3-(pyren-1-yl) Prop-2-en-1-one and TCNQ. Crystal Growth and Design, 2015, 15, 4032-4038.	3.0	31
27	Synthesis, photoluminescence properties and theoretical insights on 1,3-bis(diphenyl(9-anthryl)amino)pyrazoline and 1,3-bis(diphenyl(9-anthryl)amino)pyrazole. Luminescence, 2013, 28, 628-633.	2.9	11
28	Tuning Solid-State Fluorescence of a Twisted π -Conjugated Molecule by Regulating the Arrangement of Anthracene Fluorophores. Crystal Growth and Design, 2012, 12, 5986-5993.	3.0	63
29	Experiment and Computation of Solubility and Dissolution Properties for Enalapril Maleate and Its Intermediate in Pure Solvents. Journal of Chemical & Engineering Data, 0, , .	1.9	1