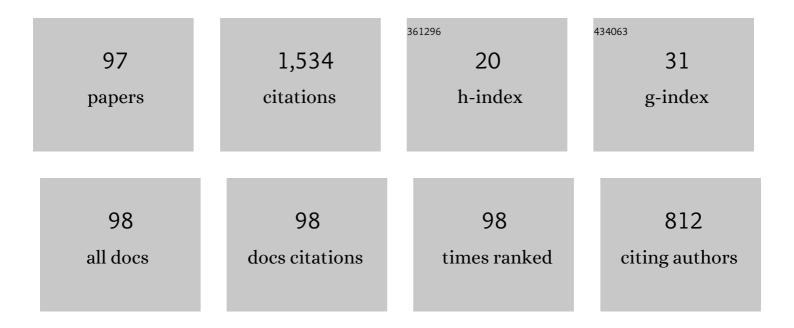
## List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Dissolution behavior, thermodynamic and kinetic analysis of malonamide by experimental measurement and molecular simulation. Chinese Journal of Chemical Engineering, 2023, 53, 260-269.	1.7	2
2	A new strategy to design isostructural salts: The case of the antitumor drug dimethylaminomicheliolide. Chinese Chemical Letters, 2023, 34, 107504.	4.8	0
3	Ultrasound assisted crystallization of cephalexin monohydrate: Nucleation mechanism and crystal habit control. Chinese Journal of Chemical Engineering, 2022, 41, 430-440.	1.7	3
4	Machine learning-based solubility prediction and methodology evaluation of active pharmaceutical ingredients in industrial crystallization. Frontiers of Chemical Science and Engineering, 2022, 16, 523-535.	2.3	10
5	Revealing the dissolution behavior of trans-p-methoxycinnamic acid in 12 organic solvents by parametric model and molecular simulation. Journal of Chemical Thermodynamics, 2022, 166, 106683.	1.0	5
6	Structural Insights into the Highly Solvating System of Axitinib via Binary and Ternary Solvates. Crystal Growth and Design, 2022, 22, 1083-1093.	1.4	5
7	Asparagine endopeptidase-targeted Ultrasound-responsive Nanobubbles Alleviate Tau Cleavage and Amyloid-β Deposition in an Alzheimer's Disease Model. Acta Biomaterialia, 2022, 141, 388-397.	4.1	15
8	Theoretical and Structural Understanding of the Different Factors Influencing the Formation of Multicomponent Crystals of 2,4-Dichlorophenoxyacetic Acid with N-heterocyclic Compounds. Crystal Growth and Design, 2022, 22, 1707-1719.	1.4	5
9	Bendable and Twistable Crystals of Flufenamic Acid Form III with Bending Mechanofluorochromism Behavior. Crystal Growth and Design, 2022, 22, 1312-1318.	1.4	17
10	Insoluble Salt of Memantine with a Unique Fluorescence Phenomenon. Molecular Pharmaceutics, 2022, , .	2.3	2
11	The heterogeneous nucleation of pimelic acid under the effect of a template: experimental research and molecular simulation. CrystEngComm, 2022, 24, 2825-2835.	1.3	0
12	Mechanical Motion and Modulation of Thermalâ€Actuation Properties in a Robust Organic Molecular Crystal Actuator. Advanced Functional Materials, 2022, 32, .	7.8	16
13	Multiple Mechanical Behaviors in One Crystal of 2,4-Dichlorophenoxyacetic Acid Form II: Thermomechanical Effect and Elastic Deformation. Crystal Growth and Design, 2022, 22, 3680-3687.	1.4	4
14	Insight into the Nucleation Mechanism of p-Methoxybenzoic Acid in Ethanol-Water System from Metastable Zone Width. Molecules, 2022, 27, 4085.	1.7	3
15	Nucleation behavior of isosorbide 5-mononitrate revealed from metastable zone widths by combining nucleation theory model and molecular simulation. Journal of Molecular Liquids, 2022, 363, 119846.	2.3	8
16	Similar but Not the Same: Difference in the Ability to Form Cocrystals between Nimesulide and the Pyridine Analogues. Crystal Growth and Design, 2021, 21, 287-296.	1.4	8
17	Optimizing the morphology of calcium <scp>d</scp> -pantothenate by controlling phase transformation processes. CrystEngComm, 2021, 23, 2162-2173.	1.3	2
18	Unraveling the Molecular Mechanisms That Influence the Color and Stability of Four Lutein Crystal Forms. Crystal Growth and Design, 2021, 21, 1762-1777.	1.4	2

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19	Structural Origins of Elastic and 2D Plastic Flexibility of Molecular Crystals Investigated with Two Polymorphs of Conformationally Rigid Coumarin. Chemistry of Materials, 2021, 33, 1053-1060.	3.2	50
20	Use of additives to regulate solute aggregation and direct conformational polymorph nucleation of pimelic acid. IUCrJ, 2021, 8, 161-167.	1.0	10
21	Intermolecular Interactions and Solubility Behavior of Multicomponent Crystal Forms of Orotic Acid: Prediction and Experiments. Crystal Growth and Design, 2021, 21, 1473-1481.	1.4	19
22	Exploring Solid Form Landscape of Anticancer Drug Dimethylaminomicheliolide Fumarate: Crystal Structures Analysis, Phase Transformation Behavior, and Physicochemical Properties Characterization. Crystal Growth and Design, 2021, 21, 2643-2652.	1.4	6
23	Preparation and Formation Mechanism of <scp>l</scp> -Valine Spherulites via Evaporation Crystallization. Industrial & Engineering Chemistry Research, 2021, 60, 6048-6058.	1.8	12
24	Tuning morphology of sulfadiazine through phase transformation of two novel solvates. Journal of Crystal Growth, 2021, 562, 126087.	0.7	3
25	Thermodynamic analysis and molecular dynamic simulation of solid-liquid phase equilibrium of imazapyr in twelve pure organic solvents. Journal of Molecular Liquids, 2021, 330, 115631.	2.3	33
26	Design of Spherical Crystallization of Active Pharmaceutical Ingredients via a Highly Efficient Strategy: From Screening to Preparation. ACS Sustainable Chemistry and Engineering, 2021, 9, 9018-9032.	3.2	21
27	Development and Structure Analysis of Crystal Forms of Apabetalone: Solvates and Polymorphs. Crystal Growth and Design, 2021, 21, 3864-3873.	1.4	13
28	Rationalizing the Formation of Belinostat Solvates with Experimental Screening and Computational Predictions. Crystal Growth and Design, 2021, 21, 4986-4996.	1.4	7
29	Template design based on molecular and crystal structure similarity to regulate conformational polymorphism nucleation: the case of α,ï‰-alkanedicarboxylic acids. IUCrJ, 2021, 8, 814-822.	1.0	6
30	Solubility measurement, correlation and computational analysis of p-Acetamidobenzoic acid in 12 pure solvents. Journal of Chemical Thermodynamics, 2021, 159, 106478.	1.0	4
31	Ultrasound-assisted solution crystallization of fotagliptin benzoate: Process intensification and crystal product optimization. Ultrasonics Sonochemistry, 2021, 76, 105634.	3.8	14
32	Enhancing continuous reactive crystallization of lithium carbonate in multistage mixed suspension mixed product removal crystallizers with pulsed ultrasound. Ultrasonics Sonochemistry, 2021, 77, 105698.	3.8	2
33	Insight into the Influential Mechanism of Polymorphic Parent Molecule with High <i>Z</i> ' on the Cocrystal Formation. Crystal Growth and Design, 2021, 21, 6385-6392.	1.4	5
34	Preparation and solid-state characterization of dapsone pharmaceutical cocrystals through the supramolecular synthon strategy. CrystEngComm, 2021, 23, 6690-6702.	1.3	4
35	Determining the solubility and understanding the solid-liquid equilibrium behavior of cyhalothric acid in eleven pure solvents. Journal of Molecular Liquids, 2020, 300, 112365.	2.3	25
36	Solubility Measurement and Data Correlation of 5,5-Dimethylhydantoin in 12 Pure Solvents at Temperatures from 283.15 to 323.15 K. Journal of Chemical & Engineering Data, 2020, 65, 814-820.	1.0	32

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37	Enhancing Stability and Formulation Capability of Fungicides by Cocrystallization through a Novel Multistep Slurry Conversion Process. Crystal Growth and Design, 2020, 20, 7356-7367.	1.4	14
38	The formation mechanism of hollow spherulites and molecular conformation of curcumin and solvate. CrystEngComm, 2020, 22, 8405-8411.	1.3	4
39	The effect of solvents on solid-liquid phase equilibrium of 1,3-Di-o-tolylguanidine. Journal of Molecular Liquids, 2020, 309, 113147.	2.3	14
40	Transformation between Two Types of Spherulitic Growth: Tuning the Morphology of Spherulitic Nitroguanidine in a Gelatin Solution. Industrial & Engineering Chemistry Research, 2020, 59, 21167-21176.	1.8	16
41	Green Mechanochemical Strategy for the Discovery and Selective Preparation of Polymorphs of Active Pharmaceutical Ingredient Î <sup>3</sup> -Aminobutyric Acid (GABA). ACS Sustainable Chemistry and Engineering, 2020, 8, 16781-16790.	3.2	14
42	Ultrasound-assisted intensified crystallization of L-glutamic acid: Crystal nucleation and polymorph transformation. Ultrasonics Sonochemistry, 2020, 68, 105227.	3.8	34
43	Additiveâ€Induced Selective Crystallization of the Elusive Formâ€II of <i>γ</i> â€Aminobutyric Acid. Chemical Engineering and Technology, 2020, 43, 1137-1143.	0.9	2
44	Reply to the â€ <sup>~</sup> Comment on "Polymorphism of levofloxacin: structure, properties and phase transformationâ€â€™ by Tejender S. Thakur, <i>CrystEngComm</i> , 2020, <b>22</b> , DOI: 10.1039/C9CE01400D. CrystEngComm, 2020, 22, 1889-1891.	1.3	1
45	Solubility Measurement and Correlation of Ceftiofur Sodium Trihydrate in Four Binary Solvent Mixtures. Journal of Chemical & Engineering Data, 2020, 65, 916-922.	1.0	3
46	Correlation and Thermodynamic Analysis of Solubility of Mesotrione in Pure Solvents. Journal of Chemical & Engineering Data, 2020, 65, 877-884.	1.0	15
47	Recent Progress in Continuous Crystallization of Pharmaceutical Products: Precise Preparation and Control. Organic Process Research and Development, 2020, 24, 1785-1801.	1.3	57
48	Solubility Determination and Thermodynamic Correlation of 2-Benzimidazolone in Twelve Pure Solvents from 283.15 to 323.15 K. Journal of Chemical & Engineering Data, 2020, 65, 2838-2845.	1.0	12
49	A tolbutamide–metformin salt based on antidiabetic drug combinations: synthesis, crystal structure analysis and pharmaceutical properties. Acta Crystallographica Section C, Structural Chemistry, 2019, 75, 1250-1258.	0.2	12
50	Drug–drug salts of mefenamic acidolfenamic acid and piperazine to improve physicochemical properties for potential veterinary use. CrystEngComm, 2019, 21, 5284-5291.	1.3	10
51	Crystal Structure, Stability and Desolvation of the Solvates of Sorafenib Tosylate. Crystals, 2019, 9, 367.	1.0	20
52	Role of solvent properties and composition on the solid-liquid equilibrium of trifloxystrobin and thermodynamic analysis. Journal of Molecular Liquids, 2019, 294, 111566.	2.3	6
53	Measurement and Correlation of the Solubility of Aminocaproic Acid in Some Pure and Binary Solvents. Journal of Chemical & Engineering Data, 2019, 64, 5312-5323.	1.0	7
54	Solubility and Data Correlation of β-Arbutin in Different Monosolvents from 283.15 to 323.15 K. Journal of Chemical & Engineering Data, 2019, 64, 5688-5697.	1.0	13

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55	Polymorphism of levofloxacin: structure, properties and phase transformation. CrystEngComm, 2019, 21, 6196-6207.	1.3	20
56	Versatile solid forms of boscalid: insight into the crystal structures and phase transformations. CrystEngComm, 2019, 21, 6838-6849.	1.3	8
57	Uncover the effect of solvent on dissolution behavior of dimethylaminomicheliolide fumarate salt. Journal of Molecular Liquids, 2019, 293, 111448.	2.3	6
58	Solubility Measurement and Correlation of Probenecid in 12 Pure Organic Solvents and Thermodynamic Properties of Mixing of Solutions. Journal of Chemical & Engineering Data, 2019, 64, 624-631.	1.0	28
59	Surprising Effect of Carbon Chain Length on Inducing Ability of Additives: Elusive Form-II of γ-Aminobutyric Acid (GABA) Induced by Sodium Carboxylate Additives. Crystal Growth and Design, 2019, 19, 3825-3833.	1.4	13
60	Influence of the Solvent Content on the Phase Transformation of Sulfadiazine N â€Methyl Pyrrolidone Solvate. Chemical Engineering and Technology, 2019, 42, 1435-1445.	0.9	4
61	Polymorphism and molecular conformations of nicosulfuron: structure, properties and desolvation process. CrystEngComm, 2019, 21, 2790-2798.	1.3	15
62	Solubility measurement, correlation and mixing thermodynamics properties of dapsone in twelve mono solvents. Journal of Molecular Liquids, 2019, 280, 175-181.	2.3	36
63	Solubility Determination and Correlation of Glibenclamide in 11 Monosolvents and (Acetone +) Tj ETQq1 1 0.784 2019, 64, 189-201.	314 rgBT 1.0	Overlock 10 23
64	Solid-liquid equilibrium behavior and thermodynamic analysis of dipyridamole in pure and binary solvents from 293.15†K to 328.15†K. Journal of Molecular Liquids, 2019, 275, 8-17.	2.3	47
65	Spherical Crystallization and the Mechanism of Clopidogrel Hydrogen Sulfate. Chemical Engineering and Technology, 2018, 41, 1259-1265.	0.9	15
66	Design and mechanism of the formation of spherical KCl particles using cooling crystallization without additives. Powder Technology, 2018, 329, 455-462.	2.1	32
67	Caking of crystals: Characterization, mechanisms and prevention. Powder Technology, 2018, 337, 51-67.	2.1	49
68	Crystal Structures and Phase Behavior of Sulfadiazine and a Method for the Preparation of Aggregates with Good Performance. Chemical Engineering and Technology, 2018, 41, 532-540.	0.9	8
69	Crystal morphology optimization of thiamine hydrochloride in solvent system: Experimental and molecular dynamics simulation studies. Journal of Crystal Growth, 2018, 481, 48-55.	0.7	17
70	Thermodynamic Study of Solubility for Imatinib Mesylate in Nine Monosolvents and Two Binary Solvent Mixtures from 278.15 to 318.15 K. Journal of Chemical & Engineering Data, 2018, 63, 4114-4127.	1.0	13
71	Measurement and Correlation of the Solubility of Pyrimethanil in Seven Monosolvents and Two Different Binary Mixed Solvents. Journal of Chemical & Engineering Data, 2018, 63, 2804-2812.	1.0	5
72	Temperature and solvent dependent thermodynamic behavior of tetrabromobisphenol A. Journal of Molecular Liquids, 2017, 241, 150-162.	2.3	17

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73	Polymorphs of daidzein and intermolecular interaction effect on solution crystallization. CrystEngComm, 2017, 19, 7146-7153.	1.3	15
74	Polymorph Control by Investigating the Effects of Solvent and Supersaturation on Clopidogrel Hydrogen Sulfate in Reactive Crystallization. Crystal Growth and Design, 2017, 17, 6123-6131.	1.4	24
75	Two novel cocrystals of lamotrigine with isomeric bipyridines and in situ monitoring of the cocrystallization. European Journal of Pharmaceutical Sciences, 2017, 110, 19-25.	1.9	20
76	Solvent-Mediated Nonoriented Self-Aggregation Transformation: A Case Study of Gabapentin. Crystal Growth and Design, 2017, 17, 4207-4216.	1.4	13
77	Solubility Correlation and Thermodynamic Analysis of Sorafenib Free Base and Sorafenib Tosylate in Monosolvents and Binary Solvent Mixtures. Journal of Chemical & Engineering Data, 2017, 62, 259-267.	1.0	67
78	Industrial Crystallization in China. Chemical Engineering and Technology, 2016, 39, 807-814.	0.9	12
79	Agglomeration Mechanism of Azithromycin Dihydrate in Acetone–Water Mixtures and Optimization of the Powder Properties. Industrial & Engineering Chemistry Research, 2016, 55, 4905-4910.	1.8	14
80	Solvent penetration mediated phase transformation for the preparation of aggregated particles with well-defined shape. CrystEngComm, 2016, 18, 9223-9226.	1.3	13
81	Nucleation behavior of eszopiclone-butyl acetate solutions from metastable zone widths. Chemical Engineering Science, 2016, 155, 248-257.	1.9	53
82	Oiling out and Polymorphism Control of Pyraclostrobin in Cooling Crystallization. Industrial & Engineering Chemistry Research, 2016, 55, 11631-11637.	1.8	27
83	Determination and correlation of solubility and solution thermodynamics of oxiracetam in three (alcohol + water) binary solvents. Journal of Chemical Thermodynamics, 2016, 96, 12-23.	1.0	31
84	Correlation and thermodynamic analysis of solubility of diphenhydramine hydrochloride in pure and binary solvents. Journal of Chemical Thermodynamics, 2016, 93, 132-142.	1.0	58
85	Size Control of Atorvastatin Calcium Particles Based on Spherical Agglomeration. Chemical Engineering and Technology, 2015, 38, 1081-1087.	0.9	19
86	Caking and adhesion free energy of maltitol: Studying of mechanism in adhesion process. Powder Technology, 2015, 272, 235-240.	2.1	17
87	Measurement and Correlation of the Solubility of Penicillin V Potassium in Ethanol + Water and 1-Butyl Alcohol + Water Systems. Journal of Chemical & Engineering Data, 2015, 60, 112-117.	1.0	33
88	Solubility of androstenedione in lower alcohols. Fluid Phase Equilibria, 2014, 363, 86-96.	1.4	45
89	The dehydration behavior and non-isothermal dehydration kinetics of donepezil hydrochloride monohydrate (Form I). Frontiers of Chemical Science and Engineering, 2014, 8, 55-63.	2.3	9
90	Solubility of Ibuprofen Sodium Dihydrate in Acetone + Water Mixtures: Experimental Measurement and Thermodynamic Modeling. Journal of Chemical & Engineering Data, 2014, 59, 3415-3421.	1.0	14

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91	Determination of the Solubility, Dissolution Enthalpy and Entropy of Donepezil Hydrochloride Polymorphic Form III in Different Solvents. Journal of Solution Chemistry, 2013, 42, 841-848.	0.6	7
92	Transformations among the New Solid-State Forms of Clindamycin Phosphate. Organic Process Research and Development, 2013, 17, 1445-1450.	1.3	8
93	In Situ Monitoring of the Solvent-Mediated Transformation of Cefadroxil DMF Solvate into Monohydrate. Organic Process Research and Development, 2013, 17, 1110-1116.	1.3	14
94	Determination of Solubility and Induction Time of Ceftazidime. Journal of Chemical & Engineering Data, 2013, 58, 176-182.	1.0	21
95	Solution Thermodynamic Analysis of <i>p</i> -(Aminomethyl)benzoic Acid in Four Binary Solvents from 288.15 to 328.15 K. Journal of Chemical & Engineering Data, 0, , .	1.0	Ο
96	Intermolecular Interactions and Solubility Behavior of Multicomponent Crystal Forms of 2,4-dichlorophenoxyacetic acid: Design, Structure Analysis, and Solid-State Characterization. CrystEngComm, 0, , .	1.3	12
97	Insights into the Role of Dipentaerythritol in the Thermodynamics and Nucleation Behavior of a Pentaerythritol–Water System. Crystal Growth and Design, 0, , .	1.4	5