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List of Publications by Year in descending order

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Version: 2024-02-01

41
papers

916
citations

567144

15
h-index

454834

30
g-index

41
all docs

41
docs citations

41
times ranked

1216
citing authors

#	ARTICLE	IF	CITATIONS
1	The Solid State Landscape of the Sildenafil Drug. <i>Journal of Pharmaceutical Sciences</i> , 2022, 111, 1104-1109.	1.6	6
2	Synthesis and Characterization of a New Norfloxacin/Resorcinol Cocrystal with Enhanced Solubility and Dissolution Profile. <i>Pharmaceutics</i> , 2022, 14, 49.	2.0	16
3	Polymorphism in the 1/1 Pterostilbene/Picolinic Acid Cocrystal. <i>Crystal Growth and Design</i> , 2022, 22, 590-597.	1.4	10
4	Revision of the Crystal Structure of the Orthorhombic Polymorph of Oxyma: On the Importance of π -Hole Interactions and Their Interplay with H-Bonds. <i>Crystals</i> , 2022, 12, 823.	1.0	1
5	A New and Highly Stable Cocrystal of Vitamin D3 for Use in Enhanced Food Supplements. <i>Crystal Growth and Design</i> , 2021, 21, 1418-1423.	1.4	11
6	A Novel, Extremely Bioavailable Cocrystal of Pterostilbene. <i>Crystal Growth and Design</i> , 2021, 21, 2315-2323.	1.4	22
7	Novel Polymorphic Cocrystals of the Non-Steroidal Anti-Inflammatory Drug Niflumic Acid: Expanding the Pharmaceutical Landscape. <i>Pharmaceutics</i> , 2021, 13, 2140.	2.0	9
8	New Cocrystal of Ubiquinol with High Stability to Oxidation. <i>Crystal Growth and Design</i> , 2020, 20, 5583-5588.	1.4	16
9	DFT Analysis of Uncommon π -H-Bond Array Interaction in a New Pterostilbene/Theophylline Cocrystal. <i>Crystal Growth and Design</i> , 2020, 20, 6691-6698.	1.4	19
10	Sildenafil-Resorcinol Cocrystal: XRPD Structure and DFT Calculations. <i>Crystals</i> , 2020, 10, 1126.	1.0	15
11	Crystal engineering of nutraceutical phytosterols: new cocrystal solid solutions. <i>CrystEngComm</i> , 2020, 22, 4210-4214.	1.3	4
12	Morphotropism and π -Quasi-Isostructurality in the Three High Z' Concomitant Polymorphs of Efinaconazole. <i>Crystal Growth and Design</i> , 2020, 20, 4238-4242.	1.4	11
13	A late appearing polymorph of nutraceutical pterostilbene. <i>CrystEngComm</i> , 2020, 22, 4680-4684.	1.3	10
14	Potentiometric CheqSol and standardized shake-flask solubility methods are complimentary tools in physicochemical profiling. <i>European Journal of Pharmaceutical Sciences</i> , 2020, 148, 105305.	1.9	2
15	H-Bonded anion-anion complexes in fentanyl citrate polymorphs and solvates. <i>Chemical Communications</i> , 2019, 55, 115-118.	2.2	26
16	Hydrogen Bond Polarization Overcomes Unfavorable Packing in the Most Stable High Z' Polymorph of Pterostilbene. <i>Crystal Growth and Design</i> , 2019, 19, 2552-2556.	1.4	10
17	Derisking Development by a Cocrystallization Screen of a Novel Selective Inhaled JAK-STAT inhibitor. <i>Crystal Growth and Design</i> , 2019, 19, 403-414.	1.4	2
18	Property prediction and pharmacokinetic evaluation of mixed stoichiometry cocrystals of zafirlukast, a drug delivery case study. <i>CrystEngComm</i> , 2018, 20, 1346-1351.	1.3	6

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19	Assembling the Puzzle of Apixaban Solid Forms. <i>Molecular Pharmaceutics</i> , 2018, 15, 1909-1916.	2.3	3
20	Combined Virtual/Experimental Multicomponent Solid Forms Screening of Sildenafil: New Salts, Cocrystals, and Hybrid Saltâ€“Cocrystals. <i>Crystal Growth and Design</i> , 2018, 18, 7618-7627.	1.4	35
21	Hydrogen bonding <i>versus</i> π -interactions: their key competition in sildenafil solvates. <i>CrystEngComm</i> , 2018, 20, 4526-4530.	1.3	9
22	Polymorphism of Sildenafil: A New Metastable Desolvate. <i>Crystal Growth and Design</i> , 2018, 18, 3740-3746.	1.4	17
23	Solubility-pH profiles of a free base and its salt: sibutramine as a case study. <i>ADMET and DMPK</i> , 2017, 5, 253-256.	1.1	2
24	The Ca^{2+} -EDTA chelation as standard reaction to validate Isothermal Titration Calorimeter measurements (ITC). <i>Talanta</i> , 2016, 154, 354-359.	2.9	22
25	Polymorphism of Cocrystals: The Promiscuous Behavior of Agomelatine. <i>Crystal Growth and Design</i> , 2016, 16, 1063-1070.	1.4	38
26	Two New Polymorphic Cocrystals of Zafirlukast: Preparation, Crystal Structure, and Stability Relations. <i>Crystal Growth and Design</i> , 2015, 15, 4162-4169.	1.4	9
27	Expanding the Crystal Form Landscape of the Antiviral Drug Adefovir Dipivoxil. <i>Crystal Growth and Design</i> , 2015, 15, 475-484.	1.4	9
28	Water wires in the nanoporous form II of carbamazepine: a single-crystal X-ray diffraction analysis. <i>CrystEngComm</i> , 2013, 15, 845-847.	1.3	12
29	A cocrystal is the key intermediates for the production of a new polymorph of Vorinostat. <i>CrystEngComm</i> , 2012, 14, 362-365.	1.3	10
30	Cooperative induction in double H-bonding donor/acceptor compounds: Chains vs. ribbons. <i>CrystEngComm</i> , 2012, 14, 5745.	1.3	22
31	Cooperativity in Solid-State Squaramides. <i>Crystal Growth and Design</i> , 2011, 11, 3725-3730.	1.4	17
32	Revisiting the Solid State of Norfloxacin. <i>Crystal Growth and Design</i> , 2010, 10, 2948-2953.	1.4	18
33	Oxyma: An Efficient Additive for Peptide Synthesis to Replace the Benzotriazoleâ€“Based HOBT and HOAT with a Lower Risk of Explosion ^[1] . <i>Chemistry - A European Journal</i> , 2009, 15, 9394-9403.	1.7	326
34	New polymorphic hydrogen bonding donorâ€“acceptor system with two temperature coincident solidâ€“solid transitions. <i>CrystEngComm</i> , 2009, 11, 52-54.	1.3	27
35	Ziprasidone malate, a new trimorphic salt with improved aqueous solubility. <i>CrystEngComm</i> , 2009, 11, 791.	1.3	21
36	Safety Evaluation of an Unexpected Incident with a Nitro Compound. <i>Organic Process Research and Development</i> , 2007, 11, 1131-1134.	1.3	17

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37	Polymorphism of (S)-triphenylglycol: kinetic dependent transformation of a new multipolymorphic system. <i>Chemical Communications</i> , 2007, , 3538.	2.2	4
38	A new polymorph of Norfloxacin. <i>Journal of Thermal Analysis and Calorimetry</i> , 2007, 89, 687-692.	2.0	46
39	Polymorphism of Norfloxacin: Evidence of the Enantiotropic Relationship between Polymorphs A and B. <i>Crystal Growth and Design</i> , 2006, 6, 1463-1467.	1.4	52
40	Combined crystallographic and computational investigation of the solvent disorder present in a new tipiracil hydrochloride methanol solvate hydrate. <i>CrystEngComm</i> , 0, , .	1.3	1
41	Static discrete disorder in the crystal structure of iododiflunisal: on the importance of hydrogen bond, halogen bond and π -stacking interactions. <i>CrystEngComm</i> , 0, , .	1.3	3