

Jonas Nyman

List of Publications by Year in descending order

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Version: 2024-02-01

15
papers

1,326
citations

759233

12
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940533

16
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docs citations

16
times ranked

1365
citing authors

#	ARTICLE	IF	CITATIONS
1	The Twelfth Solved Structure of ROY: Single Crystals of Y04 Grown from Melt Microdroplets. <i>Crystal Growth and Design</i> , 2020, 20, 7093-7097.	3.0	43
2	Systematic Finite-Temperature Reduction of Crystal Energy Landscapes. <i>Crystal Growth and Design</i> , 2020, 20, 6847-6862.	3.0	21
3	A Prolific Solvate Former, Galunisertib, under the Pressure of Crystal Structure Prediction, Produces Ten Diverse Polymorphs. <i>Journal of the American Chemical Society</i> , 2019, 141, 13887-13897.	13.7	109
4	Accuracy and reproducibility in crystal structure prediction: the curious case of ROY. <i>CrystEngComm</i> , 2019, 21, 2080-2088.	2.6	55
5	The PO13 crystal structure of ROY. <i>CrystEngComm</i> , 2019, 21, 1363-1368.	2.6	40
6	Inconvenient Truths about Solid Form Landscapes Revealed in the Polymorphs and Hydrates of Gandotinib. <i>Crystal Growth and Design</i> , 2019, 19, 2947-2962.	3.0	32
7	Crystal structure prediction is changing from basic science to applied technology. <i>Faraday Discussions</i> , 2018, 211, 459-476.	3.2	58
8	Structure searching methods: general discussion. <i>Faraday Discussions</i> , 2018, 211, 133-180.	3.2	3
9	Crystal structure evaluation: calculating relative stabilities and other criteria: general discussion. <i>Faraday Discussions</i> , 2018, 211, 325-381.	3.2	7
10	Applications of crystal structure prediction “ organic molecular structures: general discussion. <i>Faraday Discussions</i> , 2018, 211, 493-539.	3.2	8
11	Clathrate Structure Determination by Combining Crystal Structure Prediction with Computational and Experimental ¹²⁹ Xe NMR Spectroscopy. <i>Chemistry - A European Journal</i> , 2017, 23, 5258-5269.	3.3	18
12	Accurate force fields and methods for modelling organic molecular crystals at finite temperatures. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 15828-15837.	2.8	81
13	Report on the sixth blind test of organic crystal structure prediction methods. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2016, 72, 439-459.	1.1	445
14	Modelling temperature-dependent properties of polymorphic organic molecular crystals. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 31132-31143.	2.8	81
15	Static and lattice vibrational energy differences between polymorphs. <i>CrystEngComm</i> , 2015, 17, 5154-5165.	2.6	323