Stephan X M Boerrigter

List of Publications by Year
in descending order

12	688	933447 10	1199594 12
papers	citations	h-index	g-index

12	12	12	949
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Significant progress in predicting the crystal structures of small organic molecules â€“ a report on the fourth blind test. Acta Crystallographica Section B: Structural Science, 2009, 65, 107-125.	1.8	371
2	Cocrystals of nutraceuticalp-coumaric acid with caffeine and theophylline: polymorphism and solid-state stability explored in detail using their crystal graphs. CrystEngComm, 2011, 13, 611-619.	2.6	75
3	Investigation of the Milling-Induced Thermal Behavior of Crystalline and Amorphous Griseofulvin. Pharmaceutical Research, 2010, 27, 1377-1389.	3.5	56
4	The PO13 crystal structure of ROY. CrystEngComm, 2019, 21, 1363-1368.	2.6	40
5	Formation and Solid-State Characterization of a Salt-Induced Metastable Polymorph of Flufenamic Acid. Crystal Growth and Design, 2008, 8, 91-97.	3.0	39
6	New Metastable Packing Polymorph of Donepezil Grown on Stable Polymorph Substrates. Crystal Growth and Design, 2016, 16, 2552-2560.	3.0	29
7	Epitaxy of a Structurally Related Compound on the (100) Faces of Flufenamic Acid Form I and III Single Crystals. Crystal Growth and Design, 2010, 10, 518-527.	3.0	26
8	Organic vapor sorption method of isostructural solvates and polymorph of tenofovir disoproxil fumarate. European Journal of Pharmaceutical Sciences, 2013, 50, 253-262.	4.0	16
9	Investigating the effect of dehydration conditions on the compactability of glucose. International Journal of Pharmaceutics, 2011, 406, 55-61.	5.2	14
10	Anisotropic crystal deformation measurements determined using powder X-ray diffraction and a new in situ compression stage. International Journal of Pharmaceutics, 2011, 418, 199-206.	5.2	12
11	Application of errorâ€ranked singular value decomposition for the determination of potentialâ€derived atomicâ€centered point charges. Journal of Computational Chemistry, 2009, 30, 733-742.	3.3	6
12	Coreâ€“shell potentialâ€derived point charges. Journal of Computational Chemistry, 2012, 33, 950-957.	3.3	4