

# Yannick Jeanvoine

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/3079872/publications.pdf>

Version: 2024-02-01

8

papers

174

citations

1163117

8

h-index

1588992

8

g-index

8

all docs

8

docs citations

8

times ranked

269

citing authors

#	ARTICLE		IF	CITATIONS
1	Co <sup>2+</sup> -Binding Cysteine and Selenocysteine: A DFT Study. <i>Journal of Physical Chemistry A</i> , 2006, 110, 9727-9735.		2.5	41
2	SYNTHESIS OF FORMAMIDE AND RELATED ORGANIC SPECIES IN THE INTERSTELLAR MEDIUM VIA CHEMICAL DYNAMICS SIMULATIONS. <i>Astrophysical Journal</i> , 2016, 826, 107.		4.5	24
3	Hydration properties of Cm(iii) and Th(iv) combining coordination free energy profiles with electronic structure analysis. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 5824.		2.8	21
4	The formation of urea in space. II. MP2 versus PM6 dynamics in determining bimolecular reaction products. <i>Theoretical Chemistry Accounts</i> , 2019, 138, 1.		1.4	21
5	Mn <sup>2+</sup> , Fe <sup>2+</sup> , Co <sup>2+</sup> , Ni <sup>2+</sup> , Cu <sup>2+</sup> , and Zn <sup>2+</sup> -Binding Chalcogen- $\cdots$ -Chalcogen Bridges: A Compared MP2 and B3LYP Study. <i>Journal of Physical Chemistry A</i> , 2009, 113, 7878-7887.		2.5	20
6	Gas Phase Synthesis of Protonated Glycine by Chemical Dynamics Simulations. <i>Journal of Physical Chemistry A</i> , 2018, 122, 869-877.		2.5	19
7	The formation of urea in space. <i>Astronomy and Astrophysics</i> , 2018, 610, A26.		5.1	16
8	Cu <sup>2+</sup> binding chalcogen- $\cdots$ -chalcogen bridges: A problematic case for DFT. <i>Computational and Theoretical Chemistry</i> , 2010, 954, 7-15.		1.5	12