

James P Olivier

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

14
papers

9,011
citations

12
h-index

14
g-index

14
ext. papers

11,327
ext. citations

4.1
avg. IF

6.35
L-index

#	Paper	IF	Citations
14	Physisorption of gases, with special reference to the evaluation of surface area and pore size distribution (IUPAC Technical Report). <i>Pure and Applied Chemistry</i> , 2015 , 87, 1051-1069	2.1	7465
13	2D-NLDFT adsorption models for carbon slit-shaped pores with surface energetical heterogeneity and geometrical corrugation. <i>Carbon</i> , 2013 , 55, 70-80	10.4	346
12	Carbon slit pore model incorporating surface energetical heterogeneity and geometrical corrugation. <i>Adsorption</i> , 2013 , 19, 777-783	2.6	223
11	Using a New Finite Slit Pore Model for NLDFT Analysis of Carbon Pore Structure. <i>Adsorption Science and Technology</i> , 2011 , 29, 769-780	3.6	22
10	A Simple Two-Dimensional NLDFT Model of Gas Adsorption in Finite Carbon Pores. Application to Pore Structure Analysis. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 19382-19385	3.8	140
9	The Surface Heterogeneity of Carbon and Its Assessment 2008 , 147-166		5
8	An overview of physical adsorption methods for the characterization of finely divided and porous materials and their application to fluid cracking catalysts. <i>Studies in Surface Science and Catalysis</i> , 2004 , 1-33	1.8	2
7	Surface area and microporosity of pillared rectorite catalysts from a hybrid density functional theory method. <i>Microporous and Mesoporous Materials</i> , 2003 , 57, 291-296	5.3	36
6	Determination of Pore Size Distribution, Surface Area, and Acidity in Fluid Cracking Catalysts (FCCs) from Nonlocal Density Functional Theoretical Models of Adsorption and from Microcalorimetry Methods. <i>Journal of Physical Chemistry B</i> , 2003 , 107, 4128-4136	3.4	40
5	Surface Area and Microporosity of a Pillared Interlayered Clay (PILC) from a Hybrid Density Functional Theory (DFT) Method. <i>Journal of Physical Chemistry B</i> , 2001 , 105, 623-629	3.4	52
4	A new method for the accurate pore size analysis of MCM-41 and other silica based mesoporous materials. <i>Studies in Surface Science and Catalysis</i> , 2000 , 71-80	1.8	43
3	Improving the models used for calculating the size distribution of micropore volume of activated carbons from adsorption data. <i>Carbon</i> , 1998 , 36, 1469-1472	10.4	255
2	The Determination of Surface Energetic Heterogeneity Using Model Isotherms Calculated by Density Functional Theory. <i>Kluwer International Series in Engineering and Computer Science</i> , 1996 , 699-707		13
1	Modeling physical adsorption on porous and nonporous solids using density functional theory. <i>Journal of Porous Materials</i> , 1995 , 2, 9-17	2.4	369