

Jurgen Schnitker

List of Publications by Year in descending order

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16
papers

1,624
citations

567281

15
h-index

940533

16
g-index

16
all docs

16
docs citations

16
times ranked

798
citing authors

#	ARTICLE	IF	CITATIONS
1	Response to "Comment on 'An electron-water pseudopotential for condensed phase simulation'" [J. Chem. Phys. 131 , 037101 (2009)]. Journal of Chemical Physics, 2009, 131, .	3.0	15
2	Stress relaxation and misfit dislocation nucleation in the growth of misfitting films: A molecular dynamics simulation study. Journal of Applied Physics, 1998, 83, 217-227.	2.5	116
3	The reaction field method in molecular dynamics simulations of point-polarizable water models. Molecular Physics, 1996, 88, 1089-1108.	1.7	11
4	How the unit cell surface charge distribution affects the energetics of ion-solvent interactions in simulations. Journal of Chemical Physics, 1994, 101, 5024-5031.	3.0	65
5	Melting of two-dimensional colloidal crystals: A simulation study of the Yukawa system. Journal of Chemical Physics, 1994, 100, 3114-3121.	3.0	50
6	Two-dimensional melting revisited. Molecular Physics, 1993, 80, 1-24.	1.7	27
7	A comparison of classical and quantum analyses of electron localization sites in liquid water. Journal of Chemical Physics, 1992, 97, 2055-2060.	3.0	21
8	Model dependence of quantum isotope effects in liquid water. Journal of Chemical Physics, 1991, 95, 3728-3737.	3.0	92
9	Solvation dynamics of the hydrated electron: A nonadiabatic quantum simulation. Physical Review Letters, 1991, 66, 3172-3175.	7.8	235
10	Transient photophysical hole-burning spectroscopy of the hydrated electron: A quantum dynamical simulation. Journal of Chemical Physics, 1989, 90, 6916-6924.	3.0	30
11	A prioricalculation of the optical absorption spectrum of the hydrated electron. Physical Review Letters, 1988, 60, 456-459.	7.8	173
12	The hydrated electron: quantum simulation of structure, spectroscopy, and dynamics. The Journal of Physical Chemistry, 1988, 92, 4277-4285.	2.9	305
13	An electron-water pseudopotential for condensed phase simulation. Journal of Chemical Physics, 1987, 86, 3462-3470.	3.0	186
14	Quantum simulation study of the hydrated electron. Journal of Chemical Physics, 1987, 86, 3471-3485.	3.0	191
15	Quantum simulations of aqueous systems. Journal of Statistical Physics, 1986, 43, 949-965.	1.2	17
16	Electron localization in liquid water: A computer simulation study of microscopic trapping sites. Journal of Chemical Physics, 1986, 85, 2986-2998.	3.0	90