

Ersan Demiralp

List of Publications by Year in descending order

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papers

745
citations

516710

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23
all docs

23
docs citations

23
times ranked

675
citing authors

#	ARTICLE	IF	CITATIONS
1	Exact solutions for a Hamiltonian with the Morse potential and the Dirac delta shell interactions. <i>Molecular Physics</i> , 2009, 107, 2053-2062.	1.7	6
2	Bose-Einstein condensate in a harmonic trap decorated with Dirac δ functions. <i>Physical Review A</i> , 2007, 76, .	2.5	31
3	The Woods-Saxon potential with point interactions. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2007, 365, 55-63.	2.1	20
4	Bound state solutions of the Schrödinger equation for a δ -symmetric potential with Dirac delta functions. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2006, 359, 190-198.	2.1	14
5	Solutions of the Schrödinger equation for Dirac delta decorated linear potential. <i>Open Physics</i> , 2005, 3, .	1.7	6
6	Properties of a pseudo-Hermitian Hamiltonian for harmonic oscillator decorated with Dirac delta interactions. <i>European Physical Journal D</i> , 2005, 55, 1081-1084.	0.4	13
7	Bound states of n-dimensional harmonic oscillator decorated with Dirac delta functions. <i>Journal of Physics A</i> , 2005, 38, 4783-4793.	1.6	27
8	Structural and electronic properties of novel nanoscale $C_{4n}S_{4n}$ molecules. <i>Computational and Theoretical Chemistry</i> , 2003, 635, 125-131.	1.5	0
9	Properties of bound states of the Schrödinger equation with attractive Dirac delta potentials. <i>Journal of Physics A</i> , 2003, 36, 7449-7459.	1.6	24
10	Dynamic Charge Equilibration-Morse stretch force field: Application to energetics of pure silica zeolites. <i>Journal of Computational Chemistry</i> , 2002, 23, 1507-1514.	3.3	19
11	The MS-Q Force Field for Clay Minerals: Application to Oil Production. <i>Journal of Physical Chemistry B</i> , 2001, 105, 4122-4127.	2.6	23
12	Morse Stretch Potential Charge Equilibrium Force Field for Ceramics: Application to the Quartz-Stishovite Phase Transition and to Silica Glass. <i>Physical Review Letters</i> , 1999, 82, 1708-1711.	7.8	173
13	Computational Materials Chemistry at the Nanoscale. <i>Journal of Nanoparticle Research</i> , 1999, 1, 51-69.	1.9	23
14	Theoretical studies on VPI-5. 3.. <i>Computational Materials Science</i> , 1999, 14, 135-137.	3.0	11
15	Factors affecting molecular dynamics simulated vitreous silica structures. <i>Journal of Non-Crystalline Solids</i> , 1999, 253, 133-142.	3.1	121
16	Vibrational Analysis and Isotope Shifts of BEDT-TTF Donor for Organic Superconductors. <i>Journal of Physical Chemistry A</i> , 1998, 102, 2466-2471.	2.5	19
17	Conduction properties of the organic superconductor $(BEDT-TTF)_2Cu(NCS)_2$ based on Hubbard unrestricted-Hartree-Fock band calculations. <i>Physical Review B</i> , 1997, 56, 11907-11919.	3.2	25
18	Pressure Induced Phase Transformations in Silica. <i>Materials Research Society Symposia Proceedings</i> , 1997, 492, 287.	0.1	4

#	ARTICLE	IF	CITATIONS
19	Structures and Energetics Study of Tetrathiafulvalene-Based Donors of Organic Superconductors. Journal of Physical Chemistry A, 1997, 101, 8128-8131.	2.5	39
20	MSX Force Field and Vibrational Frequencies for BEDT-TTF (Neutral and Cation). Journal of Physical Chemistry A, 1997, 101, 1975-1981.	2.5	29
21	Electron-transfer boat-vibration mechanism for superconductivity in organic molecules based on BEDT-TTF. Journal of the American Chemical Society, 1995, 117, 8154-8158.	13.7	66
22	Prediction of new donors for organic superconductors. Synthetic Metals, 1995, 72, 297-299.	3.9	19
23	Ab Initio and Semiempirical Electronic Structural Studies on Bis(ethylenedithio)tetrathiafulvalene (BEDT-TTF or ET). The Journal of Physical Chemistry, 1994, 98, 9781-9785.	2.9	33