Masoud Aryanpour

List of Publications by Year in descending order

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840776 996975 16 699 11 15 citations h-index g-index papers 16 16 16 1242 docs citations times ranked citing authors all docs

| # | Article | IF | CITATIONS |
|----|---|-------------|-----------|
| 1 | An Analytical 2D Formulation for the Combined Cooling of PCM-Covered Cylindrical Battery Cells. Journal of Energy Engineering - ASCE, 2022, 148, . | 1.9 | 1 |
| 2 | Coupled Ionic-Electronic Equivalent Circuit to Describe Asymmetric Rise and Decay of Photovoltage Profile in Perovskite Solar Cells. Scientific Reports, 2019, 9, 11962. | 3. 3 | 31 |
| 3 | Computational investigation of gas detection and selectivity on TiS3 nanoflakes supported by experimental evidence. Physical Chemistry Chemical Physics, 2018, 20, 25458-25466. | 2.8 | 3 |
| 4 | Staging and In-Plane Superstructures Formed in Layered NaMO $<$ sub $>$ 2 $<$ /sub $>$ {M = Sc, Ti, V, Cr, Mn} during Na De-Intercalation: A Computational Study. Journal of the Electrochemical Society, 2015, 162, A511-A519. | 2.9 | 13 |
| 5 | Computational Study of Site-Specific Correlations among Oxygen Reduction Intermediates on Pd3Y (111). Journal of Advanced Catalysis Science and Technology, 2014, 1, 1-9. | 1.0 | 1 |
| 6 | Activity Descriptor for Catalytic Reactions on Doped Cerium Oxide. ACS Catalysis, 2013, 3, 1253-1262. | 11.2 | 39 |
| 7 | ATR–FTIR and Density Functional Theory Study of the Structures, Energetics, and Vibrational Spectra of Phosphate Adsorbed onto Goethite. Langmuir, 2012, 28, 14573-14587. | 3.5 | 142 |
| 8 | Quantum mechanical calculations on FeOH nanoparticles. Geoderma, 2012, 189-190, 236-242. | 5.1 | 10 |
| 9 | Differential Pair Distribution Function Study of the Structure of Arsenate Adsorbed on Nanocrystalline Î ³ -Alumina. Environmental Science & Technology, 2011, 45, 9687-9692. | 10.0 | 66 |
| 10 | Aluminum coprecipitates with Fe (hydr)oxides: Does isomorphous substitution of Al3+ for Fe3+ in goethite occur?. Geochimica Et Cosmochimica Acta, 2011, 75, 4667-4683. | 3.9 | 54 |
| 11 | Development of a Reactive Force Field for Ironâ^'Oxyhydroxide Systems. Journal of Physical Chemistry A, 2010, 114, 6298-6307. | 2.5 | 199 |
| 12 | Tungsten-Doped Titanium Dioxide in the Rutile Structure: Theoretical Considerations. Chemistry of Materials, 2009, 21, 1627-1635. | 6.7 | 32 |
| 13 | Mechanism of Molecular Oxygen Reduction at the Cathode of a PEM Fuel Cell: Non-Electrochemical Reactions on Catalytic Pt Particles. Journal of Physical Chemistry C, 2008, 112, 8464-8475. | 3.1 | 48 |
| 14 | First-Principles Analysis of Oxygen-Containing Adsorbates Formed from the Electrochemical Discharge of Water on Pt(111). Journal of Physical Chemistry C, 2008, 112, 9760-9768. | 3.1 | 45 |
| 15 | An algorithm for mass matrix calculation of internally constrained molecular geometries. Journal of Chemical Physics, 2008, 128, 044113. | 3.0 | O |
| 16 | Convergent Iterative Constrained Variation Algorithm for Calculation of Electron-Transfer Transition States. Journal of the Electrochemical Society, 2006, 153, E52. | 2.9 | 15 |