

Javad Beheshtian

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8048988/publications.pdf>

Version: 2024-02-01

102
papers

5,131
citations

70961

41
h-index

88477

70
g-index

103
all docs

103
docs citations

103
times ranked

3294
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 1 | Detection of phosgene by Sc-doped BN nanotubes: A DFT study. <i>Sensors and Actuators B: Chemical</i> , 2012, 171-172, 846-852. | 4.0 | 292 |
| 2 | Sensing behavior of Al and Si doped BC ₃ graphenes to formaldehyde. <i>Sensors and Actuators B: Chemical</i> , 2013, 181, 829-834. | 4.0 | 188 |
| 3 | A comparative study on the B ₁₂ N ₁₂ , Al ₁₂ N ₁₂ , B ₁₂ P ₁₂ and Al ₁₂ P ₁₂ fullerene-like cages. <i>Journal of Molecular Modeling</i> , 2012, 18, 2653-2658. | 0.8 | 160 |
| 4 | Microporous titania-silica nanocomposite catalyst-adsorbent for ultra-deep oxidative desulfurization. <i>Applied Catalysis B: Environmental</i> , 2016, 180, 65-77. | 10.8 | 153 |
| 5 | Theoretical study of CO adsorption on the surface of BN, AlN, BP and AlP nanotubes. <i>Surface Science</i> , 2012, 606, 981-985. | 0.8 | 152 |
| 6 | Van der Waals pressure and its effect on trapped interlayer molecules. <i>Nature Communications</i> , 2016, 7, 12168. | 5.8 | 137 |
| 7 | Selective function of Al ₁₂ N ₁₂ nano-cage towards NO and CO molecules. <i>Computational Materials Science</i> , 2012, 62, 71-74. | 1.4 | 136 |
| 8 | A DFT study on the functionalization of a BN nanosheet with PCX, (PC=phenyl carbamate, X=OCH ₃ , CH ₃), <i>Tj ETQq 0 0 rgBT /Overlock</i> | 3.1 | 136 |
| 9 | B ₁₂ N ₁₂ Nano-cage as Potential Sensor for NO ₂ Detection. <i>Chinese Journal of Chemical Physics</i> , 2012, 25, 60-64. | 0.6 | 126 |
| 10 | Toxic CO detection by B ₁₂ N ₁₂ nanocluster. <i>Microelectronics Journal</i> , 2011, 42, 1400-1403. | 1.1 | 124 |
| 11 | Adsorption and dissociation of Cl ₂ molecule on ZnO nanocluster. <i>Applied Surface Science</i> , 2012, 258, 8171-8176. | 3.1 | 117 |
| 12 | Highly active Fe ₂ O ₃ -doped TiO ₂ photocatalyst for degradation of trichloroethylene in air under UV and visible light irradiation: Experimental and computational studies. <i>Applied Catalysis B: Environmental</i> , 2015, 165, 209-221. | 10.8 | 117 |
| 13 | The H ₂ dissociation on the BN, AlN, BP and AlP nanotubes: a comparative study. <i>Journal of Molecular Modeling</i> , 2012, 18, 2343-2348. | 0.8 | 111 |
| 14 | Interaction of small molecules (NO, H ₂ , N ₂ , and CH ₄) with BN nanocluster surface. <i>Structural Chemistry</i> , 2012, 23, 1567-1572. | 1.0 | 103 |
| 15 | Computational study of CO and NO adsorption on magnesium oxide nanotubes. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2011, 44, 546-549. | 1.3 | 100 |
| 16 | Quantum chemical study of fluorinated AlN nano-cage. <i>Applied Surface Science</i> , 2012, 259, 631-636. | 3.1 | 97 |
| 17 | AlN nanotube as a potential electronic sensor for nitrogen dioxide. <i>Microelectronics Journal</i> , 2012, 43, 452-455. | 1.1 | 96 |
| 18 | Theoretical study of hydrogen adsorption on the B ₁₂ P ₁₂ fullerene-like nanocluster. <i>Computational Materials Science</i> , 2012, 54, 115-118. | 1.4 | 95 |

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 19 | A first-principles study of H ₂ S adsorption and dissociation on the AlN nanotube. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2012, 44, 1963-1968. | 1.3 | 92 |
| 20 | Chemisorption of NH ₃ at the open ends of boron nitride nanotubes: a DFT study. <i>Structural Chemistry</i> , 2011, 22, 183-188. | 1.0 | 88 |
| 21 | DFT study on the functionalization of a BN nanotube with sulfamide. <i>Applied Surface Science</i> , 2013, 266, 182-187. | 3.1 | 87 |
| 22 | Functionalization of BN nanosheet with N ₂ H ₄ may be feasible in the presence of Stone-Wales defect. <i>Structural Chemistry</i> , 2013, 24, 1565-1570. | 1.0 | 86 |
| 23 | Interaction of NH ₃ with aluminum nitride nanotube: Electrostatic vs. covalent. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2011, 43, 1717-1719. | 1.3 | 84 |
| 24 | Experimental and Theoretical Study of Enhanced Photocatalytic Activity of Mg-Doped ZnO NPs and ZnO/rGO Nanocomposites. <i>Chemistry - an Asian Journal</i> , 2018, 13, 194-203. | 1.7 | 83 |
| 25 | Carbon nitride nanotube as a sensor for alkali and alkaline earth cations. <i>Applied Surface Science</i> , 2013, 264, 699-706. | 3.1 | 82 |
| 26 | Electronic sensor for sulfide dioxide based on AlN nanotubes: a computational study. <i>Journal of Molecular Modeling</i> , 2012, 18, 4745-4750. | 0.8 | 80 |
| 27 | Ab initio study of NH ₃ and H ₂ O adsorption on pristine and Na-doped MgO nanotubes. <i>Structural Chemistry</i> , 2013, 24, 165-170. | 1.0 | 80 |
| 28 | Density-functional calculations of HCN adsorption on the pristine and Si-doped graphynes. <i>Structural Chemistry</i> , 2014, 25, 1-7. | 1.0 | 79 |
| 29 | A theoretical study of CO adsorption on aluminum nitride nanotubes. <i>Structural Chemistry</i> , 2012, 23, 653-657. | 1.0 | 77 |
| 30 | Benchmarking of ONIOM method for the study of NH ₃ dissociation at open ends of BNNTs. <i>Journal of Molecular Modeling</i> , 2012, 18, 1729-1734. | 0.8 | 75 |
| 31 | Ammonia monitoring by carbon nitride nanotubes: A density functional study. <i>Thin Solid Films</i> , 2013, 534, 650-654. | 0.8 | 75 |
| 32 | Theoretical investigation of C ₆₀ fullerene functionalization with tetrazine. <i>Computational and Theoretical Chemistry</i> , 2012, 992, 164-167. | 1.1 | 73 |
| 33 | Highly sensitive and selective ethanol and acetone gas sensors by adding some dopants (Mn, Fe, Co, Ni) onto hexagonal ZnO plates. <i>RSC Advances</i> , 2016, 6, 7838-7845. | 1.7 | 73 |
| 34 | The effect of surface curvature of aluminum nitride nanotubes on the adsorption of NH ₃ . <i>Structural Chemistry</i> , 2011, 22, 1261-1265. | 1.0 | 72 |
| 35 | Hydrogen dissociation on diene-functionalized carbon nanotubes. <i>Journal of Molecular Modeling</i> , 2013, 19, 255-261. | 0.8 | 72 |
| 36 | Nitrate adsorption by carbon nanotubes in the vacuum and aqueous phase. <i>Monatshefte für Chemie</i> , 2012, 143, 1623-1626. | 0.9 | 68 |

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 37 | Sensing behavior of Al-rich AlN nanotube toward hydrogen cyanide. <i>Journal of Molecular Modeling</i> , 2013, 19, 2197-2203. | 0.8 | 63 |
| 38 | Carbon nanotube functionalization with carboxylic derivatives: a DFT study. <i>Journal of Molecular Modeling</i> , 2013, 19, 391-396. | 0.8 | 63 |
| 39 | Theoretical prediction of silicene as a new candidate for the anode of lithium-ion batteries. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 29689-29696. | 1.3 | 63 |
| 40 | Functionalization of [60] fullerene with butadienes: A DFT study. <i>Applied Surface Science</i> , 2012, 258, 8980-8984. | 3.1 | 59 |
| 41 | Boron Nitride Monolayer: A Strain-Tunable Nanosensor. <i>Journal of Physical Chemistry C</i> , 2013, 117, 13261-13267. | 1.5 | 45 |
| 42 | Hydrogen storage by BeO nano-cage: A DFT study. <i>Applied Surface Science</i> , 2016, 368, 76-81. | 3.1 | 44 |
| 43 | Induced polarization and electronic properties of carbon-doped boron nitride nanoribbons. <i>Physical Review B</i> , 2012, 86, . | 1.1 | 43 |
| 44 | Computational study on the characteristics of the interaction in linear urea clusters. <i>International Journal of Quantum Chemistry</i> , 2011, 111, 3184-3195. | 1.0 | 39 |
| 45 | DFT study of NH ₃ adsorption on the (5,0), (8,0), (5,5) and (6,6) single-walled carbon nanotubes. Calculated binding energies, NMR and NQR parameters. <i>Physica B: Condensed Matter</i> , 2010, 405, 1455-1460. | 1.3 | 37 |
| 46 | Formaldehyde adsorption on the interior and exterior surfaces of CN nanotubes. <i>Structural Chemistry</i> , 2013, 24, 1331-1337. | 1.0 | 36 |
| 47 | Nitrous oxide adsorption on pristine and Si-doped AlN nanotubes. <i>Journal of Molecular Modeling</i> , 2013, 19, 943-949. | 0.8 | 36 |
| 48 | Natural pigments in dye-sensitized solar cell (DSSC): a DFT-TDDFT study. <i>Journal of the Iranian Chemical Society</i> , 2019, 16, 795-805. | 1.2 | 36 |
| 49 | Electronic and optical properties of vacancy and B, N, O and F doped graphene: DFT study. <i>Opto-electronics Review</i> , 2019, 27, 130-136. | 2.4 | 35 |
| 50 | A computational study of water adsorption on boron nitride nanotube. <i>Structural Chemistry</i> , 2010, 21, 903-908. | 1.0 | 33 |
| 51 | Exohedral and endohedral adsorption of alkaline earth cations in BN nanocluster. <i>Journal of Molecular Modeling</i> , 2013, 19, 1445-1450. | 0.8 | 33 |
| 52 | Theoretical ¹⁴ N nuclear quadrupole resonance parameters for sulfa drugs: Sulfamerazine and sulfathiazole. <i>Journal of Molecular Graphics and Modelling</i> , 2008, 27, 326-331. | 1.3 | 31 |
| 53 | Arsenic interactions with a fullerene-like BN cage in the vacuum and aqueous phase. <i>Journal of Molecular Modeling</i> , 2013, 19, 833-837. | 0.8 | 31 |
| 54 | Phase transition and mechanical properties of cesium bismuth silver halide double perovskites (Cs ₂ AgBiX ₆ , X = Cl, Br, I): a DFT approach. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 5959-5968. | 1.3 | 30 |

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 55 | Reversible structural transition in nanoconfined ice. <i>Physical Review B</i> , 2017, 95, . | 1.1 | 28 |
| 56 | Preparation of uniform 2D ZnO nanostructures by the ionic liquid-assisted sonochemical method and their optical properties. <i>Ceramics International</i> , 2014, 40, 7769-7774. | 2.3 | 26 |
| 57 | Electrochemical and DFT study of an anticancer and active anthelmintic drug at carbon nanostructured modified electrode. <i>Materials Science and Engineering C</i> , 2016, 69, 1345-1353. | 3.8 | 26 |
| 58 | Preparation of Cu ₂ O nanostructures by changing reducing agent and their optical properties. <i>Materials Letters</i> , 2015, 153, 1-4. | 1.3 | 24 |
| 59 | Effects of functionalization and side defects on single-photon emission in boron nitride quantum dots. <i>Physical Review B</i> , 2017, 96, . | 1.1 | 23 |
| 60 | First-Principles Study of Water Nanotubes Captured Inside Carbon/Boron Nitride Nanotubes. <i>Langmuir</i> , 2018, 34, 11176-11187. | 1.6 | 23 |
| 61 | Graphene-silicene bilayer: A nanocapacitor with permanent dipole and piezoelectricity effect. <i>Physical Review B</i> , 2015, 92, . | 1.1 | 20 |
| 62 | Ultrasonic route synthesis, characterization and electrochemical study of graphene oxide and reduced graphene oxide. <i>Research on Chemical Intermediates</i> , 2019, 45, 487-505. | 1.3 | 20 |
| 63 | Co-adsorption of CO molecules at the open ends of MgO nanotubes. <i>Structural Chemistry</i> , 2012, 23, 1981-1986. | 1.0 | 19 |
| 64 | Electro-Optical Properties of Monolayer and Bilayer Pentagonal BN: First Principles Study. <i>Nanomaterials</i> , 2020, 10, 440. | 1.9 | 19 |
| 65 | Doping effect on the adsorption of NH ₃ molecule onto graphene quantum dot: From the physisorption to the chemisorption. <i>Journal of Applied Physics</i> , 2013, 114, . | 1.1 | 18 |
| 66 | Spiral graphone and one-sided fluorographene nanoribbons. <i>Physical Review B</i> , 2013, 87, . | 1.1 | 17 |
| 67 | Selective separation behavior of graphene flakes in interaction with halide anions in the presence of an external electric field. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 7293-7299. | 1.3 | 16 |
| 68 | Experimental and theoretical study of CO adsorption on the surface of single phase hexagonally plate ZnO. <i>Applied Surface Science</i> , 2014, 315, 8-15. | 3.1 | 14 |
| 69 | The Alkali Metal Interactions with MgO Nanotubes. <i>Bulletin of the Korean Chemical Society</i> , 2012, 33, 1925-1928. | 1.0 | 14 |
| 70 | DFT study of NH ₃ (H ₂ O) _{n=0,1,2,3} complex adsorption on the (8,0) single-walled carbon nanotube. <i>Computational Materials Science</i> , 2010, 48, 655-657. | 1.4 | 12 |
| 71 | Computational study of ammonia adsorption on the perfect and rippled graphene sheet. <i>Physica B: Condensed Matter</i> , 2013, 429, 52-56. | 1.3 | 12 |
| 72 | Hydrogenated $\hat{\Gamma}$ -graphene as an ultraviolet optomechanical sensor. <i>RSC Advances</i> , 2020, 10, 26197-26211. | 1.7 | 12 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 73 | Adsorption of Na, Mg, and Al atoms on BN nanotubes. <i>Thin Solid Films</i> , 2012, 526, 139-142. | 0.8 | 11 |
| 74 | Effect of nitrogen doping on electronic and optical properties of ZnO sheet: DFT+U study. <i>Computational Condensed Matter</i> , 2018, 15, 1-6. | 0.9 | 11 |
| 75 | Theoretical study on the functionalization of BC ₂ N nanotube with amino groups. <i>Journal of Molecular Modeling</i> , 2013, 19, 2211-2216. | 0.8 | 10 |
| 76 | Application of hexa- <i>peri</i> -hexabenzocoronene nanographene and its B, N, and Bn doped forms in Na-ion batteries: A density functional theory study. <i>Thin Solid Films</i> , 2020, 704, 137979. | 0.8 | 10 |
| 77 | Synthesis of undoped and Fe nanoparticles doped SnO ₂ nanostructure: study of structural, optical and electrocatalytic properties. <i>Journal of Materials Science: Materials in Electronics</i> , 2017, 28, 7568-7574. | 1.1 | 9 |
| 78 | The influence of Stone-Wales defects in nanographene on the performance of Na-ion batteries. <i>Journal of Molecular Graphics and Modelling</i> , 2020, 98, 107578. | 1.3 | 9 |
| 79 | A computational study on the novel defects of graphene quantum dot as a promising anode material for sodium ion battery. <i>Materials Chemistry and Physics</i> , 2021, 265, 124484. | 2.0 | 9 |
| 80 | Photo-oxidation of phenylazonaphthol dyes and their reactivity analysis in the gas phase and adsorbed on cellulose fibers states using DFT and TD-DFT. <i>Dyes and Pigments</i> , 2011, 89, 16-22. | 2.0 | 8 |
| 81 | Voltammetric Sensor Based on Fe-doped ZnO and TiO ₂ Nanostructures-modified Carbon-paste Electrode for Determination of Levodopa. <i>Journal of Electronic Materials</i> , 2017, 46, 5657-5663. | 1.0 | 8 |
| 82 | Toxic CO detection by Li-encapsulated fullerene-like BeO. <i>Structural Chemistry</i> , 2018, 29, 231-241. | 1.0 | 8 |
| 83 | A density functional study of ¹⁵ N chemical shielding tensors in quinolines. <i>Chemical Physics Letters</i> , 2009, 476, 196-200. | 1.2 | 7 |
| 84 | Electronic, magnetic and optical properties of Fe-doped nano-BN sheet: DFT study. <i>Indian Journal of Physics</i> , 2021, 95, 823-831. | 0.9 | 7 |
| 85 | A DFT study on the potential application of pristine, B and N doped carbon nanocones in potassium-ion batteries. <i>Journal of Molecular Modeling</i> , 2021, 27, 168. | 0.8 | 6 |
| 86 | Fundamental mechanisms of hexagonal boron nitride sensing of dopamine, tryptophan, ascorbic acid, and uric acid by first-principles study. <i>Journal of Molecular Modeling</i> , 2022, 28, . | 0.8 | 6 |
| 87 | The electronic and optical properties of 3d transition metals doped silicene sheet: A DFT study. <i>Materials Research Express</i> , 2019, 6, 126326. | 0.8 | 5 |
| 88 | Nanoscale investigation of the influence of water on the elastic properties of Ca ²⁺ -H gel by molecular simulation. <i>Proceedings of the Institution of Mechanical Engineers, Part L: Journal of Materials: Design and Applications</i> , 2019, 233, 1295-1306. | 0.7 | 5 |
| 89 | A computational study on the BN and AlN nanocones as anode materials for K-ion batteries. <i>Applied Surface Science</i> , 2021, 544, 148793. | 3.1 | 5 |
| 90 | Electrodeposition of Co _x Ni _y V _z O _z Ternary Nanopetals on Bare and rGO-Coated Nickel Foam for High-Performance Supercapacitor Application. <i>Nanomaterials</i> , 2022, 12, 1894. | 1.9 | 5 |

| # | ARTICLE | IF | CITATIONS |
|-----|---|-----|-----------|
| 91 | Helium selectivity of H-, B-, N-, and F- doped nanoporous graphene membranes in the presence of natural gas: A density functional theory study. <i>Superlattices and Microstructures</i> , 2020, 141, 106478. | 1.4 | 4 |
| 92 | 15N CHEMICAL SHIFT CALCULATIONS AND NATURAL BONDING ORBITAL ANALYSES OF (BENZAMIDE) _n = 1 - 6 CLUSTERS. <i>Journal of Theoretical and Computational Chemistry</i> , 2009, 08, 973-982. | 1.8 | 3 |
| 93 | Synthesis, identification, crystal structure and theoretical study of a Ce(IV) complex. <i>Journal of the Iranian Chemical Society</i> , 2014, 11, 1353-1361. | 1.2 | 3 |
| 94 | Boron nitride nanochannels encapsulating a water/heavy water layer for energy applications. <i>RSC Advances</i> , 2019, 9, 5901-5907. | 1.7 | 3 |
| 95 | Theoretical investigation of azo dyes adsorbed on cellulose fibers: 1. Electronic and bonding structures. <i>Journal of the Iranian Chemical Society</i> , 2013, 10, 985-999. | 1.2 | 2 |
| 96 | Ab Initio Study of Mono-Layer Graphene as an Electronical or Optical Sensor for Detecting B, N, O and F Atoms. <i>Journal of Electronic Materials</i> , 2019, 48, 4265-4272. | 1.0 | 2 |
| 97 | Interfacial properties of water/heavy water layer encapsulate in bilayer graphene nanochannel and nanocapacitor. <i>Journal of Materials Science: Materials in Electronics</i> , 2019, 30, 11964-11975. | 1.1 | 2 |
| 98 | Theoretical investigation of azo dyes adsorbed on cellulose fibers: 2. Spectroscopic study. <i>Journal of the Iranian Chemical Society</i> , 2014, 11, 111-121. | 1.2 | 1 |
| 99 | Experimental and Theoretical Study of Porous Al ₂ O ₃ . <i>Transactions of the Indian Institute of Metals</i> , 2021, 74, 381-386. | 0.7 | 1 |
| 100 | Investigation of Interaction Between Graphene and Its Compounds as Carriers on Anti-Cancer Drug of 5-Fluorouracil. <i>Eurasian Journal of Analytical Chemistry</i> , 2018, 13, . | 0.4 | 1 |
| 101 | Effect of vacancy modification on the quantum capacitance of silicene- based electrode in efficient supercapacitors. <i>Thin Solid Films</i> , 2022, 756, 139378. | 0.8 | 1 |
| 102 | A Theoretical Model Based on Modified Fullerenes for Photodynamic Therapy of Cancer. <i>Journal of Computational and Theoretical Nanoscience</i> , 2018, 15, 147-152. | 0.4 | 0 |