## google-scholar-api

September 7, 2022

## 1 Google scholar

```
Source: https://pypi.org/project/scholarly/
conda install -c conda-forge scholarly
```

## 1.0.1 Example-1

```
[]: from scholarly import scholarly
     import json
     def print_pretty(input_dict):
         print("info=",json.dumps(input_dict, indent=4, sort_keys=True))
     # Retrieve the author's data, fill-in, and print
     # Get an iterator for the author results
     search_query = scholarly.search_author('James Hickman NIST')
     # Retrieve the first result from the iterator
     first_author_result = next(search_query)
     # Print first results
     print_pretty(first_author_result)
    info= {
        "affiliation": "NIST",
        "citedby": 178,
        "container_type": "Author",
        "email_domain": "@gmu.edu",
        "filled": [],
        "interests": [
            "Thermodynamics",
            "Material science",
            "Statistical mechanics",
            "Computational physics",
            "Solid state physics"
        "name": "James Hickman",
        "scholar_id": "5Gjl4o8AAAJ",
```

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"source": "SEARCH_AUTHOR_SNIPPETS",
        "url_picture":
    "https://scholar.google.com/citations?view_op=medium_photo&user=5Gj14o8AAAAJ"
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Ozoli\u0146\u0161 Department of Applied Physics Energy Sciences Institute"
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    can exhibit much higher strength than macroscopic objects, but this strength
    rarely approaches the maximum theoretical strength of the material. Here, we
    demonstrate that faceted single-crystalline nickel (Ni) nanoparticles exhibit an
    ultrahigh compressive strength (up to 34\u2009GPa) unprecedented for metallic
    materials. This strength matches the available estimates of Ni theoretical
    strength. Three factors are responsible for this record-high strength: the large
    Ni shear modulus, the smooth edges and corners of the nanoparticles, and the
    thin oxide layer on the particle surface. This finding is supported by molecular
    dynamics simulations that closely mimic the experimental conditions, which show
    that the mechanical failure of the strongest particles is triggered by
    homogeneous nucleation of dislocation loops inside the particle. The nucleation
    of a stable loop is \u2026",
            "author": "A Sharma and J Hickman and N Gazit and E Rabkin and Y
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            "citation": "Nature communications 9 (1), 1-9, 2018",
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    systems and validate a well-known but frequently contested equation predicting
    the mean square of such fluctuations. The simulations implement two virtual and
    one physical (natural) thermostat and examine the kinetic, potential, and total
    energy correlation functions in the time and frequency domains. The results
    clearly demonstrate the existence of quasiequilibrium states in which the system
    can be characterized by a well-defined temperature that follows the mentioned
    fluctuation equation. The emergence of such states is due to the wide separation
    of time scales between thermal relaxation by phonon scattering and slow energy
    exchanges with the thermostat. The quasiequilibrium states exist between these
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Nickel nanoparticles set a new record of strength

Temperature fluctuations in canonical systems: Insights from molecular dynamics simulations

Extra variable in grain boundary description

Disjoining potential and grain boundary premelting in binary alloys Development of a general-purpose machine-learning interatomic potential for aluminum by the physically informed neural network method

Thermal conductivity and its relation to atomic structure for symmetrical tilt grain boundaries in silicon

Energy spectrum of a Langevin oscillator

The impact of alloying on defect-free nanoparticles exhibiting softer but tougher behavior  $\ensuremath{\mathsf{S}}$ 

Coarsening of solid -Sn particles in liquid Pb-Sn alloys: Reinterpretation of experimental data in the framework of trans-interface-diffusion-controlled coarsening

Coarsening of Solid Particles in Liquid Pb-Sn Alloys: Reinterpretation of Data in Light of the TIDC Theory of Coarsening

Softer but tougher: The impact of alloying on defect-free nanoparticles Physically-Informed Artificial Neural Networks for Atomistic Modeling of

Materials

Investigations of Interface Phenomena via Atomistic Simulation

Atomistic modeling of pre-melted grain boundaries

Fairfax, Virginia 22030-4444 USA Vidvuds Ozoliņš Department of Applied Physics Energy Sciences Institute

Atomistic modeling of grain boundary melting and premelting in alloys

[]: # Which papers cited that publication?
citations = [print(citation['bib']['title']) for citation in scholarly.

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Temperature in and out of equilibrium: A review of concepts, tools and attempts Novel molecular-dynamics-based protocols for phase space sampling in complex systems

Fundamental thermal noise limits for optical microcavities

Assessment of mechanical, thermal properties and crystal shapes of monoclinic tricalcium silicate from atomistic simulations

Demystifying the success of empirical distributions in space plasmas

Thermal conductivity and its relation to atomic structure for symmetrical tilt grain boundaries in silicon

Local temperatures out of equilibrium

Deciphering the 'Elixir of Life': Dynamic perspectives into the allosteric modulation of mitochondrial ATP synthase by J147, a novel drug in the treatment of Alzheimer's ...

Energy spectrum of a Langevin oscillator

Efficient determination of solid-state phase equilibrium with the multicell Monte Carlo method

Temperature and its control in molecular dynamics simulations

Energy localization and excess fluctuations from long-range interactions in equilibrium molecular dynamics

Vibrational Behavior of Water Adsorbed on Forsterite (Mg2SiO4) Surfaces Fluctuation theorems in q-canonical ensembles

O(N) Fluctuations and Lattice Distortions in 1-Dimensional Systems

Fluctuating temperature outside superstatistics: Thermodynamics of small systems Temperature fluctuations for a platoon of vehicles in contact with a heat bath Alloy thermodynamics via the Multi-cell Monte Carlo (MC) 2 method

Atomistic investigation on the conversion of plastic work to heat in high-rate shear deformation

Fluid-like behaviour and definitions of temperature and heat in 1-dimensional systems

Disruption of equilibrium due to lack of change

Study on Nanoporous Graphene-Based Hybrid Architecture for Surface Bonding On the elaboration of the next generation of thermodynamic models of solid solutions

Computational Study on Surface Bonding Based on Nanocone Arrays

Resonant Spatial Light Modulation: Optical Programming and Sensing at the Fundamental Limit

Theoretical Study on the Structure and Dynamics of Hydrogen Hydrates

Counterfactual thermodynamics: Extracting work from a lack of macroscopic change Molecular dynamics modelling of gold atomic force microscopy tips on multilayer graphene

A NOTE ON THE CONSEQUENCES OF A HOT MITOCHONDRION: SOME RECENT DEVELOPMENTS AND OPEN QUESTIONS1

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[]: