

# Uthpala Herath

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## Summary

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- Computational materials scientist with a demonstrated history in employing High-Performance Computing (HPC) and High-Throughput Computing (HTC) to perform first principles *ab-initio* electronic structure calculations and molecular dynamics simulations for bulk/2D materials and heterostructures using beyond-DFT quantum many body Green's functions methods including Dynamical Mean Field Theory (DMFT).
- Interests lie in strongly correlated materials (SCMs) useful for semiconductors, high- $T_c$  superconductivity, magnetism, Mott transitions, spintronics, quantum and neuromorphic computing, electro-chemical devices and energy applications.
- Skilled in the design and characterization of novel materials together with collaborative code development.

## Education

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### Doctor of Philosophy (Physics)

2015–2022

Department of Physics and Astronomy, West Virginia University, USA

Thesis Topic: *“Development of computational methods for electronic structural characterization of strongly correlated materials: from different ab-initio perspectives.”*

Advisor: Prof. Aldo H. Romero

### Master of Science (Physics)

2015–2017

Department of Physics and Astronomy, West Virginia University, USA

### Bachelor of Science (Physics major and Mathematics minor)

2010–2014

Department of Physics, Faculty of Science, University of Peradeniya, Sri Lanka

## Research Experience

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### Duke University

2022 –present

Postdoctoral Associate

Blum Group

#### Electronic structure method development for hybrid perovskites

- Development of electronic structure methods to study heavy elements with relativistic effects useful for novel semiconductor and renewable energy materials.
- Implementation of Periodic GW with relativistic spin-orbit coupling (In progress).

### West Virginia University

2017 –2022

Graduate Research Assistant (Computational Materials Science)

Romero Group

#### Strongly Correlated Materials Research

- Developed the DMFTwDFT framework enabling users to study properties of strongly correlated materials using DFT and DMFT with minimal technical expertise in collaboration with Prof. Hyowon Park at the University of Illinois Chicago. Created interfaces to popular DFT codes (VASP, Quantum Espresso, Siesta) through Maximally Localized Wannier Functions (MLWF's) and tight-binding Hamiltonians. Maintain a code repository of 40+ stars, 10+ forks and a community forum of 15+ users.
- Implemented full charge self-consistent DFT+DMFT calculations within the DMFTwDFT framework for the DFT code Siesta in collaboration with one of its lead developers, Prof. Javier Junquera. Interfaced Quantum Espresso with DMFTwDFT (both natively and through Aiiida) to perform one-shot DMFT calculations.

- Performed electronic structure calculations to study the electronic properties of transition and rare-earth correlated nickelate perovskites using DMFT $\omega$ DFT.
- Carried out molecular dynamics simulations on SCMs using our in-house MDWC code.
- Investigated the effect of symmetrically induced oxygen vacancies based on site occupation disorder in strongly correlated rare-earth nickelate perovskites and identified potential metal-to-insulator transitions. Further studied the minimum energy pathways for oxygen vacancy diffusion using Nudged Elastic Band algorithm (NEB) within DMFT for these vacancies. Identified potential applications in neuromorphic computing and micro-electronics. Calculated total energies with self-energy corrections to train an atomic interaction potential based on artificial neural networks (ANN) in collaboration with Dr. Subramanian Sankaranarayanan at Argonne National Lab.
- Applied virtual crystal approximation (VCA) within DMFT to study alloying/defects in SCMs.

### **Contribution to materials research frameworks**

- Co-developed the popular electronic structure pre/post-processing code PyProcar based on the original implementation by Prof. Aldo Romero and Prof. Francisco Muñoz. Maintain a code repository of 100+ stars, 50+ forks and a community forum of 155+ users.
- Co-developed MechElastic; an open-source Python library to calculate elastic properties of materials in collaboration with Dr. Sobhit Singh. Extended the original code as a Python library. Implemented the Equation of State (EOS) analysis and Abinit interface along with pressure corrections in VASP. Maintain a code repository of 25+ stars, 10+ forks and a community forum of 25+ users.
- Co-developed MDWC; our in-house command line open-source Python program for constrained molecular dynamics simulations in collaboration with Dr. Arturo Hernandez. Contributed in debugging and post-processing tool development.
- Contributed to PyChemia; our in-house open-source Python library for materials structural search developed by Dr. Guillermo Avendano-Franco.

### **Grant writing**

- Took a leadership role in coordinating the efforts of the group in preparing and securing allocation of XSEDE computing resources in Stampede2 and Bridges super-computing clusters. (2018, 2019, 2020)
- Successfully secured grant for DOE-EPSCOR Project : “*Searching optimal strongly correlated materials for neuromorphic computing.*” (2019)

### **Mentorship**

- Mentored a post-doc in the group on the usage of the DMFT $\omega$ DFT framework to treat strongly correlated materials.

### **West Virginia University**

Graduate Research Assistant (Computational Space Physics))

2016 –2017

Tu Group

### **The effect of magnetic field line curvature scattering on the rapid loss of ring current ions**

- Performed test particle simulations to quantify cumulative FLC scattering of ring current ions. Based on these simulation results, a physical parameterization for the FLC scattering was attempted to be incorporated into the RAM-SCB model.
- Received hands-on experience with a C, C++ based in-house particle tracer code.
- Presented preliminary results at the 2017 GEM conference in Portsmouth, VA.

### **Arthur C. Clarke Institute for Modern Technologies, Sri Lanka**

Intern - Space Applications Division

Jan, 2014 –Dec, 2014

### **Identification of Blue Stragglers in the globular cluster M53 using CCD photometry**

Advisors : Mr. S. Gunasekara, Dr. T. Ranawake

- Worked in conjunction with an expert team of scientists at the Space Applications division of ACCIMT to perform an extensive study of globular cluster characteristics using Aperture Photometry and PSF (Point Spread Function) Fitting Photometry methods with the IRAF (Image Reduction and Analysis Facility) system.

- Telescopic data of the globular cluster M53 was used to construct a novel Color-Magnitude Diagram (CMD) of the cluster, which was in turn utilized to study its stellar population, mainly focusing on the Blue Straggler Star population.

## University of Peradeniya, Sri Lanka

2012–2013

Undergraduate Research Student - Department of Physics, Faculty of Science

### Nano Magnets and their Applications

Advisor : Prof. P. Samarasekara

- Investigated the traits, synthesis methods, functionalization and applications of Nano Magnets.

## Skills

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- **First principles *ab-initio* calculations :**
  - Density Functional Theory (DFT) codes - *FHI-aims, VASP, Siesta, Abinit, Elk, Quantum Espresso for DFT, DFT+U, Hybrid functionals, GW calculations*
  - Electronic, vibrational, magnetic, mechanical, elastic, thermal and optical properties of bulk material and heterostructures.
- **Materials Research Methods and Tools :**
  - DMFTwDFT - *An open-source framework to perform Dynamical Mean Field Theory (DMFT) calculations to study strongly correlated materials (SCMs), Mott transitions, mass renormalization effects, spectral functions. (in-house code)*
  - Wannier90 - *Basis transformation through Maximally Localized Wannier Functions.*
  - Impurity Solvers and Quantum Monte Carlo Methods - *Solving the Anderson Impurity model with the CTQMC method.*
  - Tightbinding Hamiltonians - *Employing Tightbinding Hamiltonians to study strongly correlated material.*
  - PyProcar - *A robust, open-source Python library used for pre- and post-processing of the electronic structure data coming from DFT calculations. (in-house code)*
  - PyChemia - *An open-source Python library for materials structural search. (in-house code)*
  - MechElastic - *An open-source Python library to calculate elastic properties of materials. (in-house code)*
  - AFLOW - *A materials database and search framework.*
  - VTST Tools - *A code for finding saddle points and evaluating transition state theory (TST) rate constants.*
  - DiSPy - *A utility for applying the distortion symmetry method (DSM) to the calculation of minimum energy pathways using the nudged elastic band (NEB) algorithm.*
  - SOD (Site Occupation Disorder) - *A package of tools for the computer modelling of periodic systems with site disorder, using the supercell ensemble method.*
  - Molecular Dynamics Simulations - *MDWC (A command line open-source Python program for constrained molecular dynamics simulations. - in-house code), VMD, qAgate, Jmol.*
  - Pymatgen (Python Materials Genomics) - *A robust, open-source Python library for materials analysis.*
  - phonopy - *An open source package for phonon calculations at harmonic and quasi-harmonic levels.*
  - Graphical tools - *VESTA (Visualization for Electronic Structure Analysis), gnuplot, Xmgrace, xcrysden.*
  - Vacancy, defect and alloying methods - *Virtual Crystal Approximation (VCA), Coherent Potential Approximation (CPA), SPRKKR.*
- **Programming languages :** Python, Fortran, C, C++, Bash, Matlab
- **Computational methods :** High-performance computing (HPC), High-throughput computing (HTC), Parallel computing with OpenMP and MPI, Regression analysis, Differential solvers and Integrators, Numerical analysis algorithms, Object-oriented and Procedural programming paradigms
- **Libraries :** BLAS, LAPACK, FFTW, GSL
- **Graphics and Publishing :** Latex, Markdown, Adobe Photoshop, Adobe Premiere, Adobe After Effects, Adobe Illustrator
- **Supercomputing clusters experience:** NERSC Perlmutter and Cori, WVU Thorny Flat and Spruce Knob, XSEDE Bridges2 and Stampede2, LANL HPC
- **Languages :** English (Full professional proficiency), Sinhala (Native or bilingual proficiency), Tamil (Elementary proficiency)
- **Soft skills :** Creative problem solving, Leadership, Organization, Communication, Teamwork, Flexibility, Networking

## Scholarships and Awards

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- **Ovshinsky Travel Award for the APS March Meeting 2021** 2021  
*Awarded by the American Physical Society to fund registration for the APS March Meeting in March, 2021.*
- **Robert T. Bruhn Physics Research Award at West Virginia University** 2020  
*Awarded by the Department of Physics and Astronomy to extend support to the research effort of a graduate or undergraduate student in the department in nanotechnology and material science.*

- **Office of the Provost Graduate Student Travel Award** 2018, 2019  
Awarded by the Office of the Provost, West Virginia University to fund travel to research conferences.
- **Eberly College of Arts and Sciences Graduate Student Travel Award** 2018, 2019  
Awarded by the Eberly College of Arts and Sciences, West Virginia University to fund travel to research conferences.

## Publications

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1. **Herath, U.**, Singh, V., Bhat, S. S., Park, H. & Romero, A. H. *Stability and diffusion of oxygen vacancies in  $\text{LaNiO}_3$ : a DMFT study* 2022.
2. Singh, V., **Herath, U.**, Wah, B., Liao, X., Romero, A. H. & Park, H. DMFTwDFT: An open-source code combining Dynamical Mean Field Theory with various density functional theory packages. *Computer Physics Communications* **261**, 107778 (2021).
3. Singh, S., Lang, L., Dovale-Farelo, V., **Herath, U.**, Tavadze, P., Coudert, F.-X. & Romero, A. H. MechElastic: A Python library for analysis of mechanical and elastic properties of bulk and 2D materials. *Computer Physics Communications* **267**, 108068 (2021).
4. **Herath, U.**, Tavadze, P., He, X., Bousquet, E., Singh, S., Muñoz, F. & Romero, A. H. PyProcar: A Python library for electronic structure pre/post-processing. *Computer Physics Communications* **251**, 107080 (2020).
5. Herath, A. & **Herath, U.** Developing an Expert System for Plant Pest Diagnosis. *Annals of the Sri Lanka Department of Agriculture* **15**, 381 (2012).

**Please visit my Google Scholar profile for an updated list of publications.**

<https://scholar.google.com/citations?user=m6VPFYoAAAAJ&hl=en&authuser=1>

## Publications in progress

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6. **Herath, U.**, Singh, V., Wah, B., Park, H. & Romero, A. H. A study of strongly correlated alloys through virtual crystal approximation within dynamical mean field theory.
7. Wah, B., Singh, V., **Herath, U.**, Romero, A. H. & Park, H. Force calculation within the DMFTwDFT framework.
8. Tavadze, P., Farah, F., Lang, L., Romestan, Z., Dovale-Farelo, V., **Herath, U.**, Singh, S. & Romero, A. H. Fermi3D: Expanding PyProcar's Fermi surface implementations.

## Presentations

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- **Uthpala Herath**, Pedram Tavadze, He Xu, Eric Bousquet, Sobhit Singh, Reese Boucher, Logan Lang, Freddy Farah, Francisco Muñoz and Aldo H Romero  
"A tutorial on PyProcar: A Python library for electronic structure pre/post-processing"  
Rutgers University, June 25th, 2021 (INVITED VIRTUAL TALK)
- **Uthpala Herath**, Pedram Tavadze, He Xu, Eric Bousquet, Sobhit Singh, Reese Boucher, Logan Lang, Freddy Farah, Francisco Muñoz and Aldo H Romero  
"Recent Developments in PyProcar: A Python library for electronic structure pre/post-processing"  
APS March Meeting, Mar 15-19, 2021 (VIRTUAL TALK)
- **Uthpala Herath**, Vijay Singh, Benny Wah, Xingyu Liao, Hyowon Park and Aldo H. Romero  
"DMFTwDFT: An open-source code combining Dynamical Mean Field Theory with various Density Functional Theory packages"  
APS Mid Atlantic Section Meeting, December 4-6, 2020 (VIRTUAL TALK)
- **Uthpala Herath**, Pedram Tavadze, Xu He, Eric Bousquet, Sobhit Singh, Francisco Muñoz, and Aldo H. Romero  
"PyProcar: A Python library for electronic structure pre/post-processing"  
Carolina Science Symposium, Nov 12-13, 2020 (VIRTUAL TALK)
- **Uthpala Herath**, Pedram Tavadze, Xu He, Eric Bousquet, Sobhit Singh, Francisco Muñoz, and Aldo H. Romero  
"PyProcar: A Python library for electronic structure pre/post-processing"  
Electronic Structure Workshop, June 1-4, 2020, University of California, Merced (VIRTUAL TALK)
- **Uthpala Herath**, Hyowon Park and Aldo H. Romero  
"An ab-initio DFT+DMFT study of the effect of oxygen vacancies on structural, electronic and magnetic properties of rare-earth nickelate perovskites ( $\text{RNiO}_3$ )"  
APS March Meeting, March 4-8, 2019, Boston, MA (TALK)

- **Uthpala Herath**, Hyowon Park and Aldo H. Romero  
*“Development of computational methods for the characterization of novel strongly correlated materials”*  
 International Summer School on Computational Quantum Materials, June 2018, Sherbrooke, Québec, Canada (POSTER)
- **Uthpala Herath** and Weichao Tu  
*“The Effect of Magnetic Field Line Curvature Scattering on the Rapid Loss of Ring Current Ions”*  
 Geospace Environment Modeling (GEM) conference, July 2017, Portsmouth, VA (POSTER)

## Teaching Experience

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- **Volunteer Instructor**, Department of Physics and Astronomy at West Virginia University 2017-present  
*Occasionally covered for Dr. Aldo Romero’s (thesis advisor) graduate level Quantum Mechanics class.*
- **Graduate Teaching Assistant**, Department of Physics and Astronomy at West Virginia University 2015-2017  
*Conducted lab and office hours on electrostatics, electromagnetism and optics for STEM undergraduate students in Physics 112. Graded exams and homework.*
- **Graduate Teaching Assistant**, Department of Physics at University of Peradeniya, Sri Lanka 2014-2015  
*Conducted lab and tutorial sessions for STEM undergraduate students. Graded exams and homework.*

## Contribution to code development

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See full list on <https://uthpalaherath.github.io/projects>.

- **FHI-aims**  
*“All-electron electronic structure theory with numeric atom-centered orbitals.”*  
<https://fhi-aims.org>
- **PyProcar**  
*“PyProcar is a robust, open-source Python library used for pre- and post-processing of the electronic structure data coming from DFT calculations.”*  
<https://github.com/romerogroup/pyprocar>
- **DMFTwDFT**  
*“An open-source computational package (and a library) combining DMFT with various DFT codes interfaced through the Wannier90 package.”*  
<https://github.com/DMFTwDFT-project/DMFTwDFT>
- **MechElastic**  
*“An open-source Python library to calculate elastic properties of materials.”*  
<https://github.com/romerogroup/mechelastic>
- **PyChemia**  
*“PyChemia is an open-source Python library for materials structural search.”*  
<https://github.com/MaterialsDiscovery/PyChemia>
- **mdwc**  
*“The molecular dynamics with constraints (mdwc3) package is a command line open-source Python program for constrained molecular dynamics simulations.”*  
<https://github.com/romerogroup/mdwc3>
- **pymatgen**  
*“Pymatgen (Python Materials Genomics) is a robust, open-source Python library for materials analysis.”*  
<https://github.com/materialsproject/pymatgen>

## Workshops attended

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- **Quantum Monte Carlo Virtual Workshop 2021** : Oct 5 –November 23, 2021  
 ORNL
- **Comscope Summer School 2021** : June 21 –24, 2021  
 BNL
- **PRACE Seasonal School on Modeling Materials using HPC and AI/ML** : June 15 –18, 2021

- **Bridging first-principles calculations and effective Hamiltonians** : June 7 –16  
IIT Genova, Italy
- **20th International Workshop on Computational Physics and Materials Science: Total Energy and Force Methods**  
: Feb 23 –25, 2021  
ICTP
- **BerkeleyGW Tutorial Workshop** : Jan 4 –6, 2021
- **2nd Berkeley Excited States Conference (BESC2021)** : Jan 7 –8, 2021
- **Texas A&M University Summer School On Computational Materials Science Across Scales** : June 29 –July 10, 2020
- **School on Wannier90 v3.0: new features and applications** : March 25 –27, 2020
- **School on Stochastic Approaches to Electronic Structure Calculations** : July 29th –August 2nd, 2019  
University of Pittsburgh, PA
- **3rd Summer School on Materials Growth and Design: Exotic Magnetic States in Quantum Mechanic** : June 17 –June 22, 2018  
Johns Hopkins University, Baltimore, MD
- **International Summer workshop on Computational Quantum Materials** : May 27 –June 8, 2018  
Sherbrook, Québec, Canada
- **CISM Space Weather Summer School** : July 11 –July 22, 2016  
Boulder, CO

## Extracurricular Activities

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- Co-chair (Social Committee) - Duke University Postdoc Association (DUPA) July, 2022  
- *DUPA was established to address the needs and concerns of Duke postdocs and celebrate their accomplishments. As the co-chair of the social committee I am involved in organizing events that bring postdocs together.*
- Judge - Virtual Summer Undergraduate Research Symposium, WVU July, 2020  
- *Volunteered as a graduate student judge at the 2020 Virtual Summer Undergraduate Research Symposium organized by the Office of Undergraduate Research in the Mathematics and Physics Category*
- Coordinator - Condensed Matter Seminar, Department of Physics and Astronomy, WVU 2018–Current  
- *Coordinated the Condensed Matter Seminar sessions at the Department of Physics and Astronomy, West Virginia University.*
- Outreach Volunteer - Physics and Astronomy Graduate Student Organization, WVU (PAGSO) 2015–Current  
- *Conducted physics outreach programs to enlighten the public on science in West Virginia.*
- Member - Sri Lankan Student's Organization, WVU (SLSA) 2018–Current  
- *SLSA is a student body of Sri Lankan students at West Virginia University aiming to share the Sri Lankan cultural and heritage experience with the rest of the world.*  
- *Took a leading role in organizing a vigil in memory of the Easter Sunday terrorist attacks that occurred in Sri Lanka in April, 2019.*
- President - Astronomy Society of University of Peradeniya 2013–2014  
- *Organized "Water Rocket Challenge 2014"; the first-ever water rocket competition held at University of Peradeniya*  
- *Conducted regular astronomy lectures, planetary and deep sky observation sessions and workshops in the university as well as in local schools aiming to promote astronomy among students and the general public.*  
- *Volunteered for United Nations World Space Week- Sri Lanka (2014)*
- Junior Editor - Physical Society of University of Peradeniya 2012–2013  
- *Compiled academic articles for the "Echo" magazine which is the annual scientific journal issued by the society*
- Faculty Representative - The Ceylon Drama Society of University of Peradeniya 2011–2012  
- *Represented the Faculty of Science in the Ceylon Drama Society of the University of Peradeniya, Sri Lanka.*
- Volunteer - Let Them Smile 2010–2014

-Volunteered in “Let Them Smile”, an annual charity event organized by the undergraduates of the Faculty of Science of University of Peradeniya, Sri Lanka to facilitate the education of under privileged students in rural schools around the country.  
URL: <https://www.facebook.com/letsmile14>

- Volunteer Software Developer - Department of Agriculture, Sri Lanka 2012  
-Took part in an auxiliary research project conducted by the Sri Lanka Department of Agriculture to develop a system for plant disease diagnosis. This would allow farmers around the island to use a software framework to self-diagnose and treat plant diseases. Assisted in computational aspects of this system.

## References

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| <ul style="list-style-type: none"><li>• <b>Prof. Volker Blum</b><br/>Department of Mechanical Engineering and Materials Science<br/>Duke University<br/>USA<br/>✉ <a href="mailto:volker.blum@duke.edu">volker.blum@duke.edu</a><br/>☎ +1 (919) 660 5279</li><li>• <b>Prof. Aldo H. Romero</b><br/>Department of Physics and Astronomy<br/>West Virginia University<br/>USA<br/>✉ <a href="mailto:aldo.romero@mail.wvu.edu">aldo.romero@mail.wvu.edu</a><br/>☎ +1 (304) 293-6317</li></ul> | <ul style="list-style-type: none"><li>• <b>Prof. Hyowon Park</b><br/>Department of Physics<br/>University of Illinois Chicago<br/>USA<br/>✉ <a href="mailto:hyowon@uic.edu">hyowon@uic.edu</a><br/>☎ +1 (312) 996-8913</li><li>• <b>Prof. Francisco Muñoz</b><br/>Department of Physics<br/>University of Chile<br/>Chile<br/>✉ <a href="mailto:fvmunoz@gmail.com">fvmunoz@gmail.com</a></li></ul> |
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