

Aditya Sundar, Ph.D student, University of Michigan

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Education

- 2017 – 2022 **Ph.D., University of Michigan** in Materials Science and Engineering CGPA: 3.93/4.00
Thesis title: *Multi-scale modelling of corrosion in crystalline and amorphous oxides*
Certificate: Computational Science and Discovery
Advisor: [Liang Qi](#)
- 2017 – 2015 **M.S., Cornell University** in Materials Science and Engineering CGPA: 3.90/4.00
Thesis title: *The Structure and Chemistry of Epitaxial 2D Chalcogenides*
Advisor: [Huili Grace Xing](#)
- 2011 – 2015 **B.Tech., Indian Institute of Technology Madras**
in Metallurgical and Materials Engineering CGPA: 8.66/10.00
Thesis title: *Printed Silver Lines in MIS Solar Cells*
Advisor: [Parasuraman Swaminathan](#)
Minor degree: Management studies

Experience

- Summer 2021 **Idaho National Laboratory**, Idaho Falls, USA
Research Internship, Computational Microstructural Science division
- Density functional theory based investigation of candidate moderator and cladding materials for compact nuclear microreactors
- Summer 2013 **Larsen & Toubro**, Mumbai, India
Engineering Internship, Heavy Engineering division

Skills

- | | |
|------------------------|---|
| Methods | Electronic structure, molecular dynamics, monte carlo, potential energy landscape, computational thermodynamics, spectroscopy |
| Atomic Simulation | VASP, Quantum ESPRESSO, LAMMPS, ART Nouveau (Activation Relaxation Technique) |
| Materials Simulation | Thermo-Calc, Abaqus FEA. |
| Computation and design | Mathematica, Python (NumPy, SciPy, OpenCV), \LaTeX , Linux |
| Experiments | XRD, XPS, EXAFS, Raman, SEM, Conductive-AFM |

Research Publications

Journal Articles

- 1 Hu, Y., **Sundar, A.**, Ogata, S., & Qi, L. (2021). Screening of generalized stacking fault energies, surface energies and intrinsic ductile potency of refractory multicomponent alloys. *Acta Materialia*, 116800. [link](#).
- 2 **Sundar, A.**, Chen, G., & Qi, L. (2021b). Substitutional adsorptions of chloride at grain boundary sites on hydroxylated alumina surfaces initialize localized corrosion. *npj Materials Degradation*, 5. [link](#).

- 3 **Sundar, A., & Qi, L.** (2021). Stability of native point defects in Al_2O_3 under aqueous electrochemical conditions. *Journal of Applied Electrochemistry*. [link](#).
- 4 Lu, H., Reese, C., Jeon, S., **Sundar, A.**, Fan, Y., Rizzi, E., ... Goldman, R. (2020). Mechanisms of GaN quantum dot formation during nitridation of Ga droplets. *Applied Physics Letters*, 116(6), 062107. [link](#).
- 5 Li, W., Nomoto, K., **Sundar, A.**, Lee, K., Zhu, M., Hu, Z., ... Gao, X. et al. (2019). Realization of GaN polarons using selective-area regrowth by MBE and its breakdown mechanisms. *Japanese Journal of Applied Physics*, 58(SC), SCCD15. [link](#).
- 6 Vishwanath, S., **Sundar, A.**, Liu, X., Azcatl, A., Lochocki, E., Woll, A. R., ... Peng, X. et al. (2018). MBE growth of few-layer ZnO on 3D substrates. *Journal of Crystal Growth*, 482, 61–69. [link](#).

Conference Proceedings


- 1 Hu, Y.-J., **Sundar, A.**, Chen, G., & Qi, L. (2021). Title: Screening of generalized stacking fault energies, surface energies and intrinsic ductile potency of refractory multicomponent alloys. In *The minerals, metals & materials society 2021 annual meeting & exhibition*, Orlando, USA.
- 2 Kinzer, B., **Sundar, A.**, & Bala Chandran, R. (2021). Oxide dispersion particle clustering using phase field modelling in nickel-based superalloys. In *2021 MRS Spring Meeting & Exhibit*, Phoenix, USA.
- 3 **Sundar, A.**, Chen, G., & Qi, L. (2021a). Electronic structure mechanisms to explain the onset of Cl-induced localized corrosion in Al_2O_3 . In *The minerals, metals & materials society 2021 annual meeting & exhibition*, Orlando, USA.
- 4 **Sundar, A.**, Kinzer, B., & Bala Chandran, R. (2021). Oxidation and carburization behavior of iron- and nickel-based alloys in supercritical CO_2 environments. In *2021 MRS Spring Meeting & Exhibit*, Phoenix, USA.
- 5 **Sundar, A.**, & Qi, L. (2019). Electric field dependent ionic transport in passive corundum Al_2O_3 from DFT application to localized corrosion. In *2019 MRS Fall Meeting & Exhibit*, Boston, USA.

Research Projects


2018 – 2021 ■ Ab initio modelling of chloride induced localized corrosion in passive oxides

- Explained the semi-crystalline structure in electrochemical environments; by completely sampling the energy landscape of point defects in $\alpha\text{-Al}_2\text{O}_3$ and $\alpha\text{-Cr}_2\text{O}_3$.
- Designed and implemented kinetic Monte Carlo simulations to model electromigration; by calculating electrical field dependent ionic migration barriers.
- Identified the electronic origins of Cl adsorption at weak surface sites; from static and dynamic molecular simulations in standard aqueous conditions.
- Lower adsorption energy of Cl at grain boundary sites accelerates and sustains a catalytic Al_2O_3 dissolution reaction, compared to single crystal surfaces.
- Calculations validate the Point Defect Model and explain Cl induced localized aqueous corrosion in Al_2O_3 .

Research Projects (continued)

2019 – 2019  Thermodynamic model for GaN quantum dot formation

- Provided ab initio results to explain surface coverage dependent GaN quantum dot formation on Si. Results used to interpret morphological characteristics of MBE grown GaN quantum dots.
- Designed and calculated adsorption thermodynamics of Ga atoms to explain QD size distribution on oxidised and nitrided Si(100)– 2×2 reconstructed surfaces.

2020 – 2021  Alloy selection for high temperature heat exchanger materials

- Developed CALPHAD based thermokinetic models to study supercritical CO₂ induced oxidation-carburisation of Fe and Ni alloys; up to 1100°C and 200 bar. Generated equilibrium phase diagrams for Incoloy MA 956, Haynes 214 and AMDRY 386 under different surface activities of C and O₂.
- Recommended the use of Haynes 214 due to better carburisation resistance, compared to Incoloy MA 956. Incoloy forms higher amounts of carbide precipitates.
- Collaborated with manufacturing research teams and formulated cyclic oxidation models; to estimate component lifetimes at 700°C–1100°C using parabolic oxidation behaviour.


2015 – 2017  Structural, chemical and electronic properties of MBE grown 2D materials

- Confirmed up to 44% Te alloying in MBE MoSe_xTe_{2-x} on CaF₂ and GaAs substrates; by XRD, Raman, XPS, TEM and EXAFS examinations. Compositional variation of bandgap and excitonic states shown by absorption spectroscopy.
- Imaged crystallographic orientation, twinning and mosaicity in ultrathin 2D heterostructures (graphene-transition metal dichalcogenides) by synchrotron measurements.
- Developed Mathematica scripts to parse, analyze and visualize large datasets obtained from synchrotron based diffraction and spectroscopy experiments.
- Identified highly conductive leakage paths along MBE regrown interfaces by correlating topography and current area maps, to explain limited reverse breakdown voltage.

Coursework Projects

Fall 2020  Computer Vision

Implemented a fruit-sorting algorithm using blob detection and superpixel image segmentation methods; to sort fresh and rotten fruits.

Fall 2018  Integrated Computational Materials Engineering

Developed an Abaqus finite element model to simulate uniaxial compression of an Al-Li cylinder. Calibrated multiscale strength model with experiment to study precipitate evolution, texture and mechanical behaviour under compression.

Winter 2018  Kinetic Monte Carlo simulations of crystal growth

Developed a 2D code and Mathematica visualisations to simulate the epitaxial growth of zinc blende GaAs. Surface diffusion energies from molecular dynamics calculations were used to evaluate temperature dependent point defects and surface morphologies.

Mentoring and Leadership

- 2019 – 2021
- Guided and trained Master's degree student in the design and implementation of first principles calculations using VASP. Research topics included (1) surface stability of passive oxides and (2) electronic structure origin of Cl-induced depassivation.
 - Mentored research projects for 3 undergraduate students in the (1) design of ab-initio calculations, (2) implementation of ab-initio calculations using Quantum Espresso and (3) analysis of results. Topics covered included (1) thermodynamic stability of semiconductor polymorphs and (2) electronic stability of point defects in crystal semiconductors.
 - Lab group server & website operations: <http://cms.engin.umich.edu/>
 - Vice President, SPIC MACAY (Society for the Promotion of Indian Classical Music And Culture Amongst Youth), University of Michigan chapter. Led a team of 10-15 students in the organisation of biannual musical concerts in the university. Responsible for securing grants and event logistics.
- 2018
- Graduate Student Instructor for undergraduate materials science course. Led weekly discussions and office hours, and designed problem sets and demonstrations for a class of 30 undergraduate students.

References

Available on Request