SOLVING ATOMIC STRUCTURE USING STATISTICAL MECHANICAL SEARCHES ON X-RAY SCATTERING DERIVED POTENTIAL ENERGY SURFACES

by

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DEDICATION

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Abstract

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Introduction

This is the introduction to the thesis.

ATOMIC STRUCTURE: EXTRACTION AND APPLICATION

- 1.1 Atomistic Goals
- 1.2 Atomistic Experiments
- 1.3 Atomistic Simulations

How do I combine/separate the Ensemble, PES and PDF math/computation sections?

ATOMIC PAIR DISTRIBUTION FUNCTION:

THEORY AND COMPUTATION

2.1 Theory

Derivation

2.2 Computation

HPC and **GPUs**

GPUs and vectorization

Map from ij space to k space

Periodic Boundary Conditions

- 2.3 Experiment
- 2.4 Data Processing Workflow

MetadataStore Side Loading

CHAPTER 3

STATISTICAL MECHANICAL ENSEMBLES AND POTENTIAL ENERGY SURFACES

3.1 Ensembles

Hamiltonian Monte Carlo

No-U-Turn-Sampling

Grand Canonical Monte Carlo

Configurational Biasing

BENCHMARKING

4.1 PDF

Au55

4.2 PDF WITH ADPS

Annealing and Aggregation of 2nm

Au Nanoparticles

- 5.1 Experiments
- 5.2 Data Processing
- 5.3 Data Analysis
- 5.4 SIMULATION
- 5.5 STRUCTURAL ANALYSIS
- 5.6 Conclusions

Phase Changes and Annealing Dynamics of ${\rm Pr}_2{\rm NiO}_4 \ {\rm and} \ {\rm its} \ {\rm derivatives}$

- 6.1 Experiments
- 6.2 Data Processing
- 6.3 Data Analysis
- 6.4 SIMULATION
- 6.5 STRUCTURAL ANALYSIS
- 6.6 Conclusions

Conclusion