

# **GPU-ISLE**

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# **GPU-ISLE TEAM**



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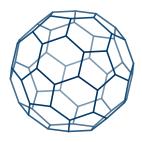


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# SIMULATING CARBON NANO SYSTEMS

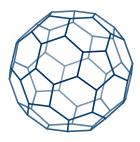
- Lattices of carbon atoms
- Various geometries
- High technical interest
  - Improved semiconductors
  - Filtering techniques
- Simulating behaviour: Quantum Field Theory
  - System description with Energy functional (Hamiltonian)
  - Generation of many states (Configurations)
  - Evaluating physical quantities





# SIMULATING CARBON NANO SYSTEMS

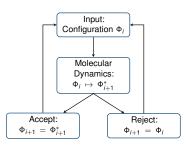
- Lattices of carbon atoms
- Various geometries
- High technical interest
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  - 2. Generation of many states (Configurations)
  - Evaluating physical quantities





### THE HYBRID MONTE CARLO ALGORITHM

- Building Markov Chain
- Iteratively updating previous configuration
- Update by Molecular Dynamics
  - Integrate Hamilton equations
    - Leapfrog
  - Requires: Matrix inversion
  - Requires: Matrix, vector operations
- Dimension  $N_{\tau} \times \mathcal{O}(N_{\sigma})$ 
  - $C_{60}: N_{\tau} \times \mathcal{O}(60)$





# **ISLE**

### **Current State**

- C++: Heavy math
- Python: Interface
- Build upon
  - Blaze
  - Lapack

#### Goal

- Port of Molecular Dynamics
  - Leapfrog
  - Matrix inverter
  - Matrix, vector operations
- Hybridization of Infrastructure (CPU,GPU)
- C++ with CUDA
- Blaze: CUDA Managed Pointer
- Utilize CUDA Math Libraries

