



# Performance Portable Assembly For Plasma Fluid Equations

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# Problem Description

## Prediction of plasma processes at moderate to high densities

- Multifluid plasma equations – Electrons, ions and neutrals
  - Six-dimensional Boltzmann equations too high of dimensionality
  - Particle methods too expensive due to plasma frequency
  - Sets of 5 or 13 coupled PDEs with stiff source terms need to be solved
- Equations need to be composable and maintainable

## Analysis Beyond Forward Simulation

- Forward solves are not enough – we want to explore complex solution spaces:
  - Simultaneous analysis and design adds requirements (typically sensitivities)
  - Do not burden analysts/physics experts with analysis algorithm requirements: i.e. programming sensitivities for implicit solvers, optimization, stability, bifurcation analysis and UQ

**Engine must be flexible, extensible, maintainable and EFFICIENT!**



# Problem Description

## Prediction of plasma processes at moderate to high densities

- Multifluid plasma equations – Electrons, ions and neutrals
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## Directed Acyclic Graph-based Assembly

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## Template-based Generic Programming

**Engine must be flexible, extensible, maintainable and EFFICIENT!**

# Electron Fluid Plasma Model

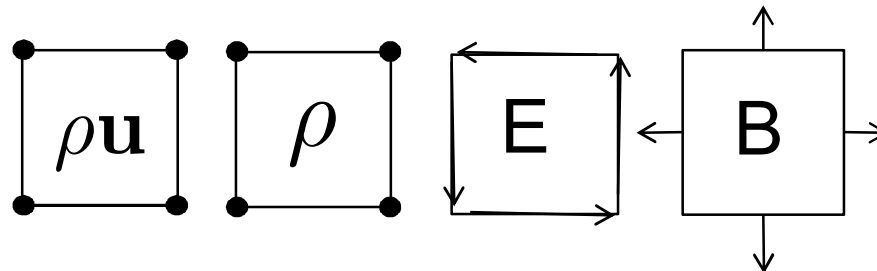
$$\frac{\partial(\epsilon \mathbf{E})}{\partial t} - \nabla \times \left( \frac{1}{\mu} \mathbf{B} \right) = -\frac{q}{m} \rho \mathbf{u}$$


$$\frac{\partial \mathbf{B}}{\partial t} + \nabla \times \mathbf{E} = \mathbf{0}$$

$$\frac{\partial(\rho \mathbf{u})}{\partial t} + \nabla \cdot (\rho \mathbf{u} \otimes \mathbf{u}) = \frac{q}{m} \rho (\mathbf{E} + \mathbf{u} \times \mathbf{B})$$

$$\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{u}) = 0$$

- Using 3D all Hex mesh
- 32 nodal + 12 edge + 6 face = 50 DOFs per element





# Drekar/Panzer: Algorithms and Software

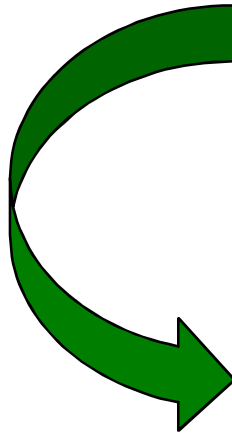
**Drekar:** Algorithms research application

- Targets coupled multi-physics
- Large scale simulation (>100k cores)
- Advanced algorithm demonstration

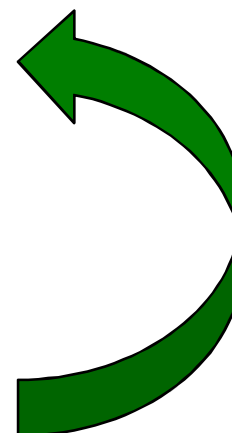
A co-dependent development relationship

**Panzer:** Multi-physics assembly engine

- Finite element focused (currently)
- Embedded analysis (AD, Sensitivities)
- Technology sharing and deployment



**Drekar drives Panzer**  
requirements and  
design goals



**Panzer provides**  
**Drekar** flexible  
infrastructure and core  
technologies

John Shadid  
Roger Pawlowski  
Eric Cyr  
Edward Phillips  
Tim Wildey

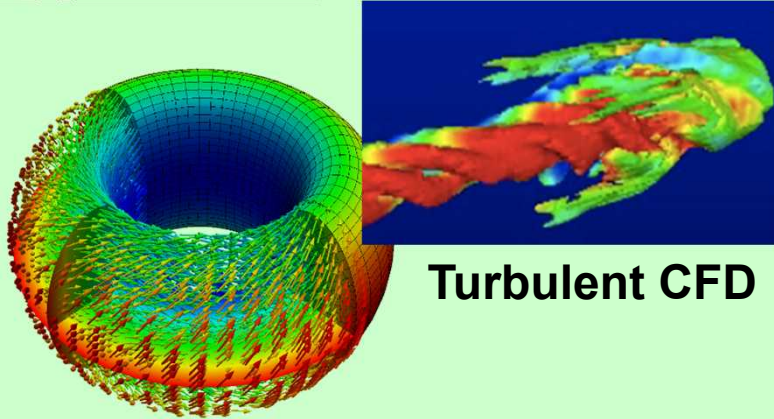
Tom Smith  
Paul Lin  
David Sondak  
Paula Weber  
Richard Kramer

Allen Robinson  
Matt Bettencourt  
Sidafa Conde  
Ben Seefeldt  
Chris Siefert

Andrew Bradley  
Greg von Winckel  
David Hensinger

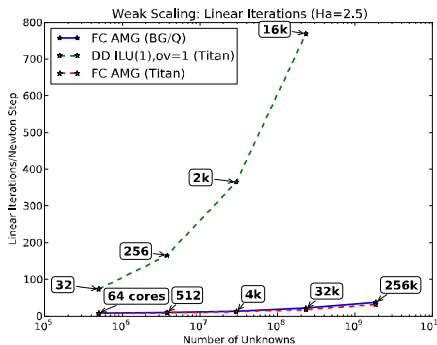
# Drekar/Panzer: Capabilities

## Applications



**Turbulent CFD**

**Magnetohydrodynamics**

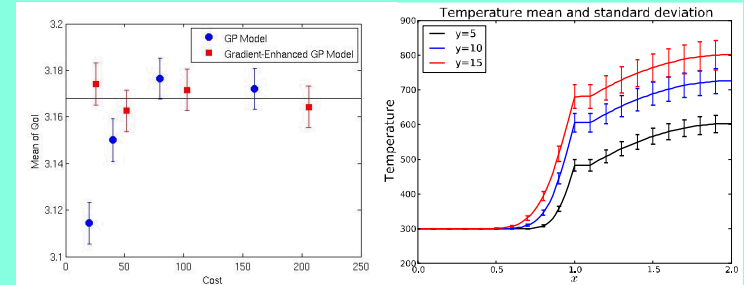


**Algebraic Multigrid  
(>100k cores)**

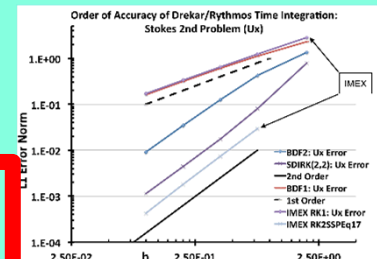
$$\mathcal{A} = \begin{bmatrix} I & \\ BF^{-1} & I \end{bmatrix} \begin{bmatrix} F & B^T \\ S & \end{bmatrix}$$
$$S = C - BF^{-1}B^T$$

**Block  
Preconditioning**

## Discretizations & Algorithms

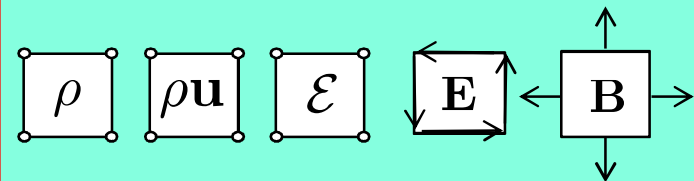


**Uncertainty Quantification**



**IMEX**

**PDE Constrained  
Optimization**

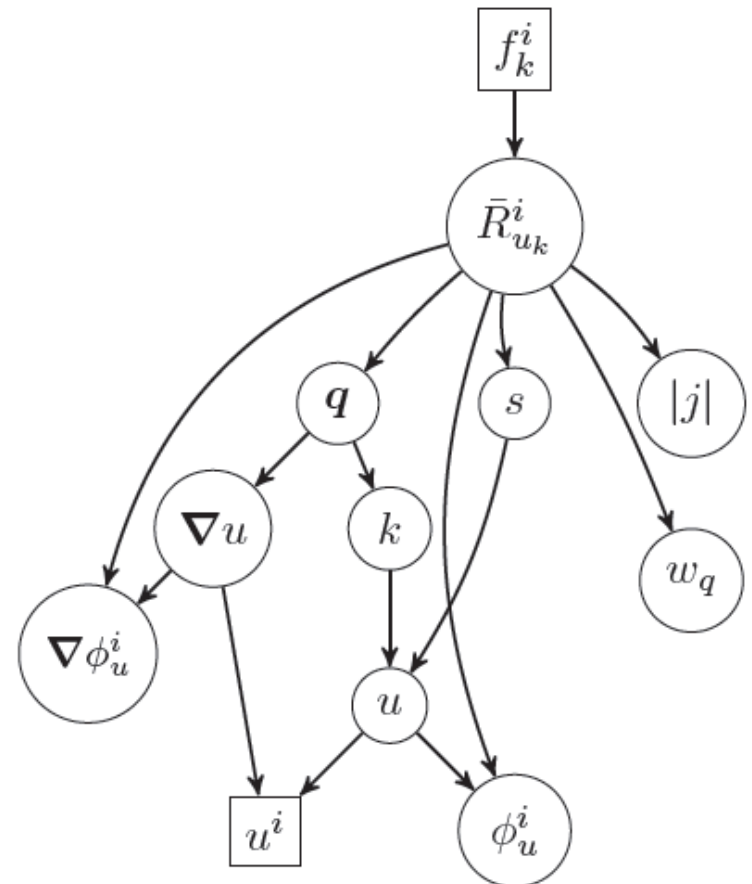


**Compatible Discretizations**

# Lightweight DAG-based Expression Evaluation

- Decompose a complex model into a graph of simple kernels (functors)
  - Decomposition is NOT unique
- Supports rapid development, separation of concerns and extensibility.
- A node in the graph evaluates one or more **fields**:
  - Declare fields to evaluate
  - Declare dependent fields
  - Function to perform evaluation
- Separation of data (Fields) and kernels (Expressions) that operate on the data
  - Fields are accessed via multidimensional array interface (shards or kokkos)

$$R_u^i = \int_{\Omega} [\phi_u^i \dot{u} - \nabla \phi_u^i \cdot \mathbf{q} + \phi_u^i s] \, d\Omega$$





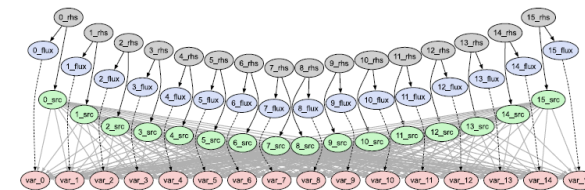
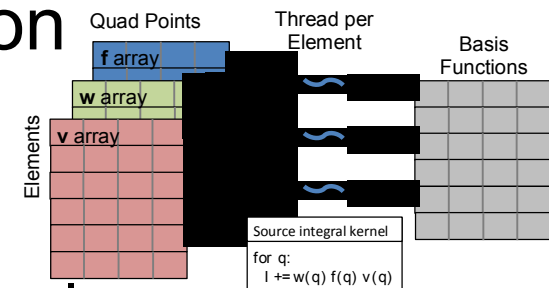
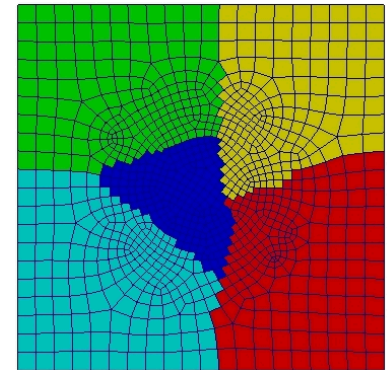
# DAG Evaluation

- The problem domain is broken into worksets
  - Smaller number of elements which fit into cache or other memory structures
  - The entire DAG is evaluated for a single workset
  - Each math kernel can be threaded over the workset elements
- Worksets are filled with a gather operation
- The workset's contributions are accrued in a matrix with a scatter operation
- All intermediate calculations are independent of if they are computing a Jacobian or residual or sensitivity



# Parallelization Strategies

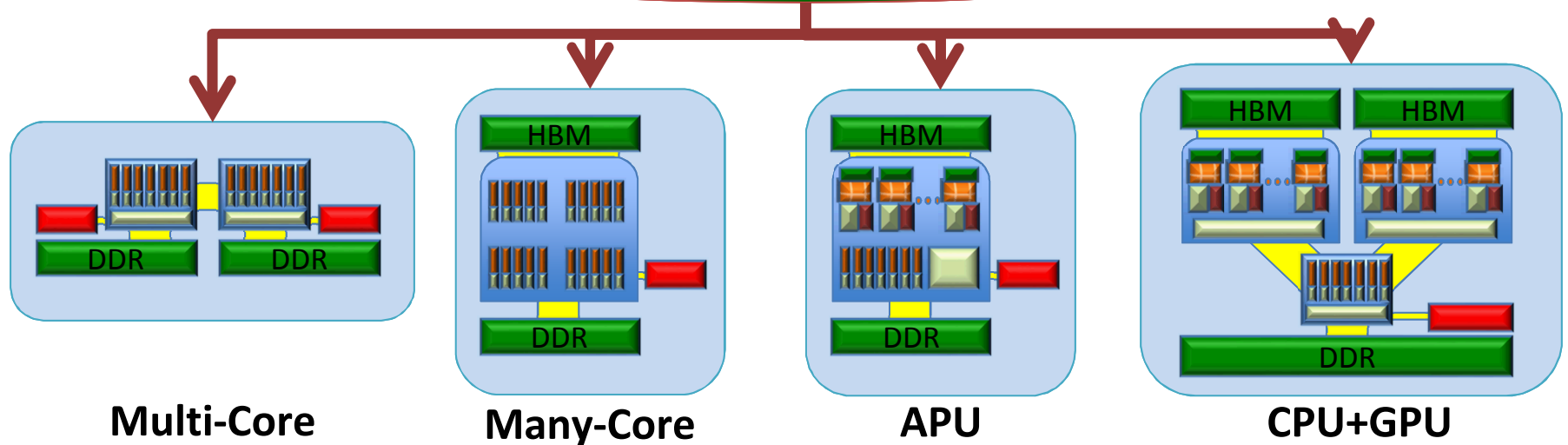
- MPI – Domain decomposition
  - This is the traditional approach
  - Not covered in this talk
- Threading – On node parallelization
  - More scalable for local cores
  - Extensible to GPUs/Cuda
  - Threads team up to accelerate a kernel
- Kernel parallelization
  - Different kernels at the same time
  - Asynchronous Many Tasking (AMT)



# What is Kokkos?



**Kokkos**  
performance portability for C++ applications



Cornerstone for performance portability across next generation HPC architectures at multiple DOE laboratories, and other organizations.



# Abstractions

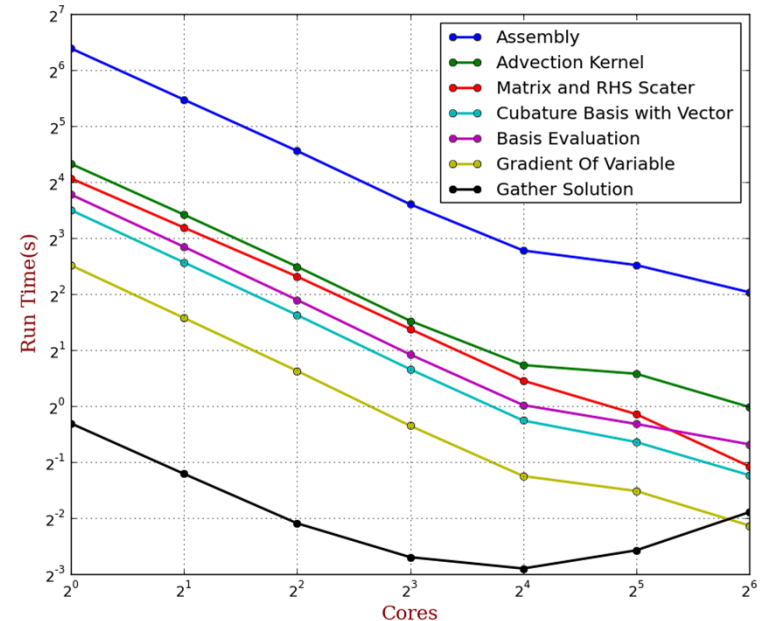
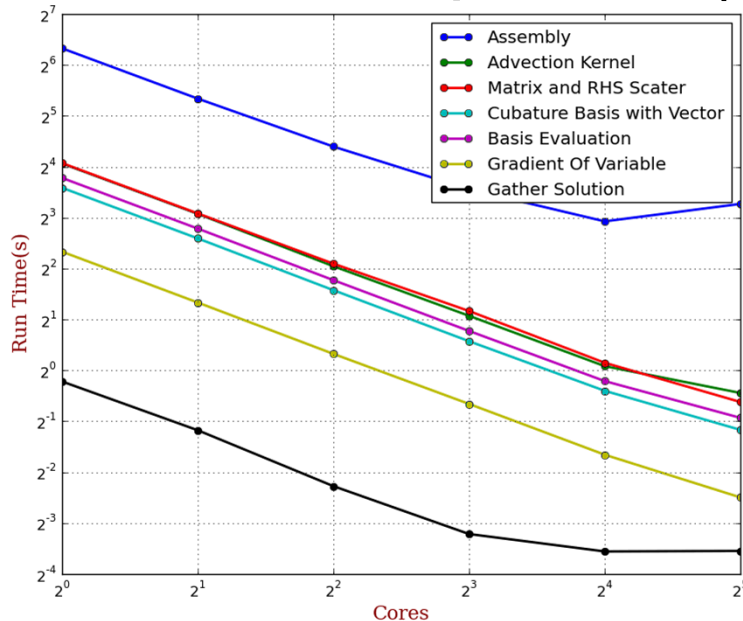
## Patterns, Policies, and Spaces

- Parallel Pattern of user's computations
  - `parallel_for`, `parallel_reduce`, `parallel_scan`, task-graph, ... (*extensible*)
- Execution Policy tells *how* user computation will execute
  - Static scheduling, dynamic scheduling, thread-teams, ... (*extensible*)
- Execution Space tells *where* computations will execute
  - Which cores, numa region, GPU, ... (*extensible*)
- Memory Space tells *where* user data resides
  - Host memory, GPU memory, high bandwidth memory, ... (*extensible*)
- Layout (policy) tells *how* user array data is laid out
  - Row-major, column-major, array-of-struct, struct-of-array ... (*extensible*)
- Differentiating: Layout and Memory Space
  - Versus other programming models (OpenMP, OpenACC, ...)
  - Critical for performance portability ...

# Threading

## Traditional Architectures

- The gather, math routines and scatter operations were all threaded
- MPI only (left) was compared to Kokkos::OpenMP (right) on 16000 elements

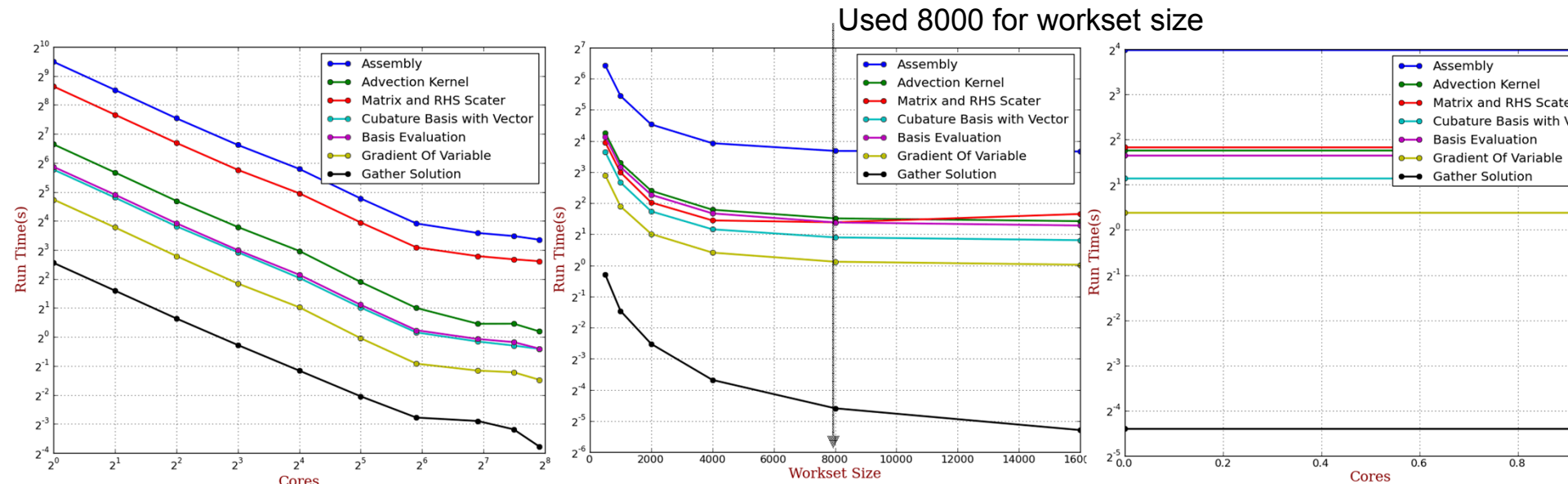


Run on a dual 16 core Haswell with hyperthreading. Left result uses one socket for 1 to 32 cores and MPI across sockets. MPI results show the average time across MPI ranks. 156k total degrees of freedom. Time is for Jacobian assembly only.

# Threading

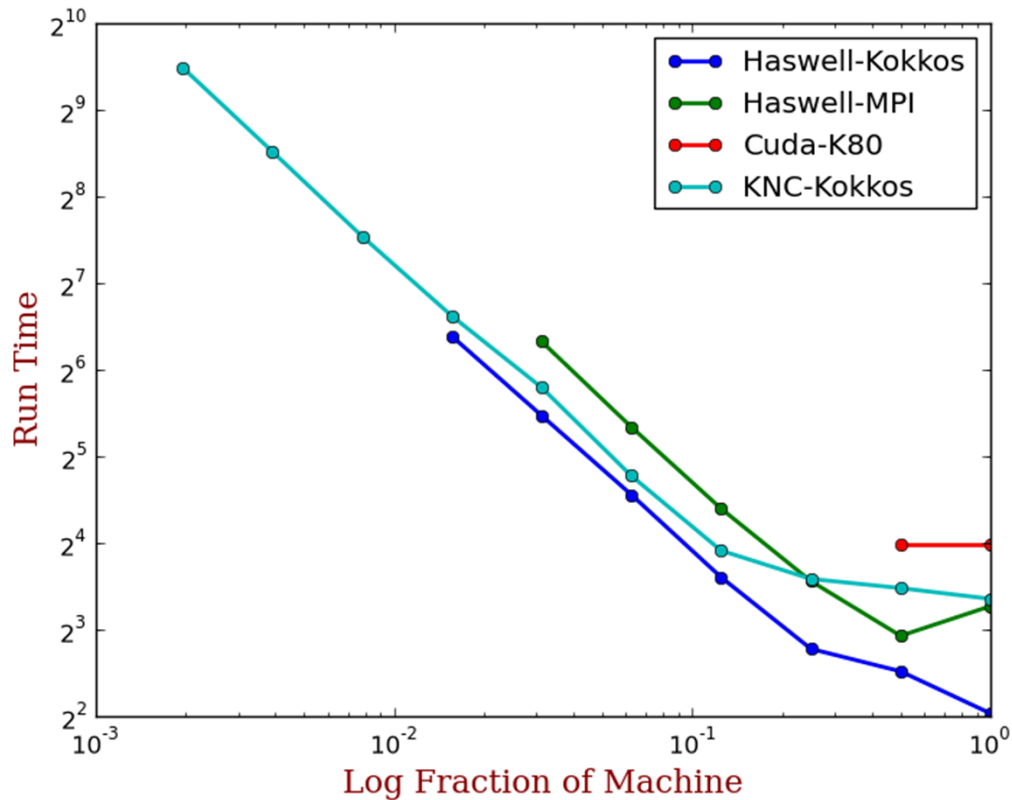
## Modern Architectures

- Kokkos allows one to migrate to new machines and backends
  - Only recompilation is required to move to Intel's Knights Corner (left) or Nvidia K80 via Cuda (right)
  - Larger worksets mean more parallelism (center)



# Summary

## Threading Results





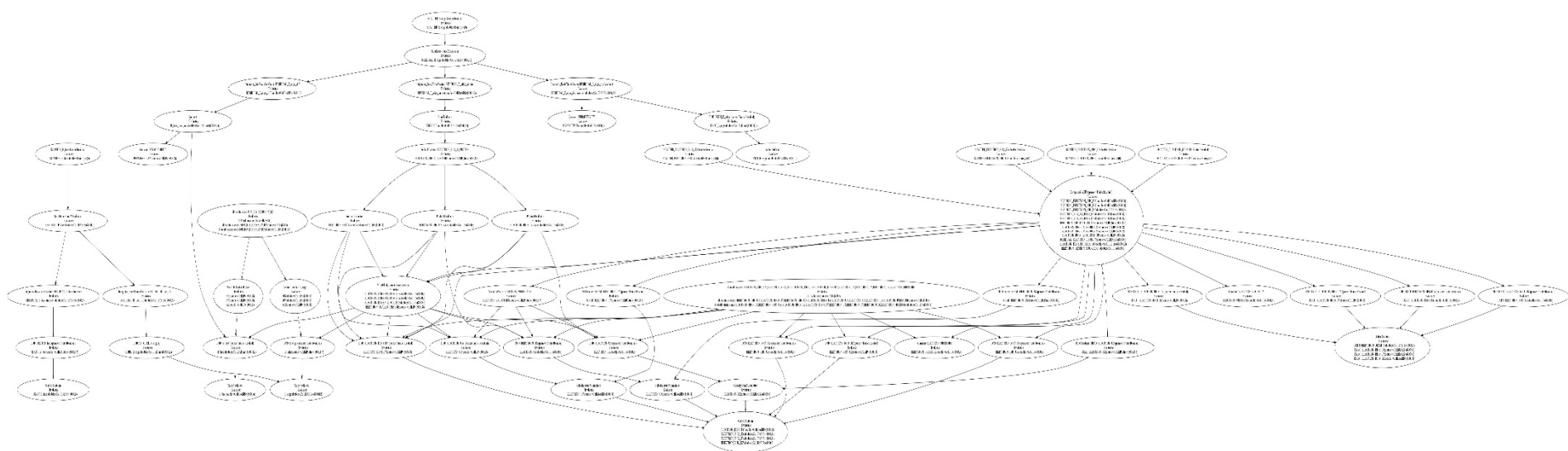
# Beyond Threading

- Threading accelerates each individual math kernel
  - Eventually there is insufficient work in a single kernel to use all the threads efficiently
- Task parallelism
  - Chunks of threads solve different kernels
  - The wider the graph, the more task parallelism
  - Asynchronous Many Task (AMT) is an active research area in scheduling these tasks
    - HPX, DHARMA, Charm++, Legion, ...

# Task Parallelism

## Single Fluid

Real DAG much more complex than the previous toy example



Theoretical speed up for this DAG

	1 Thread/Kernel	8 Threads/Kernel	16 Threads/Kernel
Jacobian	3.5	4.5	4.9
Residual	3.4	3.4	3.5





# Summary

## Next Steps

- Assembly performance can be improved with both threading and task parallelism
  - Threading shows up-to 64x speed up
  - Need to demonstrate these two approaches together
- Cuda parallelism performance lower than expected
  - Not enough parallelism per workset
    - Need bigger workset, however, limited memory
  - Need to implement hierarchical parallelism
- Kokkos makes thread parallelism easier
  - Need to incorporate Kokkos further