

Massively Parallel Monte Carlo Methods for Discrete Linear and Nonlinear Systems

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- Predictive modeling and simulation enhances engineering capability
- Modern work focused on this task leverages multiple physics simulation (CASL, NEAMS)
- New hardware drives algorithm development (petascale and exascale)
- Monte Carlo methods have the potential to provide great improvements that permit finer simulations and better mapping to future hardware
- A set of massively parallel Monte Carlo methods is proposed to advance multiple physics simulation on contemporary and future leadership class machines



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Multiple physics simulations are complicated...

- Neutronics, thermal hydraulics, computational fluid dynamics, structural mechanics, and many other physics
- Consistent models yield nonlinearities in the variables through feedback effects
- Tremendous computational resources are required with $O(1 \times 10^9)$ element meshes and $O(100,000)+$ cores used in today's simulations.

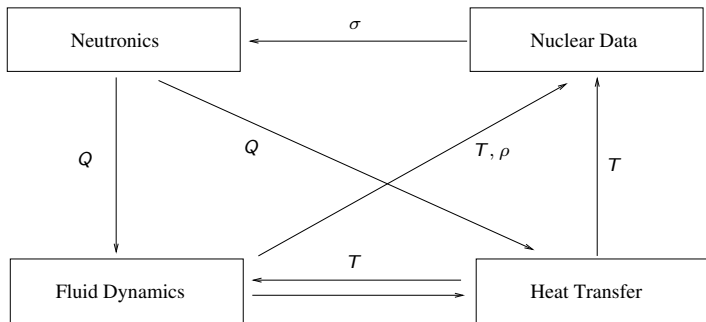


Figure: Multiphysics dependency analysis of departure from nucleate boiling.



- Modern hardware is moving in two directions: lightweight machines and heterogeneous machines characterized by low power and high concurrency.
- High concurrency and low cost units means a higher potential for both soft and hard failures.
- Monte Carlo methods bury soft failures within the variance of the tallies while hard failures are high variance events.
- New machines will also be memory restricted with a continued decrease of memory/FLOPS predicted.
- Compared to conventional methods, we aim show that Monte Carlo methods offer a memory savings.



- Parallelization of classic Monte Carlo methods.
- Parallel strategies taken from modern reactor physics methods in Monte Carlo.
- Research is required to explore how domain decomposition patterns and the discrete system properties are related.
- Research is required to explore how these methods perform in modern multiple physics simulations where strong nonlinearities are present.

We will seek solutions of the general linear problem:

$$\mathbf{Ax} = \mathbf{b} ,$$

$$\mathbf{A} \in \mathbb{R}^{N \times N}, \mathbf{A} : \mathbb{R}^N \rightarrow \mathbb{R}^N, \mathbf{x} \in \mathbb{R}^N, \mathbf{b} \in \mathbb{R}^N .$$

We will assert that \mathbf{A} is *nonsingular*. The solution is then:

$$\mathbf{x} = \mathbf{A}^{-1}\mathbf{b} .$$

We can also define the residual of the system as:

$$\mathbf{r} = \mathbf{b} - \mathbf{Ax} ,$$

where $\mathbf{r} = \mathbf{0}$ when an exact solution is found.

- General stationary methods are formed by splitting the linear operator

$$\mathbf{A} = \mathbf{M} - \mathbf{N} .$$

$$\mathbf{x} = \mathbf{M}^{-1}\mathbf{N}\mathbf{x} + \mathbf{M}^{-1}\mathbf{b} .$$

We identify $\mathbf{H} = \mathbf{M}^{-1}\mathbf{N}$ as the *iteration matrix*

$$\mathbf{x}^{k+1} = \mathbf{H}\mathbf{x}^k + \mathbf{c} .$$

- The qualities of the iteration matrix dictate convergence
- Define $\mathbf{e}^k = \mathbf{x}^k - \mathbf{x}$ as the error at the k^{th} iterate
- The iteration error is generated by a recurrence relation

$$\mathbf{e}^{k+1} = \mathbf{H}\mathbf{e}^k$$

- We diagonalize \mathbf{H} to extract its Eigenvalues

$$\|\mathbf{e}^k\|_2 = \rho(\mathbf{H})^k \|\mathbf{e}^0\|_2 ,$$

- We bound \mathbf{H} by $\rho(\mathbf{H}) < 1$ for convergence

We choose a *search subspace* \mathcal{K} and a *constraint subspace* \mathcal{L} and determine the solution to the linear problem by extracting the solution from the search subspace:

$$\tilde{\mathbf{x}} = \mathbf{x}_0 + \boldsymbol{\delta}, \quad \boldsymbol{\delta} \in \mathcal{K},$$

and by constraining it with the constraint subspace:

$$\langle \tilde{\mathbf{r}}, \mathbf{w} \rangle = 0, \quad \forall \mathbf{w} \in \mathcal{L}.$$

We can generate a more physical and geometric-based understanding of these constraints by writing the new residual as:

$$\tilde{\mathbf{r}} = \mathbf{r}_0 - \mathbf{A}\boldsymbol{\delta}$$

If $\tilde{\mathbf{r}}$ is to be orthogonal to \mathcal{L} , then $\mathbf{A}\boldsymbol{\delta}$ must be the projection of \mathbf{r}_0 onto the subspace \mathcal{L} that eliminates the components of the residual that exist in \mathcal{L} .

The Orthogonality Constraint

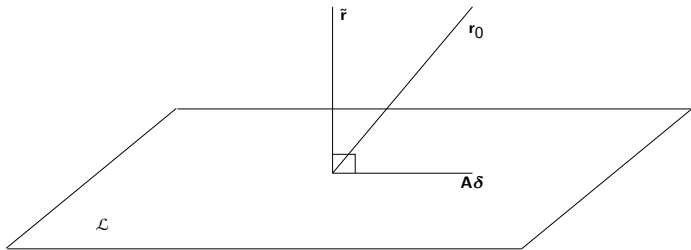


Figure: Orthogonality constraint of the new residual with respect to \mathcal{L} .

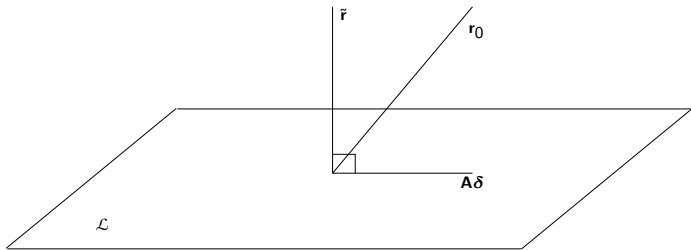


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Minimization Property

The residual of the system will always be *minimized* with respect to the constraints

$$\|\tilde{\mathbf{r}}\|_2 \leq \|\mathbf{r}_0\|_2, \quad \forall \mathbf{r}_0 \in \mathbb{R}^N,$$

Consider a matrix \mathbf{V} to form a basis of \mathcal{K} and a matrix \mathbf{W} to form a basis of \mathcal{L} .

$$\boldsymbol{\delta} = \mathbf{V}\mathbf{y}, \quad \forall \mathbf{y} \in \mathbb{R}^N.$$

From the orthogonality constraint it then follows that:

$$\mathbf{y} = (\mathbf{W}^T \mathbf{A} \mathbf{V})^{-1} \mathbf{W}^T \mathbf{r}_0,$$

We can then form an iteration sequence:

$$\mathbf{r}^k = \mathbf{b} - \mathbf{A}\mathbf{x}^k$$

$$\mathbf{y}^k = (\mathbf{W}^T \mathbf{A} \mathbf{V})^{-1} \mathbf{W}^T \mathbf{r}^k$$

$$\mathbf{x}^{k+1} = \mathbf{x}^k + \mathbf{V}\mathbf{y}^k,$$

with \mathbf{V} and \mathbf{W} updated prior to each iteration.

















Monte Carlo Solution Methods for Nonlinear Problems

















Proposed Challenge Problem



