

Computational Methods for Fusion Science 2022 IAEA Summer School

L. Chacón

May 23-27, 2022

LAUR-21-26161

Agenda

- 1. The quest for harnessing thermonuclear fusion as an energy source on Earth
- 2. Computing thermonuclear fusion: a very challenging endeavor
- 3. Computational methods for Magnetic Confinement Fusion
- 4. Computational methods for Inertial Confinement Fusion
- 5. A few examples of future research directions in Fusion Computation
- 6. Summary

Disclaimer: while I will attempt to provide as broad and as general a perspective on computing in Fusion Science as possible, my exposition will draw heavily from my own research and will reflect my personal perspective. Caveat emptor!

The quest for harnessing thermonuclear fusion as an energy source on Earth

What is thermonuclear fusion?

- " Thermonuclear Fusion" is the process whereby, upon reaching sufficiently high temperatures and densities, lighter nuclei combine to form heavier ones, converting a tiny amount of mass into a lot of energy (according to Einstein's mc2 formula)
- Fusion is the engine that drives the birth, life, and death of stars in the universe, and therefore of life on Earth
	- − In stars, matter is perfectly confined by gravity, allowing for long-term fusion energy production
	- − In the laboratory, however, one needs to figure out ways to confine matter at billions of degrees hot… this is the key challenge of harnessing fusion on Earth!

Why is thermonuclear fusion for peaceful energy production attractive?

- Fuel is hydrogen isotopes: inexhaustible, geographically distributed
- Small quantity of fuel: quick shutdown
- Inherent safety aspects
- No harmful radioactive or $CO₂$ emissions (no global warming)
- Byproduct neutrons may activate structural materials, but nextgeneration fusion reactors will be "neutron-free"

Only 100 kg deuterium (corresponding to 2800 tonnes of sea water) and 150 kg of tritium (corresponding to 10 tonnes of lithium ore) will be needed for operating a 1 GW electric power plant for one year.

What are the required conditions for thermonuclear fusion?

- One hundred million degrees!
	- − Enough kinetic energy to overcome electrostatic repulsion
- At such temperatures, matter is ionized and forms a **plasma** (4th) state of matter, most common!)
- Plasma is so hot it will instantly melt any surface: cannot meet any reactor vessel! **Confinement.**

Magnetic confinement

Θ

The toroidal geometry avoids end losses!

How to confine matter that hot? II. By compression: Inertial confinement

A taxonomy of plasmas in nature and the laboratory

Simulating thermonuclear fusion: a very challenging endeavor

The role of computing in thermonuclear fusion: Virtual experiments

- Fusion experiments are getting bigger and more expensive: does not give much room for iteration
- A successful fusion reactor should confine extremely hot matter at a sufficient density for sufficiently long: there is not much room for error!
- "Virtual experiments" are needed to find suitable operating regimes, and to guide construction, operation and optimization of future fusion reactors.
- Such "virtual experiments" are, or course, simulations. We require a predictive capability! However...

Plasma confinement (or lack thereof) results from careful interplay of physical phenomena spanning many orders of magnitude in time and space!

Challenges in thermonuclear fusion simulation: "The tyranny of scales"

Challenges in thermonuclear fusion simulation: Charge separation characteristic scales

- Plasma frequency (very fast): time scale of restoration of charge imbalance in the plasma
- Debye length (very small): length scale beyond which charge separation cannot be sustained

$$
\omega_p=\sqrt{\frac{q^2n}{\epsilon_0m}}
$$

System tends to restore neutrality but overshoots, oscillating at plasma frequency

characteristic scale the Debye length Electrons "shield" ion charge with

Challenges in thermonuclear fusion simulation: Magnetic field characteristic scales

- Gyrofrequency (very fast): time scale of particle gyration around magnetic field
- Gyroradius (very small): radius of gyration around magnetic field

Challenges in thermonuclear fusion simulation: Collision frequencies and mean-free-paths

• Collision frequency:

$$
\nu_c \propto \frac{q^{\text{-}} n}{m^{1/2} T^{3/2}}
$$

 $\overline{4}$

- Collisional mean-free-path: $\lambda_c = v_{th}/v_c$
- Collisions increase with density, decrease with temperature.
	- − Hot, low-density plasmas are weakly collisional (MFE)
	- − Warm, dense plasmas will be moderately or strongly collisional (ICF)
- Electrons are more collisional than ions by $\sqrt{m_i/m_e}$
- Collisions determine relaxation rates toward local thermal equilibrium (LTE)
- Collisionality affects momentum and energy transport, and therefore confinement.

Fusion sciences pioneered unclassified supercom

- In 1973, Dr. Alvin Trivelpiece, deputy director of the Controlled Thermonu Research (CTR) program of the Atomic Energy Commission, solicited proposals for a computing center that would aid in reaching fusion pow giving the magnetic fusion program under CTR access to computing power similar to that of the defense programs.
- CTR computing center (CTRCC) was first placed at LLNL, and in 1996 to LBL.
- CTRCC was soon renamed "National Magnetic Fusion Energy Comput Center," and in 1983 adopted its final name of National Energy Resear Supercomputing Center (NERSC) when it was open to al DOE-SC disc

https://www.nersc.gov/about/nersc-history/?start=1

 ∞

The power of algorithms! r or arg

Credit: S. Jardin, SCaLeS Report (2003)

The first-principles plasma description: Liouville's equation (an intractable problem)

• Goal: to describe particle distribution function (PDF) of N interacting particles with position **qi** and momenta **pi**:

$$
f_N = f_N(\mathbf{q}_1 \ldots \mathbf{q}_N, \mathbf{p}_1 \ldots \mathbf{p}_N, t)
$$

• PDF is governed by the (6N+1) dimensionality equation:

$$
\frac{\partial f_N}{\partial t} + \sum_{i=1}^N \frac{\mathbf{p}_i}{m} \frac{\partial f_N}{\partial \mathbf{q}_i} + \sum_{i=1}^N \mathbf{F}_i \frac{\partial f_N}{\partial \mathbf{p}_i} = 0,
$$

• Note: N is the **real** number of physical particles!

Model reduction to the tractable: BBGKY hierarchy and the Boltzmann equation

• Define s marginal PDF by integrating over $(s+1,...N)$ phase space:

$$
f_s(\mathbf{q}_1 \dots \mathbf{q}_s, \mathbf{p}_1 \dots \mathbf{p}_s, t) = \int f_N(\mathbf{q}_1 \dots \mathbf{q}_N, \mathbf{p}_1 \dots \mathbf{p}_N, t) d\mathbf{q}_{s+1} \dots d\mathbf{q}_N d\mathbf{p}_{s+1} \dots d\mathbf{p}_N
$$

• Leads to reduced equation:

$$
\frac{\partial f_s}{\partial t} + \sum_{i=1}^s \frac{\mathbf{p}_i}{m} \frac{\partial f_s}{\partial \mathbf{q}_i} - \sum_{i=1}^s \left(\sum_{j=1\neq i}^s \frac{\partial \Phi_{ij}}{\partial \mathbf{q}_i} + \frac{\partial \Phi_i^{ext}}{\partial \mathbf{q}_i} \right) \frac{\partial f_s}{\partial \mathbf{p}_i} = (N-s) \sum_{i=1}^s \int \frac{\partial \Phi_{i\,s+1}}{\partial \mathbf{q}_i} \frac{\partial f_{s+1}}{\partial \mathbf{p}_i} d\mathbf{q}_{s+1} d\mathbf{p}_{s+1}.
$$

- "Closure problem": f_s depends on f_{s+1}
- Solution: close equation at s=1, and model rhs with collisions: **Boltzmann eq.**

$$
\frac{\partial f}{\partial t} + \frac{\mathbf{p}}{m} \frac{\partial f}{\partial \mathbf{q}} + \frac{\mathbf{F}}{m} \frac{\partial f}{\partial \mathbf{p}} = \left(\frac{\partial f}{\partial t}\right)_c
$$

- Still high dimensional (6D+time)!
- 7/12/21 20 • For grazing Coulomb collisions, Boltzmann reduces to the **Vlasov-Fokker-Planck-Landau equation**: basis for all fusion modeling!

A "tractable" first-principles model: **The Vlasov-Fokker-Planck equation (+ Maxwell Eqs)**

 $\frac{E}{\beta}$ ^{*i*} ; *j*^{*i*} : *j*^{*B*} ; *j*[*]*

p

jpS(*xi xp*)

High dimensionality (3D+3V), exceedingly multiscale *^d*(**^x ^x***p*) ! *^S*(**^x ^x***p*) ; *Ep* = Â

The power of asymptotics: nested model hierarchies

- Asymptotic model reduction from the VFP equation is possible by taking advantage of time/length scale separation
- In MFE, key asymptotic parameters are:
	- $-$ Plasma frequency ω_{p} (quasineutrality, ambipolarity)
	- \sim Gyrofrequency Ω_c , the gyroradius ρ_c (magnetic field strength)
	- − Collision frequency v_c and mean-free-path λ_c (plasma collisionality)
	- − Ion/electron skin depths, $d_i e^{t} = c/\omega_{\text{size}}$ (measure scale lengths where kinetic effects are important)
- In ICF, key asymptotic parameters are:
	- − Plasma frequency
	- − Plasma collisionality and mean-free-path
	- − Plasma b (ratio of thermal-to-magnetic pressures)

Model hierarchy in MFE

Model hierarchy in MFE T_{max} space and time range scales s

Credit: Grandgirard

A taxonomy of computational methods for fusion science

Spatial discretization approaches *ing major radii at 90% flux sur*ratised above a threshold a threshold a threshold in the radial electric field *E^r* forms within the pedestal layer, bringing with it a

$$
\Phi^{n+1} = \Phi^n + \Delta t D \Delta \Phi^n
$$

Easy update (no solve), conditional stability on Δt

 \bigotimes

 $\Phi^{n+1} = \Phi^n + \Delta t D \Delta \Phi^{n+1}$

unconditionally stable in Δt ^{12/21} 28 Requires global algebraic solve (hard),

What are we looking for in a computational algorithm?

• Optimal algorithm (def): $\big|$ $\mathsf{CPU}\sim\mathsf{N}/\mathsf{R}$ ⇠ */*

$$
\mathcal{CPU} \sim \frac{N^{1+\alpha}}{n_p^{1-\beta}} \, ; \, N = \left(\frac{L}{\delta}\right)^d \, \begin{cases} \alpha \geq 0, \text{ algorithmic scalability} \\ \beta \geq 0, \text{ parallel scalability} \end{cases}
$$

- Algorithmic scalability: $\alpha \approx 0$
- Parallel scalability: $\beta \approx 0$
- Much emphasis is placed on parallel scalability, but **algorithmic scalability is also critical**: j.

 μ (weak scaling) $CDU \rightarrow p^0$ $\sim n_p$ (weak scaling), $\mathcal{CP}U\sim n_p^{\alpha+\beta}$

• Only iterative implicit algorithms may become algorithmically (and p $\sum_{\text{National Laboral L.}}$ scalable: α .

Explicit | Implicit (direct) | Implicit (iterative) | Implicit (multilevel) $\alpha = 1/d$ $\alpha = 2 - 2/d$ $\alpha > 1$ (varies) $\alpha \approx 0$

National Laborat

Computational methods for Magnetic Confinement Fusion

What is magnetic confinement fusion?

- Magnetic confinement fusion (also known as Magnetic Fusion Energy, MFE) attempts to "bottle" million-degree hot plasma using magnetic fields.
- Leading concepts are based on toroidal geometries (no end losses)
	- − Best known concept is **tokamak**, in which the plasma generates confining poloidal magnetic fields self-consistently with plasma currents.
	- − Other fusion-grade concepts include the **stellarator**, which creates confining magnetic fields with external coils.

Magnetohydrodynamics (MHD): equilibrium and stability

- MHD describes well macroscopic (bulk) plasma behavior.
- A self-sustaining long-pulse (quasi-steady-state) MFE reactor must be:
	- − In MHD equilibrium: **j**x**B**≈grad(P)
	- − MHD stable (perturbations to the equilibrium must decay, not grow exponentially)
- MHD toroidal equilibrium can be 2D or 3D, and must satisfy **j**x**B**=grad(P)
	- − 2D equilibria are computed using the Grad-Shafranov equation [reduction of **j**x**B**=g(P)]
	- − 3D equilibria requires solving full MHD equilibrium equation (e.g., in stellarator)
- MHD stability is critical for long-term reactor operation
	- − Stellarator concept does not self-generate magnetic fields, and is MHD-robust
	- − Tokamak MHD stability is more nuanced. If plasma becomes unstable in a tokamak, the plasma may terminate, causing a DISRUPTION. This should be avoided at all costs.

The MHD model

$$
\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \vec{v}) = 0,
$$
\n
$$
\frac{\partial \vec{B}}{\partial t} + \nabla \times \vec{E} = 0,
$$
\n
$$
\frac{\partial (\rho \vec{v})}{\partial t} + \left[\nabla \cdot \left[\rho \vec{v} \vec{v} - \vec{B} \vec{B} + \vec{\Pi} + \vec{\tau} \left(p + \frac{B^2}{2} \right) \right] = 0,
$$
\n
$$
\frac{\partial p_e}{\partial t} + \nabla \cdot (\vec{v} p_e) + (\gamma - 1) p_e \nabla \cdot \vec{v} = (\gamma - 1)(S - \nabla \cdot \vec{q}).
$$
\n
$$
\vec{v} = \frac{m_i \vec{v}_i + m_e \vec{v}_e}{m_i + m_e} \approx \vec{v}_i \; ; \; \vec{v}_e = \vec{v}_i - d_i \vec{\dot{f}}
$$
\n
$$
\text{Ohm's Law}: \vec{E} = -\vec{v} \times \vec{B} + \eta \vec{j} - \frac{d_i}{\rho} (\vec{j} \times \vec{B} - \nabla p_e - \nabla \cdot \vec{\Pi} \cdot e) - \frac{d_e^2 d\vec{v}_e}{d_i \; dt}
$$

MHD equilibrium [jxB=grad(P)]: Grad-Shafranov equation (F)]. Grau-Shairanov equation

- Postulate B-field as: ϵ
- Find poloidal flux $\Psi(x)$ for a given toroidal magnetic field F=R B_0 =F(Y) and pressure profile p(Y) such that:

$$
\mathbf{j} \times \mathbf{B} = \nabla p \Rightarrow \nabla \cdot \left(\frac{\nabla \Psi}{R^2}\right) = -\frac{dp}{d\Psi} - \frac{F}{R^2}\frac{dF}{d\Psi}
$$

- Highly nonlinear equation! Can be very difficult to solve.
	- − Discretized using FV, FD, FE, spectral methods.
	- − Requires nonlinear iteration
	- − Input functions p(Y), F(Y) can be eliminated by adding more physics (for instance, loop voltage and resistive decay), or can be provided in alternative forms
	- − Many available codes: EFIT, TEQ, CORSICA, CHEASE
- Codes that solve 3D MHD equilibria (stellarator) also exist: PIES (3D), VMEC (3D)

MHD stability: Methods

- After an equilibrium is found, it is important to determine whether it is MHD stable or unstable
- The question of stability is a tiered one: ideal stability (without dissipation), resistive stability, two-fluid stability, kinetic stability….
- There are many specialized tools to determine MHD stability, including some beyond-MHD effects:
	- − DCON (extended Newcomb's criterion)
	- − PEST (MHD energy principle)
	- − MARS (spectral)
	- − ELITE (edge localized modes)
	- − …
- Initial-value MHD computations by full-fledged MHD codes are also used to study MHD stability

MHD initial-value simulations of MFE: Methods

- MHD is a hyperbolic PDE system, supporting a variety of fast waves (fast and slow magnetosonic, shear Alfven).
- Spatial discretizations: FV, FD, FE, spectral,…
- Temporal discretizations: semi-implicit, fully implicit
	- − MFE benefits from quiescent plasmas, and therefore MHD simulations may need to cover a very long-time span
	- − Resolving fast timescales is impractical: implicit timestepping (Δt ω_{MHD} >> 1)
	- − Many MHD codes are available for MFE (NIMROD, M3D-C1, PIXIE3D, JOREK, SpeCyl, HiFi,…). All of them feature some level of time-implicitness.
	- − Key algorithmic requirement: SCALABILITY
- Achieving algorithmic and parallel scalability in implicit MHD codes is difficult
	- − Algorithmic scalability: CPU ~ O(N log(N)), N: number of degrees of freedom
	- − Parallel scalability: CPU ~ 1/N_p, N_p: number of processors
	- − We need **both**!
MHD: Impact of algorithms

- Impact both from spatial and temporal discretization improvements
	- − High-order, mesh adaptivity, etc.
- Time-implicitness is key
- Suitable linear and nonlinear solvers to invert associated algebraic system of equations is also important for scalability
- MHD algorithms remain an active area of research

Magnetic Fusion Energy: "Effective speed" increases
came from both faster hardware and improved algorithms

Credit: S. Jardin, SCaLeS Report (2003) $_{\rm 7/12/21}$ nnstanding 137

MHD: Implicit timestepping algorithms Why are they so difficult to scale up with problem size?

• Implicit timestepping requires an algebraic (often nonlinear) solve:

$$
G(U^{n+1}) = U^{n+1} - U^n - \Delta t F(U^{n+1}) = 0 \; ; \; U = [\rho, \mathbf{v}, p]
$$

• G(U) is generally nonlinear, and requires iteration, e.g. Newton-Raphson

$$
\left.\frac{\partial G}{\partial U}\right|_k \delta U_k = - G(U_k) \;\; ; \;\; U_{k+1} = U_k + \delta U_k
$$

- Jacobian matrix $J_k = \partial G / \partial U_k$ is a very large, sparse, ill-conditioned matrix
	- − Direct methods are prohibitive [e.g., Gaussian elimination, CPU~ O(N7/3) in 3D!]
	- − Requires iterative methods, but typically # iterations grows with the condition number of the matrix: not scalable!
- Solution: Multigrid-preconditioned Krylov methods

Krylov methods: a primer

- Krylov methods attempt to find the solution of $Ax=b$ as a series: $\mathbf{x} = \sum a_i \mathbf{d}_i$
- Here, **d**_i are A-conjugate vectors, satisfying orthogonality property:

$$
\mathbf{d}_iA\mathbf{d}_j=\delta_{ij}\Rightarrow a_i=\mathbf{d}_i^T\mathbf{b}
$$

- Problem solved! Right? No... one needs to find conjugate vectors!
- Krylov methods build conjugate basis iteratively, and orthogonalize along the way (e.g., Gram-Schmidt, QR factorization, etc.):

$$
\{\mathbf{r}, A\mathbf{r}, A^2\mathbf{r}, ...\} \to \{\mathbf{d}_1, \mathbf{d}_2, \mathbf{d}_3, ...\}
$$

- − All that is required to form basis is to multiply matrix A times a (given) vector once per iteration!
- *•* iteration!
 • For Jacobian system (Newton), matrix-vector product can be performed without ever building and storing the Jacobian matrix! (Gateaux derivative) *k* \mathbf{r} *n* be performed witl

$$
\left(\frac{\partial \vec{G}}{\partial \vec{x}}\right)_k \vec{y} = J_k \vec{y} = \lim_{\epsilon \to 0} \frac{\vec{G}(\vec{x}_k + \epsilon \vec{y}) - \vec{G}(\vec{x}_k)}{\epsilon}
$$

Jacobian-free! No need to form and store Jacobian matrix.

Preconditioning Krylov methods

- Krylov methods are much faster than other iterative methods (e.g., Jacobi, Gauss-Seidel), but still not optimal (work scales as $N^{1+\alpha}$, α >1).
- **But they can be PRECONDITIONED so that work scales as ~ N!**
- Preconditioning: rewrite linear system as:

 $AP^{-1}Px = b \Rightarrow (AP^{-1})y = b$; $x = P^{-1}y$

 P^{-1} is the preconditioner. If $P^{-1} \sim A^{-1}$, then $(A P^{-1}) \sim I$, very fast convergence!

- Matrix-vector multiplication feature of Krylov methods allow seamless implementation of preconditioner:
	- − z=(A P-1)v can be computed with 2 matrix-vector products: y= P-1v, z = Ay
- Bleeding-edge research in iterative methods is in the development of effective preconditioners
	- − Preconditioner only affects convergence, not the solution
	- − Approximations to PDE that would lead to bad solvers can be good preconditioners!
	- − Application dependent

Multigrid methods: key for algorithmic scalability (work~N)

- MG employs a divide-and-conquer approach to attack error components in the solution Preconditioning: multilevel (multigrid) methods *•* MG employs a divide-and-conquer approach to attack error components in the solution.
	- − Oscillatory components of the error are "EASY" to deal with (if a SMOOTHER exists) Oscillatory components of the error are "EASY" to deal with (if a SMOOTHER exists)
	- − Smooth components are DIFFICULT Smooth components are DIFFICULT.
- MG idea: coarsen recursively to make "smooth" components become oscillatory

- SMOOTHER is KEY component of MG
	- − Smoother is "easy" to find for parabolic and elliptic problems, but hard for hyperbolic ones • **•** • Smoother is "coopy" to find for nore *•* Smoothers for stiff hyperbolic equations are hard to formulate.
- MHD is hyperbolic: PARABOLIZATION of implicit MHD

Parabolization of stiff hyperbolic systems: Physics-based multigrid preconditioning Parabolization of stiff hyperbolic sy! Parabolization and block factorization (Schur complement) Parabolization and block factorization (Schur complement) **Physics-based multigrid preconditioning •** Parabolization is a natural consequence of implicit timestep in the sequence of implicit timestep

- Parabolization enables development of effective preconditioners for stiff hyperbolic PDEs *a*
∂∂*t*ioners for sti donore for our
- DES
- Parabolization exploits structure of implicit discretization

$$
\partial_t u = \frac{1}{\epsilon} \partial_x v \ , \ \partial_t v = \frac{1}{\epsilon} \partial_x u.
$$
\n
$$
u^{n+1} = u^n + \frac{\Delta t}{\epsilon} \partial_x v^{n+1}, v^{n+1} = v^n + \frac{\Delta t}{\epsilon} \partial_x u^{n+1}.
$$
\n
$$
u^{n+1} = u^n + \frac{\Delta t}{\epsilon} \partial_x v^n
$$

- ⁻ Parabolized systems are suitable for modern multilevel solvers (multigrid), which can be optimal $\text{[CPU} \sim O(\text{N} \text{ log}(\text{N}))$ **banable for modern mannover corvero** (mannyma*r, whi*len ca
| $f(x \text{ modern multiplication of columns } (multi) \text{ which can be$ ² *L I* #
- Connection between parabolization and block-factorization (Schur complement): *I e z*(שוα *uog(ג)*
µen parabolizatio *e* α ^{*n*} and blocl = " *I UD*¹ 2 # " *^D*¹ *UD*¹ ² *L* 0 # " *^I* ⁰ ² *L I* # *•* PARABOLIZATION via BLOCK FACTORIZATION (Schur complement): *L D*² 0 *I* 0 *D*² *D*¹

$$
\left[\begin{array}{cc} D_1 & U \\ L & D_2 \end{array}\right] = \left[\begin{array}{cc} I & U D_2^{-1} \\ 0 & I \end{array}\right] \left[\begin{array}{cc} D_1 - U D_2^{-1} L & 0 \\ 0 & D_2 \end{array}\right] \left[\begin{array}{cc} I & 0 \\ D_2^{-1} L & I \end{array}\right] \qquad \qquad D_1 - U D_2^{-1} L = \left[I - \left(\frac{\Delta t}{\epsilon}\right)^2 \partial_{xx}\right]
$$

• Provides path for generalization of parabolization strategy to complex stiff hyperbolic PDEs. " *^D*¹ *^U* **d**
<u>d</u>**1000** µzation or

Jacobian-free Newton-Krylov + Parabolization-MG delivers near-optimal scalability (parallel and algorithmic)

- Hall MHD example using GEM challenge problem (magnetic reconnection)
- Fixed implicit timestep, weak-scalability study

Application: double tearing mode in ITER (PIXIE3D)

Tokamak Disruption Simulation SciDAC Project (PI: X. Tang, LANL)

Beyond MHD: GK and micro-turbulence in tokamaks

- For MHD-stable magnetic-field tokamak configurations, small-scale (kinetic) instabilities develop that lead to micro-turbulence (both electrostatic and electromagnetic)
- Micro-turbulence has huge impact on particle and energy confinement in tokamaks
- Requires a kinetic description: Gyrokinetics
	- − Enforces quasineutrality
	- − Exploits that gyrofrequency is very fast, and gyrophase angle ignorable
	- − Does NOT assume gyroradius is too small (turbulence can in fact develop spatial scales comparable to ion and electron gyroradii)
- Gyrokinetics is one of the most successful asymptotic models in MFE
	- − Spatial discretization: both particle-in-cell and Eulerian.
	- − Temporal discretization: semi-implicit.
	- Fully implicit methods are being actively developed.

$$
\omega_{turb} \sim \omega_{*i} \sim (k_{\theta}\rho_i)\frac{v_{th}}{L_p} \sim 10^5 s^{-1} \ll \omega_{ci} = \frac{eB}{m_i} \sim 10^8 s^{-1}
$$

GK: an algorithmic revolution

- Goal: analytically remove fast plasma and gyro-frequencies (asymptotic model)
- Results in dimensionality reduction (5D instead of 6D)
- "delta-f" representations focus the numerical representation on deviations from Maxwellian (local thermal equilibrium)

Figure from SCaLeS report, Volume 2 (2004), c/o S. Jardin

GK equations Electromagnetic Gyrokinetics *f*^{*x*} *d*₃

• Transport equation 1 0 R³ $\overline{}$ II al t equation

$$
\frac{\partial f_s}{\partial t} + \dot{\mathbf{X}} \cdot \frac{\partial f_s}{\partial \mathbf{X}} + \dot{v}_{\parallel} \frac{\partial f_s}{\partial v_{\parallel}} = 0 \t f_s(\mathbf{X}, \mu, v_{\parallel}, t) : \mathbb{R}^5 \times \mathbb{R} \to \mathbb{R}^+
$$

$$
\dot{\mathbf{X}} = \frac{1}{D} \left[v_{\parallel} \left(\hat{\mathbf{b}}_0 + \delta \hat{\mathbf{b}} \right) + \frac{m_s}{q_s B_0} v_{\parallel}^2 \nabla \times \hat{\mathbf{b}}_0 - \frac{m_s}{q_s B_0} \hat{\mathbf{b}}_0 \times \left(\frac{q_s}{m_s} \langle \mathbf{E} \rangle - \mu \nabla B_0 \right) \right]
$$

$$
\dot{v}_{\parallel} = \frac{q_s}{m_s D} \left[\left(\hat{\mathbf{b}}_0 + \delta \hat{\mathbf{b}} \right) + \frac{m_s}{q_s B_0} v_{\parallel} \nabla \times \hat{\mathbf{b}}_0 \right] \cdot \left(\langle \mathbf{E} \rangle - \frac{m_s}{q_s} \mu \nabla B_0 \right)
$$

$$
D = 1 + \frac{m_s}{q_s B_0} v_{\parallel} \mathbf{b}_0 \cdot \nabla \times \mathbf{b}_0
$$

• Field equations

• Field equations
\n
$$
\bar{n}_i(\mathbf{x}) = \int_{-\infty}^{\infty} \int_0^{\infty} \langle f_i(\mathbf{X}, v_{\parallel}, \mu) \delta(\mathbf{X} - \mathbf{x} + \rho_i) \rangle B d\mu dv_{\parallel}
$$
\n
$$
-\frac{e n_0 m_i}{q_i B_0^2} \nabla_{\perp}^2 \phi = q_i \bar{n}_i - en_e
$$
\n
$$
n_e(\mathbf{x}) = \int_{-\infty}^{\infty} \int_0^{\infty} \langle f_e(\mathbf{X}, v_{\parallel}, \mu) \delta(\mathbf{X} - \mathbf{x}) \rangle B d\mu dv_{\parallel}
$$
\n
$$
-\frac{1}{\mu_0} \nabla_{\perp}^2 A_{\parallel} = j_{\parallel i} + j_{\parallel e}
$$
\n
$$
j_{\parallel s}(\mathbf{x}) = q_s \int_{-\infty}^{\infty} \int_0^{\infty} v_{\parallel} \langle f_s(\mathbf{X}, v_{\parallel}, \mu) \delta(\mathbf{X} - \mathbf{x}) \rangle B d\mu dv_{\parallel}
$$

The gyro-kinetic model: methods

- Spatial discretizations (configuration space): FD, FE, spectral…
- Velocity-space discretizations:
	- − Mesh
	- − Particles (PIC)
- Temporal discretization: semi-implicit
	- − Accuracy issues (cancellation problem)
	- − Ongoing research on fully implicit methods to resolve
- GK codes can effectively use the largest HPC computers on Earth!

A GK turbulence simulation of the tokamak edge (XGC)

High-fidelity Boundary Plasma Simulation SciDAC Project (PI: CS Chang, PPPL)

Computational methods for Inertial Confinement Fusion

What is Inertial Confinement Fusion?

News flash: the National Ignition Facility (NIF) ignited a fusion-grade ICF capsule in Aug. 2021!

- Laser input: 1.9 MJ
- Neutron yield: a few x 10^{17} (ignition threshold at $3x10^{17}$)
- Energy absorbed by capsule: ~250 KJ
- Fusion energy produced: 1.3 MJ (5x energy absorbed!)
- At threshold of engineering "ignition".

A VERY BIG DEAL!!!!!!

• However... it was not expected! Simulations remain critical for predictability!

Radiation hydrodynamics ("rad-hydro"): the workhorse model in ICF

• Simplest form: Euler + radiation transport (gray, multigroup)

$$
\frac{\partial}{\partial t}\rho + \nabla \cdot (\rho \mathbf{v}) = 0
$$
\n
$$
\frac{\partial}{\partial t}\rho \mathbf{v} + \nabla \cdot (\rho \mathbf{v} \otimes \mathbf{v}) + \nabla p = + \rho (\mathbf{a}_{ext} + \mathbf{a}_{rad})
$$
\n
$$
\frac{\partial}{\partial t}e + \nabla \cdot [(e + p) \mathbf{v}] = + \rho \mathbf{v} \cdot (\mathbf{a}_{ext} + \mathbf{a}_{rad})
$$
\n
$$
- \kappa_{\text{P}} \rho c (a_{\text{R}} T^4 - E)
$$
\n
$$
\frac{\partial}{\partial t}E + \nabla \cdot \mathbf{F} = \kappa_{\text{P}} \rho c (a_{\text{R}} T^4 - E).
$$
\n(4)

Radiation hydrodynamics ("rad-hydro"): assumptions and limitations

- Quasineutrality
- Electrostatic limit (magnetic fields ordered out due to large plasma β)
- High collisionality (local thermal equilibrium)
- Single fluid
- Cannot account for multiple species, or deviations from thermal equilibrium

Radiation hydrodynamics ("rad-hydro"): methods

- Spatial discretization:
	- − Hydro: Eulerian or Lagrangian (moving mesh)
	- − Radiation: Monte Carlo (particles), discrete ordinates (Sn). Typically considers multiple energy groups (i.e., photon frequencies)
- Temporal discretization:
	- − Hydro: explicit
	- − Radiation (stiff): fully implicit, semi-implicit (some parts implicit, others explicit).
		- Much recent work in multiscale methods (so-called high-order/low-order, HOLO), where moment descriptions ("radiation-diffusion" models) are used to accelerate kinetic solvers.

Beyond rad-hydro: evidence that more fidelity is needed

Post-simulation analysis indicate weakly collisional regimes are present

normalized ion mean-free-paths (λii 〈Z〉

4

). Symbols indicate time of

Beyond rad-hydro: evidence that more fidelity is needed Deyong rag-nygro. et performance, with little impact from hydrodynamic insta-

Simulations overpredict compression and yield performance approaches 10 because of the full single-

Beyond rad-hydro: evidence that more fidelity is needed

Evidence of importance of electromagnetic effects

[1] Li, C. K., et al. "Study Plasma Stagnation in Laser-Driven Hohlraums." *Bulletin of the American Physical Society* 62 (2017).

Beyond rad-hydro: what will it take?

- Presently, 3D-3V VFP+Maxwell solvers are out of reach
- Focus on 1D-2V geometries (planar, spherical symmetry)
- Consider suitable asymptotic limits for Maxwell equations:
	- − Electrostatic approximation (exact in 1D spherical, $\beta \approx 10^{3}$ -10⁴ in Omega)
	- σ Quasineutrality: $\rho = \sum q_i n_i = 0$
	- $-$ Ambipolarity: $j = \sum q_i n_i v_i = 0$ (in 1D)
	- − Eliminates fastest time scales (plasma frequency) and smallest length scales (Debye length)
- Consider fluid electrons:
	- − Rigorous fluid model for multiple kinetic ion species, including thermal and friction forces (Simakov et al, PoP 2014)
	- − However, it eliminates non-local heat transport effects (important; need kinetic electrons for this)
- Ions remain fully kinetic, allow for multiple species

Taitano et al., CPC **258** (2021); JCP **365** (2018); JCP **318** (2016); JCP **297** (2015)

Model equations: fully kinetic ions + fluid electrons

Vlasov-Fokker-Planck for ion species

Fluid electrons

$$
\frac{3}{2}\partial_t (n_e T_e) + \frac{5}{2}\partial_x (u_e n_e T_e) - u_e \partial_x (n_e T_e) - \partial_x \kappa_e \partial_x T_e = \sum_{\alpha} C_{e\alpha}
$$

$$
n_e = -q_e^{-1} \sum_{\alpha} q_{\alpha} n_{\alpha} \qquad u_e = -q_e^{-1} n_e^{-1} \sum_{\alpha \neq e} q_{\alpha} n_{\alpha} u_{\alpha}
$$

Electric field model: e pressure, friction, thermal forces

$$
E = -\frac{\nabla p_e + \sum_i \mathbf{F}_{ie}}{en_e} = -\frac{\nabla p_e}{en_e} - \frac{\alpha_0 (Z_{eff}) m_e}{e} \sum_i \nu_{ei} (\mathbf{V}_e - \mathbf{V}_i) - \frac{\beta_0 (Z_{eff})}{e} \nabla T_e
$$

Simakov and Molvig, PoP **21** (2014)

Beyond rad-hydro: naïve algorithms will not work, even with asymptotic models

- Mesh requirements:
	- − Intra species v_{th.max} /v_{th.min}~100
	- − Inter species ($v_{th,α}$ / $v_{th,β}$)_{max}~30
	- $N_v \sim [10(v_{th\,max}/v_{th\,min})x(v_{th\,\alpha}/v_{th\,\beta})]^2$ ~10⁹
	- $N_r \sim 10^{3} 10^{4}$
	- − **N=NrNv~1012-1013 unknowns in 1D2V!**
- Timestep requirements: $\Delta t_{exp}^{coll} \sim \frac{1}{10} \left(\frac{\Delta v}{v_{min}} \right)^2 \nu_{coll}^{-1} \sim 10^{-9} ns$
	- $-$ t_{sim}=10 ns
	- − **Nt=1010 time steps**
- Beyond exascale (>10²³ FLOPS)!

Beyond rad-hydro: algorithmic innovation is the solution!

- Fully nonlinearly time-implicit (Δt >> τ_{col})
	- Iterate solution to convergence
	- Use fluid models to accelerate kinetic solution
- Optimal, adaptive grid in phase space
	- Adaptivity in velocity space based on shift and normalization to thermal speed
	- Moving radial mesh in physical space to follow capsule implosion

• Fully conservative (mass, momentum, and energy) and asymptotic preserving (able to capture LTE solution in strongly collisional regimes)

Beyond rad-hydro: Why is strict conservation critical?

With energy conservation Without energy conservation

Beyond rad-hydro: algorithms enable hybrid VFP-fluid simulations of entire capsule implosions

- Mesh requirements:
	- − v-space adaptivity with v_{th} normalization and u_{||} shift, N_v~10⁴-10⁵
	- − Moving mesh in physical space, N_r~10²
	- − Second-order accurate phase-space discretization
	- − **N=NvNr~106~107** (vs. 1012 with static mesh)
- Timestep requirements:
	- − Optimal O(N_v) implicit nonlinear algorithms
	- − Second-order-accurate timestepping
	- $\Delta t_{\text{imp}} = \Delta t_{\text{str}} \sim 10^{-3} \text{ ns}$
	- − **Nt~103-104** (vs. 1010 with explicit methods)
- Terascale-ready! (10¹² FLOPS, any reasonable cluster)
	- − Currently taking a few hours on 400 cores for full capsule implosion simulations!

An exploding-pusher VFP full capsule implosion simulation

Beyond rad-hydro: VFP modeling successfully predicts experimental trends!

DTn and DDn YOC in 50:50 DT exploding pushers

Future research directions in Computational Methods for Fusion of personal interest

Trends and directions in fusion computation

- There are several drivers of innovation in fusion computation:
	- − Drive towards whole-device modeling in ICF and MFE
	- − Drive towards higher simulation fidelity via model integration (e.g., MHD+GK, hybrid fluid-kinetic, etc.)
	- − Drive towards exascale computing (1018 FLOPS!)
- There are many efforts around the world responding to these trends
	- − Algorithms remain a key enabling technology for the simulations of the future, in the development of multiscale numerical formulations, spatial discretization and adaptivity, or in temporal integration via advanced (scalable) solvers
- I will comment next on a few directions of particular personal interest.

Towards exascale with fully implicit, adaptive MHD solvers

- We have been exploring the use of MHD physics-based preconditioning in combination with exascale-ready libraries for spatial discretization (MFEM) and solvers (PETSc, Trilinos)
- These libraries offer tremendous flexibility in discretization and solver choices, and offer state-of-the-art adaptive mesh refinement capabilities
- Physics-based preconditioning is discretization agnostic, so can be readily implemented with any discretization strategy.
- We have demonstrated the capability with system-scale simulations of magnetic reconnection using realistic values of resistivity and viscosity

 $F_{\rm in}$ $I_{\rm out}$ \sim $\frac{1}{2}$ $\frac{1}{2}$ Simulation of island coalescence in 2D with fully implicit S_{S} and AMD in MEEM with $\frac{1}{2}$ $\frac{1}{2}$ = 10-6. Moob defe is solver and AMR in MFEM with v = η = 10⁻⁶. Mesh dofs is 0.21% of an equivalent uniform mesh.

 τ i.e., the flow velocity, i.e., the plasmoids appearing on the top part of the current sheet move upward shee Tang et al, JCP, 110967 (2022)

Fully implicit GK electromagnetic PIC solvers the convergence of our current implementation seems to require *^vte^t*

- Semi-implicit GK PIC algorithms are known to suffer from two numerical problems:
	- − Cancellation errors: arise from lack of cancellation of skin currents represented on both the mesh and the particles (needed for numerical stability)
	- − Finite-grid instabilities: due to aliasing errors arising from particles living in the continuum, while fields live on a discrete mesh
- Both these issues can be eliminated with fully implicit methods

Sturdevant et al., PoP, 072505 (2021) Sturdevant el al., JCP, submitted (2022)

E ully implicit algorithms in VCC in to the cancellation problem, and in possible economic economic economic economic economic economic economic

Toward 6D simulations of MFE devices

- GK has limitations due to embedding of asymptotic approximations:
	- − Break down in the presence of strong plasma gradients, such as in high-confinement conditions (due to pressure pedestal)
- Improving fidelity needs multiscale asymptoticpreserving integrators that can seamlessly transition between GK and full descriptions (5D to 6D)
- Will require specialized algorithms and solvers:
	- Fully implicit timestepping with strong conservation properties (for stability and accuracy)
	- § Asymptotic-preserving particle orbit integrators that capture orbit without following gyromotion
	- Use of nested model hierarchy for full algorithmic acceleration (e.g, use MHD to accelerate a fully kinetic simulation).

1. Ricketson et al, in preparation 2. Chen et al., JCP (submitted) $\frac{L}{2}$

7/12/21 72
Multi-D hybrid kinetic/fluid modeling of ICF hohlraums

- Hohlraums perform critical energy conversion process from laser to X-rays
- Hohlraum environment is rarefied (or $vacuum) \Rightarrow$ kinetic effects are important, cannot be modelled with rad-hydro
	- − Plasma expansion into vacuum
	- − Multiple ion species
	- − Beam interpenetration
- Large electromagnetic fields have been measured in hohlraums, cannot be neglected
- Hohlraum modeling uncertainty is preventing progress in ICF implosion optimization towards ignition

• At LANL, we have started developing the first multi-D hybrid kinetic ion/fluid electron code for hohlraum modeling.

Summary

Summary

- Thermonuclear fusion poses great challenges to the computational physicist
- Challenges have been met by ingenuity in developing an asymptotic model hierarchy, as well as solvers and algorithms
- Fusion has been a pioneer in HPC, driving the creation of the first unclassified computer systems (now NERSC)
- Fusion has also driven significant algorithmic innovation, and has been a pioneer in the use of modern discretizations and implicit timestepping schemes.
- Both MFE and ICF are pushing the computational frontier towards higher fidelity (kinetic) simulations, by leveraging model nesting, scalable solvers, and smart algorithms.
- These computational capabilities will continue to inform future iterations of "virtual experiments", with the goal of harnessing fusion energy on Earth.