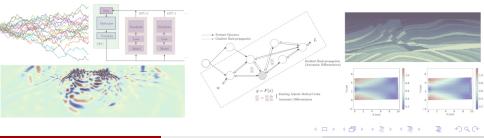
ADCME MPI: Distributed Machine Learning for Computational Engineering

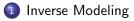
Kailai Xu and Eric Darve https://github.com/kailaix/ADCME.jl



ADCME-MPI

Distributed ML for Scientific Computing

Outline



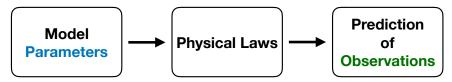
Automatic Differentiation

Distributed Computing for Computational Engineering

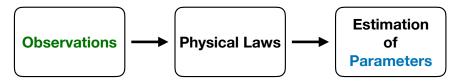
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Forward Problem



Inverse Problem



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Inverse Modeling

We can formulate inverse modeling as a PDE-constrained optimization problem

$$\min_{\theta} L_h(u_h) \quad \text{s.t. } F_h(\theta, u_h) = 0$$

- The loss function L_h measures the discrepancy between the prediction u_h and the observation u_{obs} , e.g., $L_h(u_h) = ||u_h u_{obs}||_2^2$.
- θ is the model parameter to be calibrated.
- The physics constraints F_h(θ, u_h) = 0 are described by a system of partial differential equations or differential algebraic equations (DAEs); e.g.,

$$F_h(\theta, u_h) = \mathsf{A}(\theta)u_h - f_h = 0$$

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$$\min_{\mathbf{f}} L_h(u_h) \quad \text{s.t. } F_h(\mathbf{f}, u_h) = 0$$

What if the unknown is a function instead of a set of parameters?

- Koopman operator in dynamical systems.
- Constitutive relations in solid mechanics.
- Turbulent closure relations in fluid mechanics.

• ...

The candidate solution space is infinite dimensional.

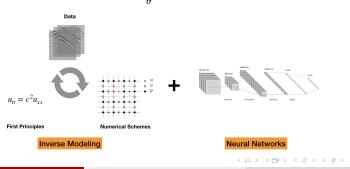
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Machine Learning for Computational Engineering

 $\min_{\theta} L_h(u_h) \quad \text{s.t.} \quad F_h(NN_{\theta}, u_h) = 0 \quad \leftarrow \text{ Solved numerically}$

- Use a deep neural network to approximate the (high dimensional) unknown function;
- Solve u_h from the physical constraint using a numerical PDE solver;
- Apply an unconstrained optimizer to the reduced problem

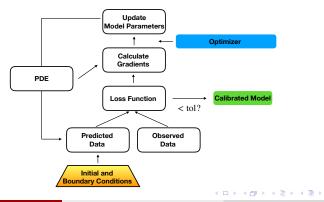


$$\min_{\theta} L_h(\underline{u}_h(\theta))$$

Gradient Based Optimization

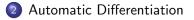
 $\min_{\theta} L_h(u_h) \quad \text{s.t. } F_h(\theta, u_h) = 0 \quad \Leftrightarrow \quad \min_{\theta} L_h(u_h(\theta))$ • We can now apply a gradient-based optimization method if we can calculate a descent direction g^k

$$\theta^{k+1} \leftarrow \theta^k - \alpha g^k$$



Outline





Distributed Computing for Computational Engineering

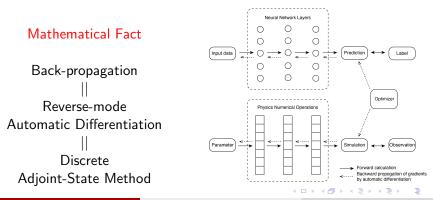
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Automatic Differentiation

The fact that bridges the technical gap between machine learning and inverse modeling:

 Deep learning (and many other machine learning techniques) and numerical schemes share the same computational model: composition of individual operators.

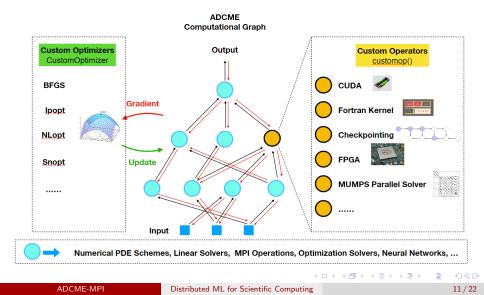


Computational Graph for Numerical Schemes

- To leverage automatic differentiation for inverse modeling, we need to express the numerical schemes in the "AD language": computational graph.
- No matter how complicated a numerical scheme is, it can be decomposed into a collection of operators that are interlinked via state variable dependencies.

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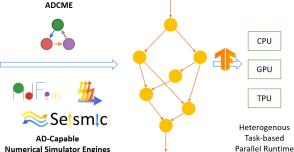
ADCME: Computational-Graph-based Numerical Simulation



How ADCME works

 ADCME translates your high level numerical simulation codes to computational graph and then the computations are delegated to a heterogeneous task-based parallel computing environment through TensorFlow runtime.

div $\sigma(u) = f(x)$ $x \in \Omega$ $\sigma(u) = C\varepsilon(u)$ $u(x) = u_0(x)$ $x \in \Gamma_u$ $\sigma(x)n(x) = t(x)$ $x \in \Gamma_n$ mmesh = Mesh(50, 50, 1/50, degree=2) left = bcmode((x,y)->xcle-5, mmesh) right = bcedge((x1,y1,x2,y2)->(x1>0.049-1e-5) && (x2>0.049-1e-5), mmesh) t1 = eval f on boundary edge((x,y)->1.0e-4, right, mmesh) t2 = eval_f_on_boundary_edge((x,y)->0.0, right, mmesh) rhs - compute fem traction term(t1, t2, right, mmesh) x = gauss_nodes(mmesh) E = abs(fc(x, [20, 20, 20, 1]))>squeeze) # E = constant(eval f on gauss pts(f, mmesh)) D = compute_plane_stress_matrix(E, nu*ones(get_ngauss(mmesh))) K = compute fem stiffness matrix(D, mmesh) bdval = [eval_f_on_boundary_node((x,y)->0.0, left, mmesh); eval f on boundary node((x,y)->0.0, left, mmesh)] DOF = [left:left .+ mmesh.ndof] K, rhs = impose Dirichlet boundary conditions(K, rhs, DOF, bdyal)



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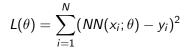


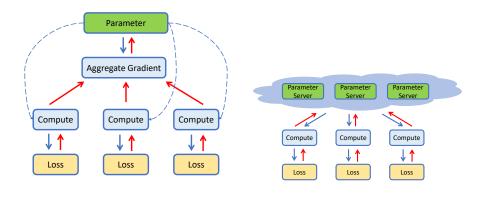
Oistributed Computing for Computational Engineering

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Common Distributed Computing Patterns in DL





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Distributed Computing in ML for Computational Engineering

Consider a time-dependent PDE, where the state variable

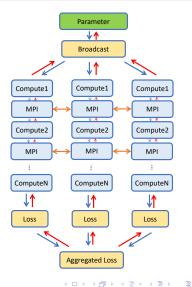
$$u_k = [u_k^{(1)} \ u_k^{(2)} \ \cdots \ u_k^{(P)}]$$

is stored on *P* machines. Each time step requires a distributed numerical solver.

$$\min_{\theta} L(u_n)$$

s.t. $A(\theta)u_2 = h(u_1; \theta) + g$
 $A(\theta)u_3 = h(u_2; \theta) + g$
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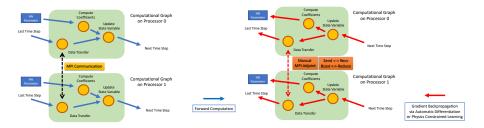
$$A(\theta)u_n = h(u_{n-1};\theta) + g$$



ADCME-MPI

ADCME-MPI abstracts distributed computing as a node in the computational graph. The ADCME-MPI model is **transparent**.

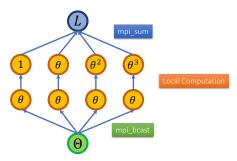
- ADCME takes responsibility for MPI communication and gradient back-propagation across clusters;
- users can adapt their single processor codes to a distributed computing environment with little efforts.



Example

Consider a simple function:

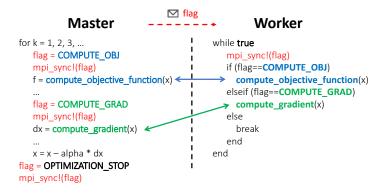
$$L(\theta) = 1 + \theta + \theta^2 + \theta^3$$



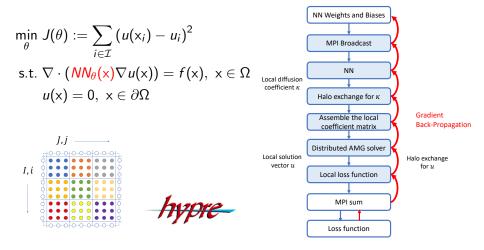
using ADCME mpi_init() # initialize MPI theta0 = placeholder(1.0)theta = mpi_bcast(theta0) l = theta^mpi_rank() $L = mpi_sum(1)$ g = gradients(L, theta0) # initialize a Session sess = Session(); init(sess) $L_value = run(sess, L)$ g_value = run(sess, g) mpi_finalize() # finalize MPI

Distributed Optimization

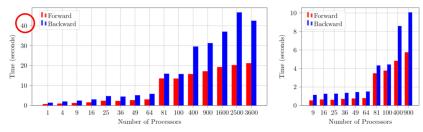
In the ADCME-MPI, we can convert a serial optimizer to a distributed optimizer by inserting some communication codes:



Benchmarks

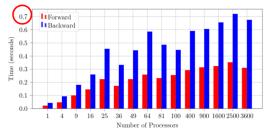


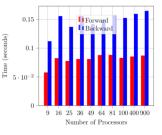
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Total Time

ADCME Overhead = Total Time – Hyper Time





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1 Core per MPI processor

4 Core per MPI processor

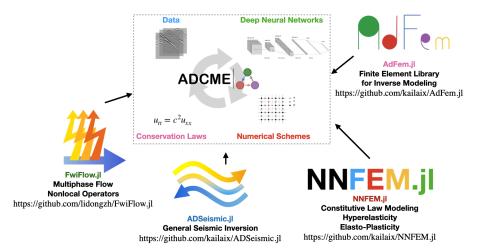
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For more technical details, benchmarks, or use cases:

- AAAI Conference Paper: ADCME MPI: Distributed Machine Learning for Computational Engineering
- Full paper: Distributed Machine Learning for Computational Engineering using MPI https://arxiv.org/pdf/2011.01349.pdf
- Software documentation: https://kailaix.github.io/ADCME.jl/dev/

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A General Approach to Inverse Modeling



Distributed ML for Scientific Computing