Meta-Programming and Hybrid Parallel Strategies for Solving PDEs: An FDM and PINN Comparison ^{1 2}

Seminar Presentation III

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full docs: https://livihai.com/html/index.html

²repository: https://github.com/livihai-official/Final-Project

Outline

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 - Related Work
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 - General Form
 - Thermal Conduction Systems
- Methodology
 - N-Dimension Matrix
 - Parallelization of N-dimension Arrays

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 - PINN Model
- S Experiments
 - On Single Node
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 - On Different Dimensions
 - With PINN on accuracy
- 6 Discussion

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Kolmogorov PDEs

Solving
$$u(x, T)$$
, for $\mathbb{R}^1 \ni T > 0$, $x \in \mathbb{R}^d$, $t \in [0, T]$, $u(t, x) = u \in \mathbb{R}^1$, $\mu(x) \in \mathbb{R}^d$, $\sigma(x) \in \mathbb{R}^{d \times d}$,
$$u_t = \frac{1}{2} \operatorname{Trace}_{\mathbb{R}^d} \left[\sigma(x) \left[\sigma(x) \right]^* \operatorname{Hess}_x u \right] + \langle \mu(x), \nabla_x u \rangle_{\mathbb{R}^d}$$
(1)

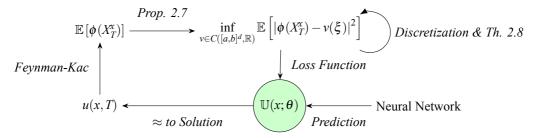


Figure: Deep Neural Network (DNN) Methodology of Solving Kolmogorov PDEs [FIRST]

Introduction

Recap - Physics Informed Nerual Network

General Form of PDEs

- -u(t,x) denotes with the target function, $x \in \mathbb{R}^d$.
- $-\Gamma[\cdot;\lambda]$ is a non-linear operator parameterized by λ .

$$u_t(t,x) + \Gamma[u;\lambda] = 0 \tag{2}$$

Define f(t,x) to be given by

$$f(t,x) = u_t(t,x) + \Gamma[u;\lambda]$$
(3)

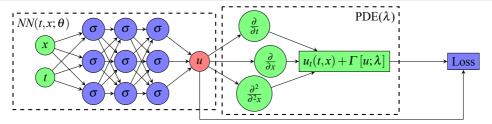


Figure: PINN, with 3 fully connected hidden layers

Introduction Recap - Conclusion

Comparing With Finite Difference Time Domain Method (FDTD)

- Deep Neural Network [FIRST]
 - Gives lower quality approximations.
 - Takes longer time to train.
 - Possible to solve high dimension PDEs
- Physics Informed Neural Network
 - Gives higher quality approximations.
 - Takes longer time to train.
 - Has more flexible way to get results.
 - Possible to solve high dimension PDEs

Objectives

This project focused on following objectives:

- Implement FDTD and PINN in C++/C.
- Implement FDTD hybrid parallel version using MPI/OpenMP.
- Implement PINN GPU parallel version using Libtorch/CUDA.
- Evaluate the efficiency and accuracy of FDTDs and PINNs.

Challanges

However, there were many obstacles including:

- O Portability.
- Overlapping Communication and Computation.
- Unnecessary intra-node communication
- Communication Overhead.
- Scalability Issues.
- Memory Management.

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General Form of problem

The PDE parametrized by number λ and an operator $\mathcal{N}[\cdot; \lambda]$, and assume the variable x is a 2D or 3D spatio-vector which is written in

$$\begin{cases} \frac{\partial u}{\partial t}(t, \vec{x}) + \mathcal{N}[u; \lambda] = 0\\ u(0, \vec{x}) = \varphi(\vec{x}) \end{cases}$$
(4)

where φ is the initial condition, and $\vec{x} \in \Omega, t \in [0, +\infty)$.

Boundary Conditions

The Dirichlet and Von Neurmann boundary conditions are formed as

$$\begin{cases} u(t,\vec{x}) = g(t,\vec{x}) \\ \frac{\partial u}{\partial \vec{n}} = g(t,\vec{x}) \end{cases}$$
 (5)

where \vec{n} is the normal vector on $\overline{\Omega}$ the boundary of domain Ω .

Thermal Conduction Systems Heat Equation 2D

The function

$$u(t,x,y) = x + y - xy, \,\forall \alpha \in \mathbb{R}^1$$
 (6)

is the solution of 2D Heat Equation 7 below

$$\frac{\partial u}{\partial t} = \alpha \left(\frac{\partial u^2}{\partial^2 x} + \frac{\partial u^2}{\partial^2 y} \right) \qquad (x, y) \in \Omega, t \in [0, +\infty)$$

$$u(0, x, y) = \varphi(x, y) = 0 \qquad (x, y) \in \Omega \qquad (7)$$

$$\begin{cases} y, \ x = 0, y \in (0, 1) \end{cases}$$

$$u(t,x,y) = g(x,y) = \begin{cases} y, & x = 0, y \in (0,1) \\ 1, & x = 1, y \in (0,1) \\ x, & y = 0, x \in (0,1) \\ 1, & y = 1, x \in (0,1) \end{cases}$$
 $t \in [0, +\infty)$

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An FDM and PINN Comparison

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Thermal Conduction Systems Heat Equation 3D

The function

$$u(t,x,y,z) = x + y + z - 2xy - 2xz - 2yz + 4xyz, \forall \alpha \in \mathbb{R}^1$$
(8)

is the solution of 3D Heat Equation 9 below

$$\frac{\partial u}{\partial t} = \alpha \left(\frac{\partial u^2}{\partial x^2} + \frac{\partial u^2}{\partial y^2} + \frac{\partial u^2}{\partial z^2} \right) \qquad (x, y, z) \in \Omega, t \in [0, +\infty)$$

$$u(0, x, y, z) = \varphi(x, y, z) = 0 \qquad (x, y, z) \in \Omega \qquad (9)$$

$$u(t,x,y,z) = g(x,y,z) = \begin{cases} y+z-2yz, & x=0, \\ 1-y-z+2yz, & x=1, \\ x+z-2xz, & y=0, \\ 1-x-z+2xz, & y=1, \\ x+y-2xy, & z=0, \\ 1-x-y+2xy, & z=1 \end{cases}$$
 $t \in [0,+\infty)$

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N-dimension Matrix Challanges

STL provides containers std::array and std::vector for creating one-dimension array. There are two way for dealing with high-dimension data:

- Nesting the one dimension arrays or vectors.
- Hierarchy approach, designing derived classes of one-dimension array base class.

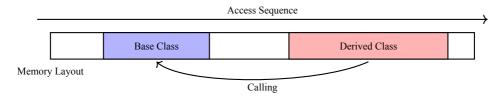


Figure: Derived Class calling members in Base class, timing is not predictable.

- Nesting multi-dimension array has non-contiguous memory layout.
- Operived class needs more time to access members in base class.
- Or cache utilization leads to poor performance.
- MPI type create requires contiguous memory layout.

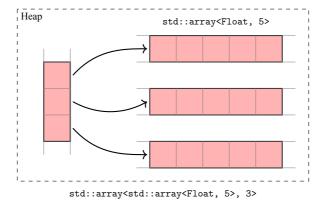


Figure: Using nested std::array<T, N> to store 2D array data.

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N-dimension Matrix Solution

- An external small __multi_array_shape object defines the routines for accessing the elements.
- Smart pointer, ensure memory's contiguous layout and safety.
- Separate into two detail and user interface objects adhering RAII rules.

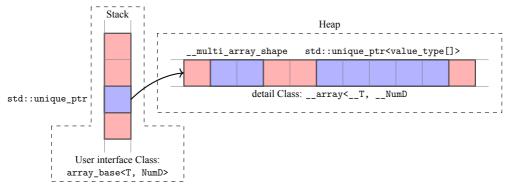


Figure: The solution of N-dimension Matrix, using detail Class, user interface Class and a shape management structure.

Parallelization of N-dimension Arrays MPI Environment

The hybrid PDE solver requires the MPI supports multi-threads on each processes.

- High-level libraries like Boost.MPI
 - have better MPI resource management and other basic communication features.
 - only provide limited useful features for latter PDE solvers.
 - 1 lead to lower performance than low-level OpenMPI.
- I developed an environment class for MPI
 - better resource management than raw MPI.
 - provides basic features exclusively for this project.

Parallelization of N-dimension Arrays MPI Topology - Challanges

Distributed N-dimension arrays are created based on MPI N-dimension Cartesian topology.

- Ghost communication is required in overlapping of MPI communication and local computation.
- 2 Parallel I/O is needed for debugging and storing results.
- **3** Topology information will be frequently used in PDE solver.

Parallelization of N-dimension Arrays MPI Topology - Solution

- An MPI Topology class defines the distribution details of N-dimension array.
- ② Using MPI_Type_create_subarray for creating Ghost MPI datatype for communication.
- Ocartesian array has members Topology class and N-dimension array class, to ensure they are closely located on memory.
- The external functions provide gather-based I/O and MPI I/O for handling different scenarios.
- User interface class unified features for both topology and array classes.
- Separate into two detail and user interface objects adhering RAII rules.

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N-dimension Boundary and Initial Conditions Challanges

- Mathematical function needs be discretized and distributed as well.
- Need to access the data.
- Initial and Dirichlet boundary conditions only applies once.
- On Neumann boundary condition participates the evolving process in FDM.

N-dimension Boundary and Initial Conditions Solution

- Use lambda function to construct classes.
- Oreating external classes for each conditions as the friend classes of PDE solver classes.
- Set Bool vectors help to determine the status of conditions and type of boundary conditions.

N-dimension PDE solvers Challanges

- PDE solver of Heat Equations in different dimension space have similar parameters and features, type-field solution lead to code redundancy.
- Applying MPI communications between local arrays, avoiding overhead.
- Three type of strategies:
 - Pure MPI parallel
 - Master-only, no overlapping hybrid parallel.
 - Master-only, communication/computation (comm./comp.) overlapping hybrid parallel.

virtual function is resolved at run-time, and only lose up to about 25% efficiency in terms of the function call mechanism,

• Create an abstract Heat base class of Heat Equation, and overriding functions in derived class on every dimension.

MPI communications are implemented in blocking and non-blocking ways using MPI_Sendrecv and MPI_Isend/MPI_Irecv.

- Three type of strategies
 - Pure MPI parallel

- Three type of strategies
 - Master-only, no overlapping hybrid parallel.

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- Three type of strategies
 - Master-only, comm./comp. overlapping hybrid parallel.

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PINN Model Challanges

For neural network implementations, there are some issues in practice: Python has many easy-use libraries such as Pytorch, Tensorflow and Caffe.

- Interpret language Python is significant slower than compile language C/C++
- Worse resource management than C/C++
- Security in parallel program.

PINN Model Solution

For getting higher performance and better safety, I chose to use

• Pytorch C++ API: Libtorch.

to reproduce the Python version of PINN.

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Single Node Scaling Tests

2D Heat Equation Single Precision

Problem Size: 512², 1024², 2048², 4096², 8192², 16384², 32768²

Non-unified Memory Access (NUMA) node Topology:

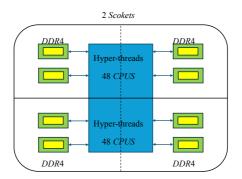
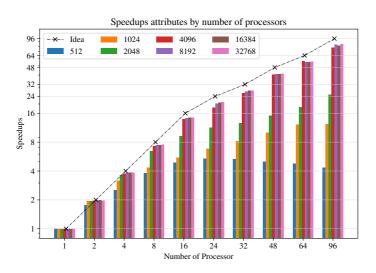
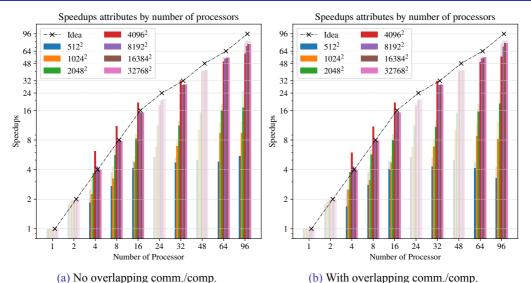


Figure: NUMA topology of single node on Cluster

Single Node Scaling Tests Strong Scaling - pure MPI



Single Node Scaling Tests Strong Scaling - MPI/OpenMP Hybrid

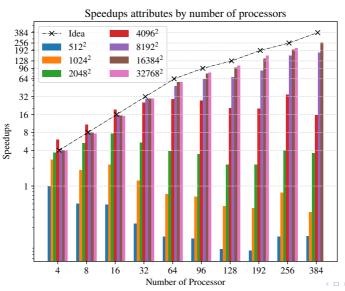


(b) With overlapping comm./comp.

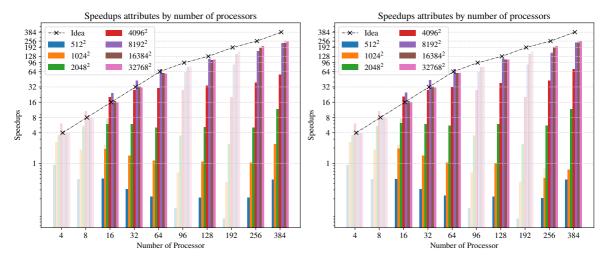
Single Node Scaling Tests Weak Scaling

Strategy	Size		C (C4)		
		4	16	64	$f_p(\%)$
Pure MPI		4.006	12.497	47.849	75.1
No Overlap	512^{2}	2.876	11.206	42.754	67.0
With Overlap		3.173	10.818	42.282	66.2
Pure MPI		3.838	9.304	33.707	53.2
No Overlap	1024^2	3.947	12.995	33.447	54.1
With Overlap		4.024	12.932	33.361	54.0
Pure MPI	re MPI	2.376	8.245	31.203	49.0
No Overlap	2048^{2}	3.874	8.972	31.510	49.8
With Overlap		3.740	8.989	31.430	49.7
Pure MPI		3.543	8.245	31.203	77.5
No Overlap	4096^{2}	3.953	13.799	49.515	78.0
With Overlap		3.948	13.800	49.989	78.7

Multi-node Scaling Tests Strong Scaling - pure MPI



Single Node Scaling Tests Strong Scaling - MPI/OpenMP Hybrid



(a) No overlapping comm./comp.

(b) With overlapping comm./comp.

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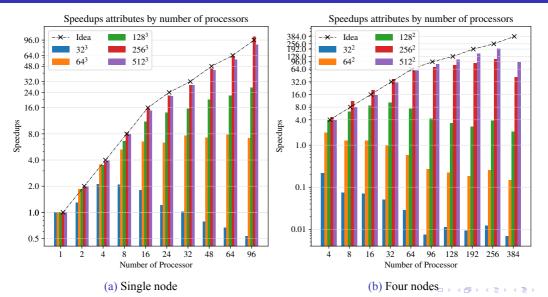
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Single Node Scaling Tests Weak Scaling

Strategy	Size		Number of CPUs			
		4	16	64	256	$f_p(\%)$
Pure MPI		3.300	10.379	26.000	124.375	48.2
No Overlap	512^{2}	-	8.094	25.931	127.648	49.3
With Overlap		-	8.386	27.126	116.951	45.5
Pure MPI		3.848	13.037	30.143	-	49.3
No Overlap	1024^2	-	13.702	44.040	-	69.8
With Overlap		-	13.834	44.090	-	69.9
Pure MPI		3.787	9.172	32.013	-	50.6
No Overlap	2048^{2}	-	13.936	34.368	-	55.7
With Overlap		-	14.274	34.414	-	55.9
Pure MPI		3.884	13.859	51.041	-	80.2
No Overlap		-	15.315	53.157	-	83.8
With Overlap		-	15.294	53.401	-	84.2

3D Space Heat Equation Strong Scaling - pure MPI



3D Space Heat Equation

Weak Scaling - MPI/OpenMP Hybrid, Single Node

3D Space Heat Equation

Weak Scaling - MPI/OpenMP Hybrid, Four Nodes

3D Space Heat Equation Weak Scaling

Memory Usage v.s. Time

Accuracy

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Closing Remarks

Thank you for your attention!

Any questions?