

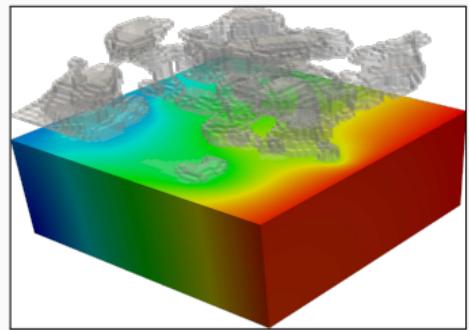
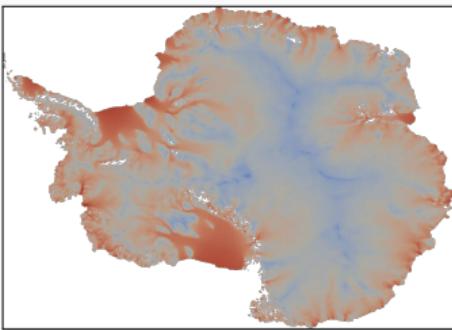
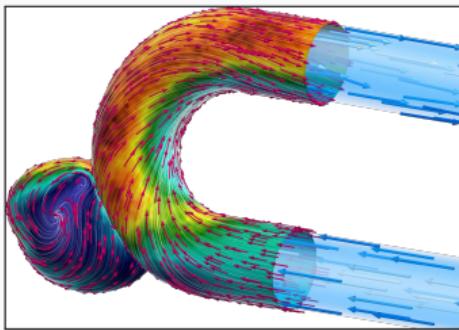
Domain decomposition for physics-informed neural networks

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Séminaire d'analyse numérique, Université de Genève, Geneva, Switzerland, February 20, 2024

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Scientific Machine Learning in Computational Science and Engineering



Numerical methods

Based on physical models

- + Robust and generalizable
- Require availability of mathematical models

Machine learning models

Driven by data

- + Do not require mathematical models
- Sensitive to data, limited extrapolation capabilities

Scientific machine learning (SciML)

Combining the strengths and compensating the weaknesses of the individual approaches:

numerical methods	improve	machine learning techniques
machine learning techniques	assist	numerical methods

Outline

1 Physics-informed machine learning & motivation

2 Deep learning-based domain decomposition method

Based on joint work with

Victorita Dolean (TU Eindhoven)

Serge Gratton and **Valentin Mercier** (IRIT Computer Science Research Institute of Toulouse)

3 Multilevel domain decomposition-based architectures for physics-informed neural networks

Based on joint work with

Victorita Dolean (University of Strathclyde, University Côte d'Azur)

Ben Moseley and **Siddhartha Mishra** (ETH Zürich)

4 Multifidelity domain decomposition-based physics-informed neural networks for time-dependent problems

Based on joint work with

Damien Beecroft (University of Washington)

Amanda A. Howard and **Panos Stinis** (Pacific Northwest National Laboratory)

Physics-informed machine learning & motivation

Artificial Neural Networks for Solving Ordinary and Partial Differential Equations

Isaac Elias Lagaris, Aristidis Likas, *Member, IEEE*, and Dimitrios I. Fotiadis

Published in **IEEE Transactions on Neural Networks, Vol. 9, No. 5, 1998.**

Approach

Solve a general differential equation subject to boundary conditions

$$G(x, \Psi(x), \nabla\Psi(x), \nabla^2\Psi(x)) = 0 \quad \text{in } \Omega$$

by solving an **optimization problem**

$$\min_{\theta} \sum_{x_i} G(x_i, \Psi_t(x_i, \theta), \nabla\Psi_t(x_i, \theta), \nabla^2\Psi_t(x_i, \theta))^2$$

where $\Psi_t(x, \theta)$ is a **trial function**, x_i sampling points inside the domain Ω and θ are **adjustable parameters**.

Construction of the trial functions

The trial functions **satisfy the boundary conditions explicitly**:

$$\Psi_t(x, \theta) = A(x) + F(x, \text{NN}(x, \theta))$$

- NN is a **feedforward neural network** with **trainable parameters** θ and input $x \in \mathbb{R}^n$
- A and F are **fixed functions**, chosen s.t.:
 - A satisfies the boundary conditions
 - F does not contribute to the boundary conditions

Neural Networks for Solving Differential Equations

Approach

Solve a general differential equation subject to boundary conditions

$$G(x, \Psi(x), \nabla\Psi(x), \nabla^2\Psi(x)) = 0 \quad \text{in } \Omega$$

by solving an **optimization problem**

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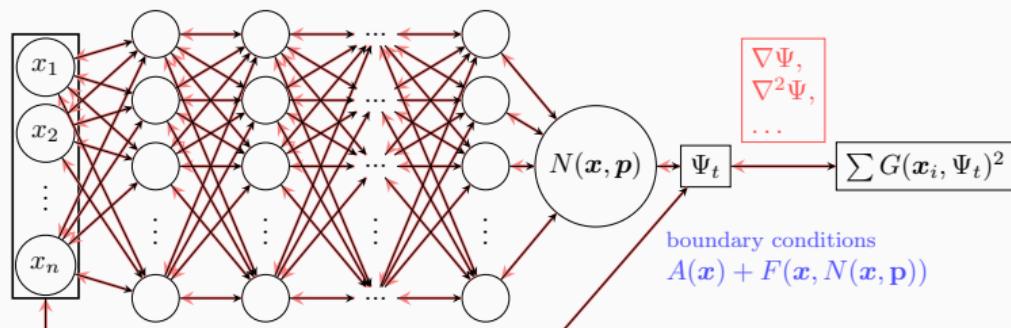
where $\Psi_t(x, \theta)$ is a **trial function**, x_i sampling points inside the domain Ω and θ are adjustable parameters.

Construction of the trial functions

The trial functions **satisfy the boundary conditions explicitly**:

$$\Psi_t(x, \theta) = A(x) + F(x, \text{NN}(x, \theta))$$

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 - F does not contribute to the boundary conditions



Lagaris et. al's Method – Motivation

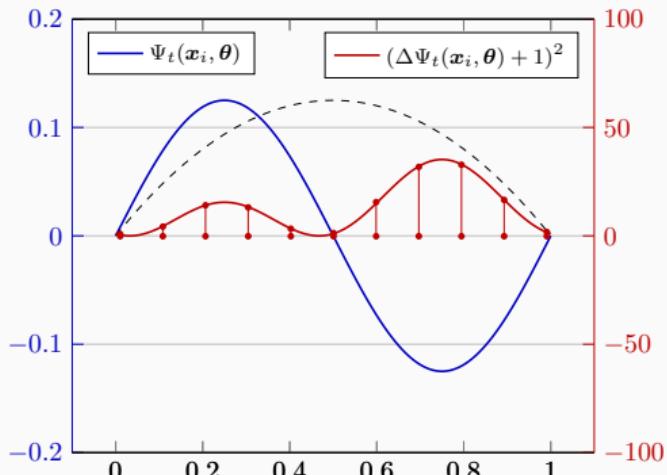
Solve the **boundary value problem**

$$\Delta \Psi_t(x, \theta) + 1 = 0 \text{ on } [0, 1],$$

$$\Psi_t(0, \theta) = \Psi_t(1, \theta) = 0,$$

via a **collocation approach**:

$$\min_{\theta} \sum_{x_i} (1 - \Delta \Psi_t(x_i, \theta))^2$$

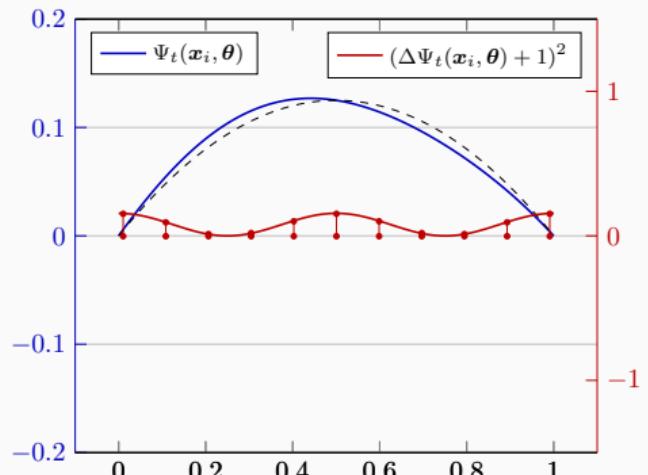


$$(\Delta \Psi_t(x_i, \theta) + 1)^2 >> 0$$

Boundary conditions

The boundary conditions can be **enforced explicitly**, for instance, via the ansatz:

$$\Psi_t(x, \theta) = \sin(\pi x) \cdot F(x, \text{NN}(x, \theta))$$



$$(\Delta \Psi_t(x_i, \theta) + 1)^2 \approx 0$$

Physics-Informed Neural Networks (PINNs)

In the **physics-informed neural network (PINN)** approach introduced by [Raissi et al. \(2019\)](#), a **neural network** is employed to **discretize a partial differential equation**

$$\mathcal{N}[u] = f, \quad \text{in } \Omega.$$

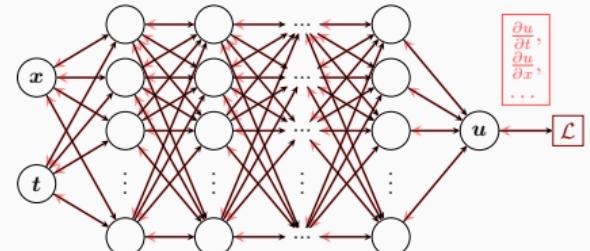
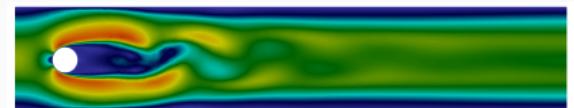
It is based on the approach by [Lagaris et al. \(1998\)](#). The main novelty of PINNs is the use of a **hybrid loss function**:

$$\mathcal{L}(\theta) = \omega_{\text{data}} \mathcal{L}_{\text{data}}(\theta) + \omega_{\text{PDE}} \mathcal{L}_{\text{PDE}}(\theta),$$

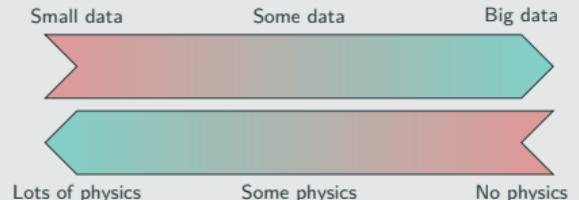
where ω_{data} and ω_{PDE} are **weights** and

$$\mathcal{L}_{\text{data}}(\theta) = \frac{1}{N_{\text{data}}} \sum_{i=1}^{N_{\text{data}}} (u(\hat{x}_i, \theta) - u_i)^2,$$

$$\mathcal{L}_{\text{PDE}}(\theta) = \frac{1}{N_{\text{PDE}}} \sum_{i=1}^{N_{\text{PDE}}} (\mathcal{N}[u](x_i, \theta) - f(x_i))^2.$$



Hybrid loss



Advantages

- "Meshfree"
- Small data
- Generalization properties
- High-dimensional problems
- Inverse and parameterized problems

Drawbacks

- Training cost and robustness
- Convergence not well-understood
- Difficulties with scalability and multi-scale problems

- Known solution values can be included in $\mathcal{L}_{\text{data}}$
- Initial and boundary conditions are also included in $\mathcal{L}_{\text{data}}$

Available Theoretical Results for PINNs – An Example

Mishra and Molinaro. *Estimates on the generalisation error of PINNs, 2022*

Estimate of the generalization error

The generalization error (or total error) satisfies

$$\mathcal{E}_G \leq C_{\text{PDE}} \mathcal{E}_{\mathcal{T}} + C_{\text{PDE}} C_{\text{quad}}^{1/p} N^{-\alpha/p}$$

where

- $\mathcal{E}_G = \mathcal{E}_G(\mathbf{X}, \theta) := \|\mathbf{u} - \mathbf{u}^*\|_V$ **general. error** (V Sobolev space, \mathbf{X} training data set)
- $\mathcal{E}_{\mathcal{T}}$ **training error** (l^p loss of the residual of the PDE)
- N **number of the training points** and α **convergence rate of the quadrature**
- C_{PDE} and C_{quad} **constants** depending on the **PDE** respectively the **quadrature** as well as on the **neural network**

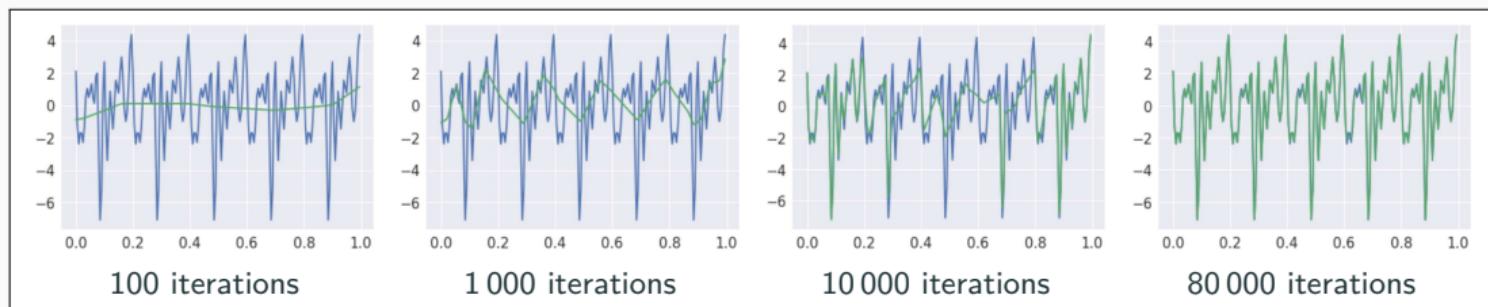
Rule of thumb:

“As long as the PINN is **trained well**, it also **generalizes well**”

Scaling Issues in Neural Network Training

Spectral bias

Neural networks prioritize learning lower frequency functions first irrespective of their amplitude.



Rahaman et al., *On the spectral bias of neural networks*, ICML (2019)

- Solving solutions on **large domains and/or with multiscale features** potentially requires **very large neural networks**.
- Training may **not sufficiently reduce the loss** or take **large numbers of iterations**.
- Significant **increase on the computational work**

Dependence on the choice of **activation functions**: Hong et al. (arXiv 2022)

Convergence analysis of PINNs via the neural tangent kernel: Wang, Yu, Perdikaris, *When and why PINNs fail to train: A neural tangent kernel perspective*, JCP (2022)

Motivation – Some Observations on the Performance of PINNs

Solve

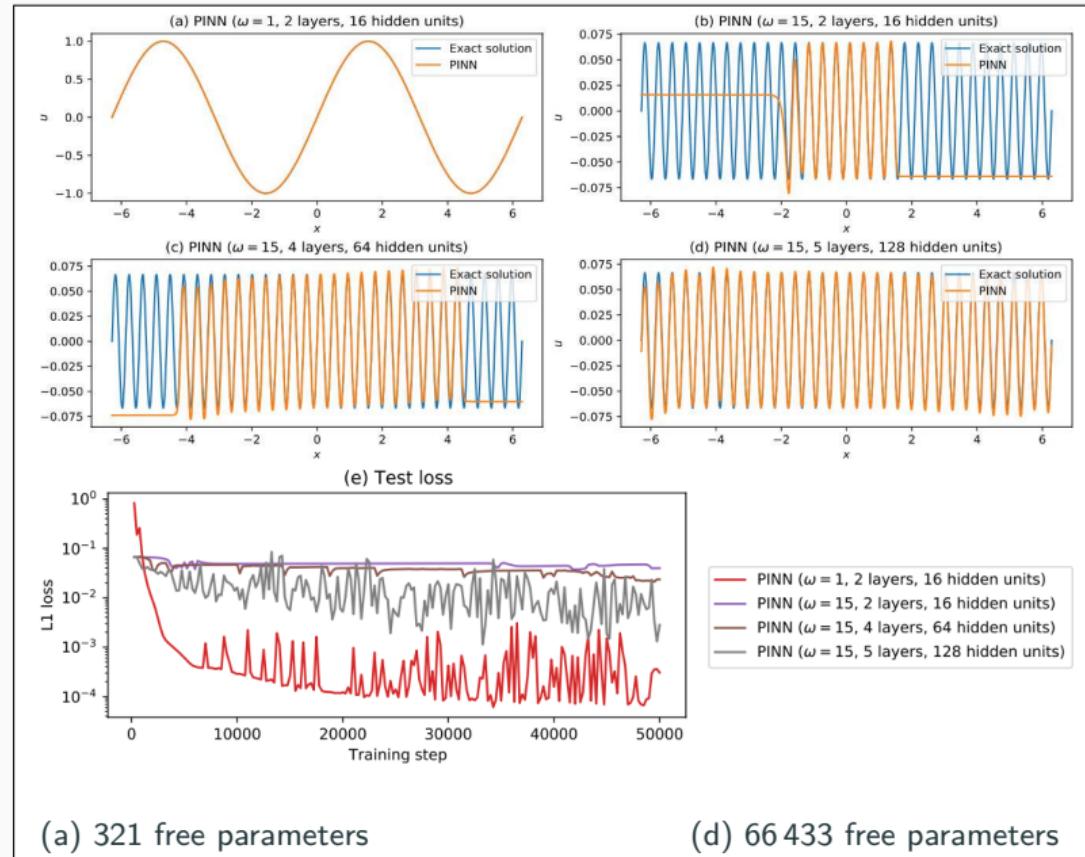
$$\begin{aligned} u' &= \cos(\omega x), \\ u(0) &= 0, \end{aligned}$$

for different values of ω
using PINNs with
varying network
capacities.

Scaling issues

- Large computational domains
- Small frequencies

Cf. Moseley, Markham, and Nissen-Meyer (2023)



Motivation – Some Observations on the Performance of PINNs

Solve

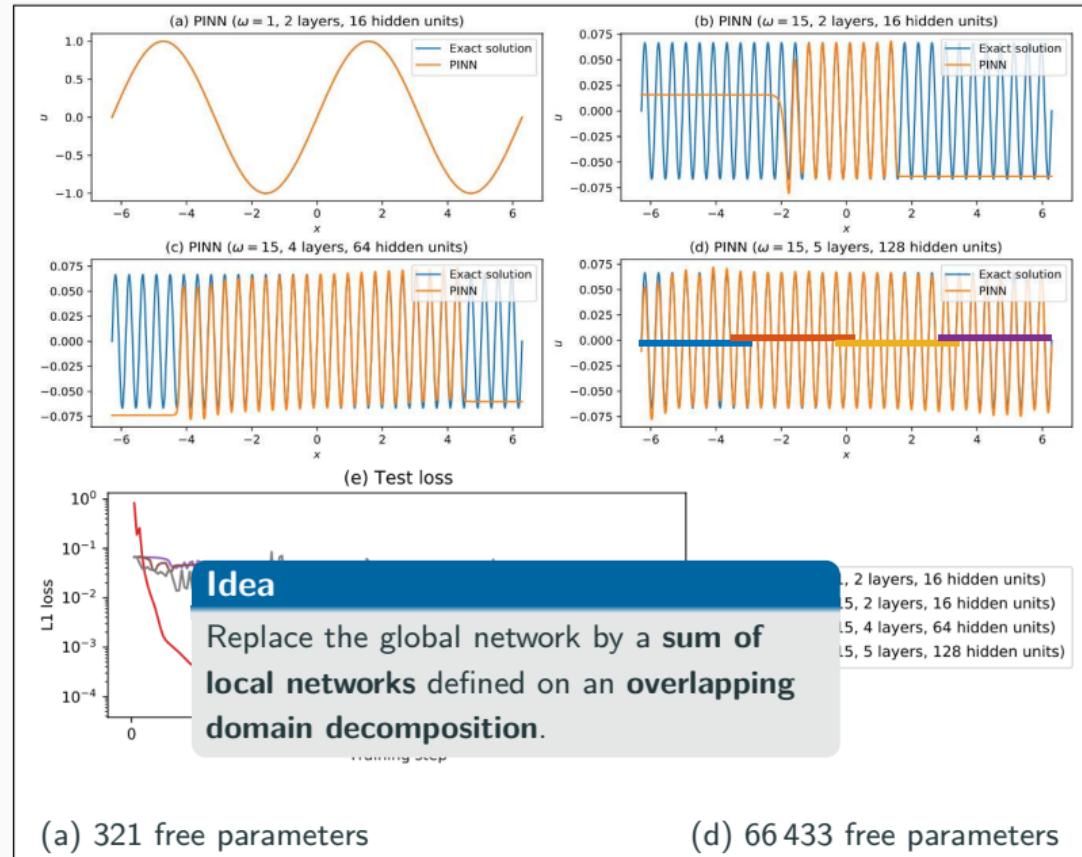
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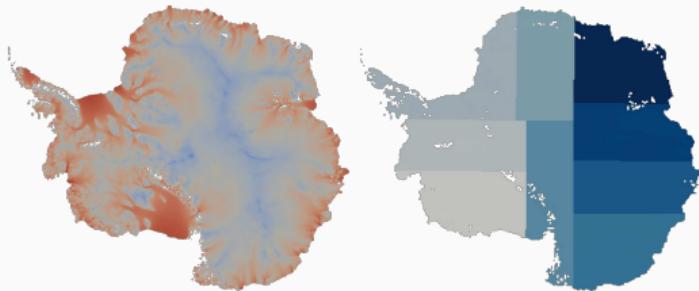
Scaling issues

- Large computational domains
- Small frequencies

Cf. Moseley, Markham, and Nissen-Meyer (2023)



Domain Decomposition Methods



Images based on Heinlein, Perego, Rajamanickam (2022)

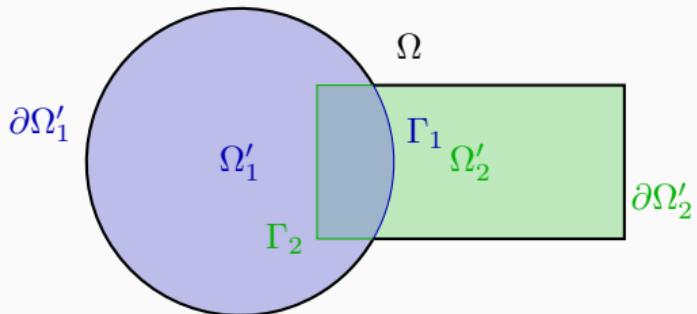
Historical remarks: The **alternating Schwarz method** is the earliest **domain decomposition method (DDM)**, which has been invented by **H. A. Schwarz** and published in **1870**:

- Schwarz used the algorithm to establish the **existence of harmonic functions** with prescribed boundary values on **regions with non-smooth boundaries**.

Idea

Decomposing a large **global problem** into smaller **local problems**:

- Better robustness** and **scalability** of numerical solvers
- Improved computational efficiency**
- Introduce **parallelism**



A non-exhaustive overview:

- cPINNs: Jagtap, Kharazmi, Karniadakis (2020)
- XPINNs: Jagtap, Karniadakis (2020)
- D3M: Li, Tang, Wu, and Liao (2019)
- DeepDDM: Li, Xiang, Xu (2020); Mercier, Gratton, Boudier (arXiv 2021); Li, Wang, Cui, Xiang, Xu (2023); Sun, Xu, Yi (arXiv 2022, arXiv 2023)
- Schwarz Domain Decomposition Algorithm for PINNs: Kim, Yang (2022, arXiv 2022)
- FBPINNs: Moseley, Markham, and Nissen-Meyer (2023); Dolean, Heinlein, Mishra, Moseley (2024, subm. 2023 / arXiv:2306.05486); Heinlein, Howard, Beecroft, Stinis (subm. 2024 / arXiv:2401.07888)

An overview of the state-of-the-art in early 2021:



A. Heinlein, A. Klawonn, M. Lanser, J. Weber

Combining machine learning and domain decomposition methods for the solution of partial differential equations — A review

GAMM-Mitteilungen. 2021.

An overview of the state-of-the-art in the end of 2023:



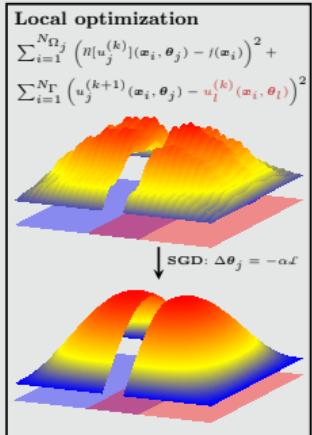
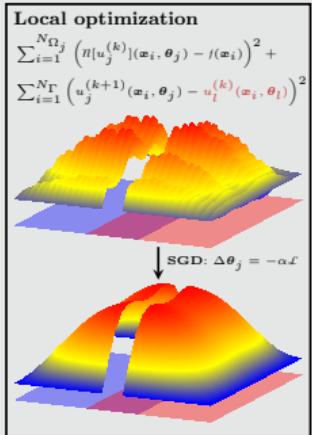
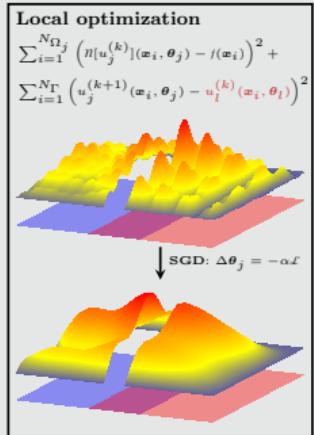
A. Klawonn, M. Lanser, J. Weber

Machine learning and domain decomposition methods – a survey

arXiv:2312.14050. 2023

Combining Schwarz Methods with Neural Network-Based Discretizations

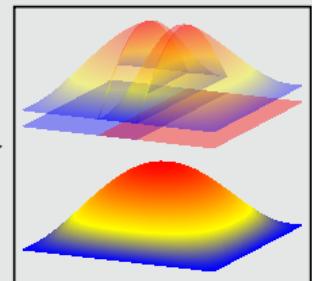
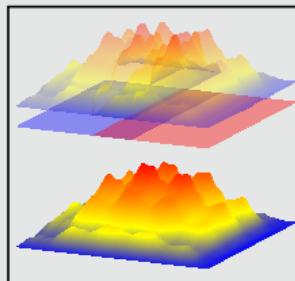
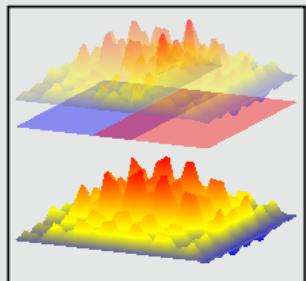
Approach 1 – Classical Schwarz iteration



Schwarz iteration

$$\begin{aligned} \Delta u_j^{(k+1)} &= f && \text{in } \Omega_j \\ u_j^{(k+1)} &= u_l^{(k)} && \text{on } \Gamma_j \end{aligned}$$

Approach 2 – Via the neural network architecture



Approach 1

**Deep learning-based domain
decomposition method**

Deep Learning-Based Domain Decomposition Method (DeepDDM)

Li, Xiang, Xu. Deep domain decomposition method: Elliptic problems. PMLR (2020)

DeepDDM for Overlapping Schwarz

In the **DeepDDM method**, we train **local networks** u_j using a **local loss function** on each subdomain Ω_j

$$\mathcal{L}_j(\theta_j) := \mathcal{L}_{\Omega_j}(\theta_j) + \mathcal{L}_{\partial\Omega_j \setminus \Gamma_j}(\theta_j) + \mathcal{L}_{\Gamma_j}(\theta_j),$$

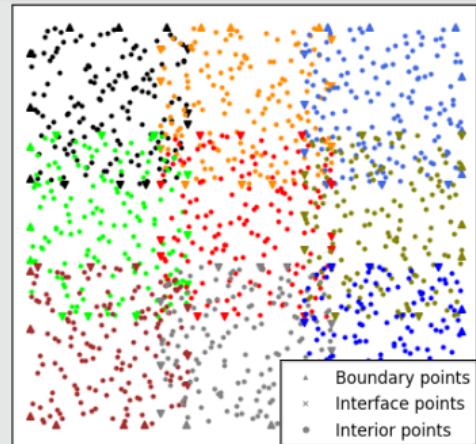
with **volume, boundary, and interface jump terms**:

$$\mathcal{L}_{\Omega_j}(\theta_j) := \frac{1}{N_{\Omega_j}} \sum_{i=1}^{N_{\Omega_j}} (n(u_j(\mathbf{x}_i, \theta_j)) - f(\mathbf{x}_i))^2$$

$$\mathcal{L}_{\partial\Omega_j \setminus \Gamma_j}(\theta_j) := \frac{1}{N_{\partial\Omega_j}} \sum_{i=1}^{N_{\partial\Omega_j}} (\mathcal{B}(u_j(\hat{\mathbf{x}}_i, \theta_j)) - g(\hat{\mathbf{x}}_i))^2$$

$$\mathcal{L}_{\Gamma_j}(\theta_j) := \frac{1}{N_{\Gamma_j}} \sum_{i=1}^{N_{\Gamma_j}} (\mathcal{D}(u_j(\tilde{\mathbf{x}}_i, \theta_j)) - \mathcal{D}(u_l(\tilde{\mathbf{x}}_i, \theta_j)))^2$$

Overl. domain decomposition



Algorithm 1: DeepDDM for Ω_j

Data: Sampling points X_j , initial network parameters θ_j^0

while convergence (local network & interface values) not reached **do**

Train local network u_j ;

Communicate & update interface values $\mathcal{D}(u_l(\tilde{\mathbf{x}}_i; \theta_j))$ from other subdomains Ω_l ;

end

Numerical Experiments

Strong scaling

Fix the problem complexity & increase the model capacity.

Optimal scaling: improving the convergence rate and/or accuracy at the same rate as the increase of model capacity.

Let first consider a **strong scaling study** for a **two-dimensional Laplacian model problem**:

$$\begin{aligned}-\Delta u &= 1 \quad \text{in } \Omega, \\ u &= 0 \quad \text{on } \partial\Omega.\end{aligned}$$

We increase the model capacity by **increasing the number of subdomains**.

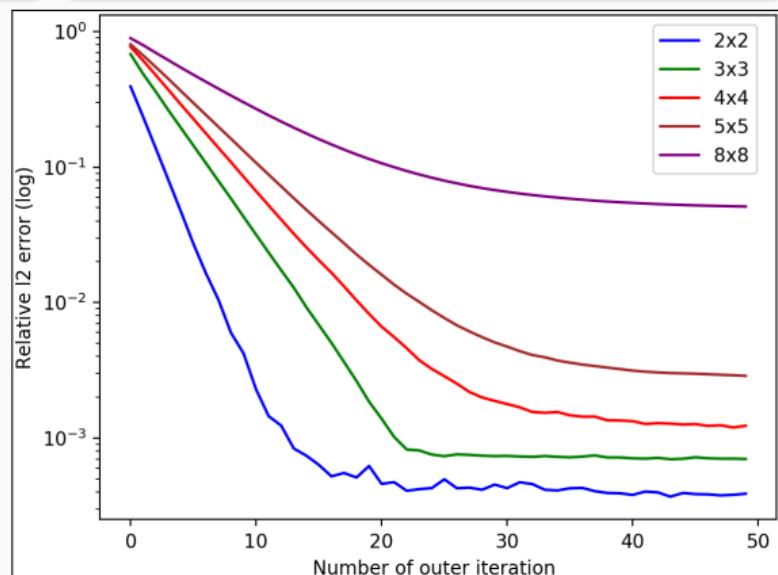
Scaling issue

We observe that the performance of the DeepDDM method deteriorates.

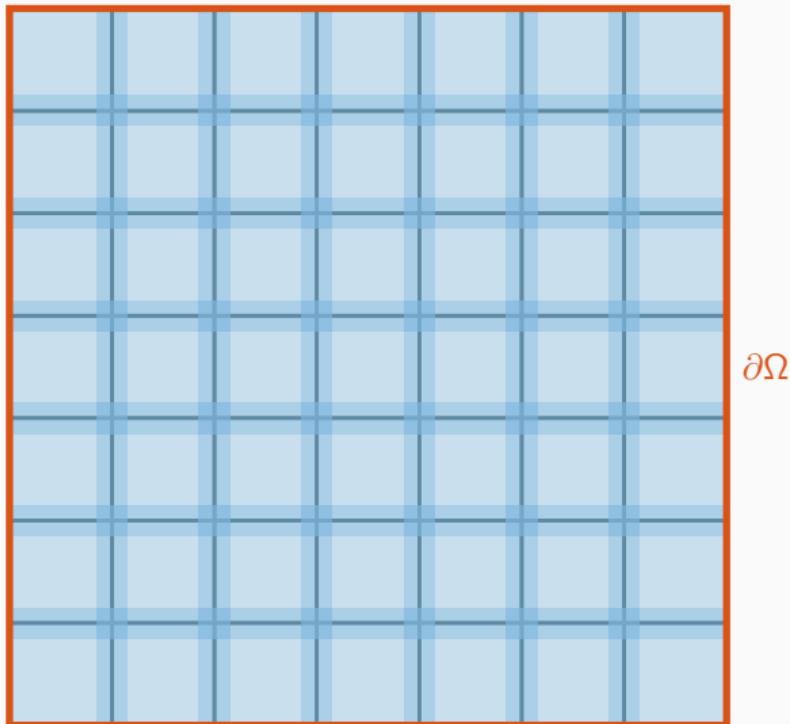
Weak scaling

Increase the problem complexity & the model capacity at the same rate.

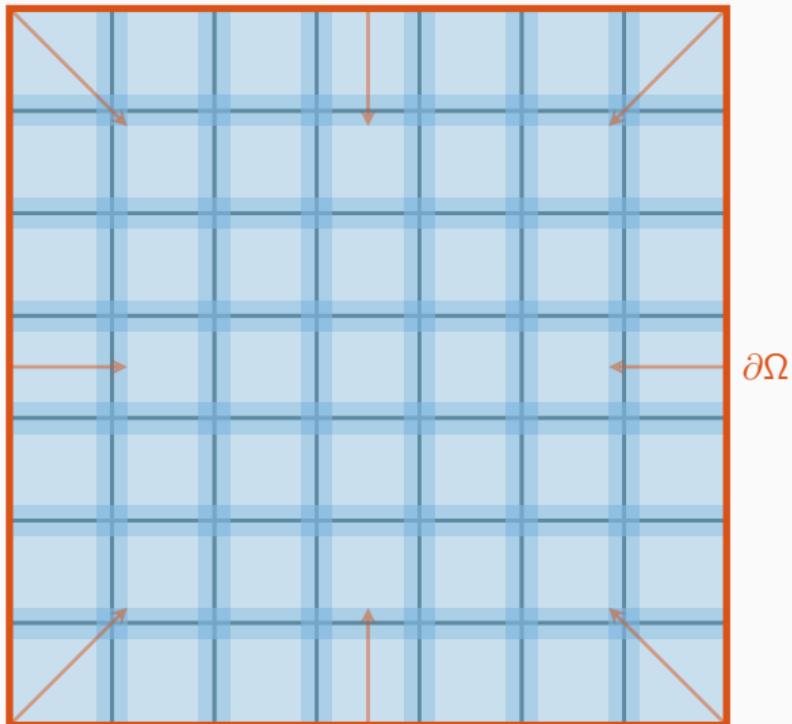
Optimal scaling: constant convergence rate and/or accuracy to stay approximately constant.



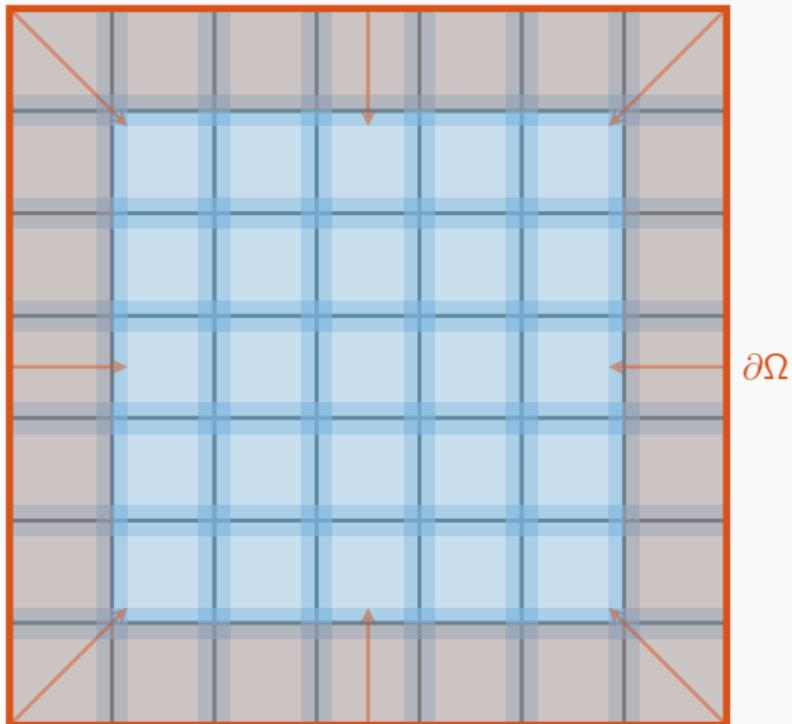
Transport of Information One-Level Overlapping Schwarz Methods



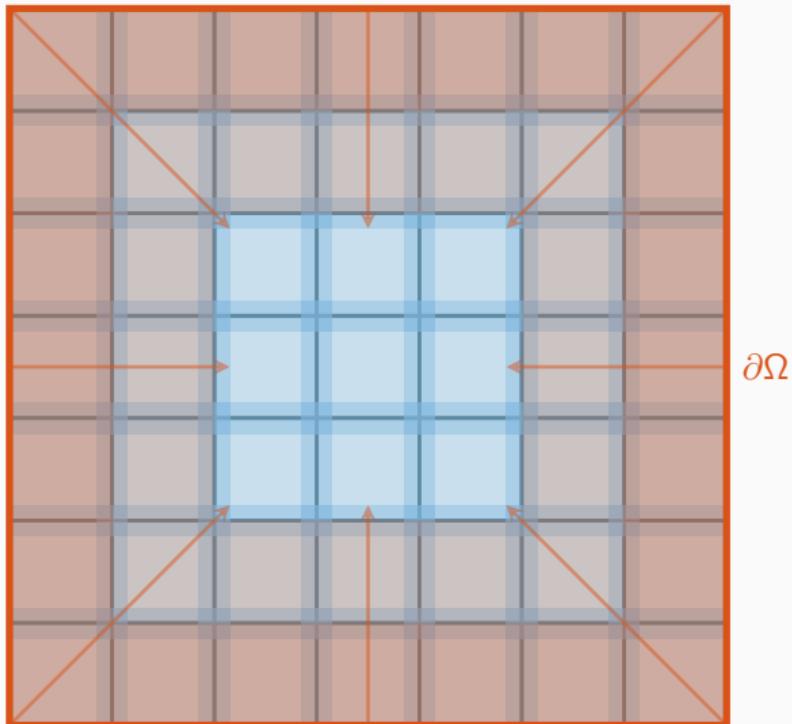
Transport of Information One-Level Overlapping Schwarz Methods



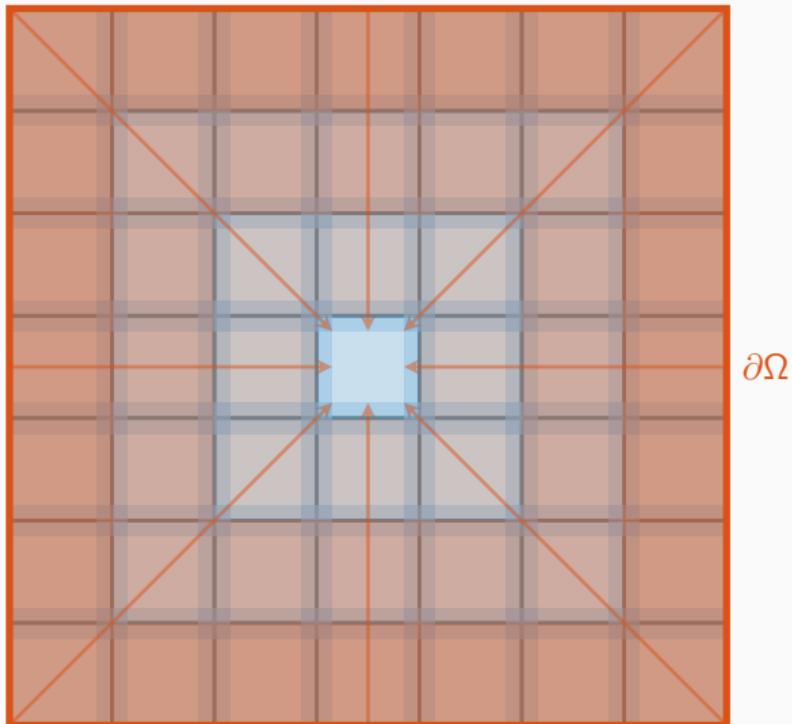
Transport of Information One-Level Overlapping Schwarz Methods



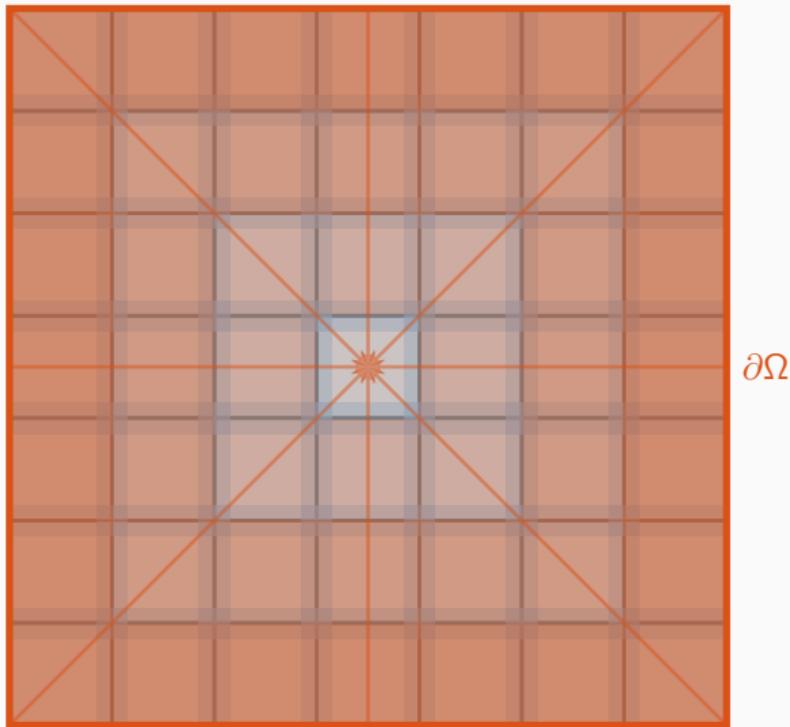
Transport of Information One-Level Overlapping Schwarz Methods



Transport of Information One-Level Overlapping Schwarz Methods



Transport of Information One-Level Overlapping Schwarz Methods



Information (in particular, boundary data) is **only exchanged via the overlapping regions**, leading to **slow convergence** → establish a faster / global transport of information.

Fast Transport of Information via a Coarse Level

Coarse space for the DeepDDM method

- Sparse sampling $\mathbf{X}_0 = \{\mathbf{x}_i^0\}_i$ over the whole domain Ω
- Train a **coarse network** (global PINN) u_0 with **additional loss term**

$$\lambda_f \frac{1}{N_0} \sum_{\mathbf{x}_i^0 \in \mathbf{X}_0} \left(u_0(\mathbf{x}_i^0) - \sum_{j=1}^J E_j(\chi_j u_j(\mathbf{x}_i^0)) \right)^2$$

for incorporating information from the first level. Here,

- E_j extension by zero outside Ω_j
- χ_j local partition of unity function
- Incorporate coarse information into the loss for the local subdomain Ω_j :

$$\frac{1}{N_{\Gamma_j}} \sum_{i=1}^{N_{\Gamma_j}} \left(\mathcal{D}(u_j(\tilde{\mathbf{x}}_i; \theta_j)) - W_j^i \right)^2$$

with $W_j^i = \mathcal{D}(\lambda_c u_l(\tilde{\mathbf{x}}_i) + (1 - \lambda_c) u_0(\tilde{\mathbf{x}}_i))$.

Algorithm 2: Two-level DeepDDM

Data: Sampling points X_j and coarse sampling points X_0 , initial network parameters θ_j^0 , weight parameters λ_f and λ_c

while convergence (local network & interface values) not reached **do**

Train local network u_j ;

Communicate & compute

$\sum_{j=1}^J E_j(\chi_j u_j(\mathbf{x}_i^0))$ for each $\mathbf{x}_i^0 \in \mathbf{X}_0$;

Train coarse network u_0 ;

Communicate & update interface values $\mathcal{D}(u_l(\tilde{\mathbf{x}}_i; \theta_j))$ from other subdomains Ω_l ;

 (**Optional**) **Update** λ_f and λ_c based on heuristic strategy;

end

2D Poisson Equation – Problem Setup

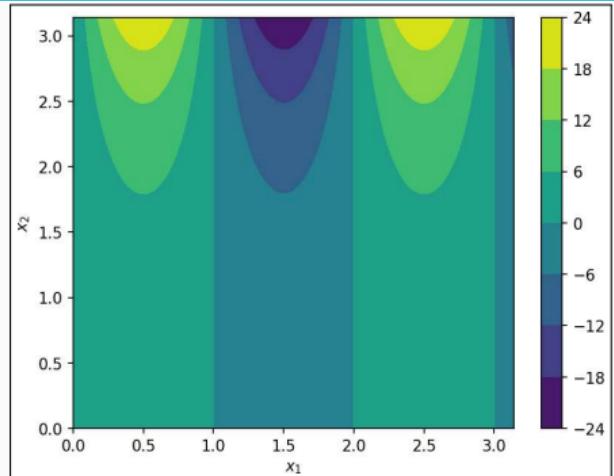
Model problem:

$$\begin{aligned}\Delta u = f &\quad \text{in } \Omega = [0, \pi] \times [0, 1], \\ u = g &\quad \text{on } \partial\Omega.\end{aligned}$$

We choose f and g such that the exact solution is

$$u(\mathbf{x}) = \sin(\alpha\pi x_1) e^{x_2},$$

where α is an integer.

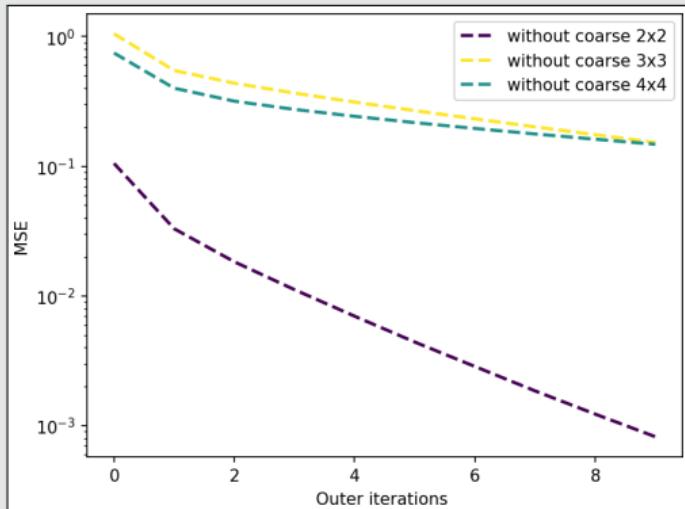


Training setup

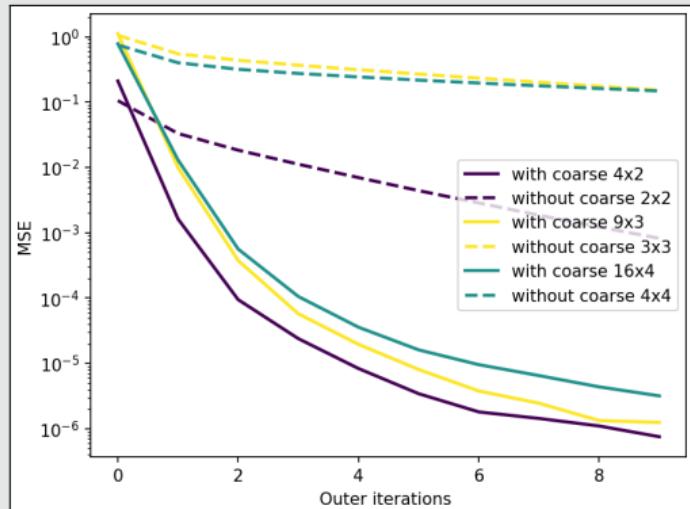
- **Strong scaling:** Latin hypercube sampling for training points with $N_\Omega = 30\,000$ and $N_{\partial\Omega} = N_\Gamma = 16\,000$.
- **Weak scaling:** Latin hypercube sampling for training points with $N_\Omega = 4\,000$ and $N_{\partial\Omega} = N_\Gamma = 1\,500$ per subdomain.
- Each network is composed of two hidden layers with 30 neurons
- Optimization of local/coarse networks: 2500 epochs using the Adam optimizer with initial learning rate $2 \cdot 10^{-4}$ and exp. decay of 0.999 every 100 epochs.
- Codes implemented in TENSORFLOW2 (v2.2.0) run on a single NVIDIA GeForce GTX 1080 Ti.
- The overlap is set to 30% of the subdomain diameter

2D Poisson Equation – Weak Scaling

One-level DeepDDM



Two-level DeepDDM



→ Adding a coarse level fixes the scaling issue.

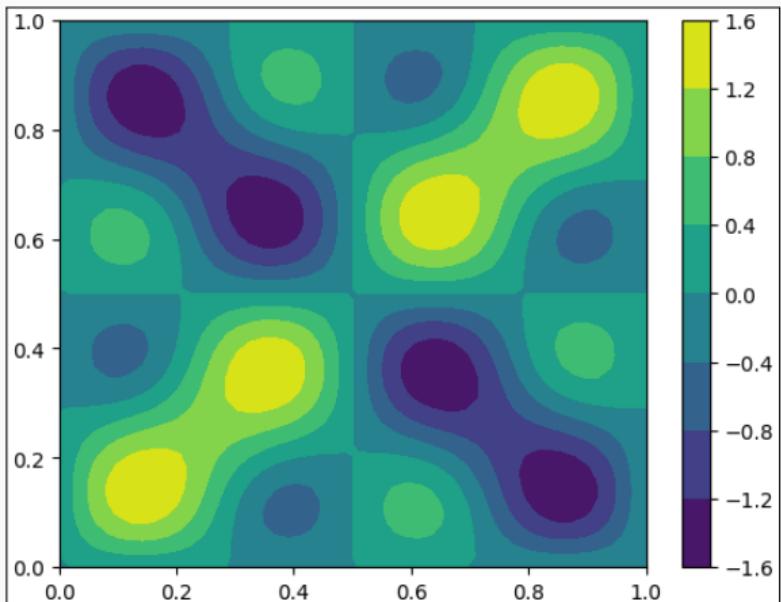
2D Poisson Equation With Variable Frequency

Model problem:

$$\begin{aligned}\Delta u = f &\quad \text{in } \Omega = [0, \pi] \times [0, 1], \\ u = g &\quad \text{on } \partial\Omega.\end{aligned}$$

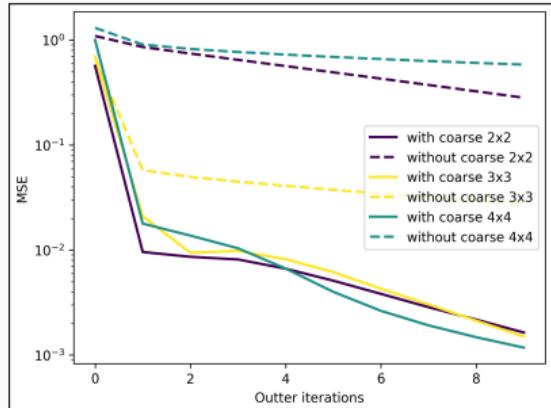
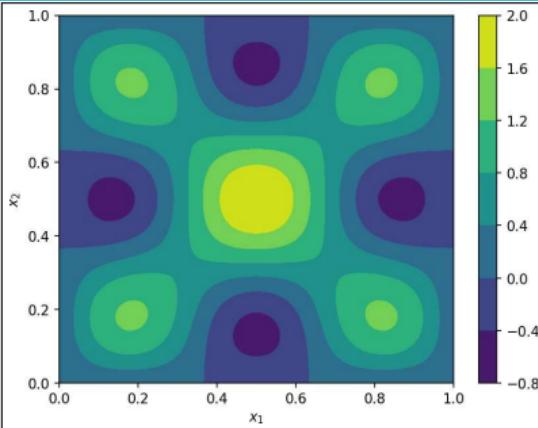
We choose f and g such that exact solution is:

$$\begin{aligned}u(\mathbf{x}) = \sin(w_1 \pi x_1) \sin(w_1 \pi x_2) \\ + \sin(w_2 \pi x_1) \sin(w_2 \pi x_2)\end{aligned}$$

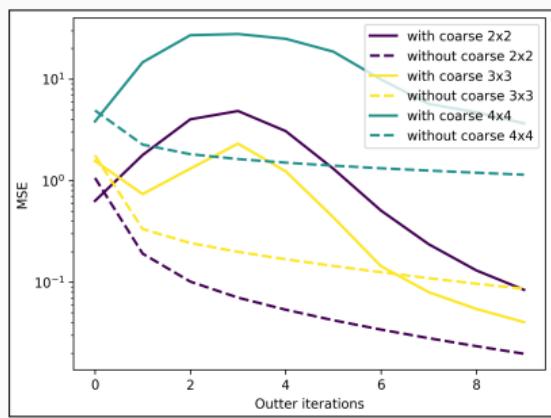
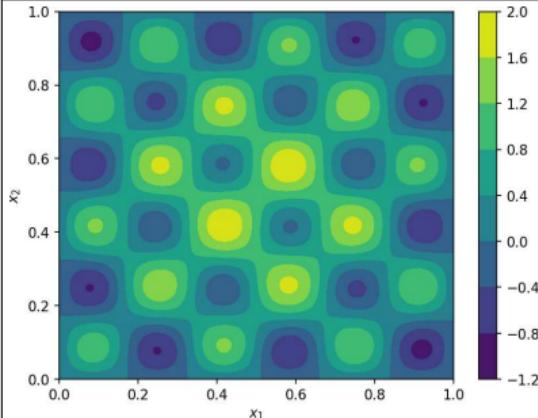


2D Poisson Equation With Variable Frequency – Weak Scaling

Low frequency test:
 $\omega_1 = 1$ and $\omega_2 = 3$

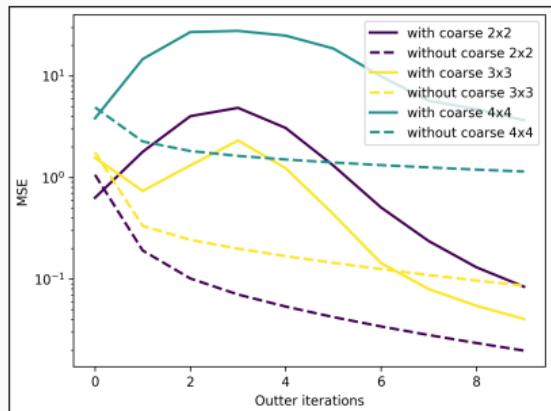
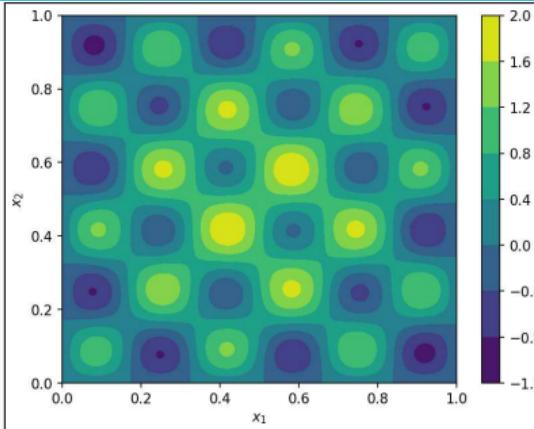


Higher frequency test:
 $\omega_1 = 1$ and $\omega_2 = 6$



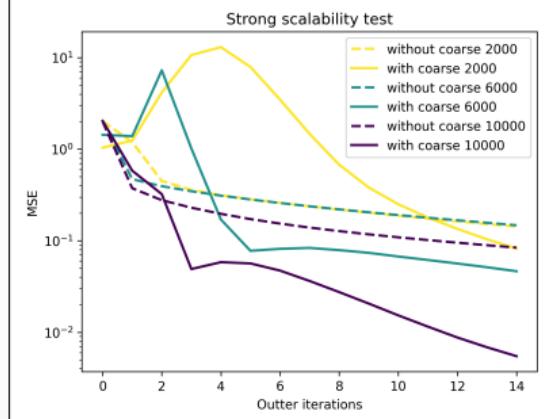
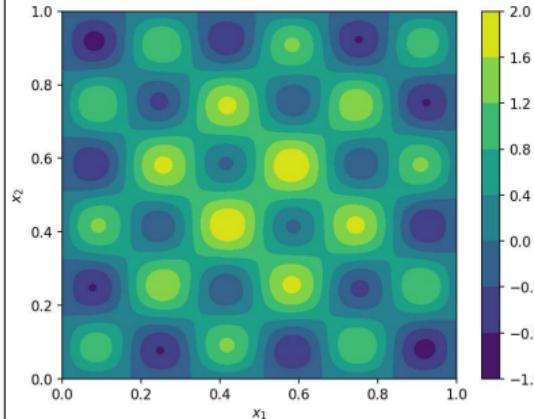
2D Poisson Equation With Variable Frequency – Weak Scaling

Higher frequency test:
 $\omega_1 = 1$ and $\omega_2 = 6$



Hyper parameter tuning

- ↑ # epochs for each sub problem
- ↑ # outer Schwarz iterations



Approach 2

**Multilevel domain decomposition-based
architectures for physics-informed neural
networks**

Finite Basis Physics-Informed Neural Networks (FBPINNs)

In the **finite basis physics informed neural network (FBPINNs) method** introduced in **Moseley, Markham, and Nissen-Meyer (2023)**, we solve the **partial differential equation**

$$\mathcal{N}[u](x) = f(x) \quad \text{in } \Omega$$

using the **PINN** approach and **hard enforcement of the boundary conditions**, similar to **Lagaris et al. (1998)**.

FBPINNs use the **network architecture**

$$u(\theta_1, \dots, \theta_J) = \mathcal{C} \sum_{j=1}^J \omega_j u_j(\theta_j)$$

and the **loss function**

$$\mathcal{L}(\theta_1, \dots, \theta_J) = \frac{1}{N} \sum_{i=1}^N \left(n[\mathcal{C} \sum_{x_i \in \Omega_j} \omega_j u_j](\mathbf{x}_i, \theta_j) - f(\mathbf{x}_i) \right)^2.$$

- **Overlapping DD:** $\Omega = \bigcup_{j=1}^J \Omega_j$
- **Window functions** ω_j with $\text{supp}(\omega_j) \subset \Omega_j$ and $\sum_{j=1}^J \omega_j \equiv 1$ on Ω

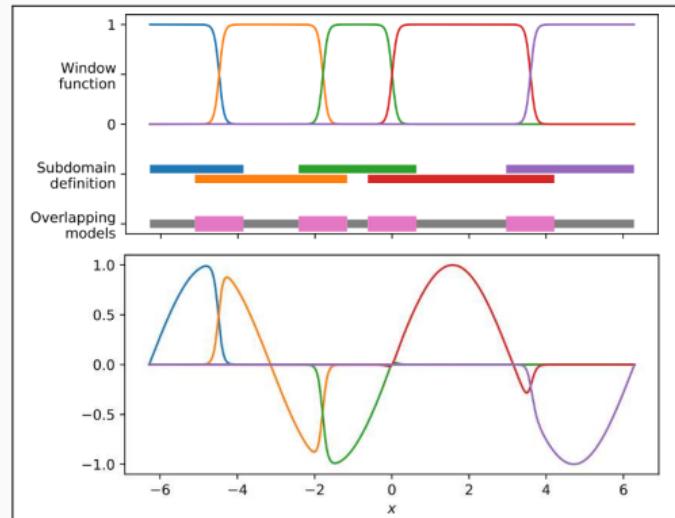
Hard enf. of boundary conditions

Loss function

$$\mathcal{L}(\theta) = \frac{1}{N} \sum_{i=1}^N (n[\mathcal{C} u](\mathbf{x}_i, \theta) - f(\mathbf{x}_i))^2,$$

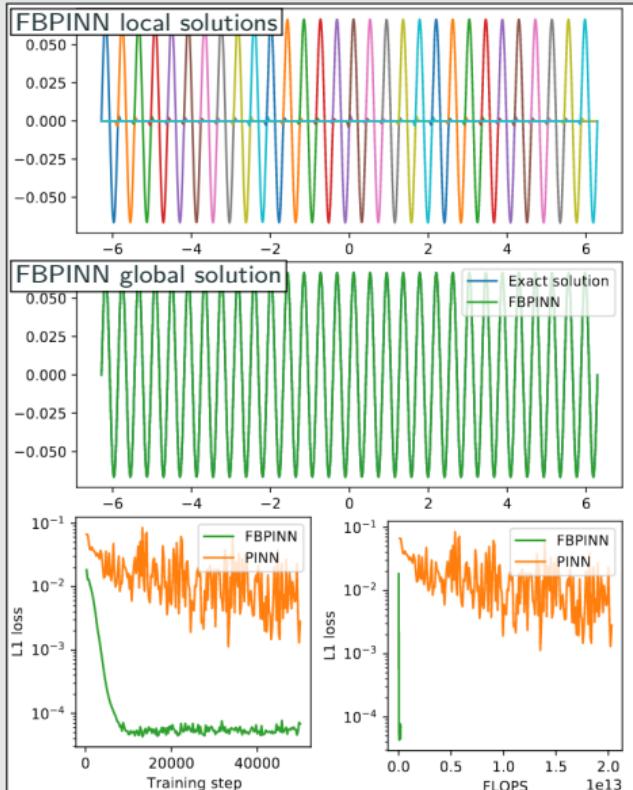
with constraining operator \mathcal{C} , which **explicitly enforces the boundary conditions**.

→ Often **improves training performance**



Numerical Results for FBPINNs

PINN vs FBPINN (Moseley et al. (2023))



Scalability of FBPINNs

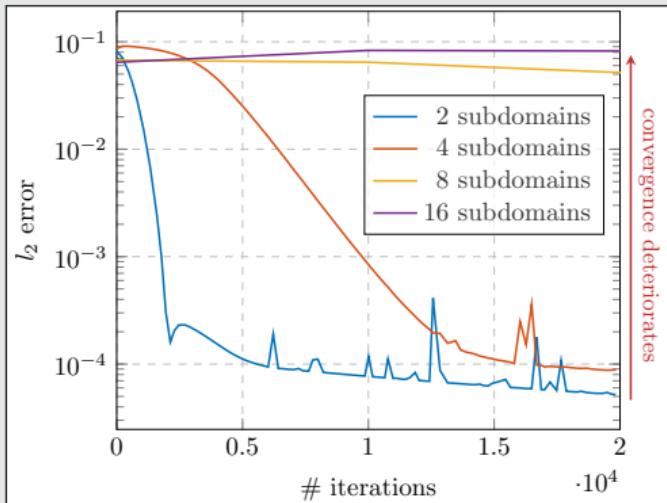
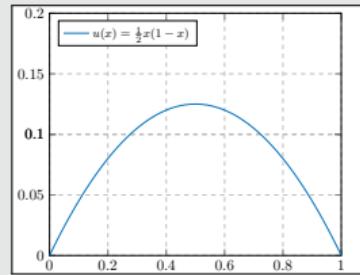
Consider the simple boundary value problem

$$-u'' = 1 \text{ in } [0, 1],$$

$$u(0) = u(1) = 0,$$

which has the solution

$$u(x) = 1/2x(1-x).$$



Multi-Level FBPINN Algorithm

We introduce a **hierarchy of L overlapping domain decompositions**

$$\Omega = \bigcup_{j=1}^{J^{(l)}} \Omega_j^{(l)}$$

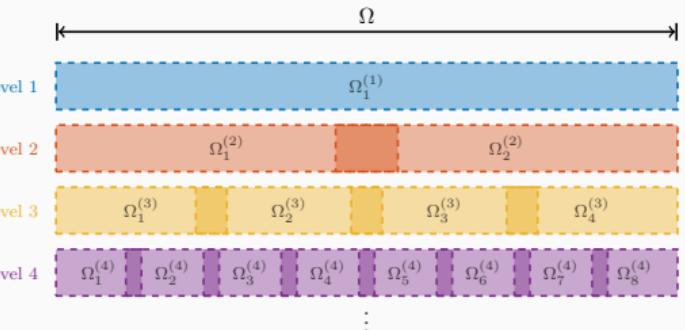
and corresponding window functions $\omega_j^{(l)}$ with

$$\text{supp}(\omega_j^{(l)}) \subset \Omega_j^{(l)} \text{ and } \sum_{j=1}^{J^{(l)}} \omega_j^{(l)} \equiv 1 \text{ on } \Omega.$$

This yields the **L -level FBPINN algorithm**:

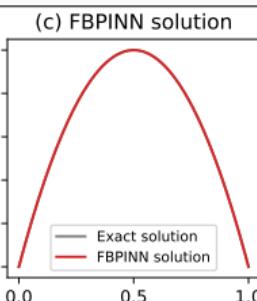
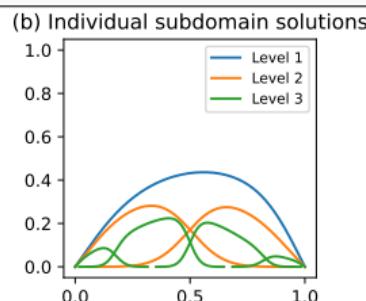
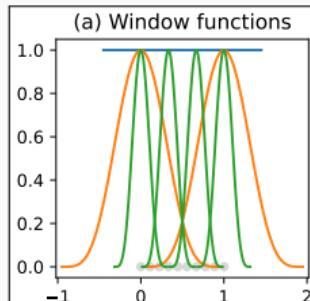
L -level network architecture

$$u(\theta_1^{(1)}, \dots, \theta_{J^{(L)}}^{(L)}) = \mathcal{C} \left(\sum_{l=1}^L \sum_{j=1}^{J^{(l)}} \omega_j^{(l)} u_j^{(l)}(\theta_j^{(l)}) \right)$$



Loss function

$$\mathcal{L} = \frac{1}{N} \sum_{i=1}^N \left(n[\mathcal{C} \sum_{x_i \in \Omega_j^{(l)}} \omega_j^{(l)} u_j^{(l)}](x_i, \theta_j^{(l)}) - f(x_i) \right)^2$$



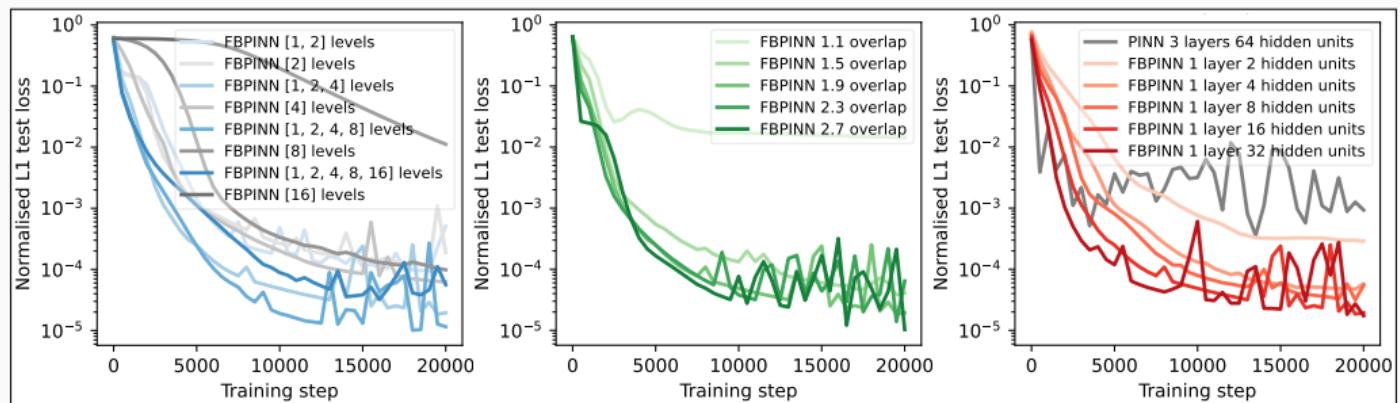
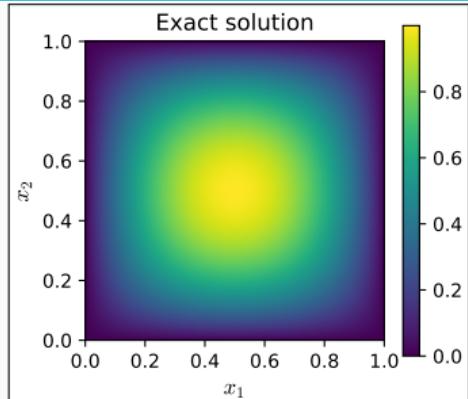
Multilevel FBPINNs – 2D Laplace

Let us consider the **simple two-dimensional boundary value problem**

$$\begin{aligned} -\Delta u &= 32(x(1-x) + y(1-y)) \quad \text{in } \Omega = [0, 1]^2, \\ u &= 0 \quad \text{on } \partial\Omega, \end{aligned}$$

which has the **solution**

$$u(x, y) = 16(x(1-x)y(1-y)).$$



Cf. Dolean, Heinlein, Mishra, Moseley (submitted 2023 / arXiv:2306.05486).

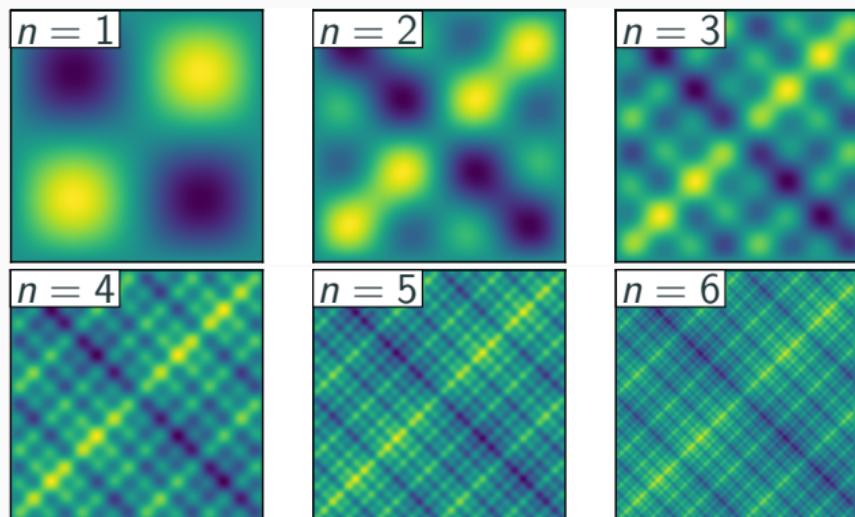
Multi-Frequency Problem

Let us now consider the **two-dimensional multi-frequency Laplace boundary value problem**

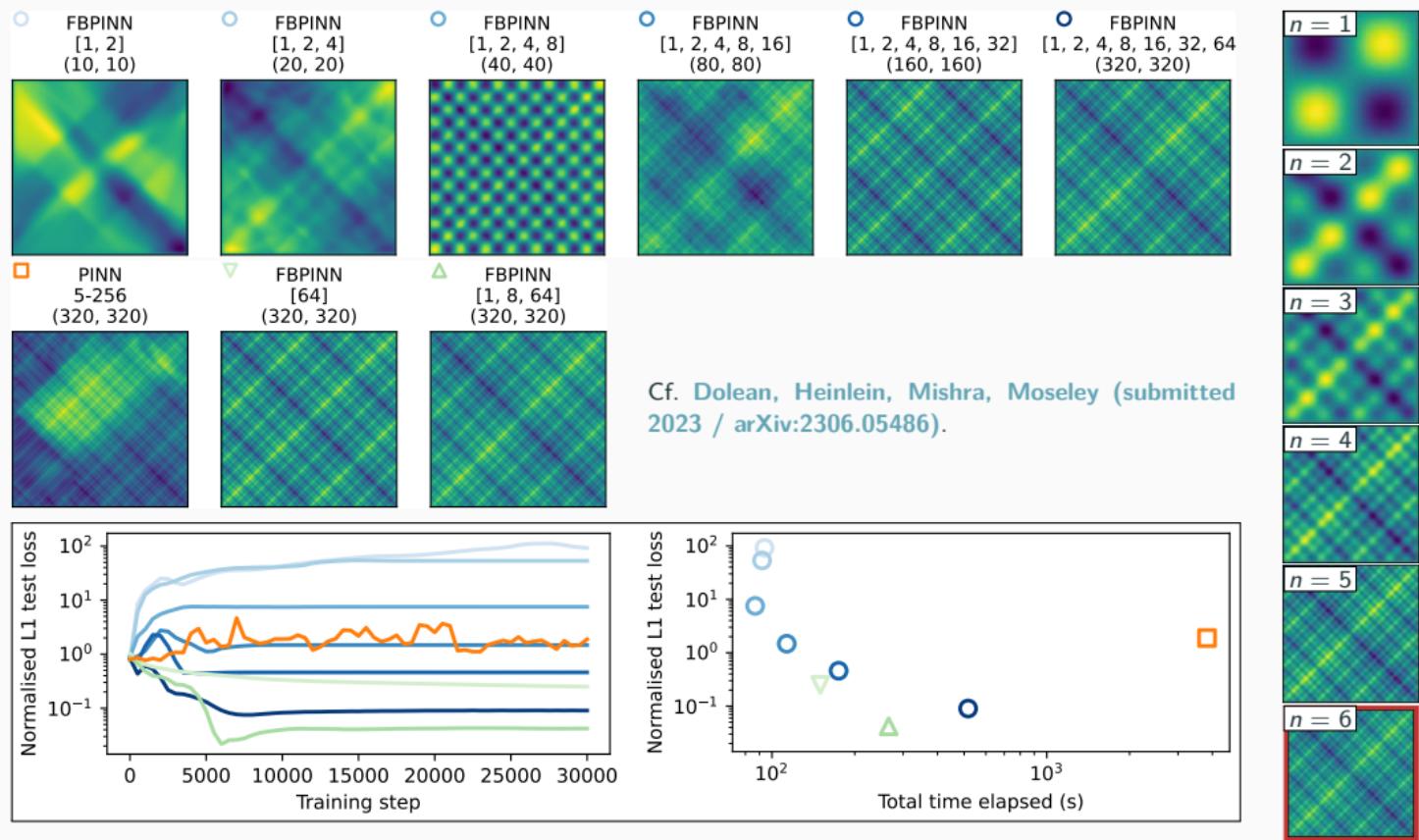
$$\begin{aligned} -\Delta u &= 2 \sum_{i=1}^n (\omega_i \pi)^2 \sin(\omega_i \pi x) \sin(\omega_i \pi y) && \text{in } \Omega = [0, 1]^2, \\ u &= 0 && \text{on } \partial\Omega, \end{aligned}$$

with $\omega_i = 2^i$.

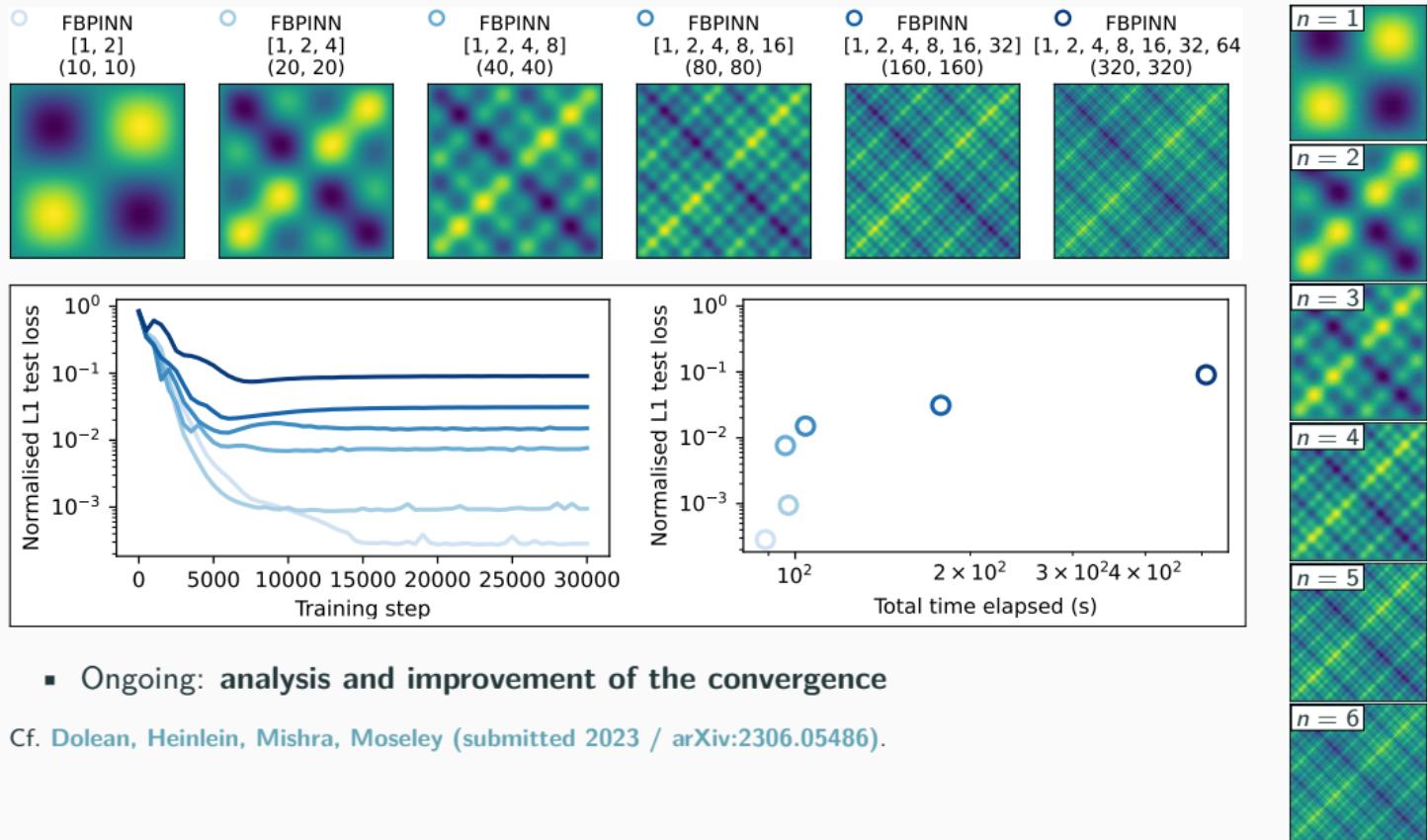
For increasing values of n , we obtain the **analytical solutions**:



Multi-Level FBPINNs for a Multi-Frequency Problem – Strong Scaling



Multi-Level FBPINNs for a Multi-Frequency Problem – Weak Scaling



- Ongoing: analysis and improvement of the convergence

Cf. [Dolean, Heinlein, Mishra, Moseley \(submitted 2023 / arXiv:2306.05486\)](#).

Helmholtz Problem

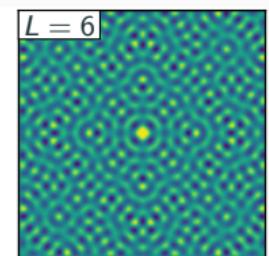
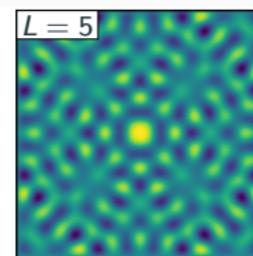
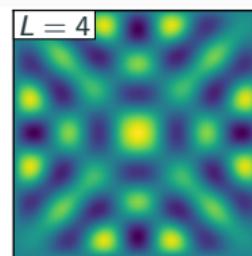
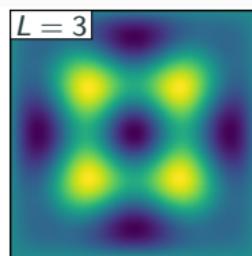
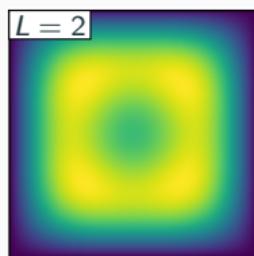
Finally, let us consider the **two-dimensional Helmholtz boundary value problem**

$$\Delta u - k^2 u = f \quad \text{in } \Omega = [0, 1]^2,$$

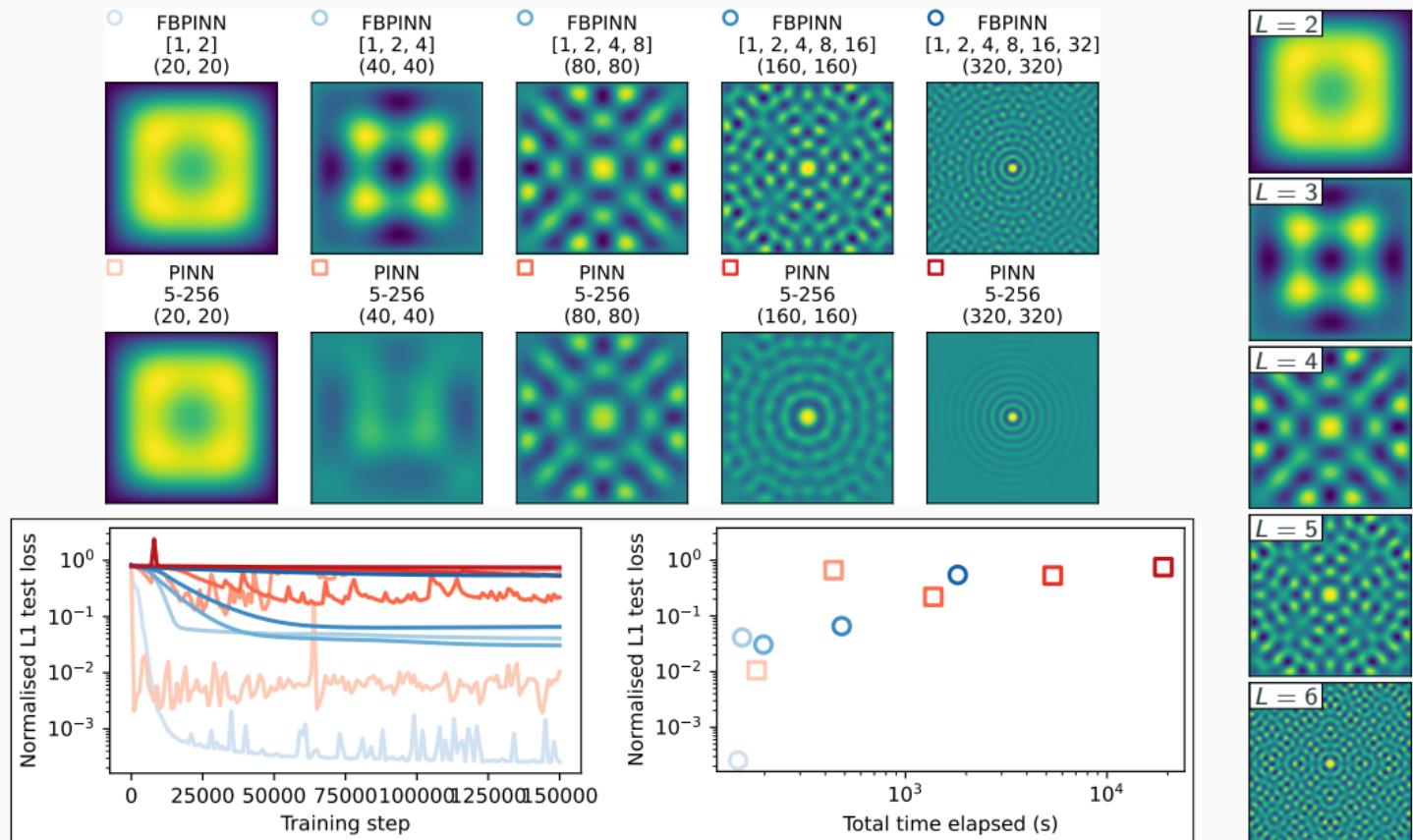
$$u = 0 \quad \text{on } \partial\Omega,$$

$$f(x) = e^{-\frac{1}{2}(\|x-0.5\|/\sigma)^2}.$$

With $k = 2^L \pi / 1.6$ and $\sigma = 0.8 / 2^L$, we obtain the **solutions**:



Multi-Level FBPINNs for the Helmholtz Problem – Weak Scaling

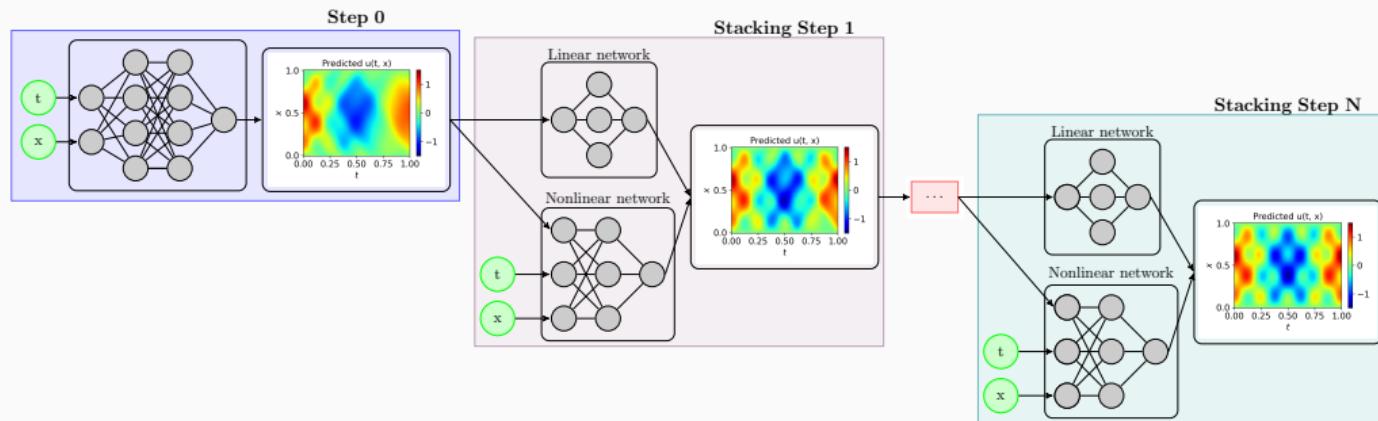


Multifidelity domain decomposition-based physics-informed neural networks for time-dependent problems

Stacking Multifidelity PINNs

In the **stacking multifidelity PINNs approach** introduced in [Howard, Murphy, Ahmed, Stinis \(arXiv 2023\)](#), multiple PINNs are trained in a recursive way. In each step, a model u^{MF} is trained as a corrector for the previous model u^{SF} :

$$u^{MF}(\mathbf{x}, \theta^{MF}) = (1 - |\alpha|) u_{\text{linear}}^{MF}(\mathbf{x}, \theta^{MF}, u^{SF}) + |\alpha| u_{\text{nonlinear}}^{MF}(\mathbf{x}, \theta^{MF}, u^{SF})$$



Related works (non-exhaustive list)

- **Cokriging & multifidelity Gaussian process regression:** E.g., [Wackernagel \(1995\)](#); [Perdikaris et al. \(2017\)](#); [Babaei et al. \(2020\)](#)
- **Multifidelity PINNs & DeepONet:** [Meng and Karniadakis \(2020\)](#); [Howard, Fu, and Stinis \(arXiv 2023\)](#); [Howard, Perego, Karniadakis, Stinis \(2023\)](#); [Murphy, Ahmed, Stinis \(arXiv 2023\)](#)

Stacking Multifidelity FBPINNs

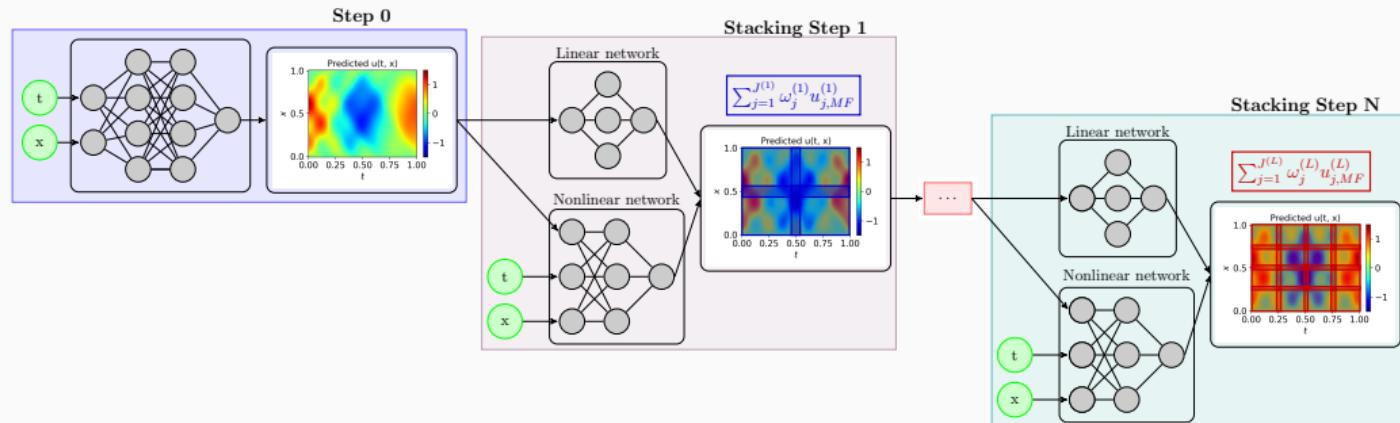
In Heinlein, Howard, Beecroft, and Stinis (subm. 2024 / arXiv:2401.07888), we **combine stacking multifidelity PINNs with FBPINNs** by using an FBPINN model in each stacking step. The resulting **stacking multifidelity FBPINN** in step I / on level I reads

$$u^{(I)}(\mathbf{x}, \theta^{(I)}) = \sum_{j=1}^{J^{(I)}} \omega_j^{(I)} u_{j,MF}^{(I)}(\mathbf{x}, \theta_j^{(I)}, u^{(I-1)}),$$

where

$$u_{j,MF}^{(I)}(\mathbf{x}, \theta_j^{(I)}) = (1 - |\alpha|) u_{j,\text{linear}}^{(I)}(\mathbf{x}, \theta_j^{(I)}, u^{(I-1)}) + |\alpha| u_{j,\text{nonlinear}}^{(I)}(\mathbf{x}, \theta_j^{(I)}, u^{(I-1)}).$$

This corresponds to a **one-way sequential coupling** of the levels.

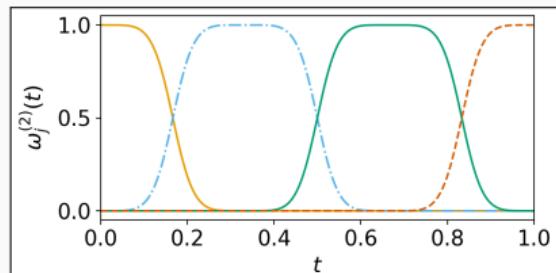


Numerical Results – Pendulum Problem

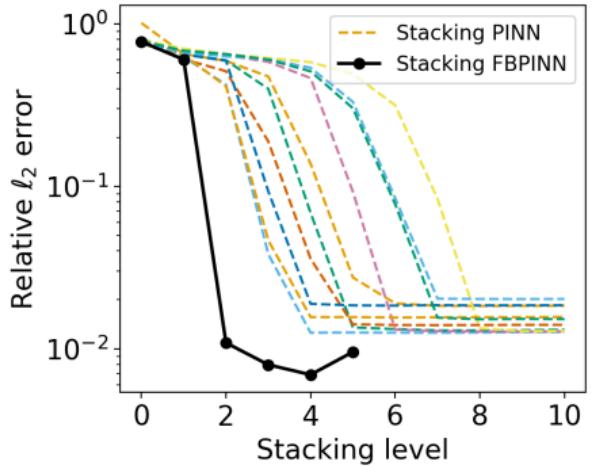
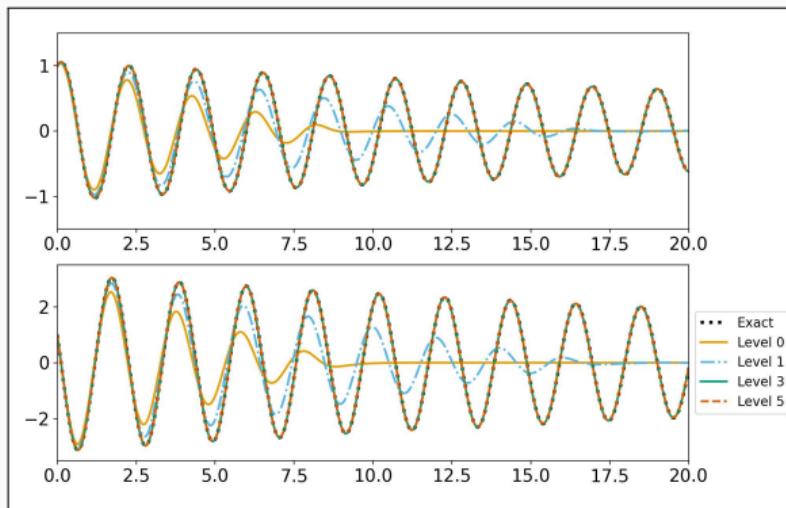
First, we consider a **pendulum problem** and compare the stacking multifidelity PINN and FBPINN approaches:

$$\begin{aligned}\frac{d\beta_1}{dt} &= \beta_2, \\ \frac{d\beta_2}{dt} &= -\frac{b}{m}\beta_2 - \frac{g}{L} \sin(\beta_1)\end{aligned}$$

with $m = L = 1$, $b = 0.05$, $g = 9.81$, and $T = 20$.



Exemplary partition of unity in time

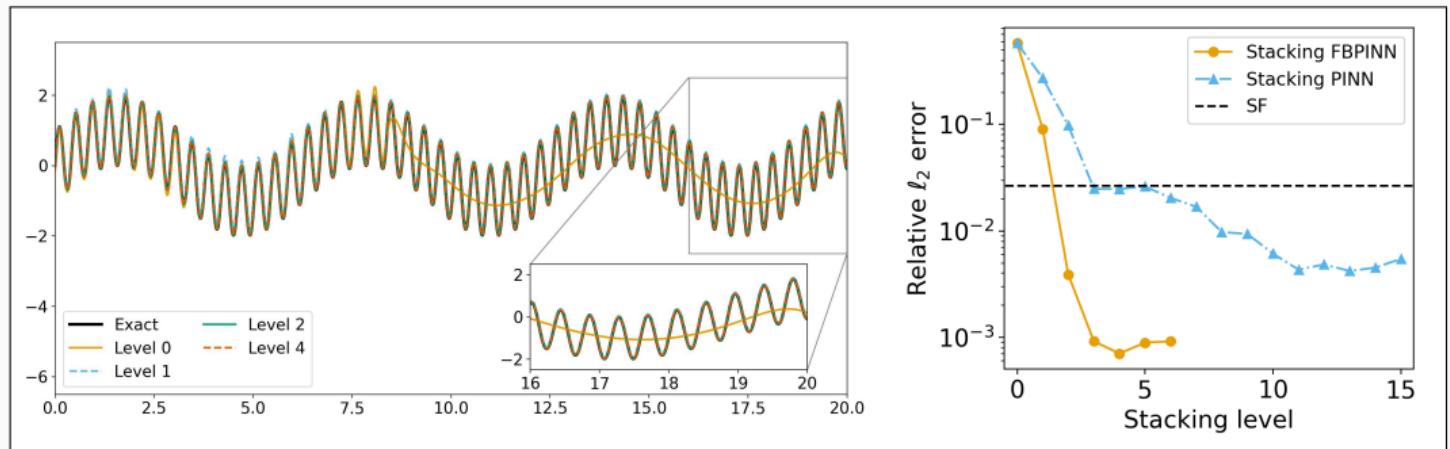


Numerical Results – Two-Frequency Problem

Second, we consider a **two-frequency problem**:

$$\begin{aligned}\frac{ds}{dx} &= \omega_1 \cos(\omega_1 x) + \omega_2 \cos(\omega_2 x), \\ s(0) &= 0,\end{aligned}$$

on domain $\Omega = [0, 20]$ with $\omega_1 = 1$ and $\omega_2 = 15$.

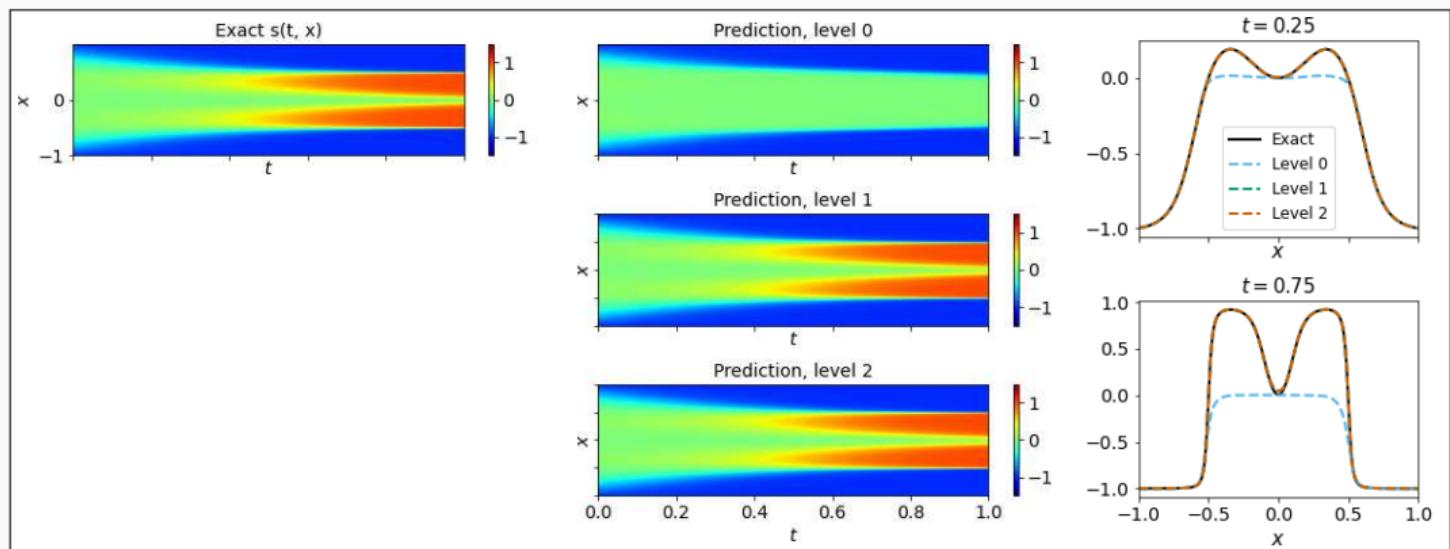


→ Due to the **multiscale structure of the problem**, the **improvements** due to the **multifidelity FBPINN approach** are **even stronger**.

Numerical Results – Allen–Cahn Equation

Finally, we consider the **Allen–Cahn equation**:

$$\begin{aligned}\dot{\vartheta}_t - 0.0001\vartheta_{xx} + 5\vartheta^3 - 5\vartheta &= 0, & t \in (0, 1], x \in [-1, 1], \\ \vartheta(x, 0) &= x^2 \cos(\pi x), & x \in [-1, 1], \\ \vartheta(x, t) &= \vartheta(-x, t), & t \in [0, 1], x = -1, x = 1, \\ \vartheta_x(x, t) &= \vartheta_x(-x, t), & t \in [0, 1], x = -1, x = 1.\end{aligned}$$



PINNs

- **Training of PINNs is often problematic** when:
 - scaling to large domains / high frequency solutions
 - multiple loss terms have to be balanced
- Convergence of PINNs has yet to be understood better

DeepDDM for PINNs

- The **DeepDDM method** is a **classical Schwarz iteration** with local PINN solver.
- **Scalability** is enabled by **adding a coarse level**.

Multilevel FBPINNs

- Schwarz domain decomposition architectures **improve the scalability of PINNs** to large domains / high frequencies, **keeping the complexity of the local networks low**
- As classical domain decomposition methods, **one-level FBPINNs** are **not scalable to large numbers of subdomains**; multilevel FBPINNs **enable scalability**.

Multifidelity stacking FBPINNs

- The combination of multifidelity stacking PINNs with FBPINNs yields **significant improvements in the accuracy and efficiency** for time-dependent problems.

Thank you for your attention!

Workshop on Computational and Mathematical Methods in Data Science 2024

Details

Date: April 24-26 2024

Location: Delft University of Technology

This workshop brings together scientists from mathematics, computer science, and application areas working on computational and mathematical methods in data science.

Confirmed invited speakers

- Christoph Brune (University of Twente)
- Victorita Dolean (TU Eindhoven)
- Thomas Richter (Otto von Guericke University Magdeburg)



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Delft

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COMiDS Workshop 2024



Workshop on Computational and Mathematical Methods in Data Science 2024
Delft University of Technology, April 25-26, 2024

About the Workshop

Welcome to the **Workshop on Computational and Mathematical Methods in Data Science 2024**. It is the 2024 edition of the annual workshop of the **GAMM Activity Group on "Computational and Mathematical Methods in Data Science" (COMiDS)** and is co-organized by the **Strategic Research Initiative "Bridging Numerical Analysis and Machine Learning"** of the **4TU Applied Mathematics Institute (AMI)**. The workshop will be hosted by **Delft University of Technology** and take place on April 25 and 26, 2024.

This workshop brings together scientists from mathematics, computer science, and application areas working on computational and mathematical methods in data science.

The meeting will be organized under the support of

- the **4TU Applied Mathematics Institute (AMI)** and
- the **TU Delft Institute for Computational Science and Engineering (ICSE)**

Registration is open!

Deadline: April 1st, 2024

More details: searhein.github.io/gamm-cominds-2024/

