

When One Level Is Not Enough

Multilevel Domain Decomposition Methods for Physics and Data-Driven Problems

Alexander Heinlein¹

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¹Delft University of Technology

Numerical Analysis and Machine Learning







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

4TU.AMI – SRI "Bridging Numerical Analysis and Machine Learning"

UNIVERSITY OF TWENTE.





Christoph Brune

Matthias Schlotthom







Alexander Heinlein

Deepesh Toshniwal



TU/e EINDHOVEN UNIVERSITY OF TECHNOLOGY





Victorita Dolean



Olga Mula



Schilders

Jemima Tabeart



Karen Veroy-Grepl



Matthias

Möller



Xiaodong Cheng

Alexander Heinlein (TU Delft)



1 Classical Schwarz Domain Decomposition Methods

2 Schwarz Domain Decomposition Preconditioners

Based on joint work with

Axel Klawonn and Jascha Knepper Mauro Perego and Siva Rajamanickam Oliver Rheinbach and Friederik Röver Olof Widlund (University of Cologne) (Sandia National Laboratories) (TU Bergakademie Freiberg) (New York University)

3 Domain Decomposition for Neural Networks

Based on joint work with

Eric Cyr Victorita Dolean Siddhartha Mishra Ben Moseley Corné Verburg (Sandia National Laboratories) (Eindhoven University of Technology) (ETH Zürich) (Imperial College London) (Delft University of Technology)

Classical Schwarz Domain Decomposition Methods

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



The Alternating Schwarz Algorithm

For the sake of simplicity, instead of the two-dimensional geometry,



0.2

0.4

0.6

0.8

1.0

$$-u'' = 1$$
, in [0,1], $u(0) = u(1) = 0$

We perform an alternating Schwarz iteration:



Figure 1: Iterate (left) and error (right) in iteration 0.

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We perform an alternating Schwarz iteration:



Figure 1: Iterate (left) and error (right) in iteration 1.

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We perform an alternating Schwarz iteration:



Figure 1: Iterate (left) and error (right) in iteration 2.

$$-u'' = 1$$
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We perform an alternating Schwarz iteration:



Figure 1: Iterate (left) and error (right) in iteration 3.

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We perform an alternating Schwarz iteration:



Figure 1: Iterate (left) and error (right) in iteration 4.

$$-u'' = 1$$
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We perform an alternating Schwarz iteration:



Figure 1: Iterate (left) and error (right) in iteration 5.

The alternating Schwarz algorithm is **sequential** because **each local boundary value problem** depends on the solution of the **previous Dirichlet problem**:

$$(D_1) \begin{cases} -\Delta u^{n+1/2} = f \quad \text{in } \Omega'_1, \\ u^{n+1/2} = \mathbf{u}^n \quad \text{on } \partial \Omega'_1 \\ u^{n+1/2} = \mathbf{u}^n \quad \text{on } \Omega \setminus \overline{\Omega'_1} \end{cases}$$
$$(D_2) \begin{cases} -\Delta u^{n+1} = f \quad \text{in } \Omega_2, \\ u^{n+1} = \mathbf{u}^{n+1/2} \quad \text{on } \partial \Omega'_2 \\ u^{n+1} = \mathbf{u}^{n+1/2} \quad \text{on } \Omega \setminus \overline{\Omega'_2} \end{cases}$$



Idea: For all red terms, we **use the values from the previous iteration**. Then, the both Dirichlet problem **can be solved at the same time**.

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The Parallel Schwarz Algorithm

The **parallel Schwarz algorithm** has been introduced by **Lions (1988)**. Here, we solve the local problems



Since u_1^n and u_2^n are both computed in the previous iteration, the problems can be solved independent of each other.

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This method is suitable for **parallel computing**!



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We perform a **parallel Schwarz iteration**:



Figure 2: Iterate (left) and error (right) in iteration 0.

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Figure 2: Iterate (left) and error (right) in iteration 1.

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Figure 2: Iterate (left) and error (right) in iteration 2.

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Figure 2: Iterate (left) and error (right) in iteration 5.

We investigate the convergence of the methods (using the alternating method as an example) depending on the **size of the overlap**:





Figure 3: Error in iteration 0.



Figure 3: Error in iteration 1.



Figure 3: Error in iteration 2.



Figure 3: Error in iteration 3.



Figure 3: Error in iteration 4.



Figure 3: Error in iteration 5.

Overlap 0.05

Overlap 0.1



Figure 3: Error in iteration 5.

 \Rightarrow A larger overlap leads to faster convergence.

Schwarz Domain Decomposition Preconditioners

Solvers for Partial Different Equations

Consider a diffusion model problem:

$$-\Delta u(x) = f \quad \text{in } \Omega = [0, 1]^2,$$
$$u = 0 \quad \text{on } \partial \Omega.$$

Discretization using finite elements yields a **sparse** system of linear equations

Ku = f.

The accuracy of the finite element solution depends on the refinement level of the mesh *h*: higher refinement \Rightarrow better accuracy.

Direct solvers

For fine meshes, solving the system using a direct solver is not feasible due to **superlinear complexity and memory cost**.

Iterative solvers

Iterative solvers are efficient for solving **sparse systems**, however, the **convergence rate depends on the condition number**: $\kappa(\mathbf{K}) = \frac{\lambda_{\max}(\mathbf{K})}{\lambda_{\min}(\mathbf{K})} \leq \frac{C}{b^2}$



Solvers for Partial Different Equations

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 \Rightarrow Introduce a preconditioner $M^{-1} \approx K^{-1}$ to improve convergence:

$$\boldsymbol{M}^{-1}\boldsymbol{K}\boldsymbol{u}=\boldsymbol{M}^{-1}\boldsymbol{f}$$



Two-Level Schwarz Preconditioners

Based on an overlapping domain decomposition, we define a one-level Schwarz operator

$$\boldsymbol{M}_{\text{OS-1}}^{-1}\boldsymbol{K} = \sum_{i=1}^{N} \boldsymbol{R}_{i}^{\top}\boldsymbol{K}_{i}^{-1}\boldsymbol{R}_{i}\boldsymbol{K}_{i}$$

where \mathbf{R}_i and \mathbf{R}_i^{\top} are restriction and prolongation operators corresponding to Ω'_i , and $\mathbf{K}_i := \mathbf{R}_i \mathbf{K} \mathbf{R}_i^{\top}$.

Condition number estimate:

$$\kappa\left(oldsymbol{M}_{\mathsf{OS-1}}^{-1}oldsymbol{K}
ight) \leq C\left(1+rac{1}{H\delta}
ight)$$

with subdomain size H and overlap width δ .














One-level Schwarz preconditioner





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Lagrangian coarse space





The two-level overlapping Schwarz operator reads

$$\boldsymbol{M}_{\text{OS-2}}^{-1}\boldsymbol{K} = \underbrace{\boldsymbol{\Phi}\boldsymbol{K}_{0}^{-1}\boldsymbol{\Phi}^{\top}\boldsymbol{K}}_{\text{coarse level - global}} + \underbrace{\sum_{i=1}^{N}\boldsymbol{R}_{i}^{\top}\boldsymbol{K}_{i}^{-1}\boldsymbol{R}_{i}\boldsymbol{K}}_{\text{first level - local}},$$

where Φ contains the coarse basis functions and $K_0 := \Phi^\top K \Phi$; cf., e.g., **Toselli, Widlund (2005)**. The construction of a Lagrangian coarse basis requires a coarse triangulation.

Condition number estimate:

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ight)$$

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FROSch (Fast and Robust Overlapping Schwarz) Framework in Trilinos





Software

- Object-oriented C++ domain decomposition solver framework with $\rm MPI\text{-}based$ distributed memory parallelization
- Part of TRILINOS with support for both parallel linear algebra packages EPETRA and TPETRA
- Node-level parallelization and performance portability on CPU and GPU architectures through KOKKOS and KOKKOSKERNELS
- Accessible through unified ${\rm TRILINOS}$ solver interface ${\rm STRATIMIKOS}$

Methodology

- Parallel scalable multi-level Schwarz domain decomposition preconditioners
- Algebraic construction based on the parallel distributed system matrix
- Extension-based coarse spaces

Team (active)

- Filipe Cumaru (TU Delft)
- Kyrill Ho (UCologne)
- Jascha Knepper (UCologne)
- Friederike Röver (TUBAF)
- Lea Saßmannshausen (UCologne)

- Alexander Heinlein (TU Delft)
- Axel Klawonn (UCologne)
- Siva Rajamanickam (SNL)
- Oliver Rheinbach (TUBAF)
- Ichitaro Yamazaki (SNL)

Overlapping domain decomposition

The overlapping subdomains are constructed by recursively adding layers of elements via the sparsity pattern of *K*.

The corresponding matrices

$$K_i = R_i K R_i^T$$

can easily be extracted from K.



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Coarse space – Example of Generalized Dryja–Smith–Widlund (GDSW)

1. Interface components



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2. Interface basis (partition of unity \times null space)

1. Interface components





For scalar elliptic problems, the null space consists only of constant functions.

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Weak Scalability up to 64 k MPI Ranks / 1.7 b Unknowns (3D Poisson; Juqueen)

GDSW vs RGDSW (reduced dimension)

Heinlein, Klawonn, Rheinbach, Widlund (2019).



Two-level vs three-level GDSW

Heinlein, Klawonn, Rheinbach, Röver (2019, 2020).



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35

0

100

1000

Cores

10000

100000

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FROSch Preconditioners for Land Ice Simulations



https://github.com/SNLComputation/Albany

The velocity of the ice sheet in Antarctica and Greenland is modeled by a **first-order-accurate Stokes approximation model**,

$$-\nabla \cdot (2\mu \dot{\epsilon}_1) + \rho g \frac{\partial s}{\partial x} = 0, \quad -\nabla \cdot (2\mu \dot{\epsilon}_2) + \rho g \frac{\partial s}{\partial y} = 0,$$



with a nonlinear viscosity model (Glen's law); cf., e.g., Blatter (1995) and Pattyn (2003).

	Ant	arctica (veloc	city)	Greenland (multiphysics vel. & temperature)			
	4 km resolu	tion, 20 layers	s, 35 m dofs	1-10 km resolution, 20 layers, 69 m dofs			
MPI ranks	avg. its	avg. setup	avg. solve	avg. its	avg. setup	avg. solve	
512	41.9 (11)	25.10 s	12.29 s	41.3 (36)	18.78 s	4.99 s	
1 024	43.3 (11)	9.18 s	5.85 s	53.0 (29)	8.68 s	4.22 s	
2 048	41.4 (11)	4.15 s	2.63 s	62.2 (86)	4.47 s	4.23 s	
4 096	41.2 (11)	1.66 s	1.49 s	68.9 (40)	2.52 s	2.86 s	
8 192	40.2 (11)	1.26 s	1.06 s	-	-	-	

Computations performed on Cori (NERSC).

Heinlein, Perego, Rajamanickam (2022)

Spectral Extension-Based Coarse Spaces for Schwarz Preconditioners

Highly heterogeneous problems ...

... appear in most areas of modern science and engineering:







Micro section of a dual-phase steel. Courtesy of J. Schröder.

Groundwater flow (SPE10); cf. Christie and Blunt (2001).

Composition of arterial walls; taken from O'Connell et al. (2008).

Spectral coarse spaces

The coarse space is **enhanced** by eigenfunctions of **local edge and face eigenvalue problems** with eigenvalues below tolerances $tol_{\mathcal{E}}$ and $tol_{\mathcal{F}}$:

$$\kappa\left(\mathbf{M}_{*}^{-1}\mathbf{K}\right) \leq C\left(1 + \frac{1}{\operatorname{tol}_{\mathcal{B}}} + \frac{1}{\operatorname{tol}_{\mathcal{F}}} + \frac{1}{\operatorname{tol}_{\mathcal{F}} \cdot \operatorname{tol}_{\mathcal{F}}}\right);$$

C does not depend on *h*, *H*, or the coefficients. OS-ACMS & adaptive GDSW (AGDSW) (Heinlein, Klawonn, Knepper, Rheinbach (2018, 2018, 2019)).

Local eigenvalue problems

Local generalized eigenvalue problems corresponding to the edges & and faces $\mathcal F$ of the domain decomposition:

$$\begin{aligned} \forall E \in \mathcal{E} : \qquad & \boldsymbol{S}_{EE} \tau_{*,E} = \lambda_{*,E} \boldsymbol{K}_{EE} \tau_{*,E}, \quad \forall \tau_{*,E} \in V_E, \\ \forall F \in \mathcal{F} : \qquad & \boldsymbol{S}_{FF} \tau_{*,F} = \lambda_{*,F} \boldsymbol{K}_{FF} \tau_{*,F}, \quad \forall \tau_{*,F} \in V_F, \end{aligned}$$

with Schur complements S_{EE} , S_{FF} with Neumann boundary conditions and submatrices K_{EE} , K_{FF} of K. We select eigenfunctions corresponding to eigenvalues below tolerances $tol_{\&}$ and $tol_{\mathcal{J}}$.

 \rightarrow The corresponding coarse basis functions are **energy-minimizing extensions** into the interior of the subdomains.



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Foam coefficient function example



Solid phase: $\alpha = 10^6$; transparent phase: $\alpha = 1$; 100 subdomains

V ₀	tol _{&}	$\mathit{tol}_{\mathcal{F}}$	it.	κ	dim V_0	dim V_0/dof
$V_{ m GDSW}$	—	—	565	$1.3 \cdot 10^{6}$	1601	0.27 %
$V_{\rm AGDSW}$	0.05	0.05	60	30.2	1 968	0.33 %
$V_{\rm OS-ACMS}$	0.001	0.001	57	30.3	690	0.12 %

Cf. Heinlein, Klawonn, Knepper, Rheinbach (2018, 2019).

Domain Decomposition for Neural Networks

Domain Decomposition for Neural Networks



A non-exhaustive literature overview:

- Machine Learning for adaptive BDDC, FETI–DP, and AGDSW: Heinlein, Klawonn, Lanser, Weber (2019, 2020, 2021, 2021, 2021, 2022); Klawonn, Lanser, Weber (2024)
- cPINNs, XPINNs: Jagtap, Kharazmi, Karniadakis (2020); Jagtap, Karniadakis (2020)
- Classical Schwarz iteration for PINNs or DeepRitz (D3M, DeepDDM, etc):: Li, Tang, Wu, and Liao . (2019); Li, Xiang, Xu (2020); Mercier, Gratton, Boudier (arXiv 2021); Dolean, Heinlein, Mercier, Gratton (subm. 2024 / arXiv:2408.12198); Li, Wang, Cui, Xiang, Xu (2023); Sun, Xu, Yi (arXiv 2022, arXiv 2023); Kim, Yang (2022, arXiv 2023)
- FBPINNs, FBKANs: Moseley, Markham, and Nissen-Meyer (2023); Dolean, Heinlein, Mishra, Moseley . (2024, 2024); Heinlein, Howard, Beecroft, Stinis (acc. 2024 / arXiv:2401.07888); Howard, Jacob, Murphy, Heinlein, Stinis (arXiv:2406.19662)
- DDMs for CNNs: Gu, Zhang, Liu, Cai (2022); Lee, Park, Lee (2022); Klawonn, Lanser, Weber (2024); . Verburg, Heinlein, Cyr (subm. 2024)

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 mid 2024:



N. Klawonn, M. Lanser, J. Weber

Machine learning and domain decomposition methods - a survey

Computational Science and Engineering. 2024

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**

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

PINNs use a hybrid loss function:

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

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

$$\begin{split} \mathcal{L}_{data}(\boldsymbol{\theta}) &= \frac{1}{N_{data}} \sum_{i=1}^{N_{data}} \left(u(\hat{\boldsymbol{x}}_i, \boldsymbol{\theta}) - u_i \right)^2, \\ \mathcal{L}_{PDE}(\boldsymbol{\theta}) &= \frac{1}{N_{PDE}} \sum_{i=1}^{N_{PDE}} \left(\mathcal{N}[u](\boldsymbol{x}_i, \boldsymbol{\theta}) - f(\boldsymbol{x}_i) \right)^2. \end{split}$$

See also Dissanayake and Phan-Thien (1994); Lagaris et al. (1998).

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



Hybrid loss



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

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Theoretical Result for PINNs

Estimate of the generalization error (Mishra and Molinaro (2022))

The generalization error (or total error) satisfies

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

- $\mathcal{E}_G = \mathcal{E}_G(\boldsymbol{X}, \boldsymbol{\theta}) \coloneqq \| \mathbf{u} \mathbf{u}^* \|_V$ general. error (V Sobolev space, \boldsymbol{X} training data set)
- δ_T training error (*I^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, quadrature, and neural network

Rule of thumb: "As long as the PINN is trained well, it also generalizes well"



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

Motivation – Some Observations on the Performance of PINNs

Solve

 $u' = \cos(\omega x),$ u(0) = 0,

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)



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



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 employ the PINN approach and hard enforcement of the boundary conditions; cf. Lagaris et al. (1998).

FBPINNs use the network architecture

$$u(\boldsymbol{\theta}_1,\ldots,\boldsymbol{\theta}_J) = C \sum_{j=1}^J \omega_j u_j(\boldsymbol{\theta}_j)$$

and the loss function

$$\mathcal{L}(\boldsymbol{\theta}_1,\ldots,\boldsymbol{\theta}_J) = \frac{1}{N} \sum_{i=1}^N \left(\mathcal{N}[\mathcal{C}\sum_{\boldsymbol{x}_i \in \Omega_j} \omega_j u_j](\boldsymbol{x}_i,\boldsymbol{\theta}_j) - f(\boldsymbol{x}_i) \right)^2.$$

Here:

- Overlapping DD: $\Omega = \bigcup_{l=1}^{J} \Omega_{j}$
- Partition of unity ω_j with $supp(\omega_j) \subset \Omega_j$ and $\sum_{j=1}^J \omega_j \equiv 1$ on Ω



Hard enf. of boundary conditions Loss function

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

with constraining operator C, which explicitly enforces the boundary conditions.

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Numerical Results for FBPINNs



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Multi-Level FBPINN Algorithm

Extension of FBPINNs to L levels; Cf. Dolean, Heinlein, Mishra, Moseley (2024).



$\label{eq:local_$

$$u\big(\theta_1^{(1)},\ldots,\theta_{j^{(L)}}^{(L)}\big) = \mathcal{C}\big(\sum_{l=1}^L\sum_{i=1}^{N^{(l)}}\omega_j^{(l)}u_j^{(l)}\big(\theta_j^{(l)}\big)\big)$$



Multi-Frequency Problem

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

$$-\Delta u = 2 \sum_{i=1}^{n} (\omega_i \pi)^2 \sin(\omega_i \pi x) \sin(\omega_i \pi y) \quad \text{in } \Omega,$$
$$u = 0 \qquad \qquad \text{on } \partial\Omega,$$

with $\omega_i = 2^i$.

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



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Multi-Level FBPINN Algorithm

Extension of FBPINNs to L levels; Cf. Dolean, Heinlein, Mishra, Moseley (2024).



L-level network architecture

$$u\big(\boldsymbol{\theta}_1^{(1)},\ldots,\boldsymbol{\theta}_{j^{(L)}}^{(L)}\big) = \mathcal{C}\big(\sum_{l=1}^L\sum_{i=1}^{N^{(l)}}\omega_j^{(l)}u_j^{(l)}\big(\boldsymbol{\theta}_j^{(l)}\big)\big)$$



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

with $\omega_i = 2^i$.

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

n = 1 n = 2 n = 3 n = 3 n = 3 n = 3 n = 3 n = 3 n = 6 n = 6

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





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Multi-Frequency Problem – What the FBPINN Learns



Cf. Dolean, Heinlein, Mishra, Moseley (2024).



Cf. Dolean, Heinlein, Mishra, Moseley (2024).

Memory Requirements for CNN Training



- As an example for a convolutional neural network (CNN), we employ the U-Net architecture introduced in Ronneberger, Fischer, and Brox (2015).
- The U-Net yields state-of-the-art accuracy in semantic image segmentation and other image-to-image tasks.

Below: memory consumption for training on a single 1024×1024 image.

12120	size	# channels		mem. feature maps		mem. weights	
lidille		input	output	# of values	MB	# of values	MB
input block	1 0 2 4	3	64	268 M	1 024.0	38 848	0.148
encoder block 1	512	64	128	167 M	704.0	221 696	0.846
encoder block 2	256	128	256	84 M	352.0	885 760	3.379
encoder block 3	128	256	512	42 M	176.0	3 540 992	13.508
encoder block 4	64	512	1024	21 M	88.0	14 159 872	54.016
decoder block 1	64	1,024	512	50 M	192.0	9 177 088	35.008
decoder block 2	128	512	256	101 M	384.0	2 294 784	8.754
decoder block 3	256	256	128	201 M	768.0	573 952	2.189
decoder block 4	512	128	64	402 M	1 536.0	143 616	0.548
output block	1 0 2 4	64	3	3.1 M	12.0	195	0.001

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Decomposing the U-Net



Cf. Verburg, Heinlein, Cyr (subm. 2024).
Decomposing the U-Net



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Decomposing the U-Net



Decomposing the U-Net



- Distribution of feature maps results in significant reduction of memory usage on a single GPU
- Moderate additional memory usage due to the communication network

Results – Synthetic Data Set





		esting	5 011 0	Subin	lages			
2	0.13	0.20	0.20	0.29	0.28	0.30	0.40	0.53
pixels 3	0.14	0.39	0.65	0.75	0.79	0.88	0.92	0.76
× 32 4	0.14	0.59	0.55	0.77	0.78	0.83	0.73	0.75
5 (32 6	0.14	0.23	0.54	0.67	0.72	0.82	0.87	0.68
mages	0.13	0.18	0.39	0.49	0.64	0.71	0.70	0.42
: subii 16	0.13	0.14	0.14	0.16	0.18	0.16	0.26	0.26
#	0	1	2	4	8	16	32	Baseline
	#	feature	e map	s com	munica	ated		

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DeepGlobe 2018 Satellite Image Data Set (Demir et al. (2018))

class	pixel count	proportion
urban	642.4M	9.35 %
agriculture	3898.0M	56.76%
rangeland	701.1M	10.21%
forest	944.4M	13.75%
water	256.9M	3.74 %
barren	421.8M	6.14 %
unknown	3.0M	0.04 %



Avoiding overfitting

The data set includes only 803 images. To avoid overfitting, we

- apply batch normalization, use random dropout layers and data augmentation, and
- initialize the encoder using the ResNet-18 (He, Zhang, Ren, and Sun (2016))



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Schwarz Domain Decomposition Preconditioners

- Numerical scalability and robust convergence for
 - heterogeneous problems
 - multiphysics problems
 - highly nonlinear problems
- ightarrow Algebraic and parallel implementation in FROSCH 🏘

Domain Decomposition for Neural Networks

- Schwarz domain decomposition architectures improve the scalability of PINNs to large domains / high frequencies, keeping the complexity of the local networks low.
- Novel DDU-Net approach decouples the training on the sub-images, allowing us to distribute the memory load among multiple GPUs. It limits communication to deepest level of the U-Net architecture using a communication network.

Thank you for your attention!