

Learning a Neural Solver for Parametric PDEs to Enhance Physics-Informed Methods

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TLDR; We propose to solve parametric PDEs using Physics-Informed methods by learning a dedicated optimizer that considerably accelerates convergence.



1. Context & problem formulation

- PINNs have demonstrated interesting performances but remain limited by training time and poor performance in parametric settings.
- We focus on solving parametric PDEs from Physical knowledge.

$$\begin{aligned}\mathcal{N}(u; \gamma) &= f \quad \text{in } \Omega, \\ \mathcal{B}(u) &= g \quad \text{on } \partial\Omega.\end{aligned}$$

Where $\gamma \in \Gamma$ are the PDE physical parameters, f are forcing terms and g can be initial and/or boundary conditions.

- We assume access to a dataset of pairs composed of the PDE data (γ, f, g) and the associated solution u on a grid.

2. Motivation

- PINNs losses are ill-conditioned and hard to optimize for traditional optimizers.
- They require extensive computational time and numerous iterations to compensate this aspect.

3. How to learn a Physics-informed solver?

- Global framework

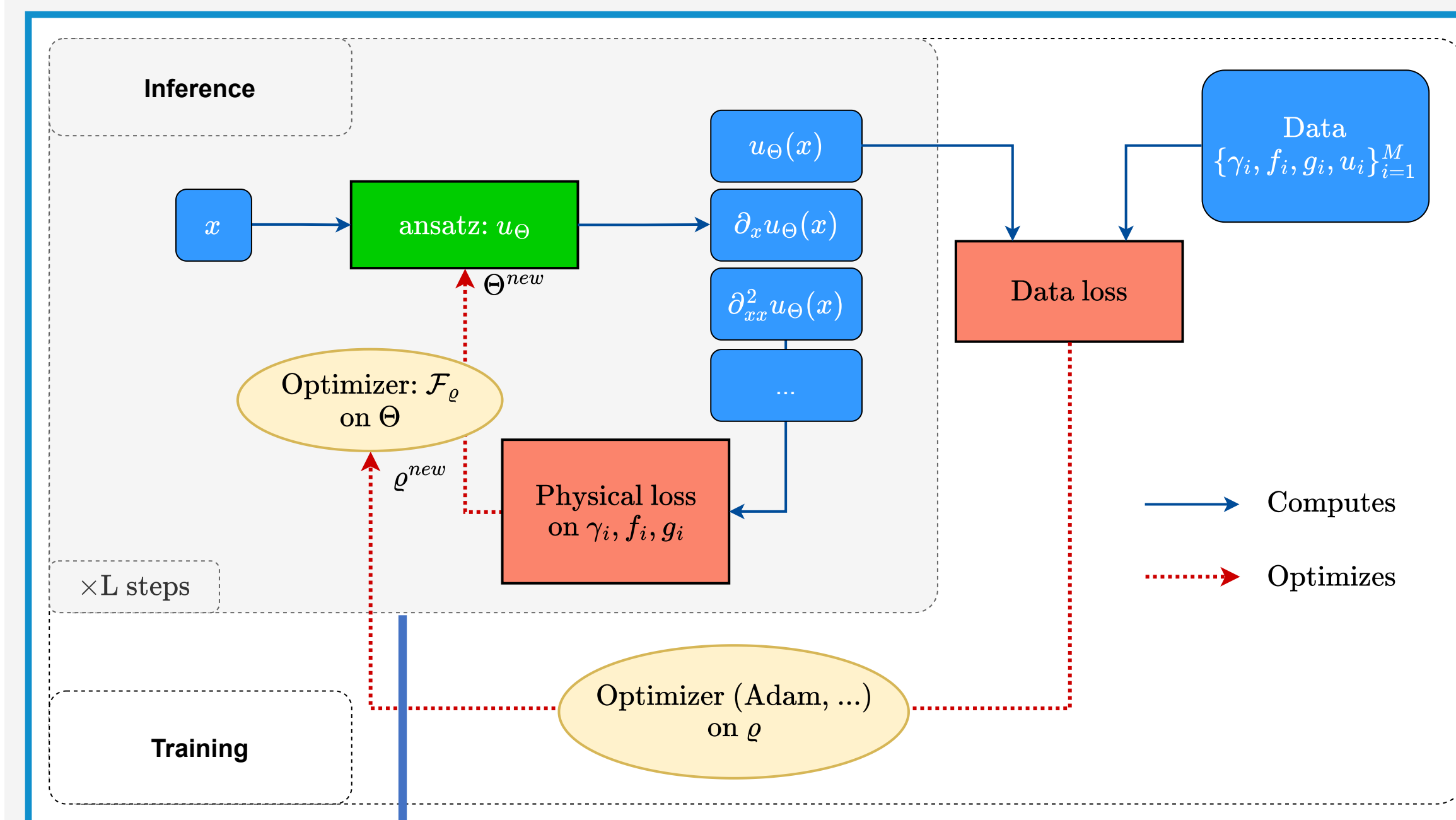


Figure 1: Optimization scheme of a physics-informed method with our framework.

Algorithm 1: Inference using the neural PDE solver.

Data: $\Theta_0 \in \mathbb{R}^n$, PDE (γ, f, g)
Result: $\Theta_L \in \mathbb{R}^n$
for $l = 0 \dots L-1$ do
 $\Theta_{l+1} = \Theta_l - \eta \mathcal{F}_\rho(\nabla \mathcal{L}_{\text{PDE}}(\Theta_l), \gamma, f, g)$
end
return Θ_L

The neural solver learns to transform the physical gradient into a more effective gradient direction that achieves fast convergence.

- Theoretical analysis in the linear case

Theorem 1. (Convergence rate in the linear case). Given a linear ansatz $u_\Theta(x) = \sum_{i=1}^N \theta_i \phi_i(x)$, assume the conditioner \mathcal{F} behaves like its linearization $P = \text{Jacobian}(\mathcal{F})$, meaning that \mathcal{F} can be replaced by P at any point. Let A be the matrix derived from the PDE loss as eq. (3) for the Poisson equation or eq. (15) in the more general case. Denote by $\kappa(A)$ the condition number of the matrix A . The number of steps $N'(\varepsilon)$ required to achieve an error $\|\Theta_l - \Theta^*\|_2 \leq \varepsilon$ satisfies:

$$N'(\varepsilon) = O\left(\kappa(PA) \ln\left(\frac{1}{\varepsilon}\right)\right), \quad (11)$$

Moreover, if \mathcal{F} minimizes $\mathcal{L}_{\text{DATA}}$ this necessarily implies $\kappa(PA) = 1 \leq \kappa(A)$. Consequently, the number of steps is effectively reduced, i.e., $N'(\varepsilon) \ll N(\varepsilon)$ with $N(\varepsilon)$ the number of steps of the vanilla PINNs.

4. Results

- Quantitative evaluation: comparison with baselines.

Table 1: Results of trained models - metrics in Relative MSE on the test set. Best performances are highlighted in bold, and second best are underlined.

	Baseline	1d		1d+time	2d	2d+time
		Helmholtz	Poisson	NLRD	Darcy-Flow	Heat
Supervised	<i>MLP + basis</i>	4.66e-2	1.50e-1	2.85e-4	3.56e-2	6.00e-1
Unsupervised	<i>PINNs+L-BFGS</i>	9.86e-1	8.83e-1	6.13e-1	9.99e-1	9.56e-1
	<i>PINNs-multi-opt</i>	8.47e-1	1.18e-1	7.57e-1	8.38e-1	6.10e-1
	<i>PPINNs</i>	9.89e-1	4.30e-2	3.94e-1	8.47e-1	1.27e-1
	<i>P2INNs</i>	9.90e-1	1.50e-1	5.69e-1	8.38e-1	1.78e-1
	<i>PO-DeepONet</i>	9.83e-1	1.43e-1	4.10e-1	8.33e-1	4.43e-1
Hybrid	<i>PI-DeepONet</i>	9.79e-1	1.20e-1	7.90e-2	2.76e-1	9.18e-1
	<i>PINO</i>	9.99e-1	<u>2.80e-3</u>	4.21e-4	1.01e-1	<u>9.09e-3</u>
Neural Solver	<i>Ours</i>	2.41e-2	5.56e-5	2.91e-4	1.87e-2	2.31e-3

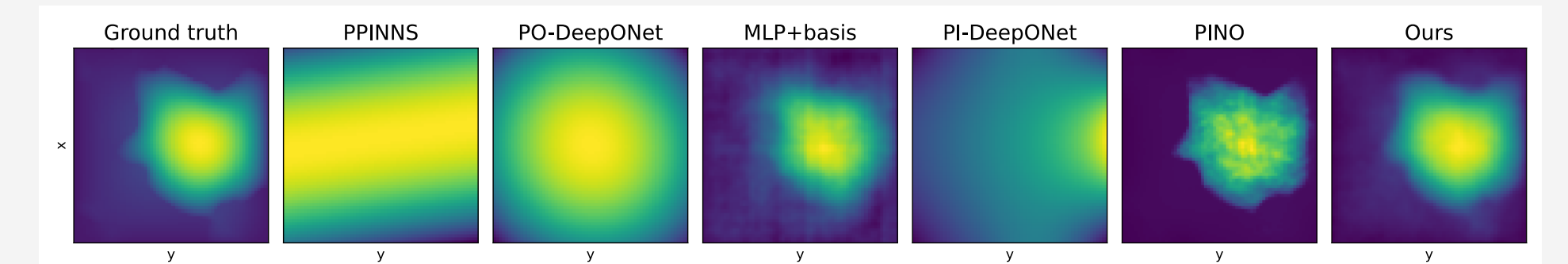


Figure 2: visual comparison of our solver's solution with baselines on the Darcy dataset.

- Solving new PDEs: comparison with optimizers.

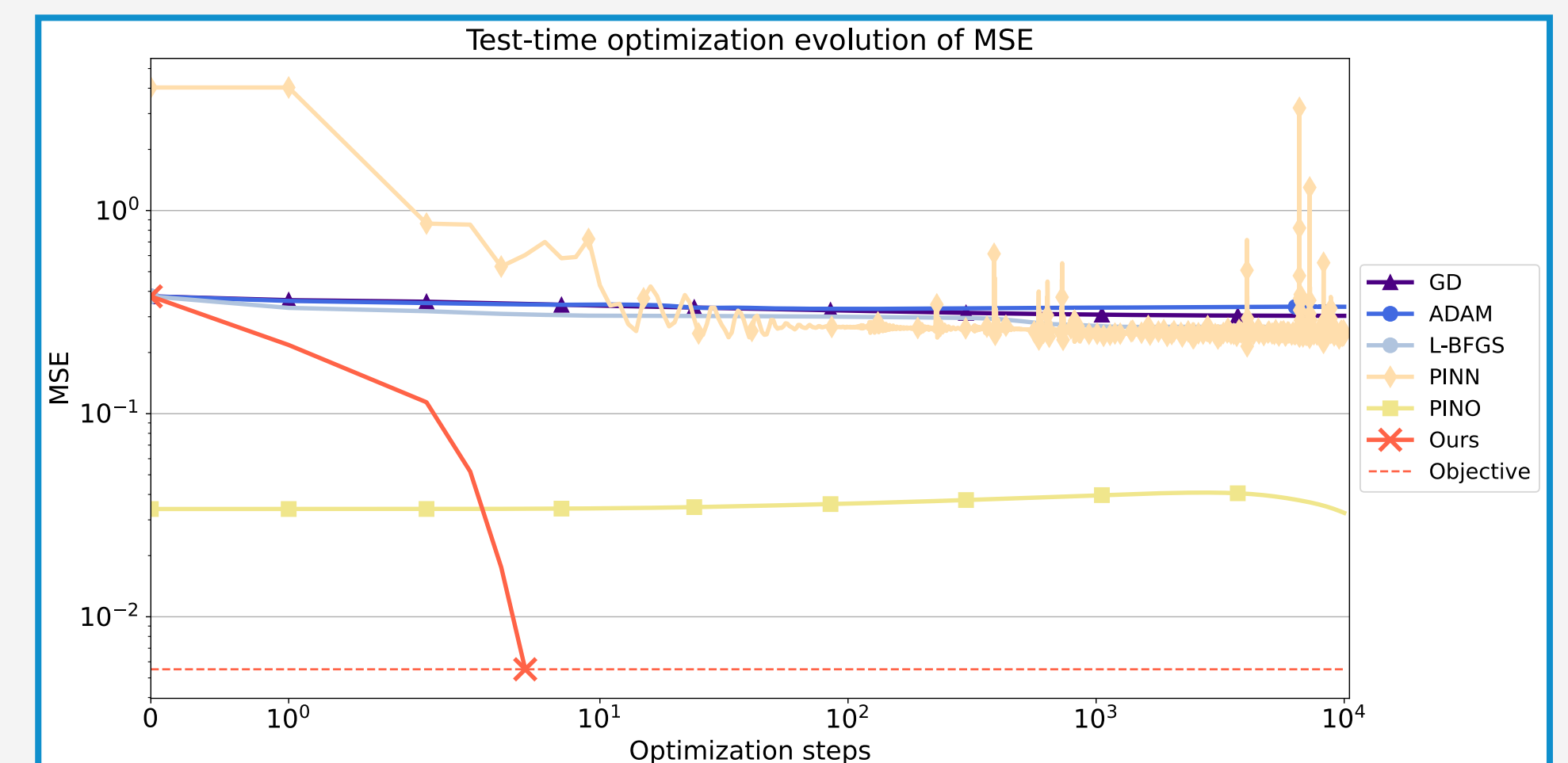


Figure 3: Test-time optimization based on the physical residual loss \mathcal{L}_{PDE} for new PDE on Darcy.

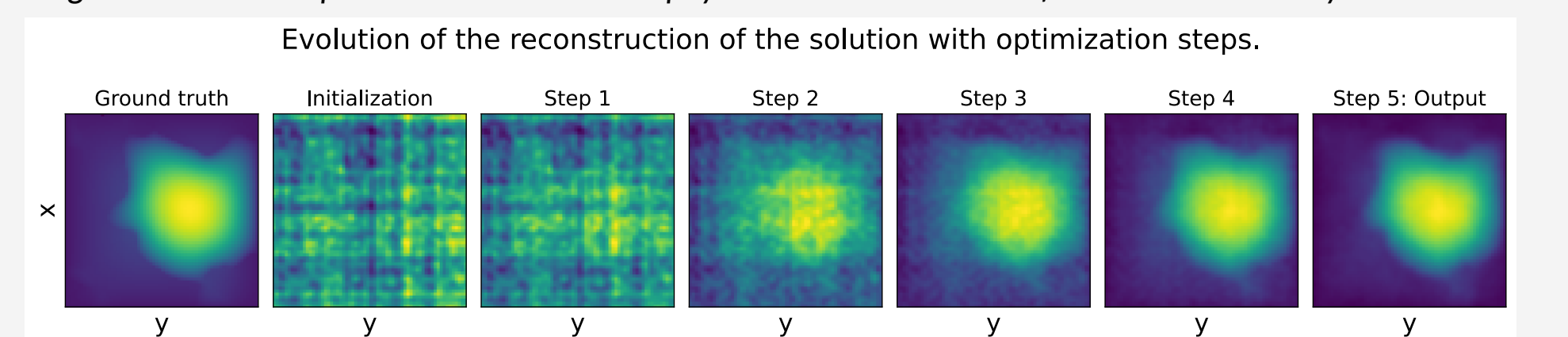


Figure 4: Visualization of the reconstruction of the solution with our method to solve a Darcy PDE.