

COMPUTATIONALLY EFFICIENT METHODS FOR UNCERTAINTY QUANTIFICATION IN SEISMIC INVERSION

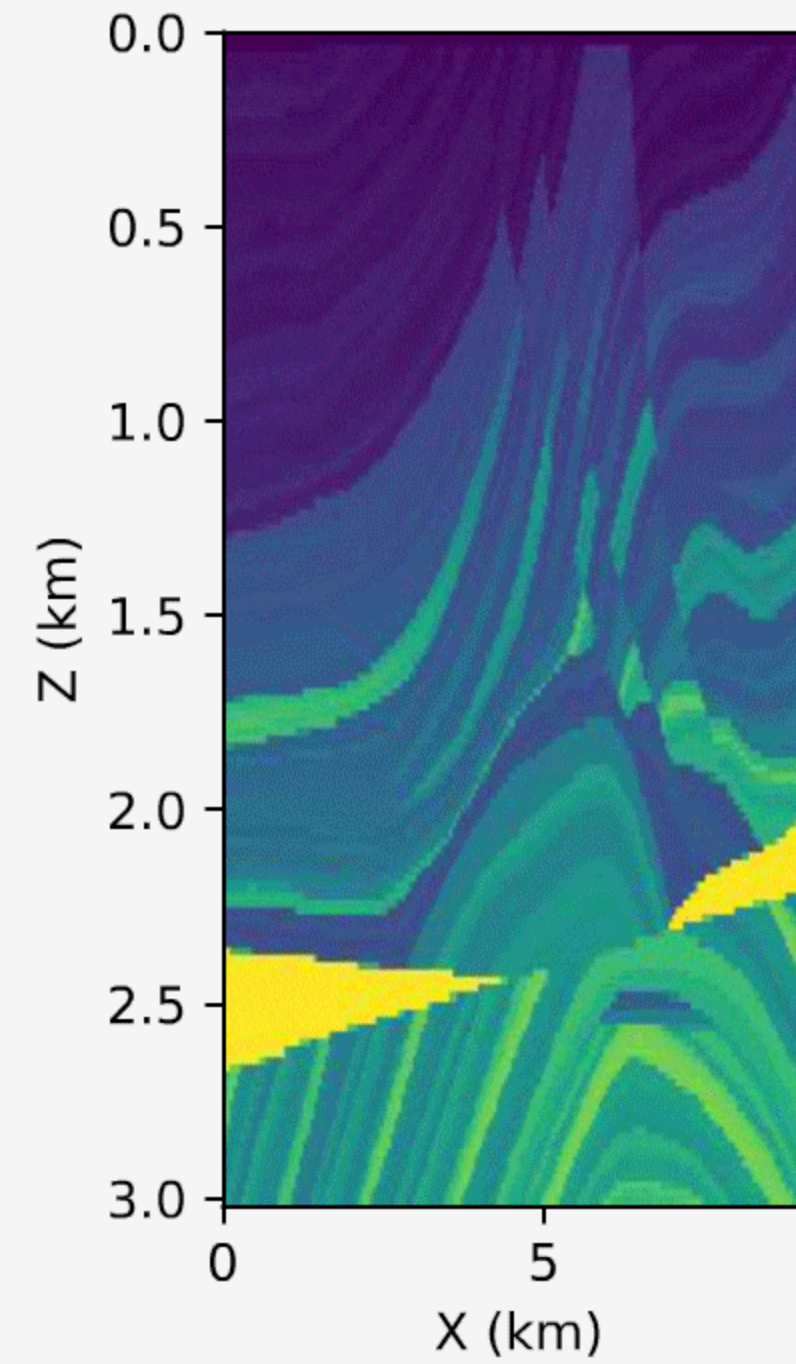
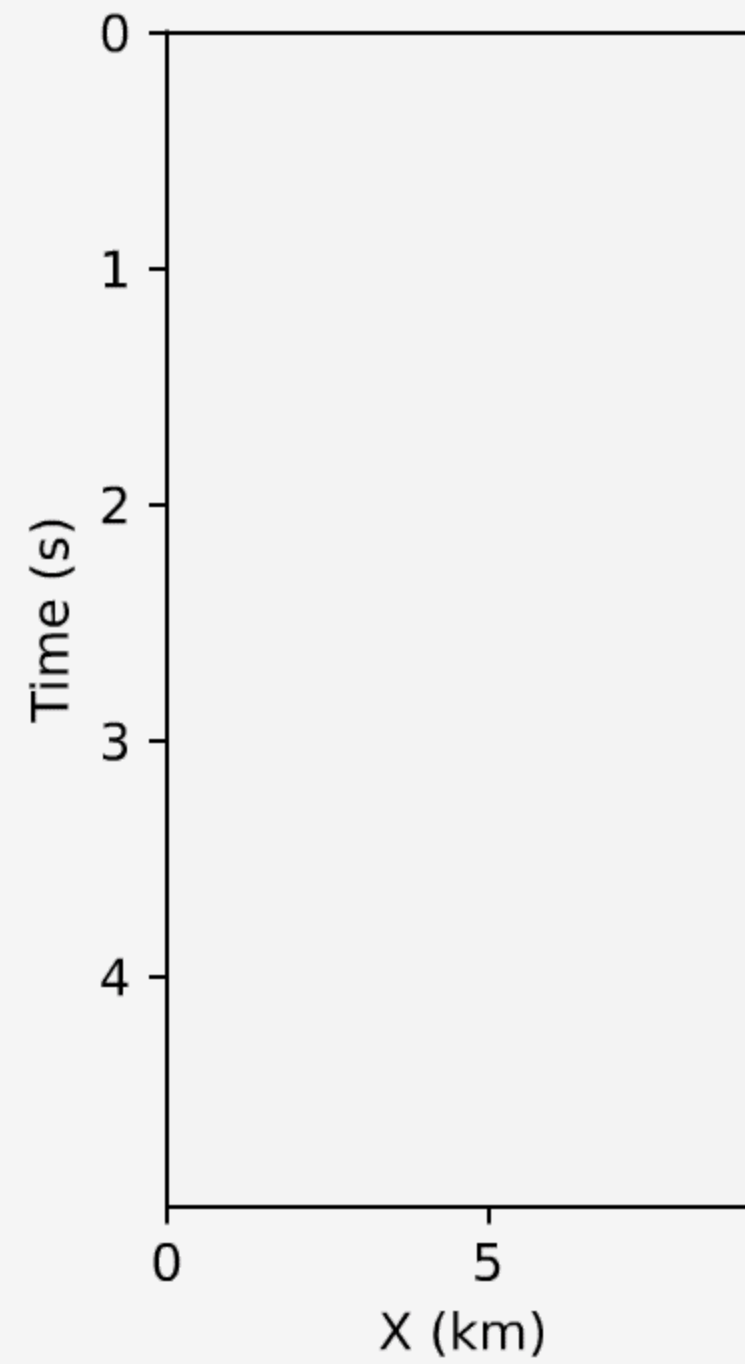
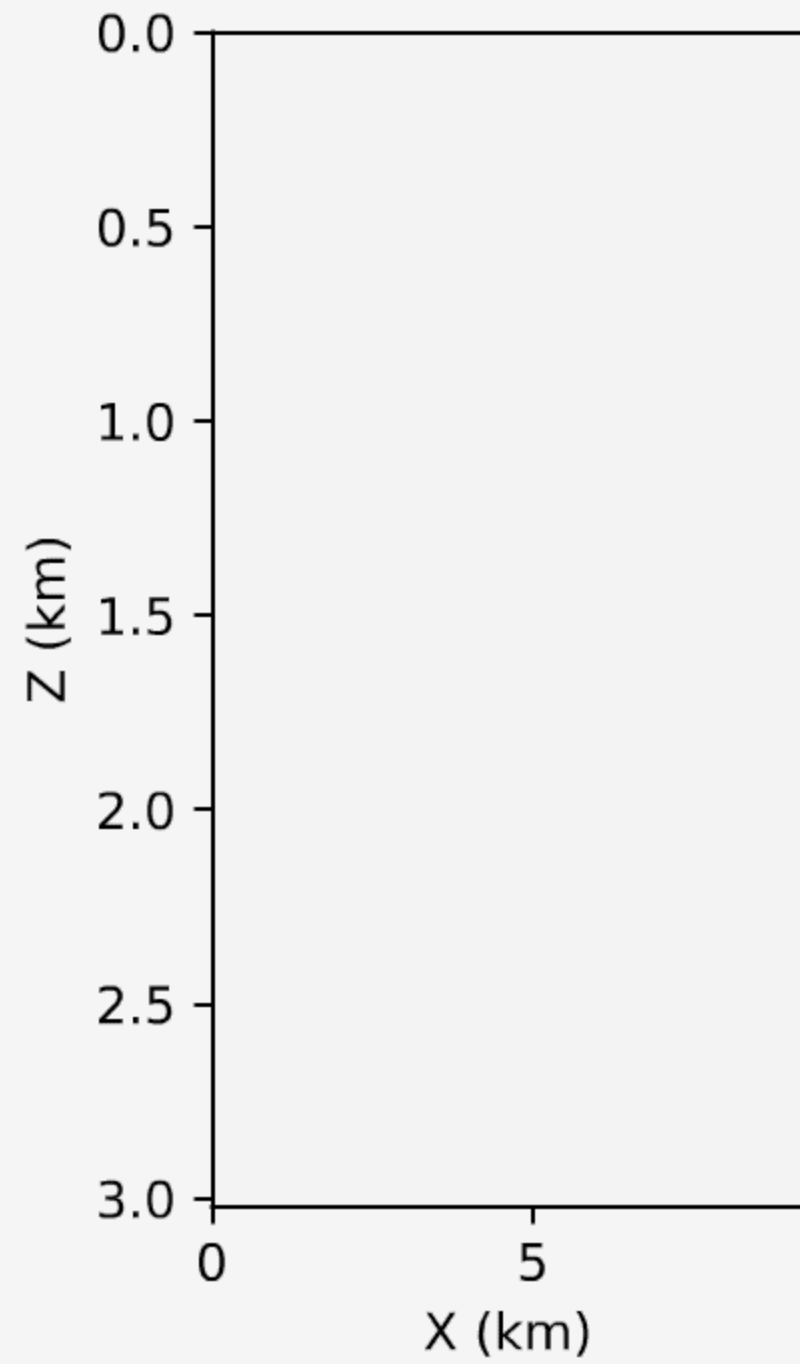


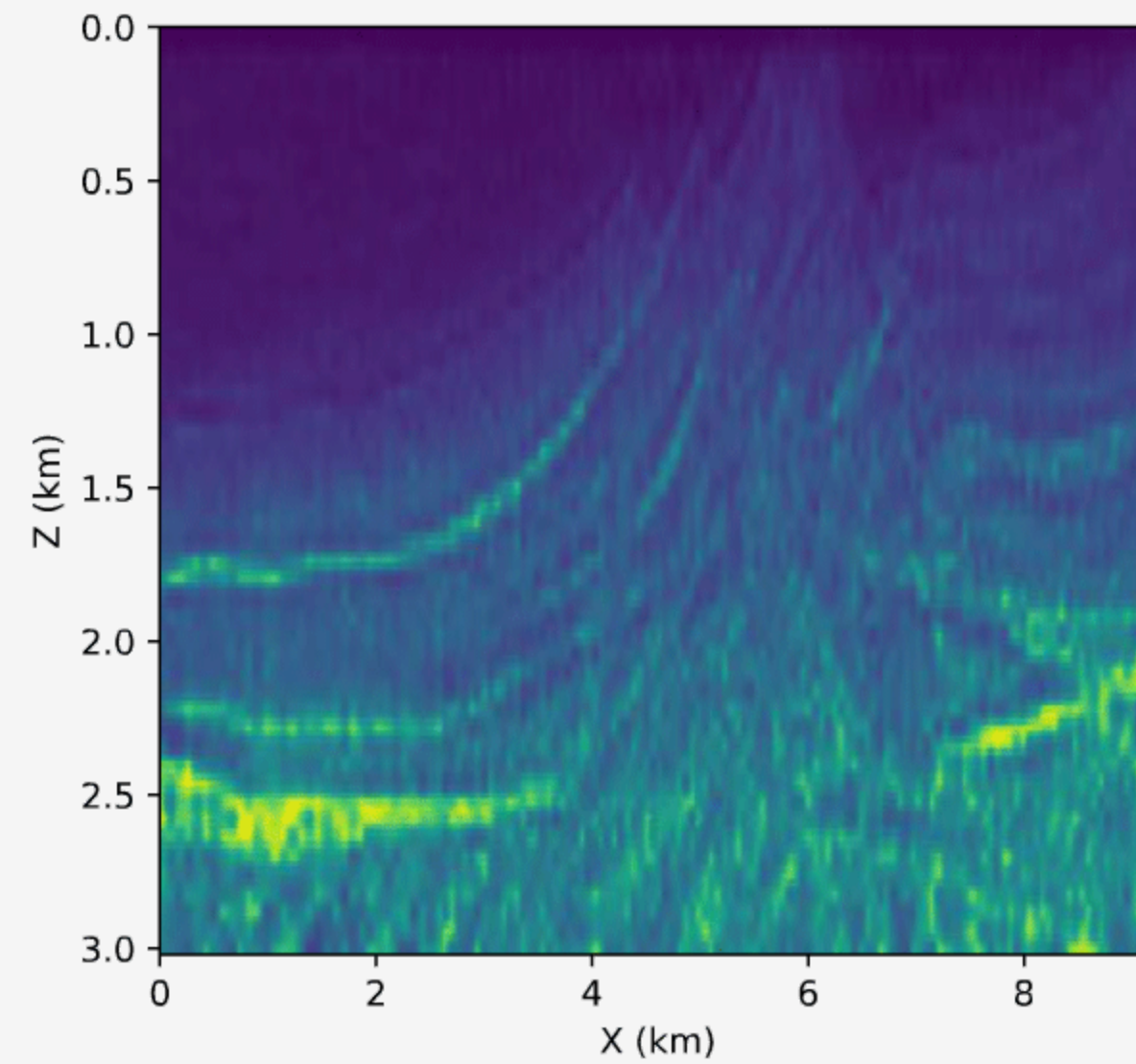
Georgia K. Stuart

The Department of Mathematical Sciences
The University of Texas at Dallas

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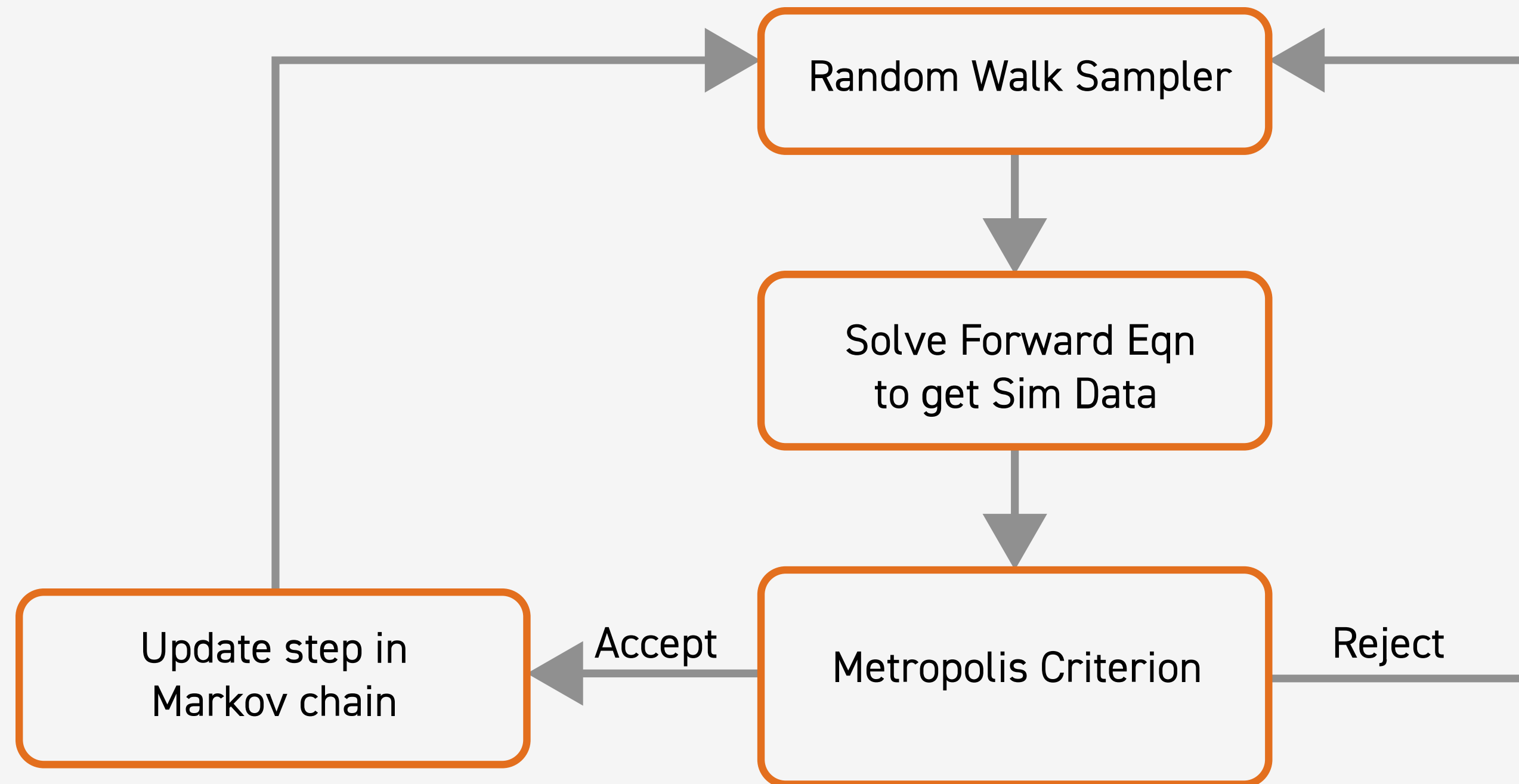












HOW CAN WE REDUCE THE
COMPUTATIONAL COST OF
MCMC METHODS FOR FWI?

- Modeling wave propagation can be **computationally expensive**.
- We use the 2D constant-density acoustic wave equation

$$\frac{1}{c^2(x, z)} \frac{\partial^2 p}{\partial t^2} - \nabla p = f$$

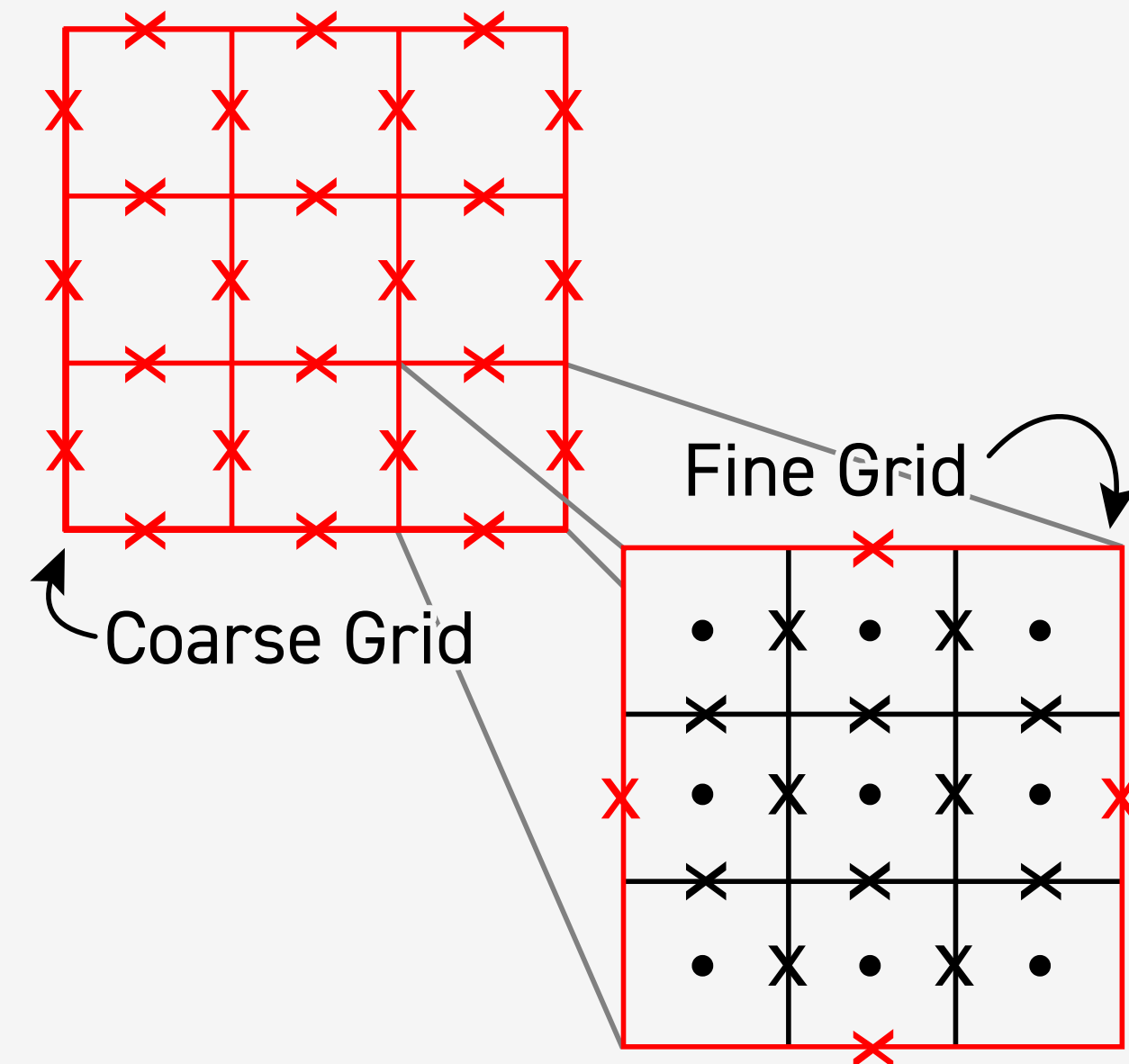
- **Operator upscaling**¹ decomposes the solution into two parts:
 1. **Fine grid** problem on independent subdomains
 2. Small **coarse grid** problem over the whole domain
- In this upscaling technique **we do NOT upscale the model**.

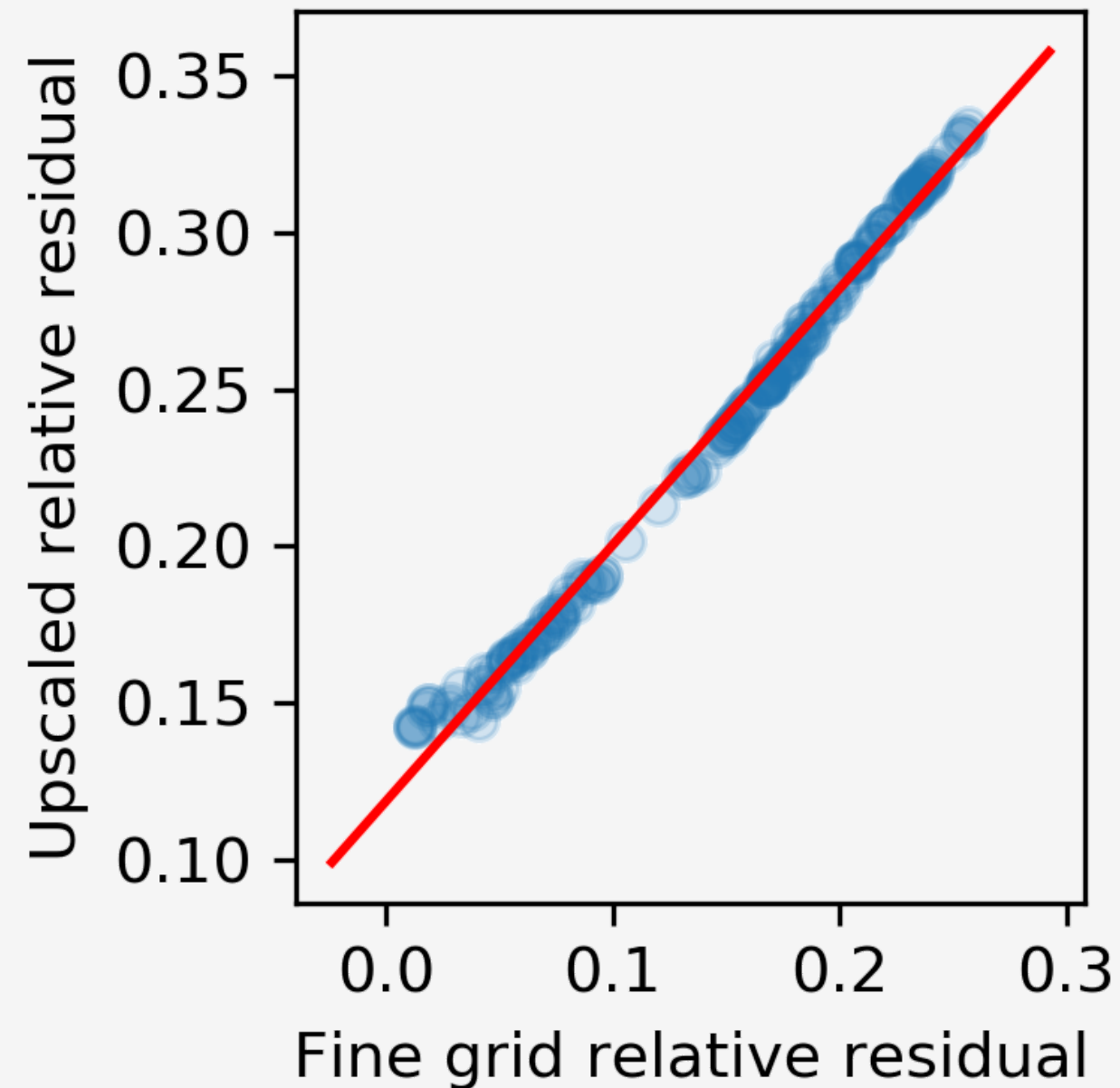
1. Write the acoustic wave equation as a system in space by introducing **acceleration, \vec{v}**

$$\vec{v} = -\nabla p$$

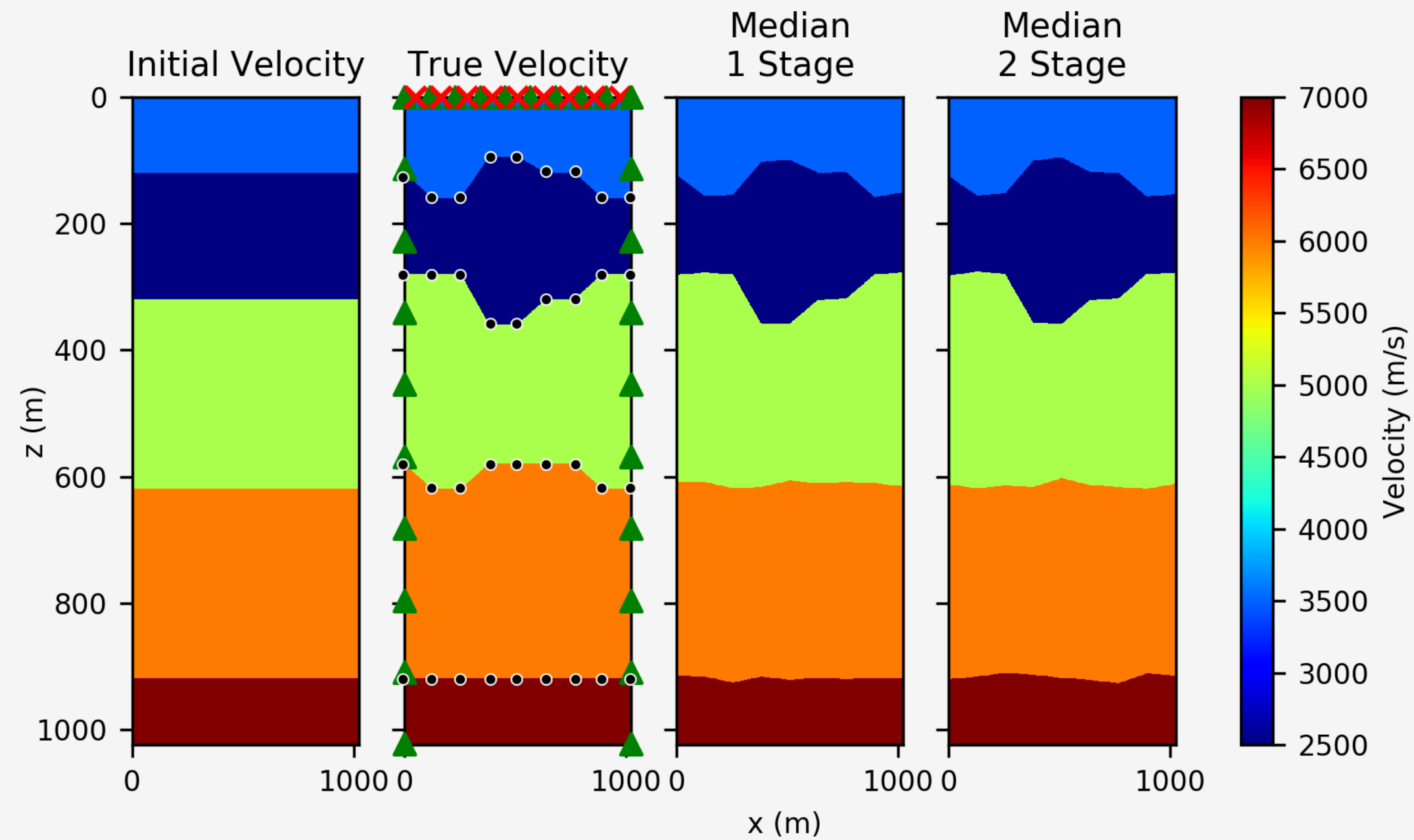
$$\frac{1}{c^2} \frac{\partial^2 p}{\partial t^2} = -\nabla \cdot \vec{v} + f$$

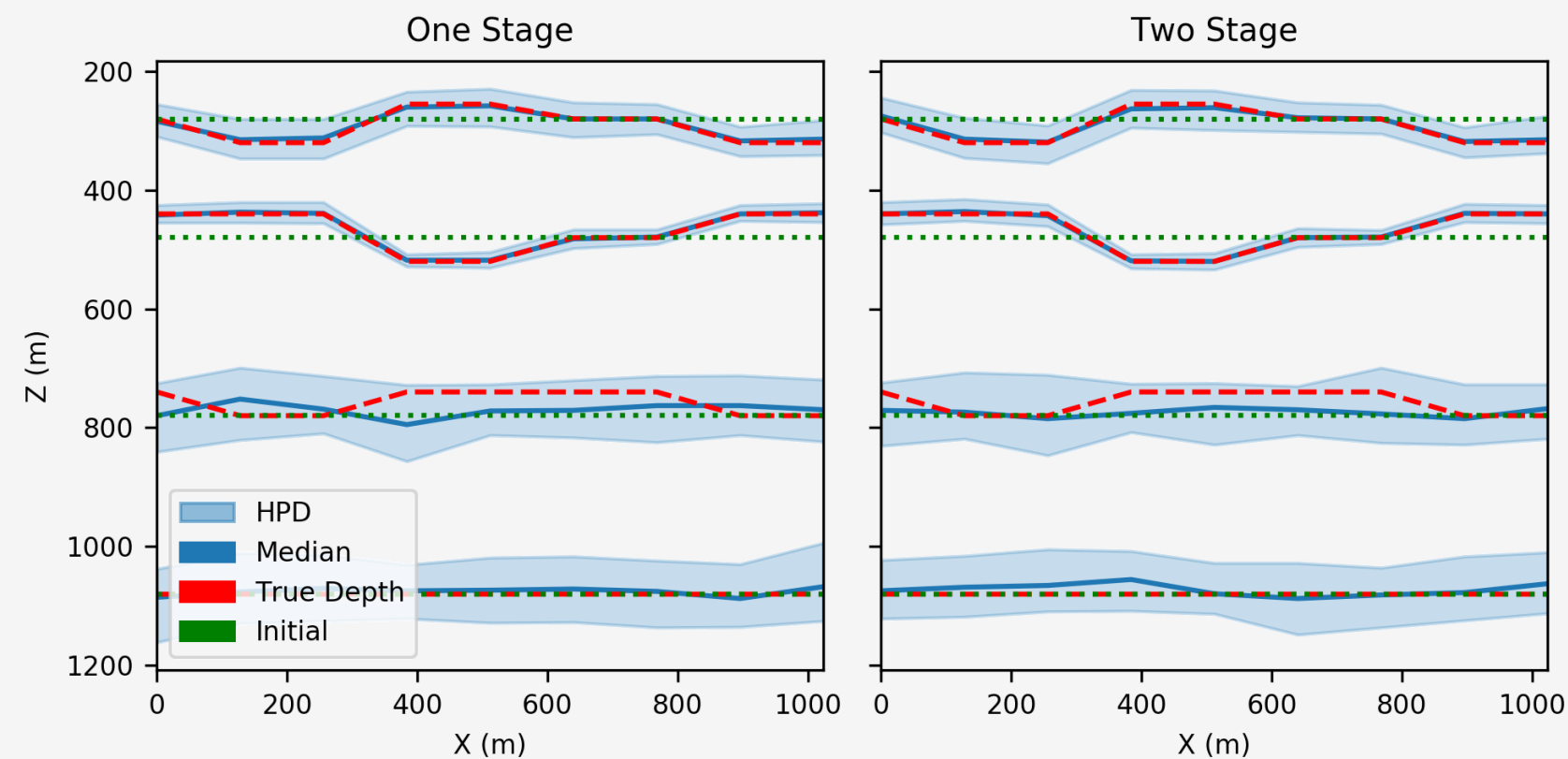
2. Solve in parallel for fine grid pressure and acceleration over each **independent** coarse block. **No communication** is required at this stage.
3. Solve for coarse grid acceleration over the whole domain.





- We see a **strong linear relationship** between the fine grid relative residuals and the upscaled relative residuals for a layered velocity model.
- This indicates that the upscaling filter is a **good surrogate** for the fine grid solver.

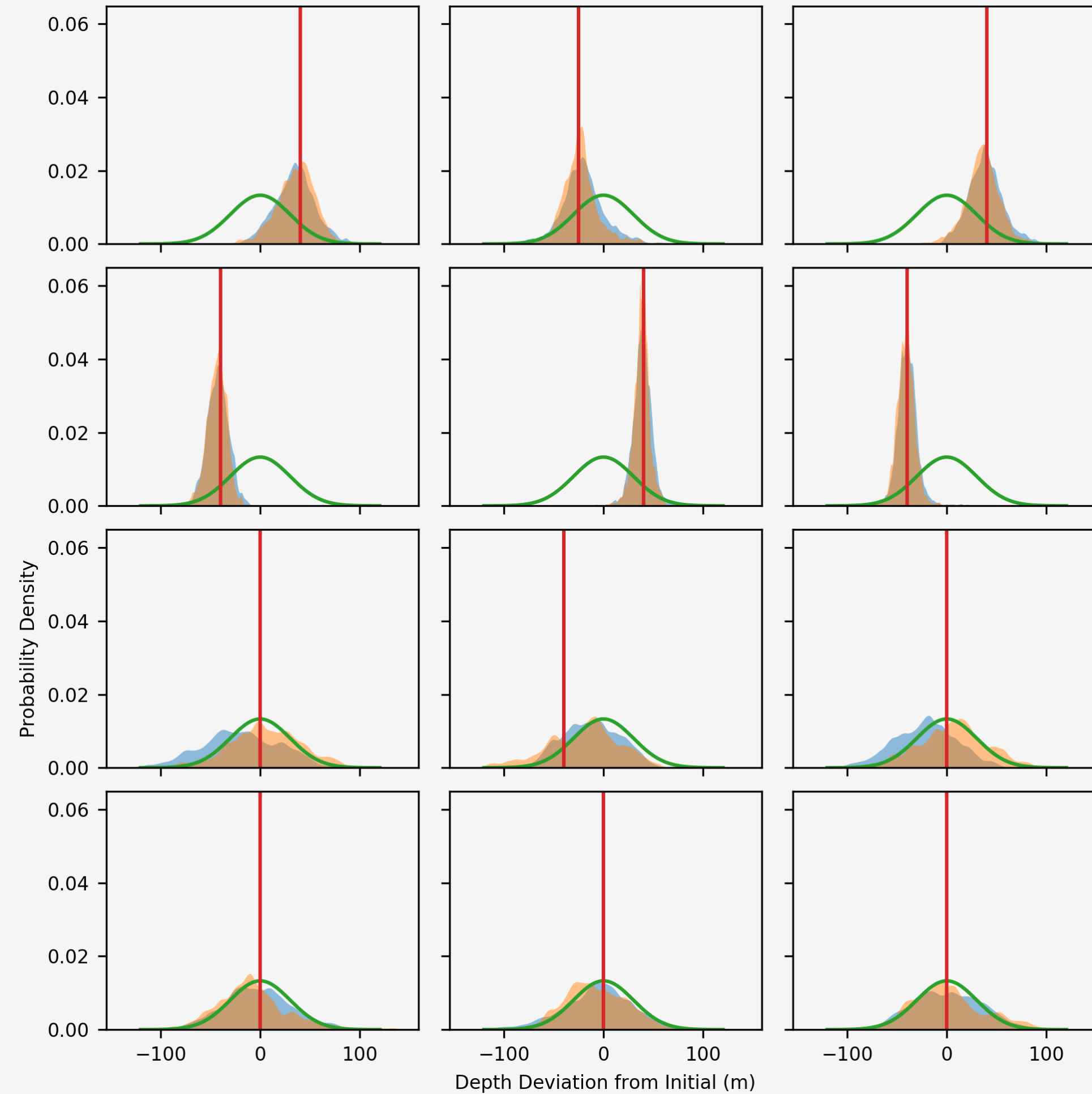


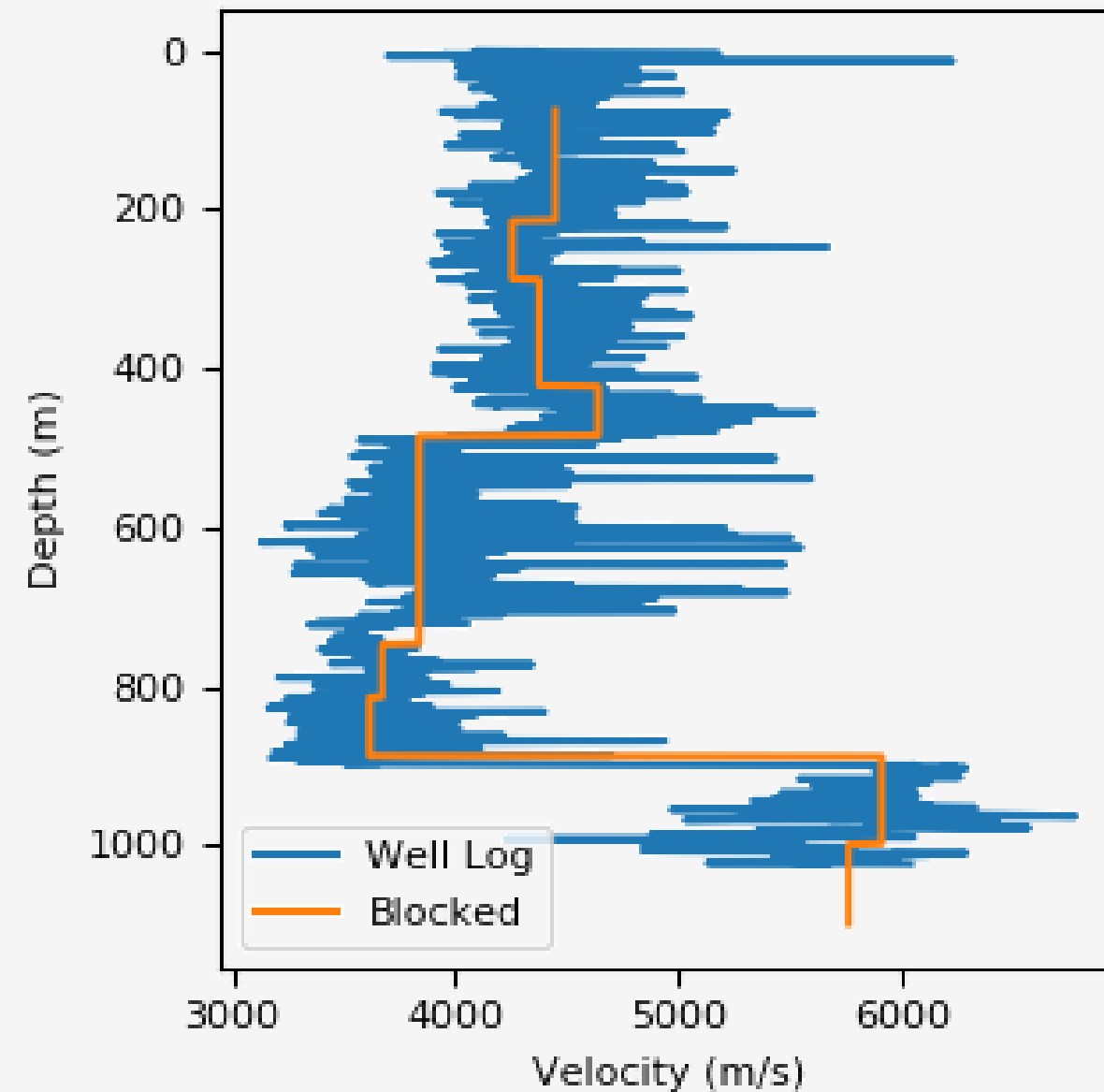


A comparison between one-stage MCMC highest posterior density (HPD) intervals and two-stage MCMC HPD intervals.

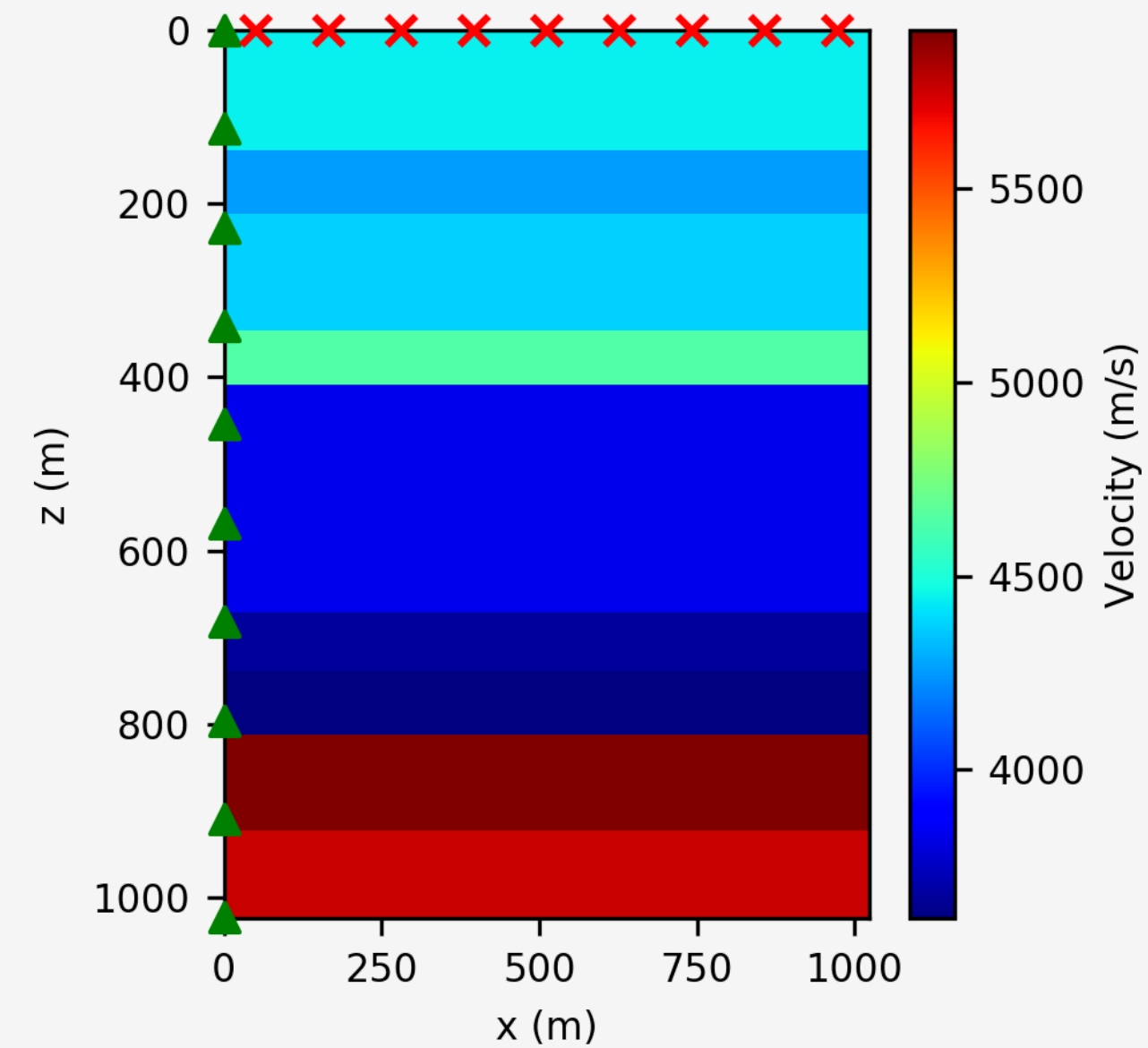
- Acceptance rate **increases** from 10% to 40%.
- Time per sample **decreases** by 22% (40% in other experiments).
- Time per rejection **decreases** by 33%.

RESULTS: TWO-STAGE MCMC WITH UPSCALING





Well log from the Midland, TX basin (blue, courtesy of [Pioneer Natural Resources](#) and 9-layer block (orange).



Flat layered experimental setup (Stuart et al. 2019a)

- In **theory**, MCMC will converge to the target distribution.
- In **practice**, methods based on random walk sampling (RWS) can handle a limited number of unknowns (**< 100 in our experience**)
- RWS produces samples that are highly **correlated**.

THE RANDOM WALK SAMPLER
PRACTICALLY LIMITS THE
NUMBER OF UNKNOWNNS WE
CAN USE

Potential
Energy

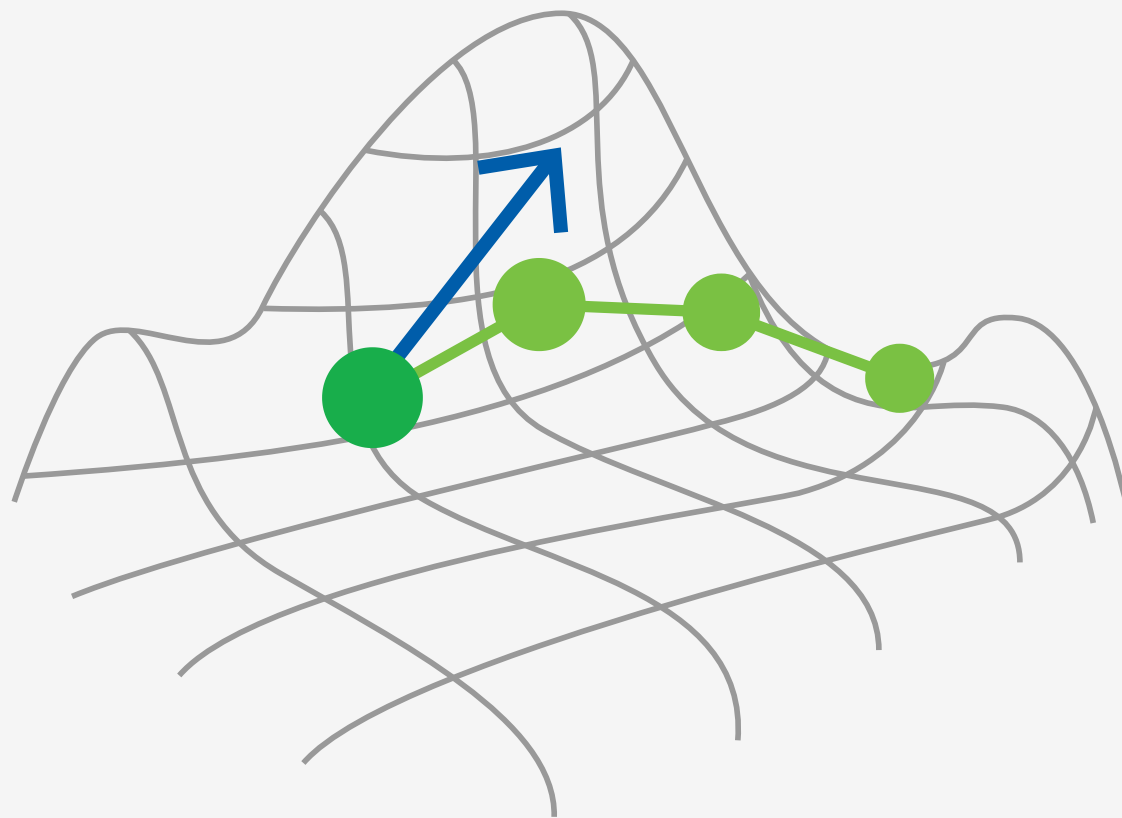
Kinetic Energy

$$H(q, p) = U(q) + K(p)$$

Hamiltonian

Position

Momentum



$$\begin{aligned}\frac{dq}{dt} &= \frac{\partial H(q, p)}{\partial p} \\ \frac{dp}{dt} &= -\frac{\partial H(q, p)}{\partial q}\end{aligned}$$

PROBLEM: GRADIENT COMPUTATION IS EXPENSIVE!



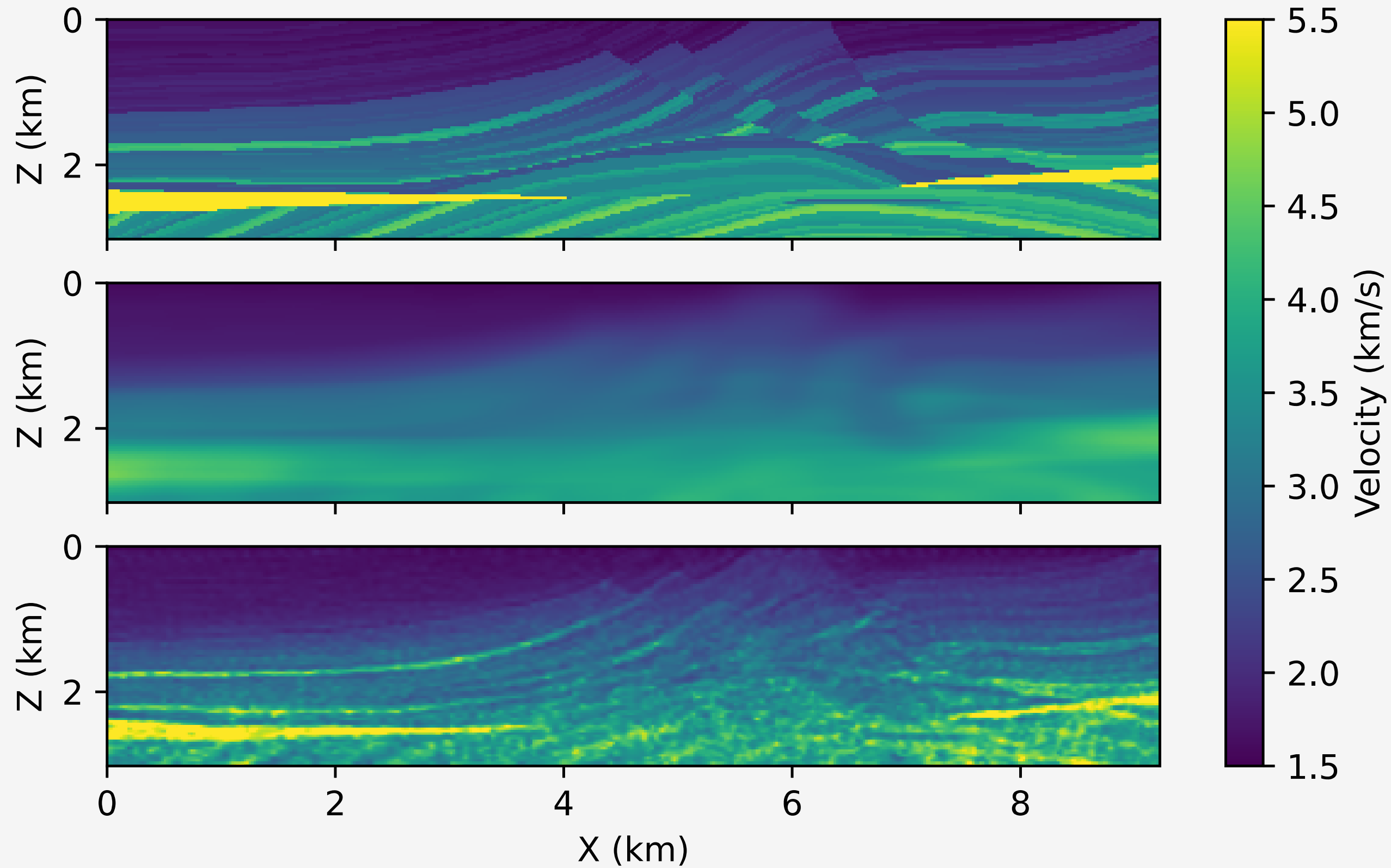
NEURAL NETWORK-ENHANCED TWO-STAGE HMC (NNHMC)

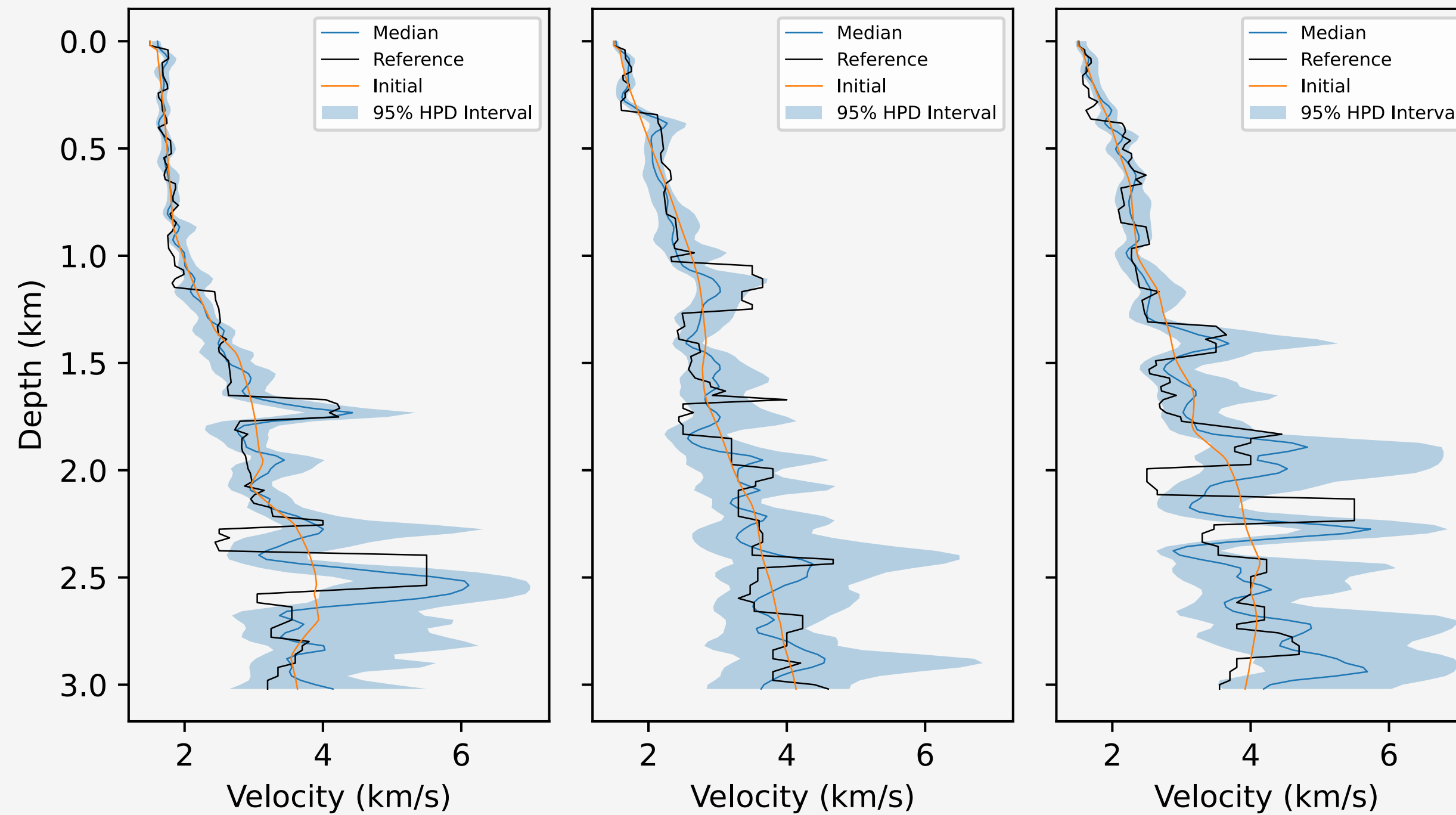


HMC REQUIRES USER-
SPECIFIED PARAMETERS TO
DISCRETIZE THE HAMILTONIAN
DYNAMICS

THE NO-U-TURN SAMPLER (NUTS)







Vertical slices with HPD intervals

Posterior Distributions

