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Geophysical Inversion using Parametric Variational Inference GeoPVI User Guide

Xuebin Zhao¹ and Andrew Curtis¹

¹ School of GeoSciences, University of Edinburgh, United Kingdom

E-mail: xuebin.zhao@ed.ac.uk, andrew.curtis@ed.ac.uk

SUMMARY

This user guide describes how to use GeoPVI, a Python package for Geophysical inversion using Parametric Variational Inference methods. In GeoPVI, a variational distribution is defined to approximate the Bayesian posterior probability distribution function (pdf) and is represented by parametric (analytic) expressions. GeoPVI provides a suite of tools to facilitate variational Bayesian inversion, allowing users to efficiently model uncertainties in their geophysical parameter estimations. It differs from other non-parametric variational methods, such as Stein variational gradient descent (SVGD), in the sense that the posterior probability distribution is parmetrised explicitly and (semi-)analytically, such that posterior probability value of any model sample can be evaluated efficiently after inversion (without invoking Bayes' rule). This parametric feature can be used to vary prior information post Bayesian inversion using a variational prior replacement method. GeoPVI currently features both mean field and full rank automatic differentiation variational inference (ADVI), physically structured variational inference (PSVI), normalising flows and boosting variational inference (BVI). Future updates will expand this package to incorporate other parametric variational methods that have been tested in geophysics.

1 INTRODUCTION

Many geophysical problems can be cast as typical inverse problems that estimate a set of latent model parameters from observed geophysical data. Given that geophysical inversion often yields non-unique solutions, it is important to find all possible parameters that fit the observed data to within data errors to estimate the corresponding uncertainties.

This can be achieved under a probabilistic framework by calculating the so-called *posterior* probability distribution function (pdf) $p(\mathbf{m}|\mathbf{d}_{obs})$ using Bayes' rule – a problem often referred to as Bayesian inference

$$p(\mathbf{m}|\mathbf{d}_{obs}) = \frac{p(\mathbf{d}_{obs}|\mathbf{m})p(\mathbf{m})}{p(\mathbf{d}_{obs})}$$
(1)

where vectors \mathbf{m} and \mathbf{d}_{obs} stand for model parameter and observed data, respectively. Term $p(\mathbf{m})$ describes the *prior* information available on \mathbf{m} , $p(\mathbf{d}_{obs}|\mathbf{m})$ is the *likelihood* function, and $p(\mathbf{d}_{obs})$ is a normalisation constant called the *evidence*.

Monte Carlo sampling methods are often used to solve such problems by generating an ensemble of posterior samples that embody the Bayesian solution. However, these methods can be computationally expensive, especially for large-scale problems that contain thousands, even millions of unknown model parameters.

Alternatively, variational inference solves Bayesian inversion problems under an optimisation framework, by seeking one optimal approximation to the true posterior pdf from a family of predefined and known distributions. Variational methods therefore convert a random sampling problem into a numerical optimisation one, while still providing fully probabilistic results. Therefore, the method can be more efficient than random sampling methods and can provide better scaling to higher dimensional problems (Bishop 2006; Blei et al. 2017; Zhang et al. 2018).

Many variational methods have been applied to geophysical inverse problems and are proved to be far more efficient than Monte Carlo methods. Currently, these methods can be categorised into two groups based on their approaches to represent an approximate posterior distribution. The first group generates a set of posterior samples to implicitly approximate statistical information of the posterior distribution, such as Stein variational gradient descent (SVGD – Zhang & Curtis 2020;

Zhang et al. 2023). The second group directly models a parametric representation (expression) of the posterior pdf (Zhao et al. 2021; Zhao & Curtis 2024b,c). This parametric (closed form) nature enables efficient evaluation of posterior probability value of any model sample once the inversion process is finished, without performing forward simulation. We can further make use of this feature to update prior information post Bayesian inversion using a variational prior replacement methodology Zhao & Curtis (2024a).

GeoPVI implements on a suite of methods that optimise a parametric posterior pdf, currently including automatic differentiation variational inference (ADVI – Kucukelbir et al. 2017), physically structured variational inference (PSVI – Zhao & Curtis 2024c), normalising flows (Rezende & Mohamed 2015), and boosting variational inference (BVI – Guo et al. 2016; Miller et al. 2017). Future updates will expand this package to incorporate other parametric variational inference methods.

2 INSTALLATION

2.1 Requirements

GeoPVI uses NumPy to perform basic numerical implementations, and uses automatic differentiation library provided by PyTorch to optimise variational parameters. SciPy is used to implement some scientific calculations, such as sparse matrix representation and inverse of large-scale matrix. In addition, multiprocessing package is used to parallelise forward simulation for different model samples. Alternatively, Dask can be used for this purpose, which provides an effective interface to use multiple CPU cores from modern high performance computing clusters. In a 2D FWI example provided below, forward simulation code is written by C and Cython. These packages can be installed using the Python's package-management system pip:

```
$ pip install -r requirements.txt
```

2.2 Installation

Once you have downloaded the GeoPVI package and installed the above required packages

```
$ cd GeoPVI
2 $ sh setup.sh install
```

We recommend to install GeoPVI in an editable mode. Alternatively, if you do not want to install the package, simply do

```
sh setup.sh
```

Then, you need to tell scripts which use the GeoPVI package where the package is. For example, simply run a script with

```
$ PYTHONPATH=/your/GeoPVI/path python fwi2d.py
```

3 IMPLEMENTATION

In variational inference, a family of distributions (called the variational family) $Q(\mathbf{m}) = \{q(\mathbf{m})\}$ is defined, from which an optimal member $q^*(\mathbf{m})$ is selected to best approximate the true posterior distribution $p(\mathbf{m}|\mathbf{d}_{obs})$. This optimal distribution can be found by maximising the following evidence lower bound (ELBO $[q(\mathbf{m})]$) (Zhao & Curtis 2024c):

$$ELBO[q(\mathbf{m})] = \mathbb{E}_{q(\mathbf{m})}[\log p(\mathbf{m}, \mathbf{d}_{obs})] - \mathbb{E}_{q(\mathbf{m})}[\log q(\mathbf{m})]$$
(2)

The calculation of the objective function in equation 2 requires to calculate $\log p(\mathbf{m}, \mathbf{d}_{obs})$ and $\log q(\mathbf{m})$, which can be constructed separately using GeoPVI.

3.1 Forward simulation

The calculation of $\mathbb{E}_{q(\mathbf{m})}[\log p(\mathbf{m}, \mathbf{d}_{obs})]$ in equation 2 requires the logarithmic joint probability value $\log p(\mathbf{m}, \mathbf{d}_{obs})$, which is the summation of the logarithmic likelihood and prior values, according to Bayes' rule (equation 1). In GeoPVI, this can be done by defining a function that takes model parameter \mathbf{m} as input and calculates $\log p(\mathbf{m}, \mathbf{d}_{obs})$. For example,

```
def log_prob(m):

# input has a shape of (num_of_samples, dim)

# requires a forward solver that takes x as input and outputs the modelled data

# then calculate the log-likelihood and log-prior probability values
```

```
logp = log_prior + log_like
return logp
```

where logp, log_like and log_prior are the logarithmic joint probability, likelihood and prior values, respectively.

Currently, GeoPVI supports Normal and Uniform prior distributions. User-specified prior distributions can also be defined using a template in *geopvi/prior.py*. To use the Normal and Uniform prior pdfs provided in GeoPVI:

```
from geopvi.prior import Uniform, Normal

prior = Uniform(lowerbound, upperbound, smooth_matrix = L)
```

This defines a Uniform prior distribution bounded by a lowerbound and upperbound. If a smoothed prior distribution is desired, then pass a smooth matrix L to smooth_matrix. Otherwise, it can be set as None to represent prior information without smoothing.

To define a Normal prior distribution:

```
prior = Normal(loc, std = None, covariance = None, scale_tril = None, precision = None)
```

where loc is the mean vector of the Normal distribution. The covariance matrix can be represented by one of the following four options: std (1D array representing standard deviations), covariance (positive-definite covariance matrix), scale_tril (lower-triangular factor – Cholesky decomposition – of the covariance matrix, with positive-valued diagonal entities), or precision (positive-definite precision matrix). Note that only one of the four options can be specified.

The log-prior probability value for a given model sample m can be calculated using GeoPVI built-in function:

```
log_prior = prior.log_prior(m)
```

If the likelihood function is assumed to be a Gaussian distribution, then it can be calculated by the misfit value between observed data and synthetic data generated by a forward modeller. Note that if the forward simulation code is written using PyTorch's built-in functions, the computational graph is already able to be recorded by PyTorch autograd package, such that the gradient of the likelihood value with respect to input sample m can be calculated directly through automatic differentiation. Otherwise, if the forward modeller itself is not auto-differentiable or rely on non-

PyTorch libraries (e.g., NumPy or C), users will need to wrap it to interface with the autograd engine. Below is an example for this purpose

```
from torch.autograd import Function
 class ForwardModel(Function):
 # This class is used to register the output of the external (non-PyTorch) code
 # into automatic differentiation engine
 @staticmethod
 def forward(ctx, input, your_external_solver):
 # forward simulation using external solver
 # returns simulation output (log-likelihood value in this case)
 # also register its gradient w.r.t. input, and save for back propagation
 output, grad = your_external_solver(input)
 ctx.save_for_backward(input, torch.tensor(grad))
 return torch.tensor(output)
 @staticmethod
 def backward(ctx, grad_output):
 # this function returns the gradient of loss function w.r.t the input tensor
 # therefore, the return shape should be the same as the shape of input tensor
 input, grad = ctx.saved_tensors
 grad_input = (grad_output[...,None] * grad)
 return grad_input, None
 def your_external_solver(m):
 # forward simulation of m
 # return log-likelihood value and its gradient w.r.t m
 \mbox{\tt\#} grad should have the same shape as that of \mbox{\tt m}
 return log_like, grad
 def log_like_for_back_propagation(m):
 # m has a shape of (num_of_samples, dim)
 # forward simulation of m: your_external_solver(m)
 log_like = ForwardModel.apply(m, your_external_solver)
return log_like
```

where your_external_solver is an external (non-PyTorch) Python function that performs forward simulation for input sample m. It returns log-likelihood value log_like and its gradient with respect to m. By calling function log_like_for_back_propagation, the calculated log-

likelihood value is registered in the automatic differentiation graph and can be back propagated during training process. More details can be found in a 2D FWI example provided below and on the PyTorch website.

GeoPVI provides two forward functions used typically in seismic imaging problems: fast marching method (FMM) for travel time tomography and 2D acoustic full waveform simulation for full waveform inversion (FWI). More details can be found in *geopvi/forward/tomo2d/posterior.py* and *geopvi/forward/fwi2d/posterior.py*.

3.2 Flows-based variational distribution

In GeoPVI, we build a parametric variational distribution $q(\mathbf{m})$ using a normalising flows model (Rezende & Mohamed 2015; Zhao et al. 2021). Normalising flows provide a flexible way to construct a parametric probability distribution by passing a simple and known initial (base) probability distribution (e.g., a standard Normal or a Uniform distribution) through a series of invertible and differentiable transforms, called the flows. To construct such a model:

```
from geopvi.vi.models import VariationalDistribution
```

Users can choose specific transforms (flows) based on problems at hand. Python script geopvi/vi/flows.py provides several commonly used flows in literature (more available flows will be implemented in future release). Below is an example to construct a flows-based model to perform automatic differentiation variational inference (ADVI) and physically structured variational inference (PSVI):

```
from geopvi.vi.flows import Linear, Real2Constr

flows = [Linear(dim, kernel = 'structured', mask = off_diag_mask)]

flows.append(Real2Constr(lower = lowerbound, upper = upperbound))

variational_pdf = VariationalDistribution(flows, base = 'Normal')
```

This initialises a variational model using one Linear flow and one Real2Constr flow, applied on a standard normal base probability distribution. Real2Constr transforms model parameters from the space of real numbers into a constrained space bounded by hyper-parameters lowerbound and upperbound. If model parameters do not have lower/upper bound, the corresponding parameter

can be set to None. Linear flow performs a linear transform on the input vector x

$$z = \mu + Lx \tag{3}$$

The variational parameters are μ and L in this case. If x follows a standard Normal distribution $\mathcal{N}(0, I)$, one single Linear flow essentially transforms $\mathcal{N}(0, I)$ into any normal distribution defined by $\mathcal{N}(\mu, LL^T)$. In Linear flow, dim is the dimensionality (number of unknown parameters to be inferred) of vector x. Parameter kernel controls different methods used to construct matrix L, and its available options include 'diagonal', 'fullrank' and 'structured', each corresponding to a Gaussian distribution with a diagonal, full, and structured (desired correlation information - see details in Zhao & Curtis 2024c) covariance matrix, respectively. These three options can be used to perform mean field ADVI, full rank ADVI, and PSVI, followed by a Real2Constr flow. If 'structured' is used, a mask defining the desired correlation structure (off_diag_mask) is required (see details in examples provided in GeoPVI). Otherwise it can be ignored safely by setting it to None.

Perform

```
x = variational_pdf.sample_from_base(100) # x has a shape of (100, dim)
multiple m, logq = variational_pdf(x)
```

which generates 100 samples x from base distribution of the flows-based model, and transforms them to get samples m distributed according to the constructed variational distribution $q(\mathbf{m})$ and obtain their corresponding logarithmic probability values $\log q(\mathbf{m})$.

After obtaining the two ingredients $\log q(\mathbf{m})$ and $\log p(\mathbf{m}, \mathbf{d}_{obs})$, we can estimate the ELBO in equation 2 using Monte Carlo integration with n random samples drawn from $q(\mathbf{m})$

$$ELBO[q(\mathbf{m})] \approx \frac{1}{n} \sum_{i} (\log p(\mathbf{m}, \mathbf{d}_{obs}) - \log q(\mathbf{m}_{i}))$$
 (4)

Define a loss function to be the negative ELBO, which can be calculated by performing

```
x = variational_pdf.sample_from_base(100) # x has a shape of (100, dim)
m, logq = variational_pdf(x)
logp = log_prob(m)
# -ELBO can be estimated using Monte Carlo integration
```

```
negative_elbo = -torch.mean(logp - logq)
```

The optimal variational distribution can be found by iteratively minimising this loss function until convergence. This training process can be accomplished by calling a Variational_Inversion class defined in GeoPVI:

```
from geopvi.vi.models import VariationalInversion
inversion = VariationalInversion(variationalDistribution = variational_pdf,
log_posterior = log_prob)
```

Then, variational inversion can be performed by:

which updates the flows-based variational model using the Adam optimizer for 1000 iterations with 10 samples per iteration for Monte Carlo integration in equation 4. This function returns the negative_elbo value within each iteration. Once the training process is finished, posterior samples can be obtained by sampling from the variational distribution

```
samples = variational_pdf.sample(1000)
```

3.3 BVI-based variational distribution

In addition to the flows-based variational model, GeoPVI also provides a second approach to build a variational distribution through boosting variational inference (BVI), in which a mixture of simple (e.g., normal) distributions is used to approximate the posterior probability distribution. BVI trains and adds one component distribution to the mixture distribution once a time using a greedy algorithm

$$q^{t}(\mathbf{m}) = (1 - w_t)q^{t-1}(\mathbf{m}) + w_t g_t(\mathbf{m})$$
(5)

where $q^t(\mathbf{m})$ is a mixture of t distributions, obtained by combining the previous mixture distribution $q^{t-1}(\mathbf{m})$, weighted by $(1-w_t)$, with the new component $g_t(\mathbf{m})$ weighted by w_t . New component distribution $g_t(\mathbf{m})$ is found by maximising a so-called residual-ELBO (RELBO):

$$RELBO[g_t(\mathbf{m})] = \mathbb{E}_{g_t(\mathbf{m})}[\log p(\mathbf{m}, \mathbf{d}_{obs})] - \mathbb{E}_{g_t(\mathbf{m})}[\log q^{t-1}(\mathbf{m})] - \lambda \mathbb{E}_{g_t(\mathbf{m})}[\log g_t(\mathbf{m})]$$
 (6)

More theoretical details can be found in Zhao & Curtis (2024b).

To run BVI,

```
from geopvi.boostingVI.component import GaussianComponent

gaussian = GaussianComponent(dim, kernel = 'diagonal', constrained = False)
```

this defines a Gaussian component distribution with a dimensionality of dim and a diagonal covariance kernel (similar to the Linear flow introduced above). In this case, for each component we need to optimise its mean and standard deviation vectors. Similarly, if the model parameter m to be inverted is defined within a constrained space, then simply set parameter constrained = True, otherwise set it as False.

A BVI model built by boosting a set of Gaussian distributions can be constructed by

```
from geopvi.boostingVI.bvi import BoostingGaussian

bvi_pdf = BoostingGaussian(componentDistribution = gaussian, log_posterior = log_prob)
```

To train this BVI model, simply run

```
result = bvi_pdf.update(ncomponent = 5, n_iter = 1000, nsample = 10,

optimizer = 'torch.optim.Adam')
```

which builds a mixture of 5 Gaussian distributions, with each updated for 1000 iterations and 10 samples per iteration to calculate the expectations in equation 6. After training, the inversion result, defined by a Python dictionary containing optimised parameters for each component, is returned. Similarly, posterior samples can be obtained by

```
samples = bvi_pdf.sample(1000)
```

4 EXAMPLES

In this section we show examples of using the GeoPVI package to solve 2D travel time tomography, and 2D and 3D full waveform inversion (FWI) problems. Note that the implementation of 3D FWI requires a suitably efficient 3D forward and adjoint wavefields solver, which is not included in the current release of GeoPVI.

4.1 2D travel time tomography

For 2D travel time tomography we use a fast marching method (FMM) to model the travel time data (Rawlinson & Sambridge 2005), which can be found in *geopvi/forward/tomo2d* folder. The implementation details of variational travel time tomography can be found in *examples/tomo2d*. To use the code, simply do

```
the contraction of the contracti
```

The input and output of the code are described below. Note that hyperparameters for building a variational model and performing optimisation are defined as a Python argparser.parse_args() object in *examples/tomo2d/tomo2d.py*, and are self-explained.

4.1.1 Input

The following input files are stored in examples/tomo2d/input.

config.ini – This file contains nuisance parameters required by the 2D FMM code.

prior.txt – This file contains the hyperparameters of the prior distribution whose name is given by the argument args.prior_param in tomo2d.py. If a Uniform distribution is used, the file contains the lower (first column) and upper (second column) bounds for each parameter. In the case of a normal distribution, the file contains the mean (first column) and standard deviation (second column) for each parameter.

sources.txt and receivers.txt – Source and receiver coordinates files. Each row contains the (x, y) location for a source (or a receiver).

traveltime.txt – Travel time data file, whose name is given by the argument args.datafile in tomo2d.py. Each row contains a travel time. The order is first source and first receiver, first source and second receiver, ..., first source and last receiver; second source and first receiver, second source and second receiver, ... and so on. If the travel time for a given source-receiver pair is not available, set it to zero.

4.1.2 output

The outputs are stored in the directory *examples/tomo2d/output*.

*parameter.npy – A NumPy NPY format file which contains the optimised variational parameters. If Linear flow is used, this output file includes the vector $\boldsymbol{\mu}$ (the first dim elements), diagonal elements of the matrix \mathbf{L} (the second dim elements) and off-diagonal elements of the matrix \mathbf{L} (the remaining elements). If mean field ADVI is performed, no off-diagonal elements is output. Parameters $\boldsymbol{\mu}$ and \mathbf{L} can then be used to define a Gaussian variational distribution $\mathcal{N}(\boldsymbol{\mu}, \mathbf{L}\mathbf{L}^T)$.

*model.pt – A PyTorch .pt format file which stores the entire variational model using Python's pickle module. This file can be used to resume training (optimisation) process from a certain checkpoint, or to load the obtained variational model for post-inversion analysis.

*samples.npy - A NumPy NPY format file which contains the posterior samples.

*loss.txt – A text file which stores the negative ELBO values during training.

4.1.3 Examples

A synthetic example described in Zhao et al. (2021) can be found in the folder examples/tomo2d.

4.2 2D FWI

In a 2D FWI example, the forward modelling of waveform data is performed using a finite difference method, and the gradient of the misfit function with respect to velocity parameters is obtained using the adjoint method. Its implementation details can be found in *geopvi/forward/fwi2d*. The example code that uses variational inference to solve this 2D FWI problem is in *examples/fwi2d/fwi2d.py*. To run the code, simply do

```
cd examples/fwi2d
sh run.sh
```

The input and output of the code are described below. Note that hyperparameters for building a variational model and performing optimisation are defined as a Python argparser.parse_args() object in *examples/fwi2d/fwi2d.py*, and are self-explained.

4.2.1 Input

The following input files are stored in examples/fwi2d/input.

config.ini – This file contains the control parameters to perform forward and adjoint simulation.

An example is shown below, in which every line beginning with '#' is a comment.

```
# nuisance parameter for 2D fwi code
2 [FWI]
3 # grid points in x direction (nx)
4 \text{ nx} = 250
5 # grid points in z direction (nz)
6 \text{ nz} = 110
7 # depth of (water) layer where velocity is set as true value during inversion
8 layer_fixed = 10
9 # true velocity value in the layer with fixed velocity value
vel_fixed = 1950
# pml points (pml0)
12 pml = 10
# Finite difference order (Lc)
_{14} Lc = 3
15 # Method to calculate Laplace operator: 0 for FDM and 1 for PSM, (laplace_slover)
16 laplace_slover = 0
# Total number of sources (ns)
18 \text{ ns} = 12
# Total time steps (nt)
20 nt = 2001
21 # Shot interval in grid points (ds)
22 ds = 20
23 # Grid number of the first shot to the left of the model (ns0)
ns0 = 15
25 # Depth of source in grid points (depths)
_{26} depths = 1
27 # Depth of receiver in grid points (depthr)
_{28} depthr = 9
29 # Receiver interval in grid points (dr)
# Time step interval of saved wavefield during forward (nt_interval)
32 nt_interval = 2
^{33} # Grid spacing in x direction (dx)
```

```
34 dx = 20.0
35 # Grid spacing in z direction (dz)
36 dz = 20.0
37 # Time step (dt)
38 dt = 0.002
39 # Donimate frequency (f0)
40 f0 = 10.0
```

prior.txt – This file contains the hyperparameters of the prior distribution whose name is given by the argument args.prior_param in fwi2d.py. If a Uniform distribution is used, the file contains the lower (first column) and upper (second column) bounds at each depth (Z direction). In the case of a normal distribution, the file contains the mean (first column) and standard deviation (second column) at each depth. The prior distribution is assumed to be laterally homogeneous.

waveform.npy – A NumPy NPY format file which contains the observed waveform data. The file name is given by the argument args.datafile in *fwi2d.py*. The data is stored in a NumPy array with a shape (ns, nr, nt) where ns is the number of sources, nr is the number of receivers and nt is the number of time points of each trace.

4.2.2 output

The outputs are stored in the directory examples/fwi2d/output.

*parameter.npy – A NumPy NPY format file which contains the optimised variational parameters. If Linear flow is used, this output file includes the vector μ (first dim elements), diagonal elements of the matrix L (second dim elements) and off-diagonal elements of the matrix L (remaining elements). If mean field ADVI is performed, no off-diagonal elements is output. Parameters μ and L can then be used to define a Gaussian variational distribution $\mathcal{N}(\mu, \mathbf{LL^T})$.

*model.pt – A PyTorch .pt format file which stores the entire variational model using Python's pickle module. This file can be used to resume training (optimisation) process from a certain checkpoint, or to load the obtained variational model for post-inversion analysis.

*samples.npy - A NumPy NPY format file which contains the posterior samples.

*loss.txt – A text file which stores the negative ELBO values during training.

4.2.3 Examples

A synthetic example that uses a part of the Marmousi model can be found in the folder *examples/fwi2d*. The details of this example are described in Zhao & Curtis (2024c).

4.3 3D FWI

In a range of geophysical applications one may need to solve a large scale inverse problem. This has become possible in the light of the development of modern high performance computation (HPC) facilities. In this section we take 3D FWI as an example to show how to use the GeoPVI package to solve large scale problems.

Throughout this suite of methods, the forward simulation can be computed independently which makes it easy to parallelise. For example, one can use the Dask package to parallelise the computation using multiple CPU cores.

```
from dask.distributed import Client
def fwi_i(model):
      # forward midelling for the ith model sample
      log_like, grad = external_fwi_code(model)
      return log_like, grad
8 def fwi(models):
      # forward modelling for all model samples
      # First, create a dask client for cros-core parallisation
      client = Client()
      # submit simulation job to dask client
      futures = []
      for i in range(models.shape[0]):
          futures.append(client.submit(fwi_i, model, pure = False))
      # gather the computational results
      results = client.gather(futures)
19
      log_like = np.zeros((models.shape[0],))
      grad = np.zeros_like(models)
21
      for i in range(models.shape[0]):
```

The function fwi can thereafter be used with the GeoPVI package to implement 3D FWI, just as displayed above. In this way the code can use many cores that are available on modern HPC facilities to provide an efficient method. Note that the code can be further improved by combining the modern Distributed Resource Management Systems and by using minibatch data sets. For a detailed example that uses the SGE system, see *geopvi/forward/fwi3d* and *examples/fwi3d/fwi3d.py*.

5 TERMS OF USE AND DISCLAIMER

To the best of our knowledge this code is correct, error-free as of June 2024. We can not guarantee its accuracy however. If after your own detailed investigation you believe there are errors/bugs in the code or examples, please contact Xuebin Zhao at xuebin.zhao@ed.ac.uk, or Andrew Curtis at andrew.curtis@ed.ac.uk.

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