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# Measuring higher mode surface wave dispersion using a transdimensional Bayesian approach

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## SUMMARY

We present a method for measuring the dispersion of higher mode surface wave phase velocities from a single seismogram using a hierarchical transdimensional Bayesian approach. The 1-D shear velocity profiles down to 800 km depth between sources and seismic stations are regarded as the controlling parameters to tune the phase velocities of fundamental and higher modes. The misfits between synthetics and real waveforms indicate whether the phase velocities are recovered well from the data. We use Monte Carlo Markov chains to approximate the posterior distribution of each model parameters and assess the uncertainties from these probability density functions. These techniques can test models of varying dimensions while being parsimonious, thereby letting the data themselves control the complexity of the solution. Another advantage is that the algorithm can decide how much data noise is needed in order to explain the data without overfitting them. The data noise can be treated as an unknown and different noise levels can be applied to the different time windows considered. The posterior noise distributions can then be used as an indicator of the quality of the waveform fit within each frequency–time window. We considered phase velocities between 50 and 200 s for each mode and performed a reliability analysis to determine which modes and periods are reliably constrained. In this paper, we first present the method and demonstrate its feasibility with synthetic tests, which show that the technique is robust. We then illustrate it with applications to real data. We applied the method to two paths sampling Australia using earthquakes at regional distances and obtained results that agree well with previous studies. The new method can be used in regional and global tomographic studies to obtain phase velocity maps and 3-D models of seismic velocities and anisotropy at depths that are not well resolved by fundamental mode surface waves or body waves.

**Key words:** Inverse theory; Probability distributions; Tomography; Statistical seismology; Surface waves and free oscillations.

## 1 INTRODUCTION

Higher mode surface waves carry unique, independent constraints on structure at greater depths than commonly used fundamental mode surface waves, and thus enhance the vertical resolution of surface wave tomographic models in the deep upper mantle, transition zone and uppermost lower mantle. However, most surface wave tomography studies are still limited to fundamental surface waves because traditional surface wave measurements can only be readily applied to fundamental modes and measuring higher modes is complicated. Only in very specific and rare cases has it been possible to isolate and measure the dispersion of the first higher mode Rayleigh wave using single mode dispersion methods (Crampin 1964; Roult & Romanowicz 1984). The difficulties in measuring higher modes

are mainly due to the fact that their group velocities overlap significantly in a broad frequency range and thus they do not appear as a clear wave train on the seismogram. When distant events are used, a small difference in group velocity can be sufficient to obtain mode separation using numerical techniques, but it is not always enough.

One way to take advantage of the information contained in higher mode surface waves is to retrieve velocity structure directly, without extracting dispersion curves first. Most of these techniques involve determining path-averaged shear wave velocity structure for multiple station–receiver pairs and combining them to obtain 3-D models of Earth's internal structure. Cara & Lévêque (1987) determined path-averaged  $V_S$  models from a single seismogram using mode branch cross-correlation functions (Lerner-Lam & Jordan 1983) as secondary observables. This method, which was designed to

minimize the dependence on the starting model, was later automated and combined with a regionalization scheme (Montagner 1986) to model 3-D velocities (Debayle 1999) and anisotropy (Maggi *et al.* 2006). Nolet (1990) developed a two-step nonlinear partitioned waveform inversion (PWI) technique that was automated by Lebedev & Nolet (2003). Using synthetic seismogram calculations, the PWI finds along-path velocity models that fit recorded waveforms and inverts the velocity models obtained for all paths to constrain 3-D structure. Li & Romanowicz (1995, 1996) were also able to obtain the averaged structure along the great circle between the source and receiver using a waveform modelling technique that included nonlinear asymptotic coupling theory and coupling across different branches.

Despite the difficulty, several attempts have been made to separate higher modes. Some methods operate in the frequency–wavenumber ( $\omega - k$ ) domain and use seismic arrays (Cara 1973, 1978, 1979; Nolet 1975). While this kind of direct measurement is fast to implement, there are a few limitations: (i) the method can only be applied in regions with dense arrays and thus cannot be applied at the global scale, (ii) it requires the usage of linear regional seismic arrays approximately aligned with the epicentre, which reduces the geographical ray coverage, and (iii) variations within an array cannot be assessed (Laske & Widmer-Shnidrig 2015).

Another method, based on mode branch stripping, was developed by van Heijst & Woodhouse (1997). The principle behind this technique is that the signals of overtone branches can be isolated by fitting the cross-correlation function of a single mode using mode branch cross-correlation functions. However, while this method does not rely on arrays and thus can be used in global tomography, it does not work well with epicentral distances shorter than  $30^\circ$  (van Heijst & Woodhouse 1999) since individual higher modes cannot be separated from others in a seismogram when the path is too short. Additionally, it might be difficult to separate one single mode from observed seismograms when that mode is contaminated by interference from other modes.

Other researchers have utilized waveform fitting techniques to extract higher mode dispersion. Stutzmann & Montagner (1993) and Stutzmann *et al.* (1994) developed a waveform fitting technique to invert phase velocity dispersion and velocity structure at depth in successive steps. Beucler *et al.* (2003) proposed a roller-coaster technique in which the phase velocity perturbation is obtained by fitting the synthetic seismogram to the real data. Yoshizawa & Kennett (2002, 2004) and Yoshizawa & Ekström (2010) used the Neighbourhood Algorithm (Sambridge 1999a) to search the model space for multimode dispersion in a nonlinear waveform inversion. Visser *et al.* (2007) developed a method similar to that of Yoshizawa & Kennett (2002) to estimate multimode dispersion curves for global tomography, though they were additionally able to obtain quantitative uncertainties on the phase speed measurements by performing a Bayesian appraisal of the models sampled (Sambridge 1999b). Uncertainties were determined from the whole ensemble of dispersion models obtained instead of taking the standard deviation of the 1000 best models as in Yoshizawa & Kennett (2002).

In our study, we developed a waveform modelling method using a reversible jump Markov Chain Monte Carlo (rj-MCMC) inversion (Bodin & Sambridge 2009) to extract higher mode dispersion curves and their uncertainties at periods between 50 and 200 s. Like the methods of Yoshizawa & Kennett (2002), Yoshizawa & Ekström (2010), Visser *et al.* (2007), and Visser (2008), our technique has the advantage of sampling the model space instead of choosing one model among many possible solutions with a strong regularization

scheme. It therefore provides more reliable posterior model uncertainties than regularized inversions. Additionally, contrary to those previous studies in which the authors adopted a fixed dimension model space chosen *a priori*, the rj-MCMC sampler offers a way to treat the dimension of the model space as variable. In addition, because we use a Hierarchical Bayes approach, data noise parameters are treated as unknowns in the inversion, which accounts for the part of the signal that we are not able to explain with our forward theory and the chosen parametrization. This allows the model to explain the data without overfitting them. In this paper, we first present the method and validate it using a synthetic test, and then show two real data examples.

## 2 Method

Our goal is to measure phase velocity dispersion for fundamental and higher mode Rayleigh waves using waveform modelling. To do this, we seek a large number of path-averaged 1-D shear-velocity models that can fit the filtered waveform, and the resulting models are employed to calculate dispersion curves. These path-specific 1-D models represent the average fundamental and higher mode dispersion curves for the chosen source–receiver path as in Yoshizawa & Kennett (2002), Yoshizawa & Ekström (2010), Visser *et al.* (2007) and Visser (2008).

### 2.1 Waveform Modelling

A synthetic seismogram ( $s$ ) can be calculated by summation of normal modes ( $m$ ) of amplitude  $A_m$  in the frequency domain ( $\omega$ ) for a 1-D model as follows (e.g. Dahlen & Tromp 1998):

$$s(\omega) = \sum_m A_m(\omega) \exp[i\omega\Delta/c_m(\omega)], \quad (1)$$

where  $c_m(\omega)$  is the phase velocity of the mode  $m$  at angular frequency  $\omega$  and  $\Delta$  is the epicentral distance. The relationship between seismograms and their corresponding velocity model is thus highly nonlinear. A fully nonlinear approach, which does not require any partial derivatives with respect to model parameters or any strong constraints on parametrization, is therefore desirable to perform waveform inversion.

Fortran code Mineos (Masters *et al.* 2011) can be used to compute synthetic seismograms for a reference 1-D model of Earth's interior by normal mode summation. The advantage of such formulation is that it builds a direct connection between seismic waveforms and phase velocities. However, the calculation of normal mode eigenfunctions and eigenfrequencies for a given mantle model is very time-consuming and thus we cannot use the fully nonlinear formulation that is eq. (1) at each iteration of an MCMC scheme. We thus decided to linearize the forward modelling problem in order to overcome the computational speed limitation. As in Yoshizawa & Kennett (2002), we obtained a first synthetic seismogram for a reference model using fully nonlinear calculations (eq. 1). We then used perturbation theory to update the seismogram for other models generated at each iteration of the Markov Chain. For a small perturbation, the change in mode eigenfrequency can be calculated assuming unperturbed eigenfunctions:

$$\delta \ln(\omega) = \int_0^a \left( \frac{\delta\alpha}{\alpha}(r) K_\alpha(r, \omega) + \frac{\delta\beta}{\beta}(r) K_\beta(r, \omega) + \frac{\delta\rho}{\rho}(r) K'_\rho(r, \omega) \right) dr + \sum_d \delta d [K_d(\omega)]_{-}^{+}, \quad (2)$$

where  $\delta \ln(\omega) = \delta \omega / \omega$ ,  $a$  is the radius of the Earth, and  $\alpha$ ,  $\beta$ ,  $\rho$  and  $d$  are  $P$ -wave velocity,  $S$ -wave velocity, density and radius of discontinuities, respectively.  $K_\alpha$ ,  $K_\beta$ ,  $K'_\rho$  and  $K_d$  are the Fréchet derivatives, which relate the change in wave velocities, density and depth of discontinuity from the reference model to changes in the eigenfrequencies. The Fréchet derivatives can be calculated for each mode using the eigenfunctions determined for the reference model (Woodhouse 1980).

In theory, all parameters ( $\delta\alpha$ ,  $\delta\beta$ ,  $\delta\rho$  and  $\delta d$ ) could be allowed to vary. However, for computational reasons, we decided to only search for shear-wave velocity models. This will unlikely affect the results since  $P$ -wave velocity and density have little influence on the phase velocity perturbation in the frequency range considered (5–20 mHz) as tested by Visser (2008). Similarly, we also neglected perturbations in the depth of mantle discontinuities, which are not well resolved by the data considered (Meier *et al.* 2009) and including them in the parametrization does not significantly affect the resulting velocity models (Weidner *et al.* 2017). Correctly accounting for the Moho depth has been shown to be important as it can have strong nonlinear effects on waveform modelling and phase velocity calculations (Montagner & Jobert 1988). We decided to account for deviations from PREM in Moho depth by adopting a reference model that includes crustal constraints from CRUST1.0 (Laske *et al.* 2013), averaged over the length of the chosen path, and calculating the corresponding shear wave sensitivity kernels. The partial derivatives employed therefore reflect departures of the Moho from PREM. Hence, eq. (2) becomes

$$\delta \ln(\omega) = \int_0^a \frac{\delta \beta}{\beta} K_\beta dr. \quad (3)$$

The updated frequency can then be converted into phase velocity for a normal mode of angular order  $l$  using (Jeans 1923; Dahlen & Tromp 1998):

$$c(\omega^*) = \frac{\omega^* a}{l + 0.5}, \quad (4)$$

where  $\omega^*$  is the updated eigenfrequency for the mode considered. In Fig. 1, we compare synthetic seismograms calculated using the normal mode summation without linearization (eq. 1) and the linearized method based on perturbation theory described above (eq. 3). The results show that the phase velocity information from the linearized perturbation method and the nonlinear formulation match well, especially in the period range we are interested in (50–200 s).

The rj-MCMC method described below performs a guided Monte Carlo sampling of the model space using the values of the misfit between the real and the synthetic seismograms. The data and the synthetics are compared in different time and frequency windows, which are chosen such that they include the fundamental and several higher modes. We used different time and frequency windows to compare the fundamental and higher mode waveforms. We employed group velocities to choose the beginning and the end of the time windows for the fundamental mode and we employed  $S$ - or  $SS$ -wave arrival times and group velocities to choose the higher mode time window as in Lebedev & Nolet (2003). For epicentral distance shorter than  $35^\circ$  the higher mode window start time is set just before the arrival of the  $S$ -wave train. For epicentral distance between  $35^\circ$  and  $70^\circ$  the start time is just after the  $S$  and before the  $SS$  wave train. Group velocities can be picked manually by inspection of the waveform or using a frequency–time analysis (Dziewonski *et al.* 1969). Those windows are summarized in Table 1. An example of time window selection on real data is shown in Fig. 2 for an event in Chiapas, Mexico and station HRV in north eastern America.

Because the amplitude of the synthetic seismograms depends on the scalar moments of the event, and because inaccurate scalar moments in catalogues might cause an increase in the data misfit, we implemented an energy equalization method (Lebedev *et al.* 2005) before calculating the misfit between observed and synthetic seismograms (defined in Section 2.2). The aim is to equalize the energy of the synthetic with that of the data by multiplying the amplitude of the synthetics by the following constant:

$$f^{\text{eq}} = \sqrt{\frac{\sum_{t=t_1}^{t=t_2} d^2(t)}{\sum_{t=t_1}^{t=t_2} s^2(t)}}, \quad (5)$$

where  $d(t)$  and  $s(t)$  are the observed and synthetic seismograms, respectively, filtered in the frequency range considered (5–20 mHz).  $t_1$  indicates the beginning of the overtone time window (window 3 defined in Table 1) and  $t_2$  is the end of the first fundamental mode time window (window 1 in Table 1).

## 2.2 Bayesian inference

In a Bayesian formulation, model parameters are described by probability density functions (PDFs), which effectively indicate the uncertainty associated with these parameters. The aim of a Bayesian inference is to quantify the *a posteriori* probability distribution of all model parameters given some prior information and the data. Bayes' theorem (Bayes 1763) gives the posterior as

$$p(\mathbf{m}|\mathbf{d}_{\text{obs}}) \propto p(\mathbf{d}_{\text{obs}}|\mathbf{m})p(\mathbf{m}), \quad (6)$$

where  $A|B$  means  $A$  given (or conditional on)  $B$ , that is, the probability of having  $A$  when  $B$  is fixed.  $\mathbf{m}$  is the vector of model parameters and  $\mathbf{d}_{\text{obs}}$  is the observed data. The term  $p(\mathbf{d}_{\text{obs}}|\mathbf{m})$  is called likelihood function, which shows the probability of observing data  $\mathbf{d}_{\text{obs}}$  given a particular model  $\mathbf{m}$ . The *a priori* probability of model,  $p(\mathbf{m})$ , contains what we assume about the model  $\mathbf{m}$  before having the observed data. Accordingly, the posterior distribution represents the probability of the model given the observed data. The parts of the model space that are more frequently required by the data than others manifest with greater posterior probability, and hence are more likely to reflect properties of the Earth (Backus 1988).

The likelihood function  $p(\mathbf{d}_{\text{obs}}|\mathbf{m})$  describes the probability of data given the current model. In this study, we compared the waveform fit in three different frequency–time windows, and we assumed the data noise in each window follows a multivariate normal distribution with zero mean and covariance matrix  $\mathbf{C}_{d_i}$ , where  $i$  is the window index. The likelihood term for each frequency–time window is thus written in this form:

$$p_i(\mathbf{d}_i|\mathbf{m}) = \frac{1}{\sqrt{(2\pi)^n |\mathbf{C}_{d_i}|}} \times \exp\left[-\frac{\Phi_i(\mathbf{m})}{2}\right], \quad (7)$$

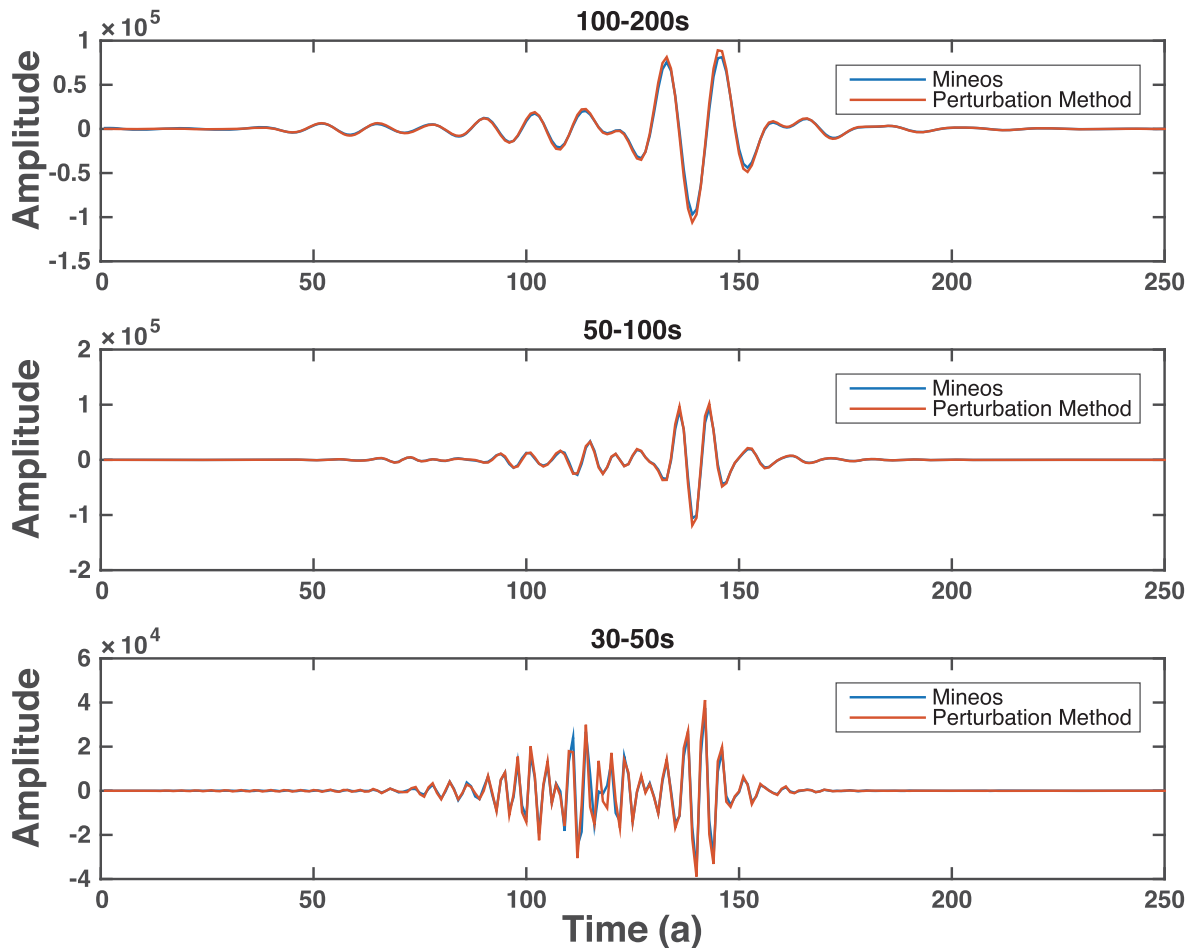
where  $d_i$  is the data in window  $i$ ,  $\Phi_i(\mathbf{m})$  is the function describing the distance between real data and synthetics predicted by the current model. The general expression for  $\Phi_i(\mathbf{m})$  considering correlated noise is (Bodin *et al.* 2012):

$$\Phi_i(\mathbf{m}) = (\mathbf{G}(\mathbf{m}) - \mathbf{d}_i)^T \mathbf{C}_{d_i}^{-1} (\mathbf{G}(\mathbf{m}) - \mathbf{d}_i), \quad (8)$$

where  $\mathbf{d}$  represents the measured data vector and  $\mathbf{G}(\mathbf{m})$  is for the synthetic data predicted by model  $\mathbf{m}$ .

For simplicity, we assumed Gaussian and uncorrelated noise, in which case the covariance matrix is diagonal. Then the equation above becomes

$$\Phi_i(\mathbf{m}) = M_i(\mathbf{m})/\sigma_i^2, \quad (9)$$



**Figure 1.** Comparison between synthetic seismograms calculated with normal mode summation using Mineos without linearization (blue) and with our linearized method (red) in three different period ranges: 100–200 s (top), 50–100 s (middle) and 30–50 s (bottom). The shear velocity model used to compute the seismograms is shown in Fig. 3.

**Table 1.** Selection of frequency–time windows. The first and second window indices correspond to the fundamental mode and the third window is for higher modes. The start and end times for the fundamental mode are determined by the group velocities  $U$  indicated in the table and the event epicentral distance  $\Delta$ . For the higher modes, the beginning of the time window is determined by the  $S$ - or  $SS$ -wave arrival times. The end of the window is determined by  $\Delta/U$ .

Window index	Frequency (mHz)	Start	End
1	5–10	$U \approx 4.45 \text{ km s}^{-1}$	$U = 2.95 \text{ km s}^{-1}$
2	10–20	$U = 4.30 \text{ km s}^{-1}$	$U = 3.20 \text{ km s}^{-1}$
3	10–20	$S$ or $SS$	$U = 4.30 \text{ km s}^{-1}$

where  $\sigma_i$  is the standard deviation of the Gaussian uncorrelated noise in the  $i$ th frequency–time window, that is, the diagonal element of the covariance matrix, and  $M_i$  is defined by

$$M_i(\mathbf{m}) = \sum_{j=1}^{L_i} (d_i^j - f^{\text{eq}} s_i^j)^2, \quad (10)$$

where  $L_i$  is the length of the selected time window,  $d_i$  is the observed data,  $s_i$  is the synthetics calculated for model  $\mathbf{m}$  and  $f^{\text{eq}}$  is the energy equalization factor defined in eq. (5). The expression for the likelihood becomes

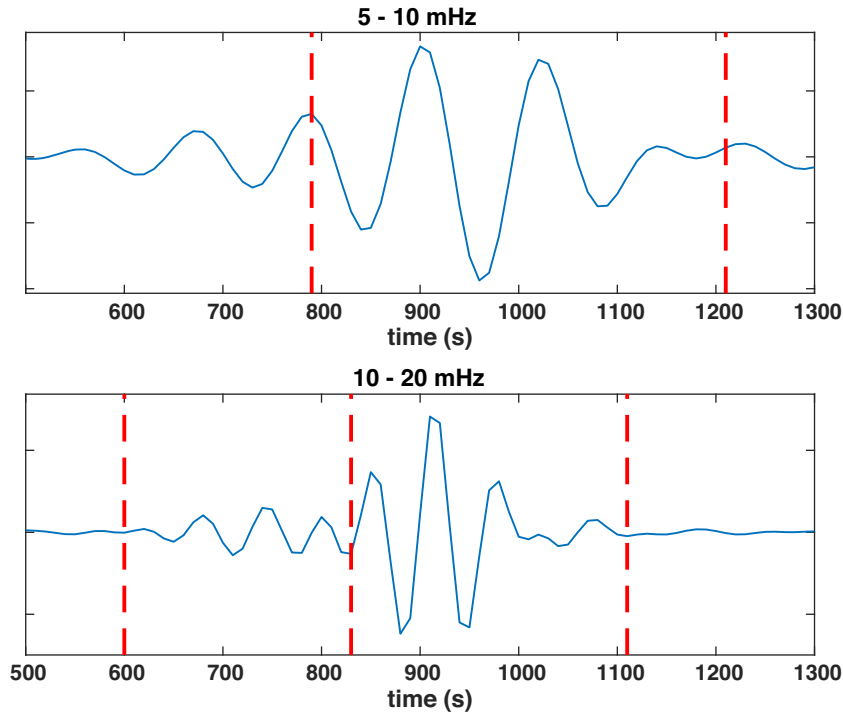
$$p(\mathbf{d}_i | \mathbf{m}) = \frac{1}{(\sqrt{2\pi}\sigma_i)^n} \times \exp \left[ -\frac{M_i(\mathbf{m})}{2\sigma_i^2} \right]. \quad (11)$$

The total likelihood function is then the product of the likelihood functions of three frequency–time windows:

$$p(\mathbf{d}_{\text{obs}} | \mathbf{m}) = p(\mathbf{d}_1 | \mathbf{m}) p(\mathbf{d}_2 | \mathbf{m}) p(\mathbf{d}_3 | \mathbf{m}). \quad (12)$$

As shown in eq. (7), the level of noise accounts for theoretical errors, that is, the part of the signal that we are not able to explain with the forward theory and with the chosen parametrization, which is a simplified and discretized representation of reality. Unlike traditional inversion methods in which the noise level is fixed at a presumed level, we use a Hierarchical Bayesian approach, treating noise parameters  $\sigma_1$ ,  $\sigma_2$  and  $\sigma_3$  as unknowns in the inversion. These noise parameters therefore determine the width of the Gaussian likelihood function and the relative weight given to different frequency–time windows during the inversion. The diagonal element  $\sigma_i$  is given by a uniform prior distribution, which is explored during the Monte





**Figure 2.** An example of our automatic time window selection for a path across North America. The waveform is for the observed data. The station, HRV, is located in northeastern America. The event was in Chiapas, Mexico and occurred on 2007 July 6. The time window in the upper panel is window 1 as defined in Table 1. The time window on the right-hand side of the lower panel is for the fundamental mode (window 2) and the time window on the left-hand side of the lower panel is the higher modes window (window 3).

Carlo search. An advantage of this approach is that the weights of the different frequency–time windows are constrained by the data, thereby avoiding an arbitrary choice of noise levels by the user and the risk of overfitting the data. In addition, windows with better waveform fit will naturally have higher weights since the relative weight between different windows is controlled by the data error levels  $\sigma_1$ ,  $\sigma_2$  and  $\sigma_3$ . We discuss the effects of data noise level in Section 3.2.

Following Bodin *et al.* (2012), a uniform prior distribution with relatively wide bounds is adopted here so that the final model will be dominated by the data rather than by prior information. When only independent parameters are considered, the prior probability distribution can be written as the product of three terms:

$$p(\mathbf{m}) = p(\mathbf{c}, \mathbf{v}|k)p(k)p(\mathbf{h}), \quad (13)$$

where  $\mathbf{c}$ ,  $\mathbf{v}$  are vectors containing the depth and velocity of each interpolation point,  $p(k)$  is the prior on the number of layers and  $p(\mathbf{h})$  is the prior on noise hyper parameters. Details of the derivation for the expression of  $p(\mathbf{m})$  can be found in Bodin *et al.* (2012). Here, we only show the expression for velocity priors at the  $i$ th interpolation point  $v_i$ :

$$p(\mathbf{v}|k) = \begin{cases} \frac{1}{\Delta v} & \text{if } V_{\min} \leq v_i \leq V_{\max} \\ 0 & \text{otherwise,} \end{cases} \quad (14)$$

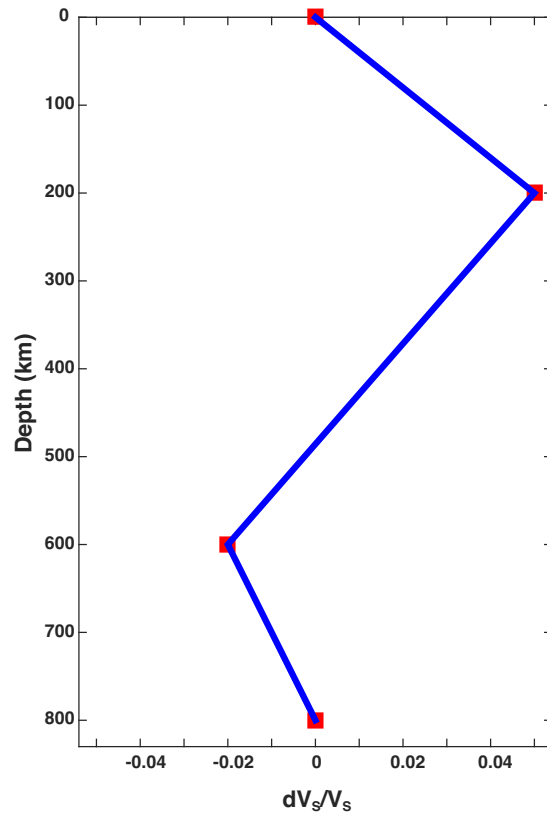
where  $V_{\max}$  and  $V_{\min}$  are the upper and lower bounds of velocity, and  $\Delta v$  is equal to  $(V_{\max} - V_{\min})$ . Any velocity that falls outside this pre-defined boundaries leads to a null prior. For our study, we chose  $\Delta v = 10$  per cent of the velocity in the reference model at a given depth.

### 2.3 The rj-MCMC method

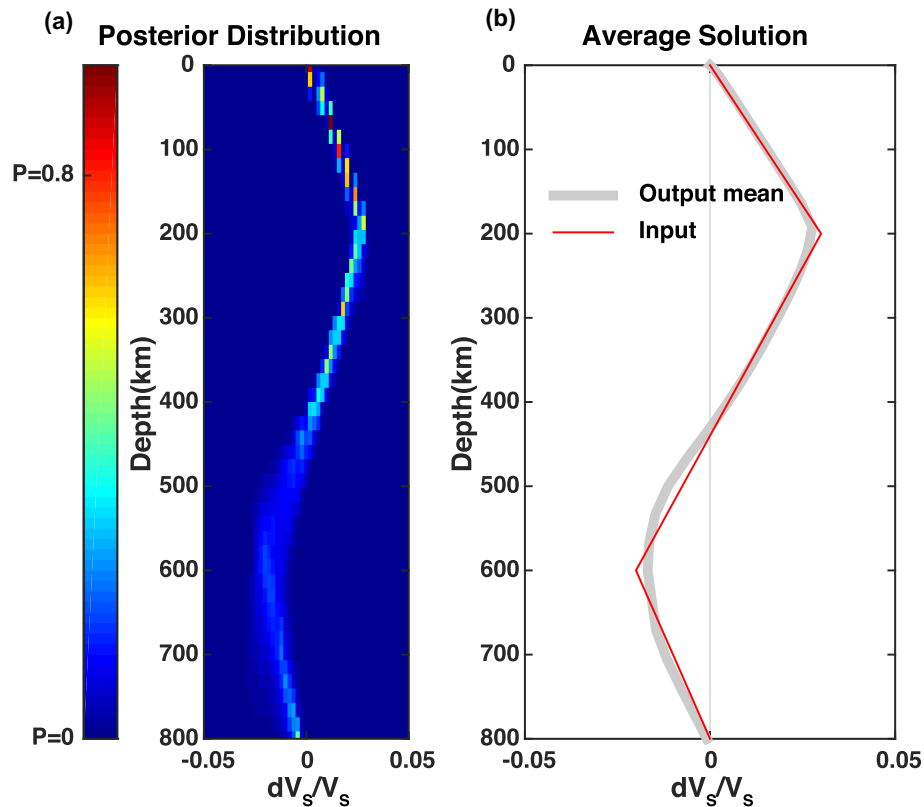
The goal of the rj-MCMC method is to generate an ensemble of earth models distributed according to a target distribution. In our case, we want to find the 1-D shear velocity models that best-fit the waveform data to represent the dispersion of multiple modes along a specific source–receiver path.

Most inversion schemes adopt a fixed dimension model space. However, we do not know the complexity of the  $V_S$  profile, that is, the dimension of the model space, *a priori*. One potential drawback of fixing the depth parametrization is that different seismograms may require different number of model parameters to improve the fit and the results could potentially depend on the chosen prior parametrization. Here, we adopted a transdimensional Bayesian inversion, which consists in a sampling-based algorithm that include the number of parameters in the set of unknowns and hence the constraints added on the uncertain parametrization or uncertain data errors can be released. The transdimensional approach lets the data themselves constrain the allowable model complexity instead of having the user choose a model parametrization (e.g. number of layers or cells) *a priori*. It is well-known, however, that increasing the complexity of the model parametrization often results in better data fit and introduces parameter trade-offs. One difficulty in inverse problems is to determine which parameters are well resolved and how much complexity is really needed by the data. To avoid this problem, the rj-MCMC sampler, first proposed by Geyer & Møller (1994), is designed to find a parsimonious solution (Malinverno 2002), that is, it naturally discourages high dimensional (many layers) models and the least complex explanation or model for an observation is preferred to avoid overfitting the data.

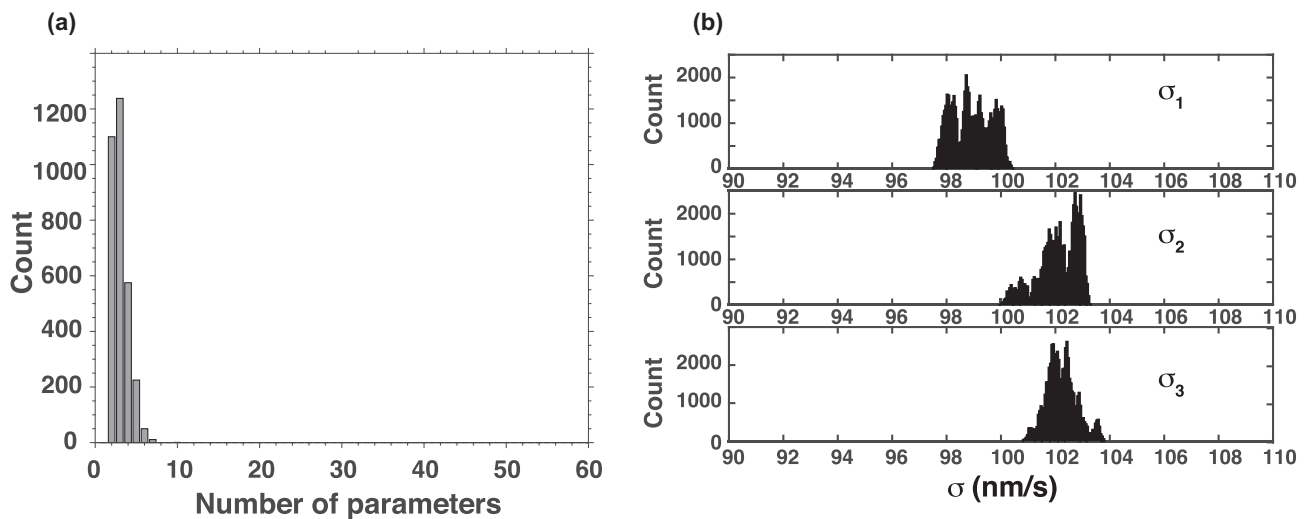
The rj-MCMC method has been explained in detail in Bodin & Sambridge (2009) and Bodin *et al.* (2012). Our method is based on



**Figure 3.** Example of model parametrization. The red squares denote the points at which the velocity model is perturbed. The complete  $dV_S/V_S$  profile is obtained by linear interpolation between those points.



**Figure 4.** (a) Colour density plot representing the ensemble of solutions obtained by inversion of the synthetic seismogram that was calculated with the input model shown in (b); the posterior mean model resulting from this distribution of solutions is also displayed in (b).



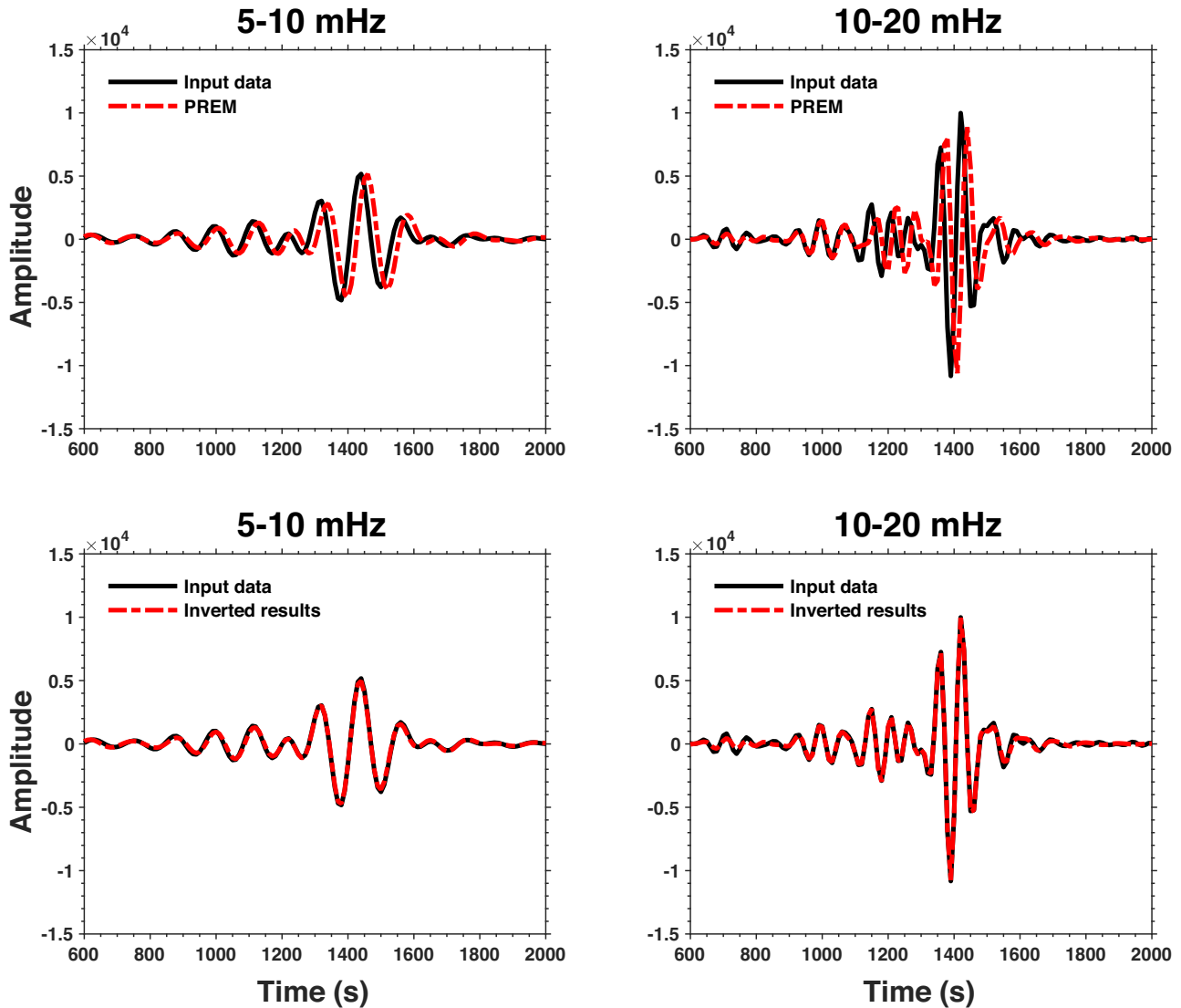
**Figure 5.** (a) Posterior distribution for the number of parameters in the synthetic test of Fig. 4; (b) Posterior distribution for the data noise level for each of the three time windows considered.

theirs except that we used a different depth parametrization. The nodes in the two different parametrizations have, nevertheless, a similar meaning and the derivations are similar to theirs. The main difference comes from the way the velocity models are constructed. In Bodin *et al.* (2012), each Voronoi nucleus defines its nearest neighbour region as one layer, which has the same velocity as the nucleus within it. By contrast, in our model the shear wave velocity perturbation at a given depth is interpolated from its two nearby nodes.

In the rj-MCMC method, every new model is generated by iteratively perturbing the last one according to some chosen proposal distribution. There are five types of perturbation based on the parametrization we chose:

- (1) Change the velocity of one interpolation point;
- (2) Birth: create a new interpolation point;
- (3) Death: remove one interpolation point at random;
- (4) Move: randomly pick one interpolation point and move it to a new depth;





**Figure 6.** Comparison between input synthetic waveform and the waveform predicted by PREM (top) and by the mean model resulting from our inversion (bottom). The seismograms shown here represent the velocity of the ground motion. The input seismogram was calculated for the true model shown in Fig. 4. Two frequency bands, as indicated in Table 1, are considered separately. The waveforms match is much improved after inversion.

**Table 2.** Source parameters from GCMT catalogue. Event 200503021042A was used for the synthetic test. 200910151211A was for the western Australia path and 201004172315A for the eastern Australia path.

ID	Date	Time	lat (°)	lon (°)	Depths (km)	$M_0$ (dyne cm)	Strike, dip, rake
200503021042A	2005/03/02	10:42:16.9	-6.54	129.99	196.1	5.73e+26	41, 88, 55
200910151211A	2009/10/15	12:11:19.8	-3.04	139.45	105.1	1.34e+25	296, 83, 145
201004172315A	2010/04/17	23:15:27.9	-6.82	147.30	62.9	2.19e+25	268, 50, 91

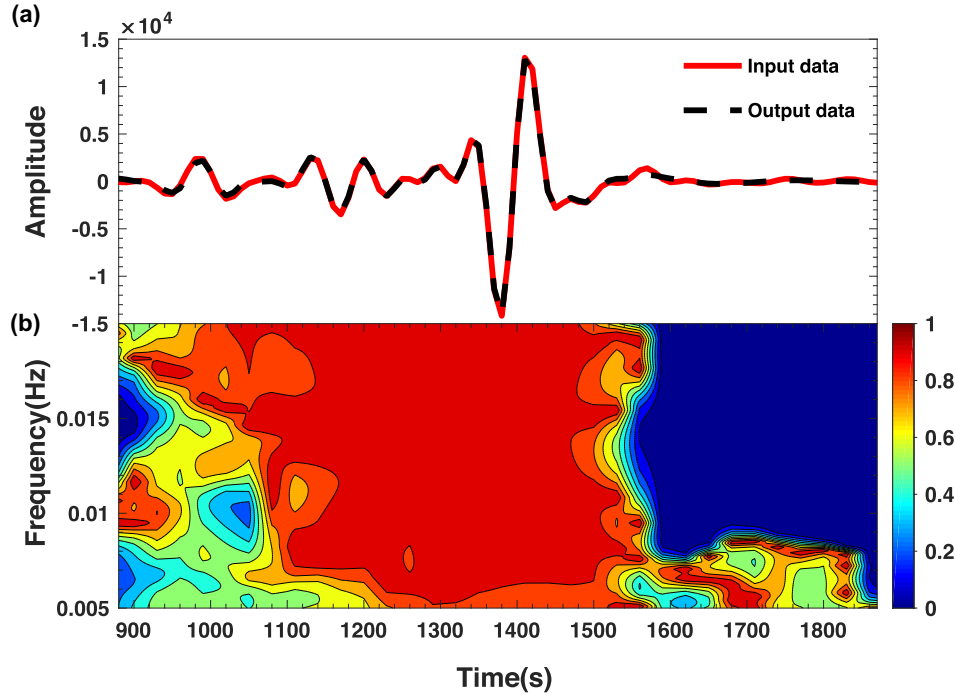
(5) Change the noise level;

Here we have employed a slightly different model proposal for Birth than Bodin *et al.* (2012) due to the different parametrization used here. The other proposals are the same. For the birth step, we added a new interpolation point at depth  $c'_{k+1}$  and then assigned a new velocity perturbation value  $v'_{k+1}$  to this new node. This is drawn from a Gaussian proposal probability density:

$$q(v'_{k+1}|v_*) = \frac{1}{\theta_{\text{birth}}\sqrt{2\pi}} \exp\left\{-\frac{(v'_{k+1} - v_*)^2}{2\theta_{\text{birth}}^2}\right\}, \quad (15)$$

where  $v_*$  is the current velocity perturbation value at the depth  $c'_{k+1}$  where birth takes place. The value of  $v_*$  can be calculated by a simple interpolation between its two nearest nodes. The standard deviation  $\theta_{\text{birth}}$  of the Gaussian distribution is a parameter to be chosen. We refer the reader to Bodin *et al.* (2012) for the proposal distributions of other types of perturbation.

New models generated from those perturbations are then randomly accepted or rejected according to the acceptance ratio. In order to converge to the target distribution  $p(m|\mathbf{d}_{\text{obs}})$ , the acceptance probability from the current model  $\mathbf{m}$  to the proposed model



**Figure 7.** (a) Comparison between the output and input synthetic data in the time domain filtered in the 50–200 s period range; (b) Spectrogram of the waveform fit  $f(\omega, t)$  as defined in eq. (19). The regions with warmer colours correspond to better waveform fit in the F–T domain.

$\mathbf{m}'$ ,  $\alpha(\mathbf{m}'|\mathbf{m})$ , has to meet the following requirement:

$$\alpha(\mathbf{m}'|\mathbf{m}) = \min \left[ 1, \frac{p(\mathbf{m}')}{p(\mathbf{m})} \cdot \frac{p(\mathbf{d}_{\text{obs}}|\mathbf{m}')}{p(\mathbf{d}_{\text{obs}}|\mathbf{m})} \cdot \frac{q(\mathbf{m}|\mathbf{m}')}{q(\mathbf{m}'|\mathbf{m})} \cdot |\mathbf{J}| \right], \quad (16)$$

where the matrix  $\mathbf{J}$  is the Jacobian of the transformation from  $\mathbf{m}$  to  $\mathbf{m}'$ .

The derivations of acceptance probability in rj-MCMC method were given by Bodin *et al.* (2012), Appendix C. Here we will only describe the ideas briefly and show the equations that differ from theirs. For moves without change of dimension, both the Jacobian term and the ratio of proposal distributions are 1 and the reader can find the same results as Bodin *et al.* (2012) by inserting eqs (11) and (13) into eq. (16). For the birth and death steps that involve a change of dimension, it can be shown that the Jacobian term is equal to 1. However, due to the different birth proposal distribution we selected, the expression for the acceptance term for the birth and death steps are different from Bodin *et al.* (2012). For the birth step:

$$\alpha(\mathbf{m}'|\mathbf{m}) = \min \left[ 1, \frac{\theta_{\text{birth}}\sqrt{2\pi}}{\Delta v} \cdot \exp \left\{ -\frac{(v'_{k+1} - v_*)^2}{2\theta_{\text{birth}}^2} \right\} \cdot \frac{p(\mathbf{d}_{\text{obs}}|\mathbf{m}')}{p(\mathbf{d}_{\text{obs}}|\mathbf{m})} \right]. \quad (17)$$

The meaning of  $v'_{k+1}$  and  $v_*$  is the same as in (15). For the death step:

$$\alpha(\mathbf{m}'|\mathbf{m}) = \min \left[ 1, \frac{\Delta v}{\theta_{\text{birth}}\sqrt{2\pi}} \cdot \exp \left\{ -\frac{(v_j - v'_i)^2}{2\theta_{\text{birth}}^2} \right\} \cdot \frac{p(\mathbf{d}_{\text{obs}}|\mathbf{m}')}{p(\mathbf{d}_{\text{obs}}|\mathbf{m})} \right], \quad (18)$$

where  $v_i$  is the velocity of the  $i$ th interpolation point to be removed at depth  $c_i$  and  $v_*$  is the velocity at depth  $c_i$  in the new structure after the removal of the  $i$ th node.

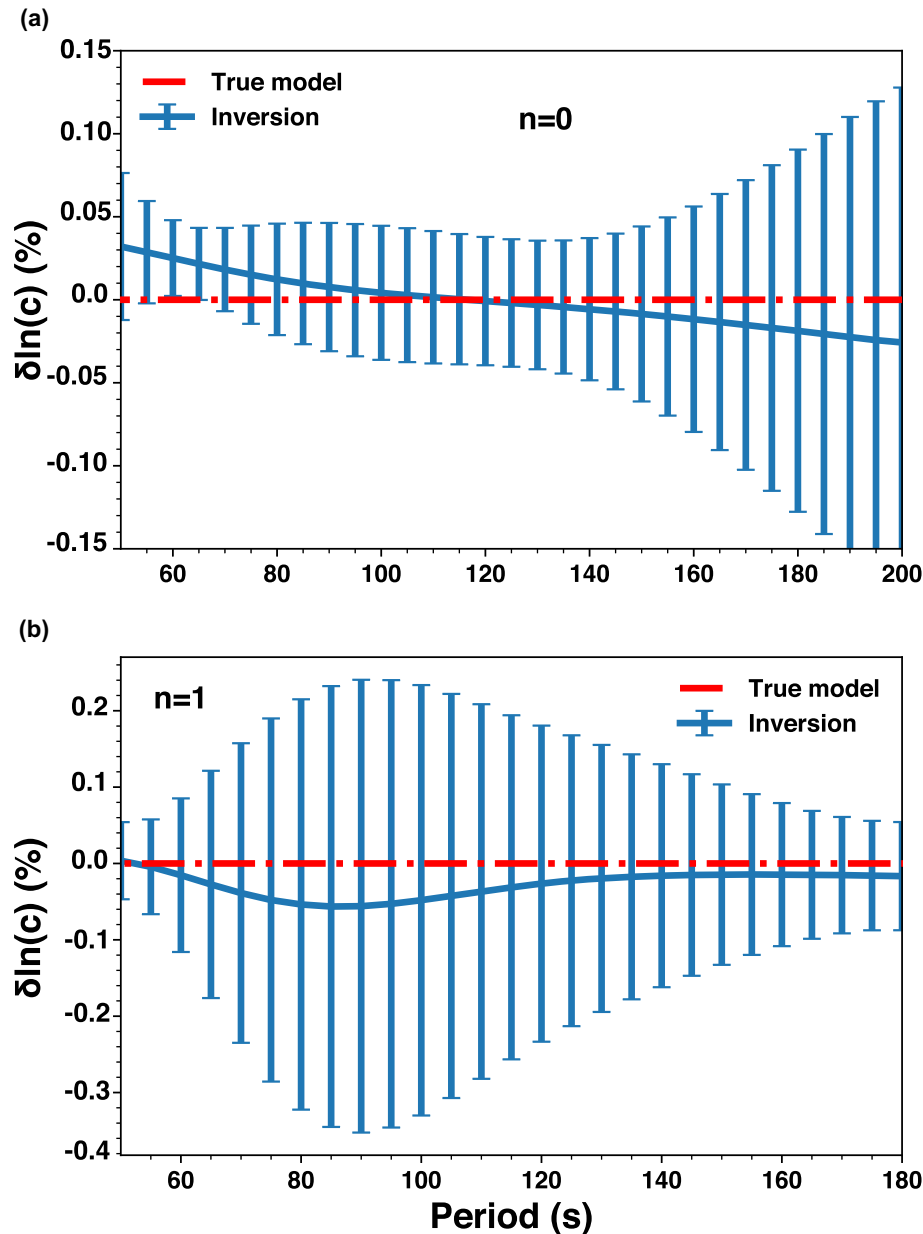
If the newly proposed model is rejected, then the last model is retained for another iteration. The Markov chain is generated via hundreds of thousands of iterations. The first part of the chain (called the burn-in period) is discarded, after which the random walk is assumed to be stationary and starts to sample the model space according to the posterior distribution  $p(\mathbf{m}|\mathbf{d}_{\text{obs}})$ . If the algorithm is run long enough, these samples should then provide a good approximation of the structure of Earth as constrained by the data.

This ensemble solution contains many models with variable parametrization and each  $V_S$  model in the ensemble corresponds to one dispersion relationship. The expected dispersion is the weighted average through the posterior distribution sampled by the rj-MCMC algorithm. One can use this average to choose one dispersion curve with proper uncertainties for interpretation purpose.

## 2.4 Forward modelling and parametrization

In this study, we seek a 1-D depth-dependent isotropic shear velocity model to represent the dispersion of multiple modes between a chosen source and receiver. The isotropic nature of the model is assumed mainly for computational reason, but it is a reasonable assumption as demonstrated by Visser (2008) who showed that the differences in the phase velocities calculated assuming isotropy or anisotropy are small. The  $V_S$  profile is described by a variable number  $k$  of interpolation points (Fig. 3). The vertical position of these points defines the depths at which  $V_S$  is perturbed and the horizontal position of these points corresponding to the amount  $V_S$  is perturbed relative to a reference model using the velocity prior described by eq. (14).

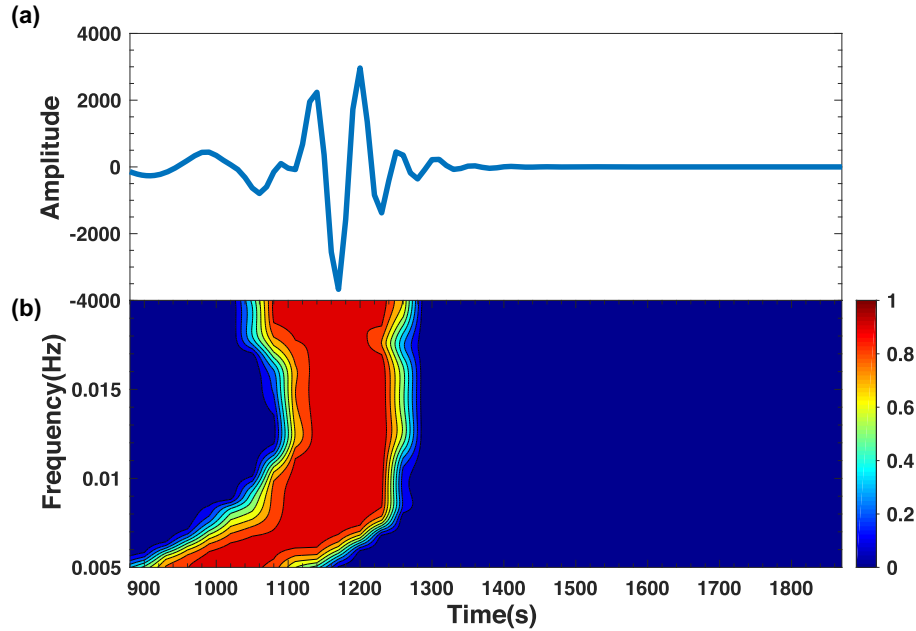
We adopted the rj-MCMC algorithm (Bodin & Sambridge 2009) to explore the model space and sample the distribution of model parameters informed by the data. The procedure of our waveform inversion can be summarized as follows:



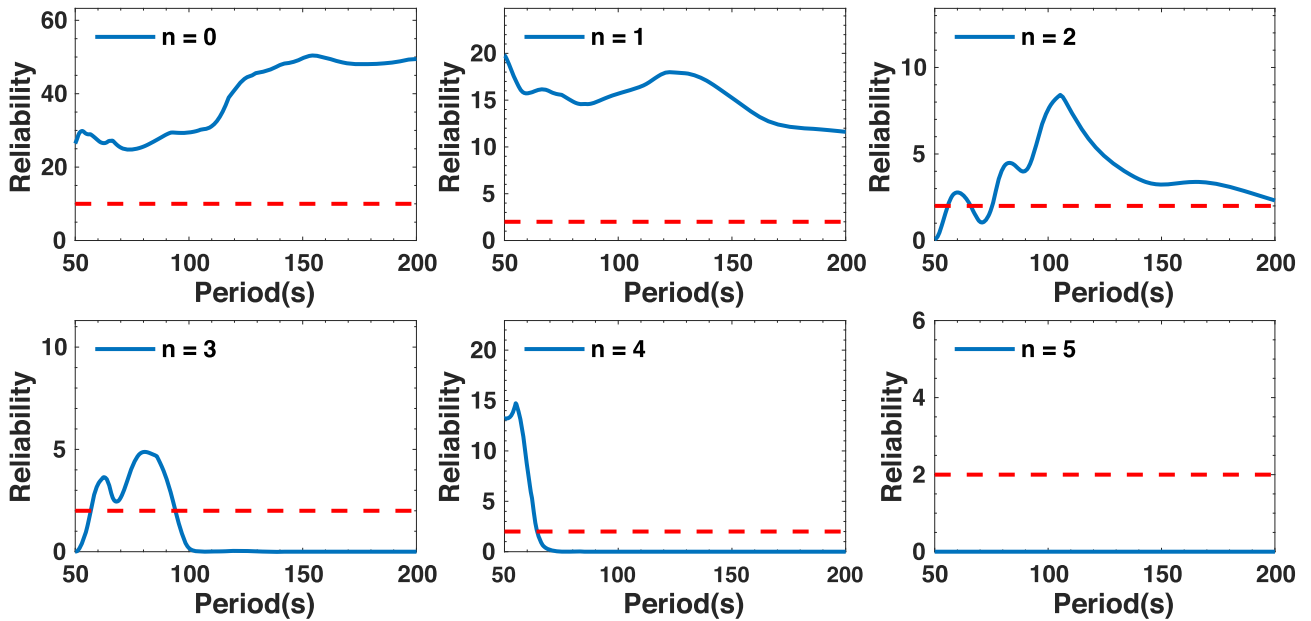
**Figure 8.** Inverted fundamental mode (a) and first overtone (b) phase velocity dispersion and uncertainties (blue) compared to the dispersion curve calculated for the true model (red). The uncertainties shown here correspond to  $2\sigma$  obtained from the posterior distribution.

- (1) Choose path-specific reference 1-D shear velocity model
- (2) Calculate normal mode eigenfrequencies for the reference model using MINEOS and compute a reference synthetic seismogram using normal-mode summation (eq. 1)
- (3) Generate path-specific 1-D shear velocity model by perturbing the reference model according to the proposal distribution of the rj-MCMC scheme;
- (4) Compute the synthetic seismogram predicted by the 1-D model using eq. (3);
- (5) Calculate the misfit between synthetic and observed waveforms in each window using the  $L_2$  norm;
- (6) Randomly accept or reject the proposed model according to the acceptance ratio;

- (7) If the model is accepted, calculate and save the phase velocities predicted by the model using eq. (4) and generate a new  $V_S$  model by perturbing the accepted model according to the proposal distribution. If the model is rejected, generate a new  $V_S$  model by perturbing the previously accepted model according to the proposal distribution;
- (8) Repeat steps (4) to (7) to generate the Markov chains;
- (9) Gather all the saved phase velocities to determine the distribution of dispersion curves obtained at each period/frequency;
- (10) Calculate the mean and standard deviation of these distributions to obtain a measured dispersion curve and uncertainties of phase velocities, respectively;
- (11) Estimate the reliability of the dispersion curves (see Section 2.5).



**Figure 9.** (a) Synthetic waveform of the first overtone calculated from the inverted 1-D  $V_S$  profile and filtered in the 50–200 s period range; (b) Spectrogram  $P_1(\omega, t)$  of the relative weight of the first overtone in the synthetic test, as defined in eq. (20). The regions with warmer colours in the figure represent larger weights in the F–T domain.



**Figure 10.** Reliability parameters as a function of frequency for the fundamental mode ( $n = 0$ ) and the first five overtones in the synthetic test. The thresholds we defined for each mode are denoted by the red dashed lines. Only periods at which the reliability parameters are above the threshold are kept as reliable inversion results.

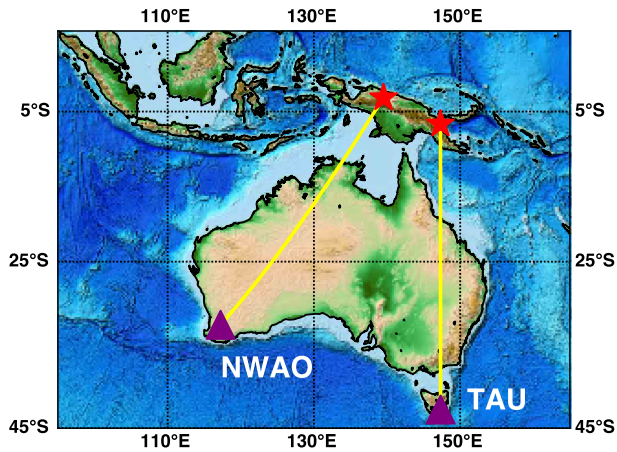
## 2.5 Reliability analysis

In principle, with the proposed method we are able to obtain the phase velocities for any mode at any period since they are calculated from the ensemble of  $V_S$  models obtained. However, in practice, only some modes at certain periods can be reliably constrained by a single seismogram. It is thus important to estimate the reliability of the measurements in each frequency band for different modes. Here, we adopted a procedure similar to Yoshizawa & Kennett (2002) and Yoshizawa & Ekström (2010) based on the frequency–time analysis method (Dziewonski *et al.* 1969).

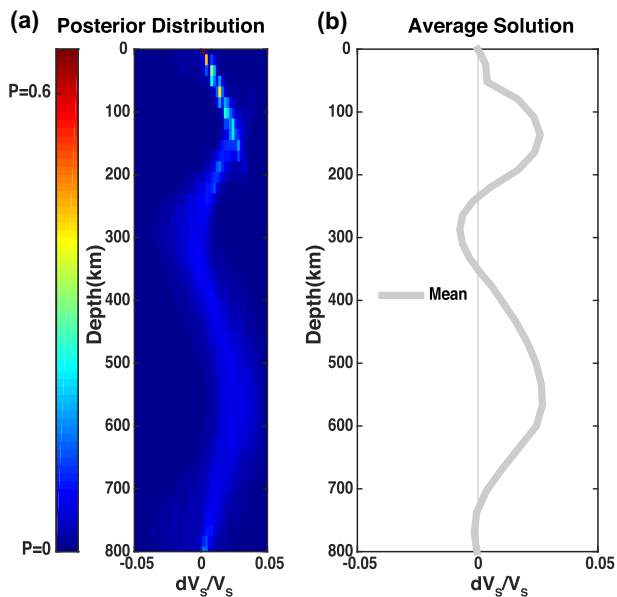
Let us first define the waveform fit  $f(\omega, t)$ :

$$f(\omega, t) = \exp \left[ -\frac{S^{\text{mis}}(\omega, t)}{S^{\text{syn}}(\omega, t)} \right], \quad (19)$$

where  $S^{\text{mis}}(\omega, t)$  represents the spectrogram of the difference between real data and the inverted synthetics.  $S^{\text{syn}}(\omega, t)$  is the spectrogram of the whole synthetic waveform. Similarly, the relative power



**Figure 11.** Locations of events (stars) and stations (triangles) used for the real data test. The yellow lines show the great circle paths from sources to receivers for western and eastern Australia.



**Figure 12.** (a) Posterior distribution of relative velocity perturbations with respect to Debayle *et al.* (2016) for the western Australia path; (b) Resulting mean  $dV_S/V_S$  model.

of the  $j$ th mode  $p_j(\omega, t)$  is defined as

$$p_j(\omega, t) = \exp \left[ - \frac{S_j^{\text{res}}(\omega, t)}{S_j^{\text{syn}}(\omega, t)} \right], \quad (20)$$

where  $S_j^{\text{syn}}(\omega, t)$  is the spectrogram of the  $j$ th mode waveform and  $S_j^{\text{res}}(\omega, t)$  is the spectrogram of the residual seismogram calculated as the difference between the full synthetic and the  $j$ th mode waveform. In practice, both the frequency  $\omega$  and time  $t$  are discrete, so the spectrograms can be expressed in matrix form. For example, the element in the  $k$ th row and  $l$ th column of the spectrogram matrix represents the amplitude of  $S(\omega_k, t_l)$ , where  $\omega_k$  is the discretized frequency and  $t_l$  is the centre of the  $l$ th time window used in the frequency–time domain analysis.

The waveform fit  $\mathbf{f}$  measures the difference between the synthetic and real data in the frequency–time domain. It is designed such that  $f(\omega, t) = 1$  at all frequencies and time when the synthetic seismograms are identical to the observed data, and  $\mathbf{f}$  decreases

when the waveform fit becomes worse. The relative power  $\mathbf{p}_j$ , which measures the relative power of the  $j$ th mode, is designed in a similar way so that  $p_j(\omega, t) = 1$  at all frequencies and time when the  $j$ th mode in the seismogram is completely dominant and there is no contribution from other modes, and  $\mathbf{p}_j$  decreases as the contribution from the  $j$ th mode becomes smaller.

For the reliability analysis, we want to consider the overall waveform fit and the relative weight of one single mode at the same time because both factors indicate how well the phase velocities at certain frequencies are constrained by the data. That is, the measurement of the  $j$ th mode becomes more reliable if the waveform fit is better, or if the relative weight of the  $j$ th mode increases. Following this criteria, we define the reliability parameter as

$$r_j(\omega) = \sum_l p_j^{kl} f^{kl}, \quad (21)$$

where  $p$  and  $f$  are defined in eqs (19) and (20). To get the reliability parameter of the  $j$ th mode  $r_j(\omega)$ , we first multiply the waveform fit by the relative weight of the  $j$ th mode in the F–T domain, and then integrate the resulting matrix with respect to time. In this way, the  $r_j(\omega)$  is a function of frequency and can be regarded as an indicator of reliability at frequency  $\omega$  for the  $j$ th mode. Examples of the reliability analysis are shown in Section 3.

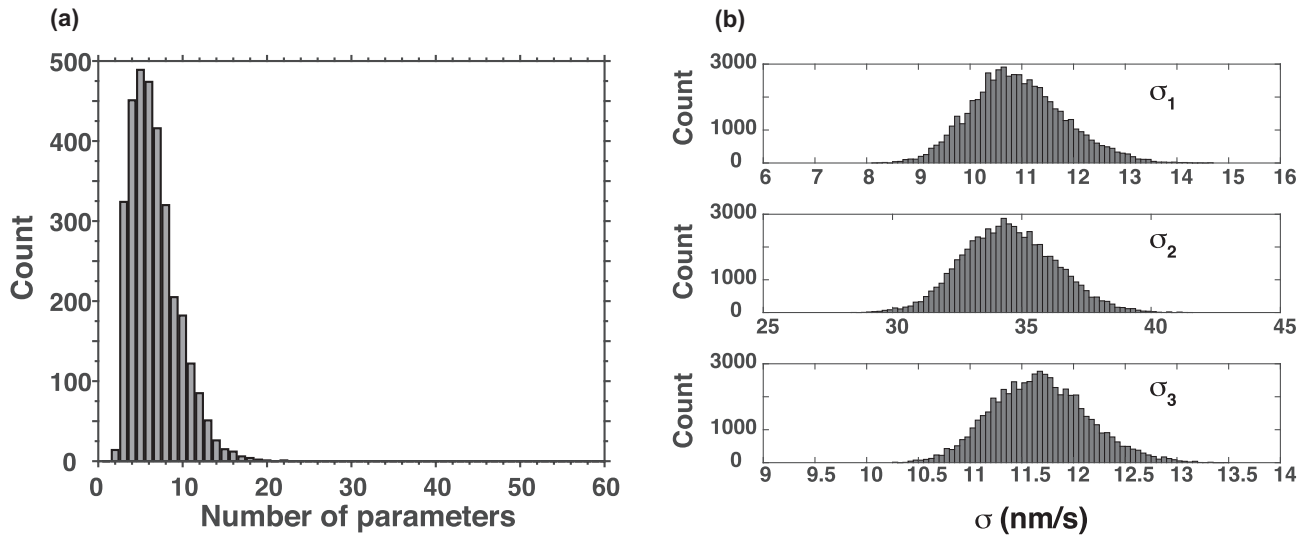
### 3 TESTS AND RESULTS

#### 3.1 Synthetic tests

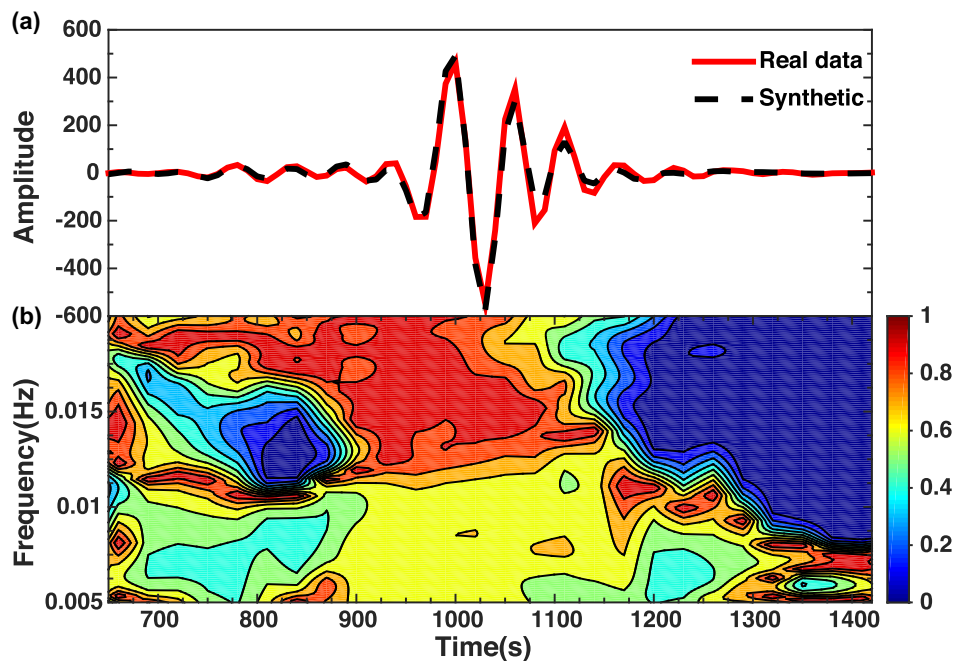
We first present a synthetic test to demonstrate that our method can recover an input shear wave velocity model well (Figs 4 and 5). A shear wave velocity profile generated by perturbing the Preliminary Reference Earth Model (PREM; Dziewonski & Anderson 1981) was used to calculate a synthetic seismogram. Fig. 4(b) displays the input perturbation  $dV_S(r)/V_S(r)$ . It is characterized by +3 per cent anomaly relative to PREM at 200 km depth and –2 per cent at 600 km depth. The synthetic seismogram (Fig. 6), representing the velocity of the ground motion, was calculated for a great circle path between the Banda Sea and Beijing, which corresponds to a distance of 5347 km. We used the fundamental mode and the first ten overtones to calculate the synthetics and to invert the synthetic data. The event source parameters were taken from the GCMT catalogue (Dziewonski *et al.* 1981; Ekström *et al.* 2012) and are shown in Table 2. Only the vertical component is considered for the waveform fitting. We generated correlated noise by bandpass-filtering a sequence of uncorrelated Gaussian (white) noise, and then added the correlated noise to the synthetic seismogram. This way, the filtered noise sequence is correlated and data points in the sequence are no longer independent. The noise level was set at 10 per cent of the average amplitude of the waveform contained in the third time window as defined in Table 1, which includes higher modes and excludes the fundamental mode.

The inversion of the synthetic waveform data was performed using eight Markov chains starting from different random models, with the length of each chain fixed at 120 000 iterations. To guarantee the convergence of the inversion when sampling the posterior distribution, the first 60 000 iterations in each chain were marked as ‘burn-in’ and we removed them from the ensemble of sampled models. The ensemble of  $dV_S/V_S$  solutions represented by a colour density plot is displayed in Fig. 4(a). The brighter colour means the algorithm spent more time in that region, and therefore it corresponds to a more likely  $V_S$  structure at that particular depth. The mean output model calculated from the ensemble of





**Figure 13.** (a) Posterior distribution of the number of velocity parameters for the western Australia path; (b) Posterior data error distribution for each of the three time windows. Note the different scales on the horizontal axes.



**Figure 14.** The same as Fig. 7 but for the western Australia path. (a) real data and synthetic seismogram filtered in the 50–200 s period range and calculated using the inverted mean model of Fig. 12; (b) F–T analysis of the misfit as defined in eq. (19).

solutions (Fig. 4b), which can be regarded as the expectation of the ‘true model’, is close to the input model. Fig. 5(a) displays the posterior distribution of the number of parameters used by the algorithm during the inversion and demonstrates the parsimony of the transdimensional framework: although we can almost always fit the data better by introducing more parameters, the rj-MCMC method tends to prevent overfitting the data and prefer models of smaller dimensions. In this case the result shows that there is a higher probability that the data can be explained by less than eight depth nodes. Fig. 5(b) represents the posterior data noise level and shows that fitting the first time window required less data noise than the other two time windows.

A comparison between the input and output seismograms is shown in Fig. 6, which shows that the predicted waveform resulting

from our inversion fits the input data much better than the waveform predicted by PREM. The misfit between the inverted waveform and the input data is represented in Fig. 7 both in the time domain and the F–T domain.

The fundamental mode and first overtone phase velocity dispersion was calculated and compared with the dispersion relationship predicted by the true model (Fig. 8). The standard deviation of our measurements at each period were calculated from the ensemble of dispersion curves obtained. We see that the true dispersion curve is within the 95 per cent likelihood contour of our inversion results at all periods.

We further applied the reliability analysis described in the previous section to our synthetic test. The relative weight of the first overtone is used as an example to demonstrate the method (Fig. 9).



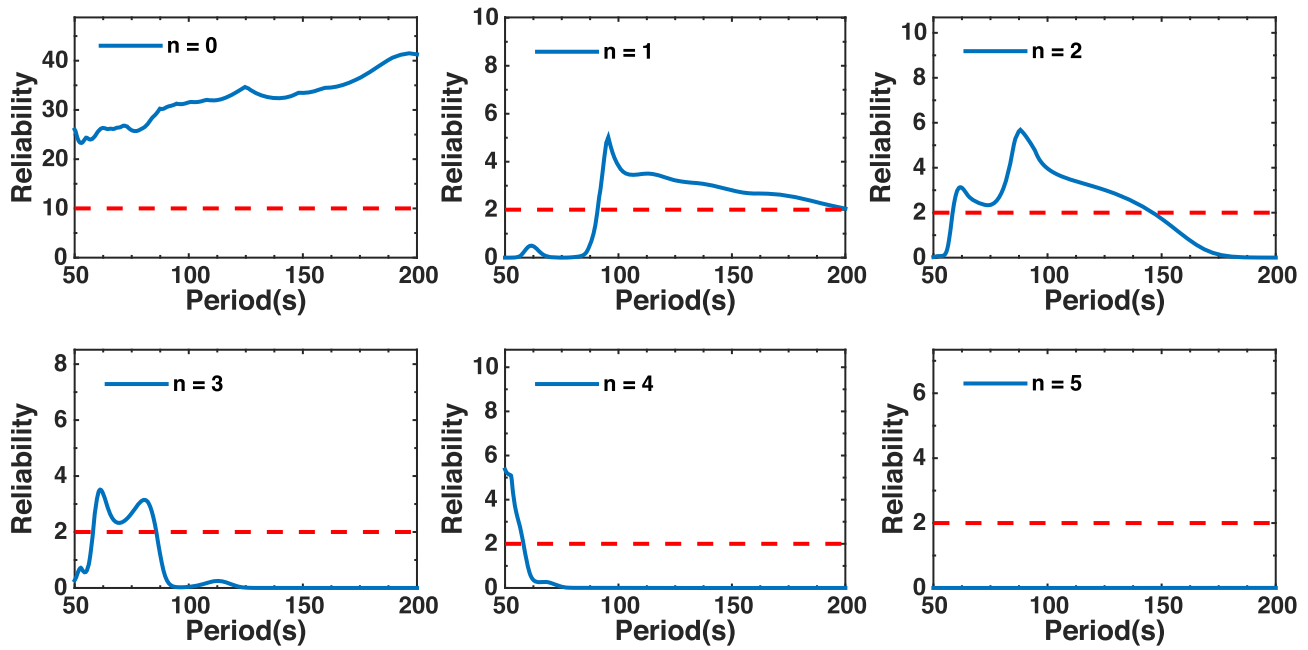


Figure 15. Reliability parameters for the western Australia path (see Fig. 10 for details).

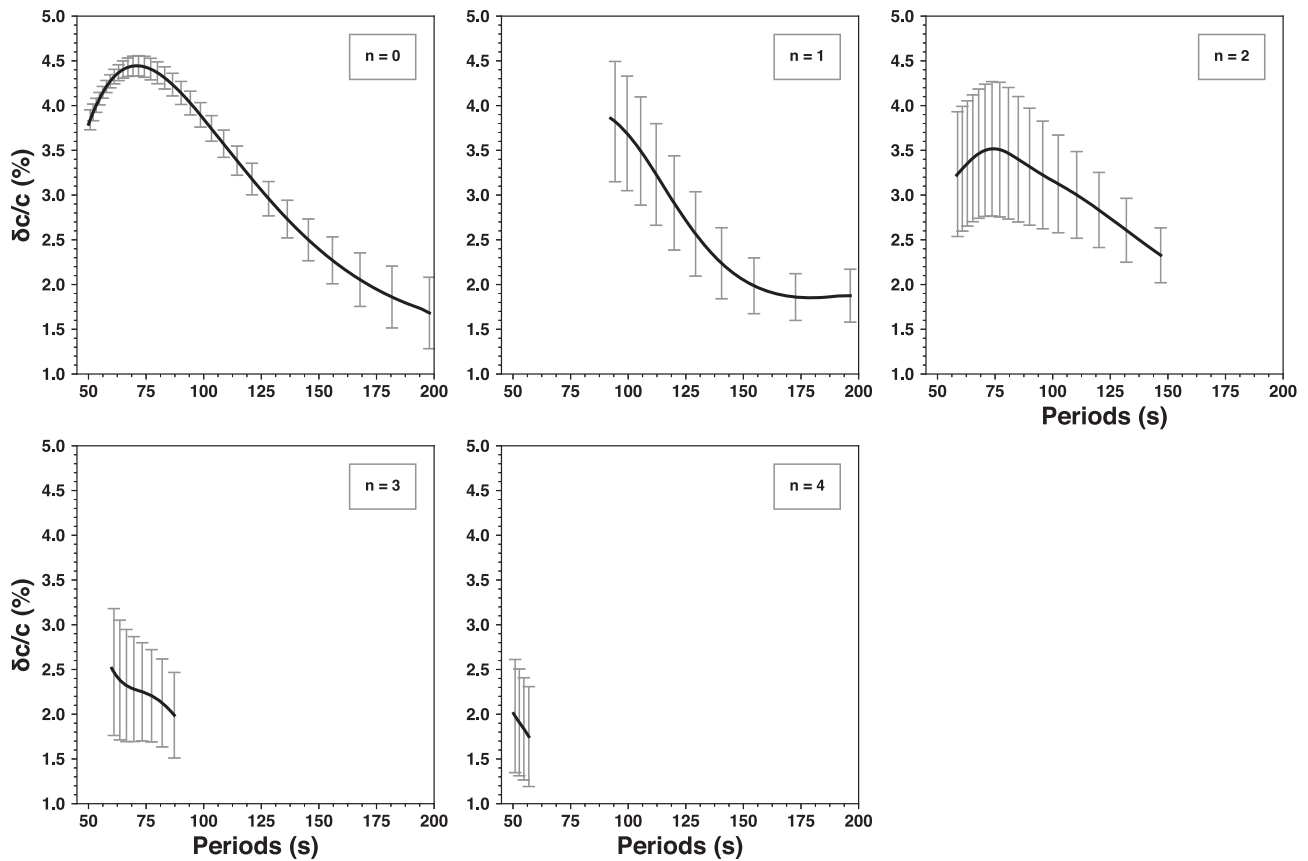


Figure 16. Measured phase velocities anomalies for the western Australia path and the modes and periods that were estimated to be reliable (Fig. 15). The measurements were converted from perturbations with respect to the average version of the 3-D reference model into perturbations with respect to PREM. Uncertainties correspond to  $2\sigma$ .

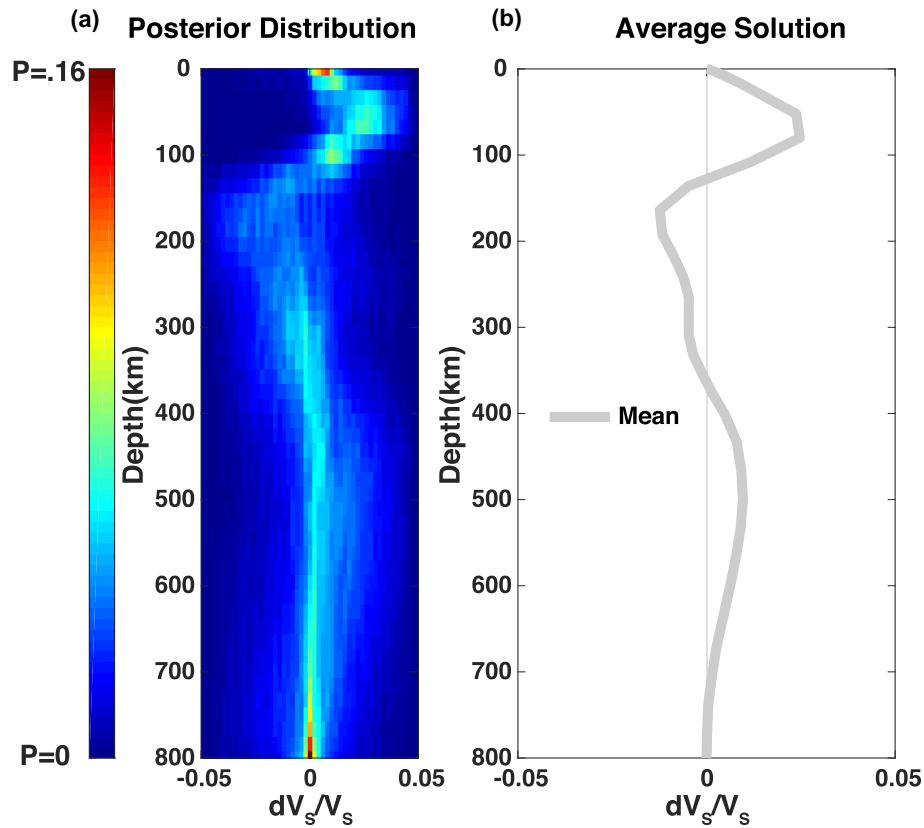


Figure 17. Posterior ensemble of solutions for the eastern Australia path. See Fig. 4 for details.

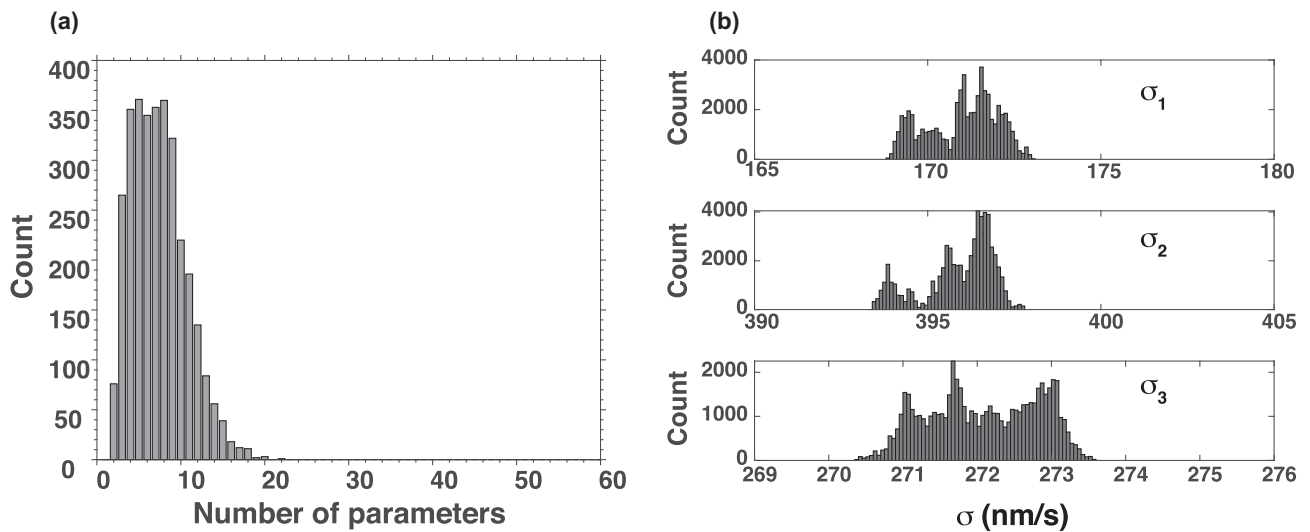
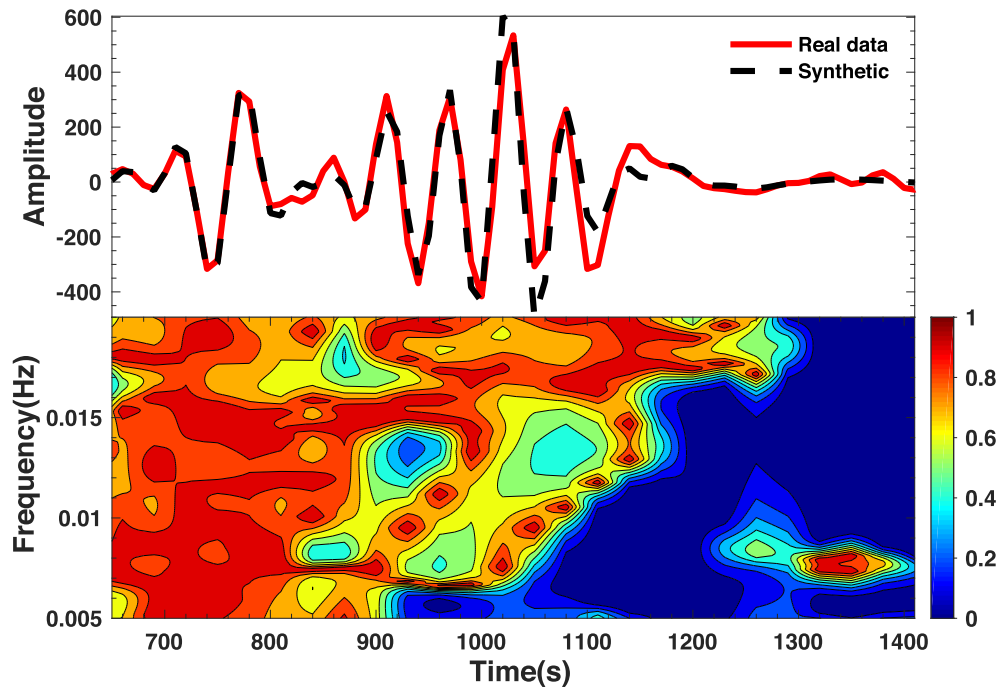


Figure 18. (a) Posterior ensemble for the number of parameters for the eastern Australia path; (b) Posterior noise distribution for each of the three time windows. Note the different horizontal scales.

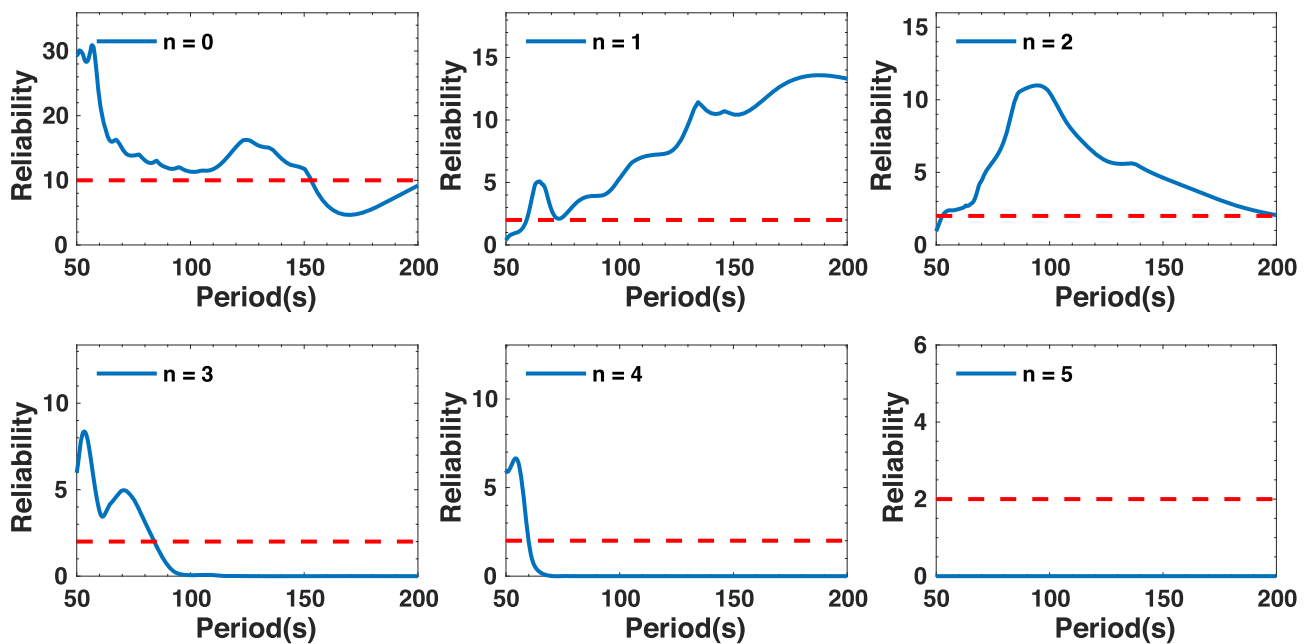
The spectrogram of misfit in Fig. 7 and the spectrograms of relative weights (like in Fig. 9) were multiplied using eq. (21) and summed along the time domain to obtain the reliability parameters as a function of frequency for each mode (Fig. 10). In order to automate the process, we arbitrarily define the threshold values at 10 for the fundamental mode and 2 for higher modes. For each mode, only periods with reliability parameters larger than the threshold will be kept. As is shown in the synthetic test results, the fundamental mode

and the first overtone have high reliability at all periods. The second overtone is also above the threshold at most periods but the third and fourth higher modes only have a short range of reliable results. The fifth overtone cannot be determined reliably in the synthetic test, likely because it was not excited well at those frequencies by the employed seismic source.

Further synthetic tests were performed to verify whether our method is valid for uncorrelated noise (Supporting Information



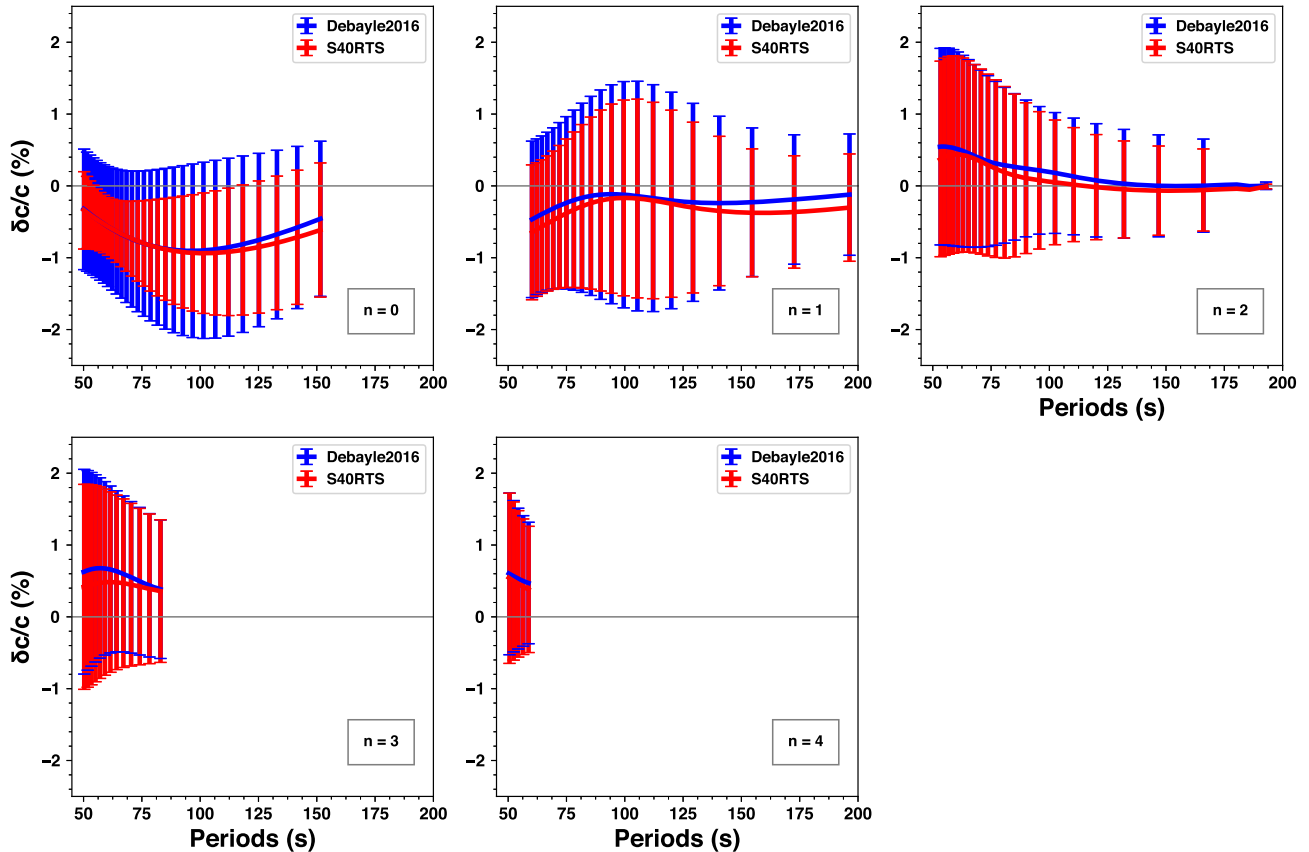
**Figure 19.** Real data and synthetic seismogram filtered in the 50–200 s period range and calculated using the inverted mean model of Fig. 17 (top); F–T analysis of the misfit as defined in eq. 19 (bottom).



**Figure 20.** Reliability test for the eastern Australia path. See Fig. 10 for details.

Fig. S1) and for fixed depth nodes, that is, when the transdimensionality aspect of the software is turned off (Supporting Information Fig. S2). This transdimensionality test showed that for a parametrization with fixed number of parameters, similar to what was implemented by others (Yoshizawa & Kennett 2002, 2004; Visser *et al.* 2007; Visser 2008; Yoshizawa & Ekström 2010), we can obtain phase velocity measurements with error bars that include the predictions of the input model, that is, the ‘true’ data. However, the mean phase velocities deviate from the true values, highlighting

the importance of reporting proper uncertainties with the measurements. We also see that the mean phase velocities do not reproduce the input data as well as when the number of parameters is allowed to vary in the inversion. This is visible when comparing Supporting Information Figs S1 and S2. This is also true for the mean velocity model, which is not as close to the input model as the mean model of Supporting Information Fig. S1. It should be noted, however, that when the dimension of the model is fixed, smoothed parametrizations such as B-splines are more commonly employed (Yoshizawa & Kennett 2002, 2004; Visser *et al.* 2007; Visser 2008; Yoshizawa



**Figure 21.** Reliable measured phase velocities expressed as perturbations with respect to PREM for the eastern Australia path and the modes and periods that were estimated to be reliable (Fig. 20). The red curve represents measurements performed using the model of Debayle *et al.* (2016) as a reference model. The blue curves are measurements performed using model S40RTS of Ritsema *et al.* (2011) as reference. The phase velocities were converted from perturbations with respect to the 3-D reference model into perturbations with respect to PREM. Uncertainties correspond to  $2\sigma$ .

& Ekström 2010), since they can more easily represent smooth 1-D velocity variations with a limited number of nodes. In addition, we note that the computational time when the transdimensionality is turned on was not very different from when it was turned off.

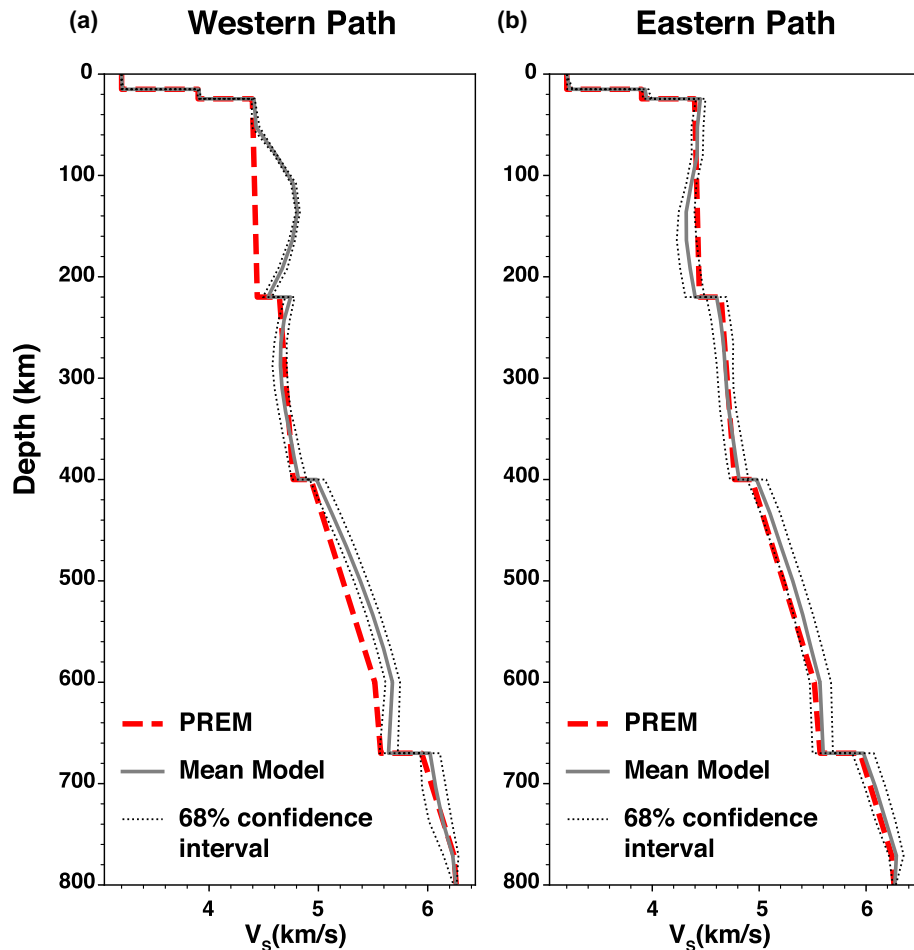
We additionally tested the effect of the data noise level by underestimating it (Supporting Information Fig. S3) and overestimating it (Supporting Information Fig. S4) compared to the input data. These figures demonstrate that the prior data noise level can have large influence on the inverted phase velocities as well as their estimated uncertainties: when the noise level is underestimated, the posterior uncertainties on the estimated phase velocities become smaller but the input data are not well represented by those uncertainties as they lie outside (or close to) two standard deviations. When the noise level is overestimated, however, the uncertainties on the phase velocities are much larger but the mean is close to the input data. We therefore argue that overestimating the noise level leads to more reliable measurements, albeit with larger uncertainties, and reduces the possibility of overfitting the data.

### 3.2 Application to real data

In this section we present the phase velocity inversion results for two paths across western and eastern Australia. Fig. 11 shows the locations of the events and stations. Both selected events have depths larger than 50 km. The source parameters of the two events are listed in Table 2. The epicentral distance for the western path is 4036 km and for the eastern path is 4012 km. Although a nonlinear inversion

method is used here to calculate phase velocities from the  $S$ -wave velocity models, it can still be preferable to use a reference model that is as close as possible to the true structure of the region studied. This is mainly because our method to calculate the synthetic seismograms is not fully nonlinear for computational reasons (see Section 2). The forward modelling part of the algorithm was linearized, and we search for model parameters  $dV_S/V_S$  between  $-5$  and  $5$  per cent around a reference model in order to reduce errors caused by the linearization. Here, we adopted the 3-D shear velocity model of Debayle *et al.* (2016). We averaged the  $V_S$  profiles extracted from the reference model along each station–event pair and used this average as a reference model for the specific 1-D path considered. The prior data noise level was a uniform distribution between 1 and 500  $\text{nm s}^{-1}$  for each time window.

The posterior distribution of  $dV_S/V_S$  models for the western Australia path is shown in Fig. 12, followed by the number of parameters and data noise distributions (Fig. 13), the waveform fit (Fig. 14), the reliability analysis (Fig. 15) and the estimated phase velocity dispersion curves (Fig. 16). The inverted  $dV_S/V_S$  structure (Fig. 12) displays a modest positive deviation from the reference model between 50 and 200 km depth as well as below 400 km depth. The uppermost positive anomaly is consistent with a thick lithosphere as expected in a cratonic area. The number of velocity parameters required by the data is around four, which demonstrates the parsimony of the algorithm, and the posterior data noise distributions have relatively narrow Gaussian distributions with a standard deviation much lower than the assumed prior noise level. Overall, the



**Figure 22.** The mean  $V_S$  solution and one standard deviation for the western (a) and Eastern (b) Australia paths compared to PREM.

fundamental mode in the first window (5–10 mHz) is the lowest, followed by the higher modes window.

The synthetic waveform calculated from the mean velocity model fits the real data well (Fig. 14), and we see that phase velocities for the fundamental mode has reliability parameters larger than the pre-defined threshold, 10, at all periods considered (Fig. 15). For the first overtone, the reliable measurements are in the period range 92–200 s, with the threshold set at 2. Similarly, the second overtone phase velocities were estimated to be reliable in the period range 58–148 s. The reliability parameters of the third overtones is between 60 and 88 s. Finally, a small portion (50–57 s) of the fourth overtone was found to be reliable.

For the eastern Australia path, we repeated the procedure described above. The results are shown in Figs 17–21. We see a strong positive  $V_S$  anomaly compared to the reference model at about 60 km depth and a negative  $V_S$  anomaly at 120–220 km depth. Our solution can be interpreted as a thin lithosphere (about 100 km thick) with slightly positive velocity anomalies and a low velocity layer that may coincide with the asthenosphere (Fig. 22b), consistent with Phanerozoic geological features and with previous studies (Yoshizawa & Kennett 2015). It should also be noted, however, that our  $V_S$  solutions are not directly comparable with 3-D models, but rather can be regarded as a 1-D approximation of the real Earth structure along the selected path. Fig. 18 shows that the number of parameters required to fit the data is between four and eight and the data noise level is clearly the lowest for the fundamental mode

filtered between 5 and 10 mHz (window 1). Overall the noise level is much higher for the eastern Australia path than for the western Australia path. With Fig. 20, we show that the fundamental mode and the first four overtones can be reliably retrieved: the fundamental mode has high reliability at 50–150 s. The first and second overtones have large reliability at longer periods (60–200 s, 54–200 s). The reliable periods for third and fourth overtones are 50–83 and 50–61 s, respectively. Measurements for the modes and periods estimated reliable are shown in Fig. 21. This figure also displays phase velocities measured using S40RTS (Ritsema *et al.* 2011) as a reference instead of the Debayle *et al.* (2016) model to test the dependence of our method upon the reference model. The results show no significant dependence on the reference model. The reliability tests were not strongly affected by the reference model either (not shown here).

The mean absolute  $V_S$  model were also plotted for paths and compared to PREM (Fig. 22a). This figure shows a clear difference between the eastern and western paths in the top 200 km, beyond which the two velocity profiles become more similar.

In this study we used the centroid moment tensor from the GCMT catalogue. We acknowledge, however, that for some paths dispersion measurements performed using waveform modelling at single stations can be affected by uncertainties in source parameters. While quality control can be applied to filter out unreliable earthquake sources by comparing source parameters from different earthquakes in the same region or by comparing different source catalogues, joint

inversions of structure and source parameters are preferable (Valentine & Trampert 2012). A future implementation of our technique will include source parameter among the unknowns in order to account for possible errors in the source parameters.

#### 4 CONCLUSION

We applied the rj-MCMC technique to the nonlinear problem of measuring fundamental and higher mode Rayleigh wave phase velocity dispersion using waveform modelling. The use of higher mode surface waves in depth inversions of seismic velocities and anisotropy is very valuable as it increases the vertical resolution of tomographic models in the upper-mantle and transition zone. The forward problem consisted in calculating synthetic seismograms by normal mode summation using computer program Mineos and a linearized approximation to the calculation of normal mode eigenfrequencies. The rj-MCMC method enabled us to find a distribution of 1-D shear velocity models that best fit the waveform and represent the dispersion of multiple modes along a specific source–receiver path. An advantage of this technique is that the algorithm can decide how much data noise is needed to fit the data without overfitting them. The posterior noise distributions can then be used as an indicator of the quality of the waveform fit within each frequency–time window. The data noise was thus treated as an unknown and different noise levels were applied to the different time windows considered. The resulting distribution of velocity models was used to calculate a posterior distribution of dispersion curves from which a mean and standard deviation were obtained for different modes. A reliability analysis was then performed following Yoshizawa & Kennett (2002) to assess which higher modes were reliably separated.

With synthetic tests, we showed that the newly developed method presented here can extract the first four overtones and the fundamental mode at most periods between 50 and 200 s, but that the third and fourth overtones can only be measured reliably at relatively short periods and in a narrower period band (50–100 s for  $n = 3$  and 50–60 s for  $n = 4$ ). Measurements performed for real data along two paths sampling western and eastern Australia showed similar results: the fundamental mode and first two overtones were measured reliably over most frequencies considered, and the third and fourth overtones could be extracted and their dispersion measured over a narrower frequency band for shorter periods. We also demonstrated that the measured phase velocities do not strongly depend on the reference model used to calculate the initial seismogram.

While computationally intensive, the technique presented here has several advantages. It allows the user to sample the model space without requiring any explicit regularization other than specifying the bounds of the model space explored. In addition, because it is transdimensional and parsimonious, it can include data noise and the number of parameters among the unknowns without overfitting the data. It therefore lets the data themselves control the complexity of the solution. The reliability of mode separation can be assessed and quantitative uncertainties on the dispersion curves of each mode can be readily obtained. This provides useful information on the data to seismologists who use these measurements to model 3-D seismic velocity and anisotropy. Our technique can also be easily adapted to include source parameters among the unknowns and perform joint inversions of model structure and source parameters to account for the effect of source uncertainties on the measured phase velocities. A future implementation of the method will also include perturbations in Moho depth as it can have an important effect on waveform modelling for paths that travel mostly across

continental areas. We also note that the phase velocities are not the only products of this technique. One could also envision using the resulting path-averaged velocity models and uncertainties and combine them to obtain a 3-D velocity model.

#### ACKNOWLEDGEMENTS

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## SUPPORTING INFORMATION

Supplementary data are available at *GJI* online.

**Figure S1.** (A)  $V_S$  posterior distribution for a synthetic test with uncorrelated noise. Both noise levels and model dimensions are variable during the inversion. The red line in panel (B) is the true model. The grey curve is the mean of the distribution of panel (A). (C) Inverted fundamental mode (top) and first overtone (bottom) phase velocity dispersion curves and uncertainties (blue) compared to the dispersion curve calculated for the true model (red). The uncertainties shown here correspond to two standard deviation obtained from the posterior distribution.

**Figure S2.**  $V_S$  posterior distribution (A) for the synthetic test with uncorrelated noise and fixed depth parametrization. We fixed the number of depth nodes to 9 and we fixed their depths at 5, 20, 80, 140, 210, 310, 430, 550 and 690 km. Only the noise level and amplitude of  $dV_S/V_S$  are allowed to vary during the inversion. The red line in panel (B) is the true model and the grey curve is the mean of the distribution of panel (A). (C) Inverted fundamental mode (top) and first overtone (bottom) phase velocity dispersion curves and uncertainties (blue) compared to the dispersion curve calculated for the true model (red). The uncertainties shown here correspond to two standard deviation obtained from the posterior distribution.

**Figure S3.**  $V_S$  posterior distribution for the synthetic test with fixed, uncorrelated and underestimated noise and fixed depth parametrization. The transdimensionality of the algorithm was turned off and the depths of the nodes were the same as in Fig. S2. We assumed the fixed noise level was underestimated in the inversion by a factor of 100 compared to the synthetic input data noise level. Only the amplitude of  $dV_S/V_S$  was allowed to vary during the inversion. The red line in panel (B) is the true model and the grey curve is the mean of the distribution of panel (A). (C) Inverted fundamental mode (top)

and first overtone (bottom) phase velocity dispersion curves and uncertainties (blue) compared to the dispersion curve calculated for the true model (red). The uncertainties shown here correspond to two standard deviation obtained from the posterior distribution.

**Figure S4.**  $V_S$  posterior distribution for the synthetic test with fixed, uncorrelated and overestimated noise and fixed depth parametrization. The transdimensionality of the algorithm was turned off and the depths of the nodes were the same as in Fig. S2. We assumed the fixed noise level was overestimated in the inversion by a factor of 10 compared to the synthetic input data noise level. Only the amplitude of  $dV_S/V_S$  was allowed to vary during the inversion. The red line

in panel (B) is the true model and the grey curve is the mean of the distribution of panel (A). (C) Inverted fundamental mode (top) and first overtone (bottom) phase velocity dispersion curves and uncertainties (blue) compared to the dispersion curve calculated for the true model (red). The uncertainties shown here correspond to two standard deviation obtained from the posterior distribution.

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