



## 1. Motivation

- Normal modes provide some of the only direct constraints on density
- There is a lack of modern computational tools for accurate low-frequency spectra
- Self-coupling codes have significant errors (Akbarashrafi et al., 2018)

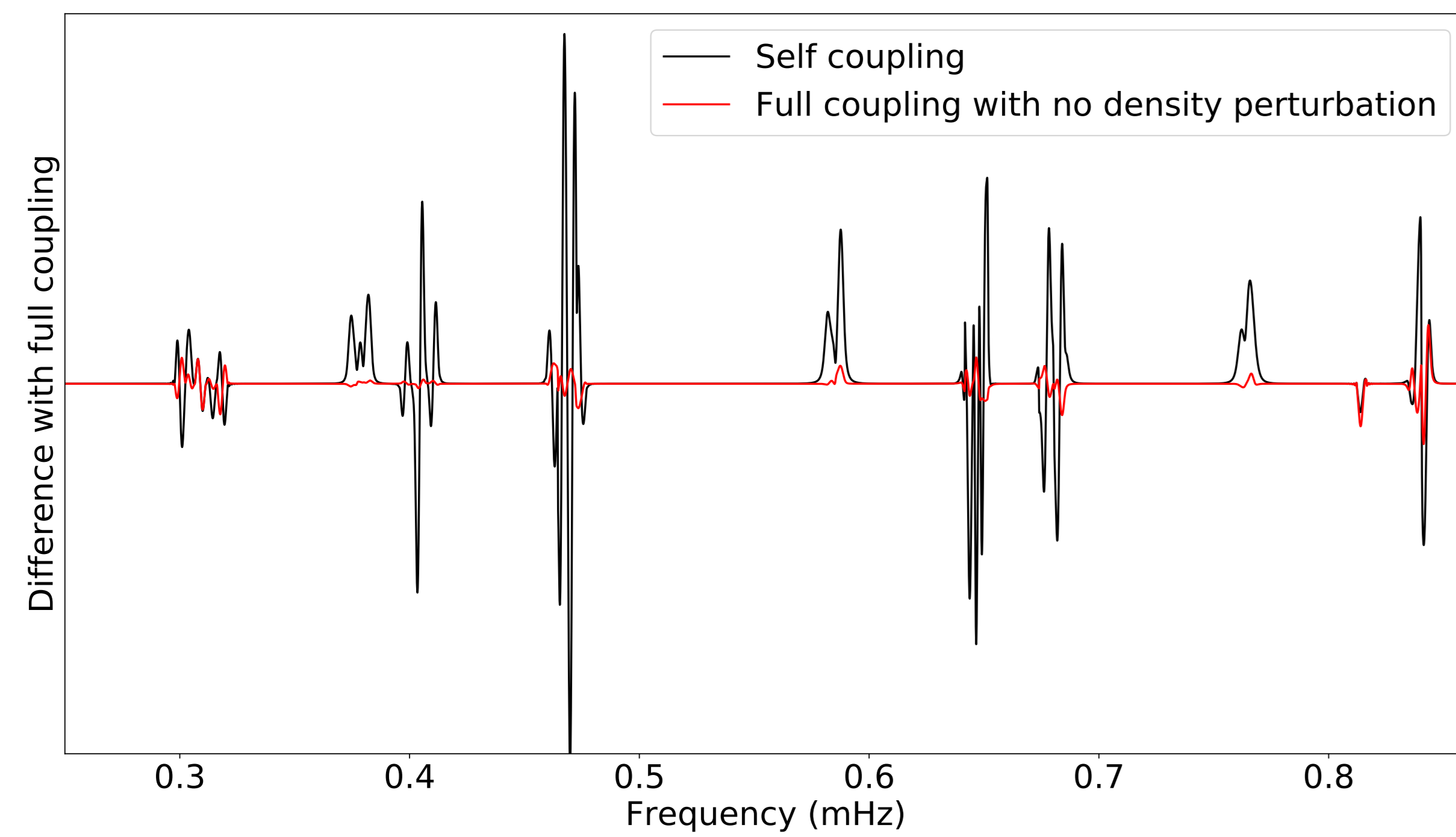


Fig. 1: Comparison of errors in self-coupling vs the size of the signal from density

- Full coupling still has approximations in both density and boundary terms (Al-Attar et al., 2018)

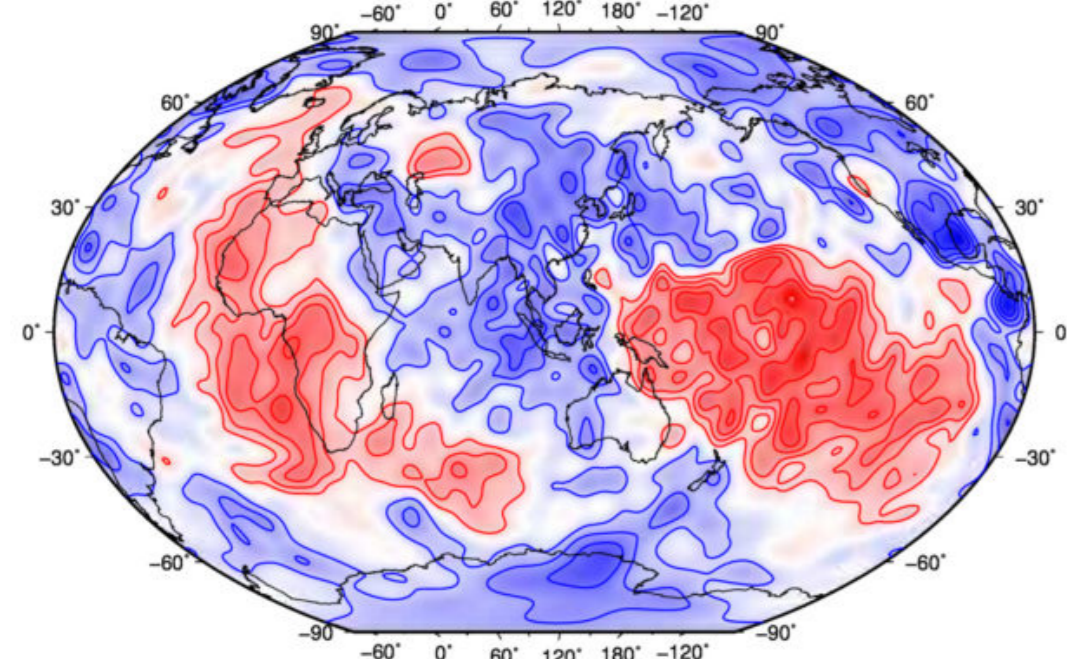


Fig. 2: Tomography model S40RTS (Ritsema et al., 2011) at the CMB

The LLSVPs are two large anomalies that consistently appear in tomography models. Their dynamic interpretation is still not fully known. Knowledge of their density from normal mode observations is important in determining their influence.

## 2. Normal mode coupling

- The elastic wave equation in the frequency domain (Dahlen & Tromp, 1998)

$$\mathcal{H}\mathbf{u} + 2i\omega\boldsymbol{\Omega} \times \mathbf{u} = \omega^2\mathbf{u} \quad (1)$$

- Expanding the wavefield using the 1D normal modes as a basis yields a matrix differential equation

$$S(\omega)\mathbf{u} = -\omega^2 T\mathbf{u} + 2i\omega W\mathbf{u} + V\mathbf{u} = \mathbf{f}(\omega) \quad (2)$$

### Calculating matrices

- Traditionally done using Wigner-Eckart theorem, e.g. for kinetic energy

$$T_{k,k'} = \sum_{st} (-1)^m \left[ \frac{(2l+1)(2s+1)(2l'+1)}{4\pi} \right]^{1/2} \begin{pmatrix} l & s & l' \\ -m & t & m' \end{pmatrix} \times \left\{ \int_0^a \delta\rho_{st} T_{\rho} r^2 dr + \sum_d d^2 \delta\tilde{d}_{st} [T_d]_{\pm}^{\pm} \right\} \quad (3)$$

- Ellipticity treated in a perturbative manner

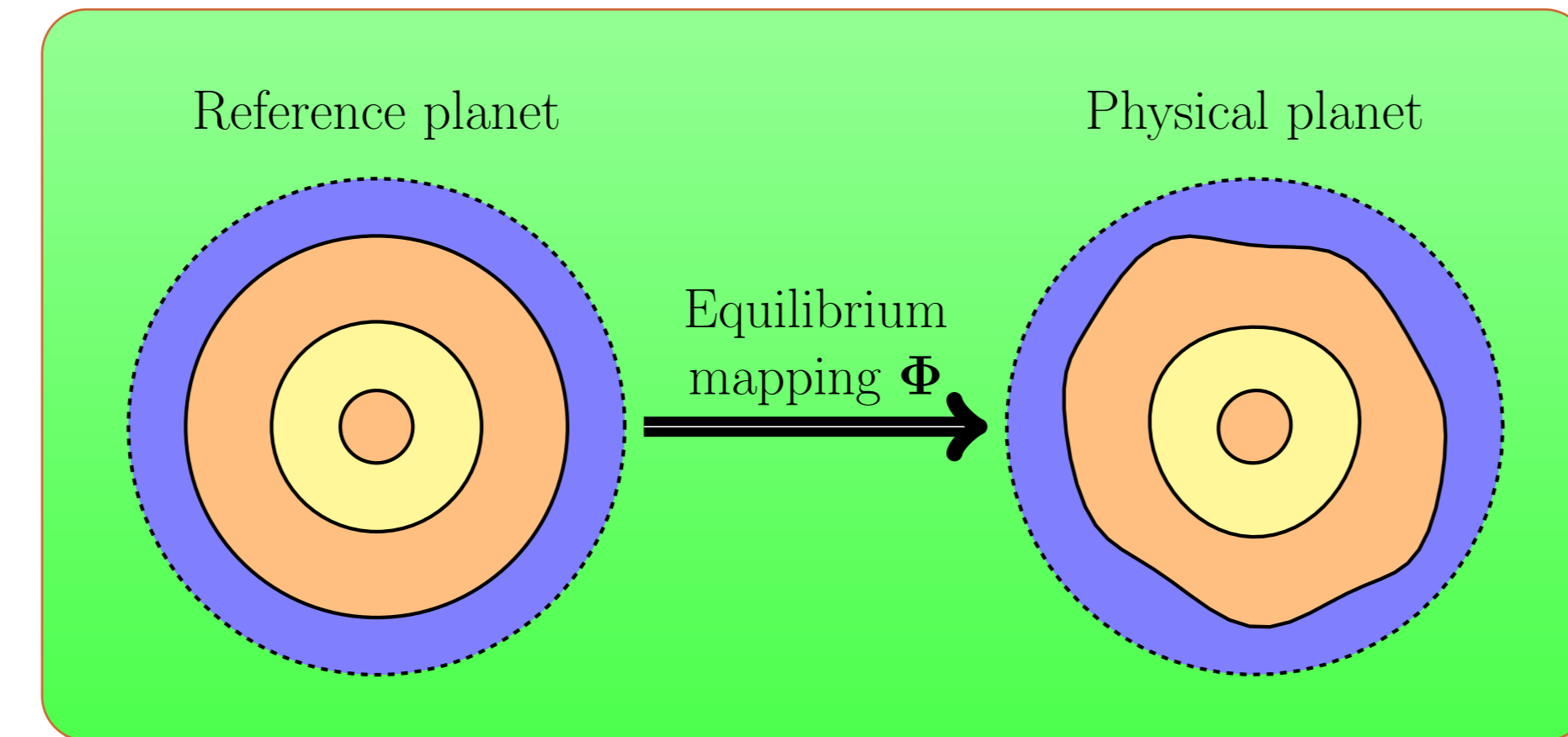
### Solving the matrix differential equation

- Non-standard eigenvalue/eigenvector problem
- Adds inaccuracy to solution for non-simple rheologies

The iterative direct solution method (IDSM) is an exact, efficient alternative suggested by Al-Attar et al. (2012), which solves for  $\mathbf{u}(\omega) = S^{-1}(\omega)\mathbf{f}(\omega)$  and performs an inverse Fourier-Laplace transform

## 3. The referential formulation

- Mapping  $\Phi$  between geometrically spherical reference planet and physical planet



- The boundary conditions on solid-fluid boundaries are complicated

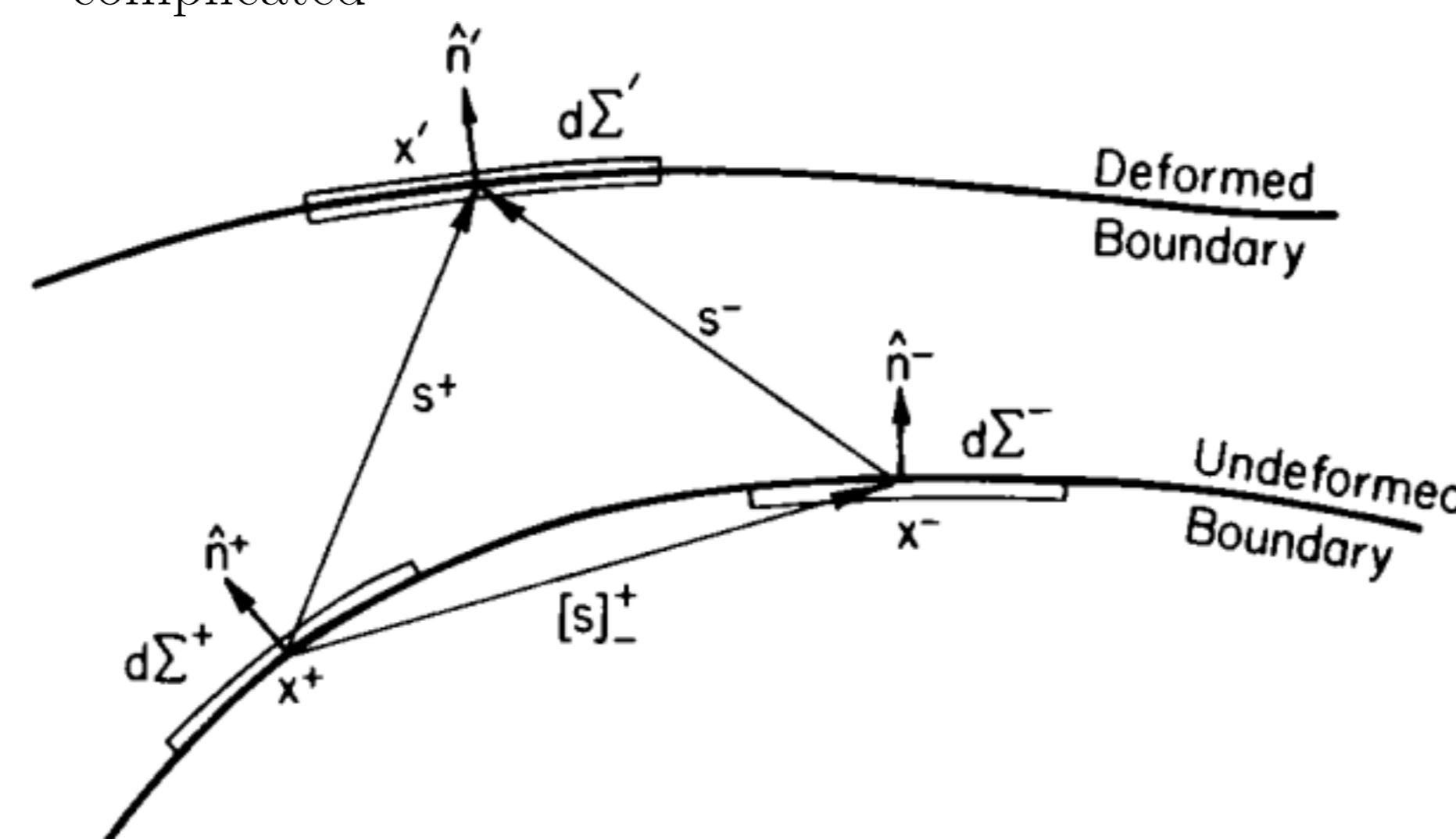


Fig. 3: Schematic of a slipping boundary from Woodhouse and Dahlen (1978)

## 5. Gravitational potential

- Weak form for the gravitational potential  $\zeta$  defined on the reference body with  $\chi$  as the test function is (Maitra & Al-Attar, 2019)

$$\mathcal{A}(\zeta, \chi) = \int_B \langle a \nabla \zeta, \nabla \chi \rangle d^3x + \sum_{lm} (l+1) b \zeta_{lm}(b) \overline{\chi_{lm}(b)} \quad (8)$$

$$= -4\pi G \int_M \rho \bar{\chi} d^3x$$

- $a = JF^{-1}F^{-T}$  where  $F_{ij} = \partial\xi_i/\partial x_j$  and  $\xi$  maps the referential to the physical body

- Separate into a radial and a spherical harmonic basis, ie

$$\zeta(r, \theta, \varphi) = \sum_{n=1}^N \sum_{l=0}^L \sum_{m=-l}^l \zeta_{lmn} h_n(r) Y_{lm}(\theta, \varphi) \quad (9)$$

- Solve without deriving the matrix  $A_{ij} = \mathcal{A}(\chi_i, \chi_j)$
- Uses generalised spherical harmonic (GSPH) transforms
- Motion is a perturbation to the equilibrium mapping

## 4. Exact matrices

- Normal modes don't satisfy tangential slip condition in referential body which is

$$\langle (\text{Def } \Phi)^{-1} \cdot (\mathbf{u}_2 - \mathbf{u}_1), \hat{\mathbf{n}} \rangle = 0 \quad (4)$$

- Require the set of functions  $\hat{\mathbf{u}}_n = (\text{Def } \Phi) \cdot \mathbf{u}_n$  where  $\mathbf{u}_n$  are the set of 1D normal modes

- Matrices are

$$T_{n'n} = \int_M \rho \langle \hat{\mathbf{u}}_n, \hat{\mathbf{u}}_{n'} \rangle d^3x \quad (5)$$

$$W_{n'n} = \int_M \rho \langle \boldsymbol{\Omega} \times \hat{\mathbf{u}}_n, \hat{\mathbf{u}}_{n'} \rangle d^3x \quad (6)$$

$$V_{n'n} = \int_M \{ \rho \langle \boldsymbol{\Omega} \times (\boldsymbol{\Omega} \times \hat{\mathbf{u}}_n) - \gamma_n, \hat{\mathbf{u}}_{n'} \rangle + \langle \boldsymbol{\Lambda} \cdot \text{Def } \hat{\mathbf{u}}_n, \text{Def } \hat{\mathbf{u}}_{n'} \rangle \} d^3x + \int_{\Sigma} \varpi \langle Q \cdot [\hat{\mathbf{u}}_n]_{\pm}, \text{Def } \hat{\mathbf{u}}_{n'} \rangle dS + \int_{\Sigma} \varpi \langle \text{Def } \hat{\mathbf{u}}_n, Q \cdot [\hat{\mathbf{u}}_{n'}]_{\pm} \rangle dS - \int_{\Sigma} \varpi \langle S \cdot [\hat{\mathbf{u}}_n]_{\pm}, [\hat{\mathbf{u}}_{n'}]_{\pm} \rangle dS \quad (7)$$

- We need to be able to solve exactly for the gravitational potential and its perturbation under motion
- The form of the matrix equations is still the exact same, ie the second part of the problem is unaffected and the IDSM can be used
- Calculation of matrix elements will be done via "spectral" method (Lognonné & Romanowicz, 1990)

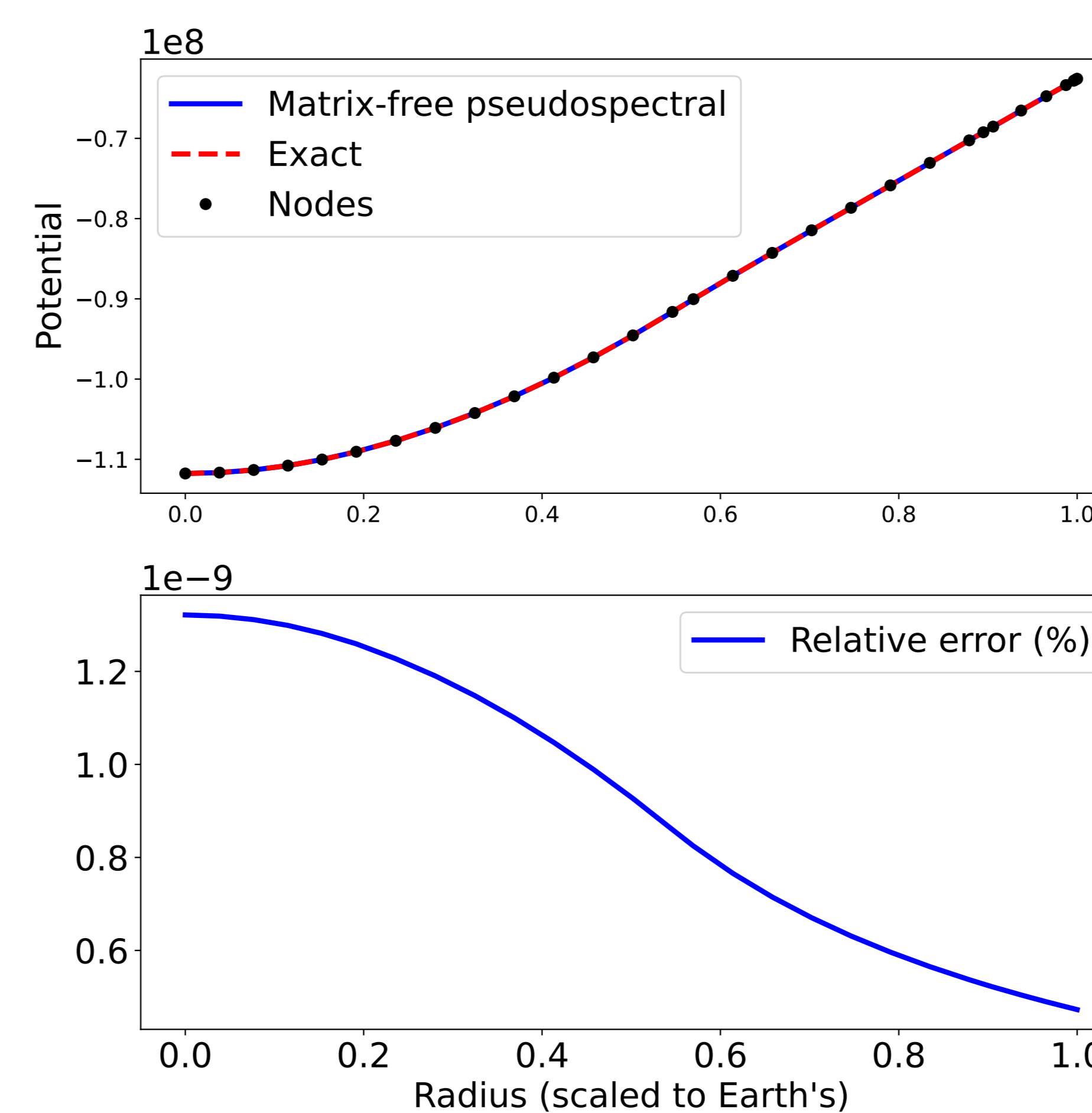
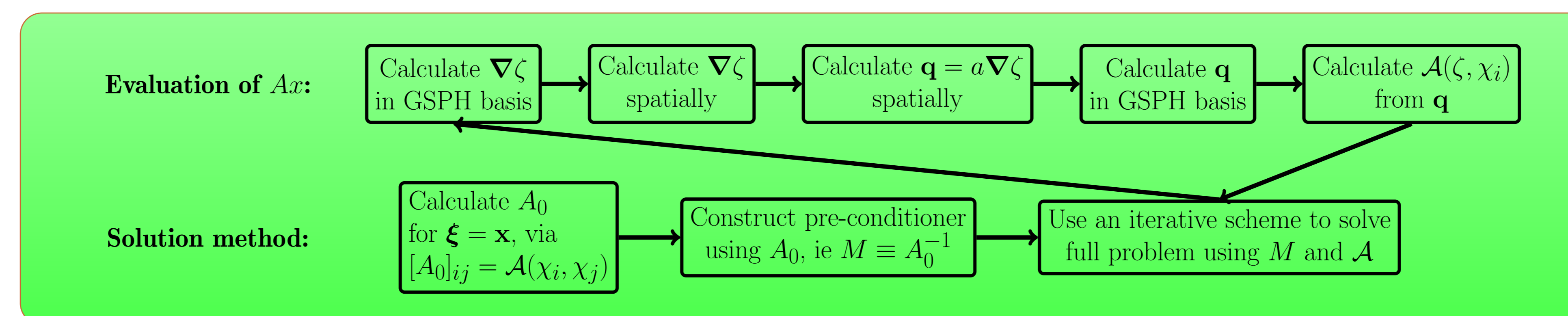


Fig. 4: Benchmark of matrix-free pseudospectral scheme against integral solution for the potential within the PREM model



## 6. Improving the IDSM

- The IDSM enables rapid, accurate computation

- Decompose  $S(\omega) = S_0(\omega)\tilde{S}(\omega)$  for which  $S \approx S_0$  and  $S_0$  is easily invertible

- Matrix-free iterative scheme ensures parallelisation performed with minimal memory usage increase via

$$S(\omega)\mathbf{u} = -\omega^2(T\mathbf{u}) + 2i\omega(W\mathbf{u}) + V\mathbf{u} \quad (10)$$

- BiCGSTAB and improved parallelisation result in significantly more rapid code

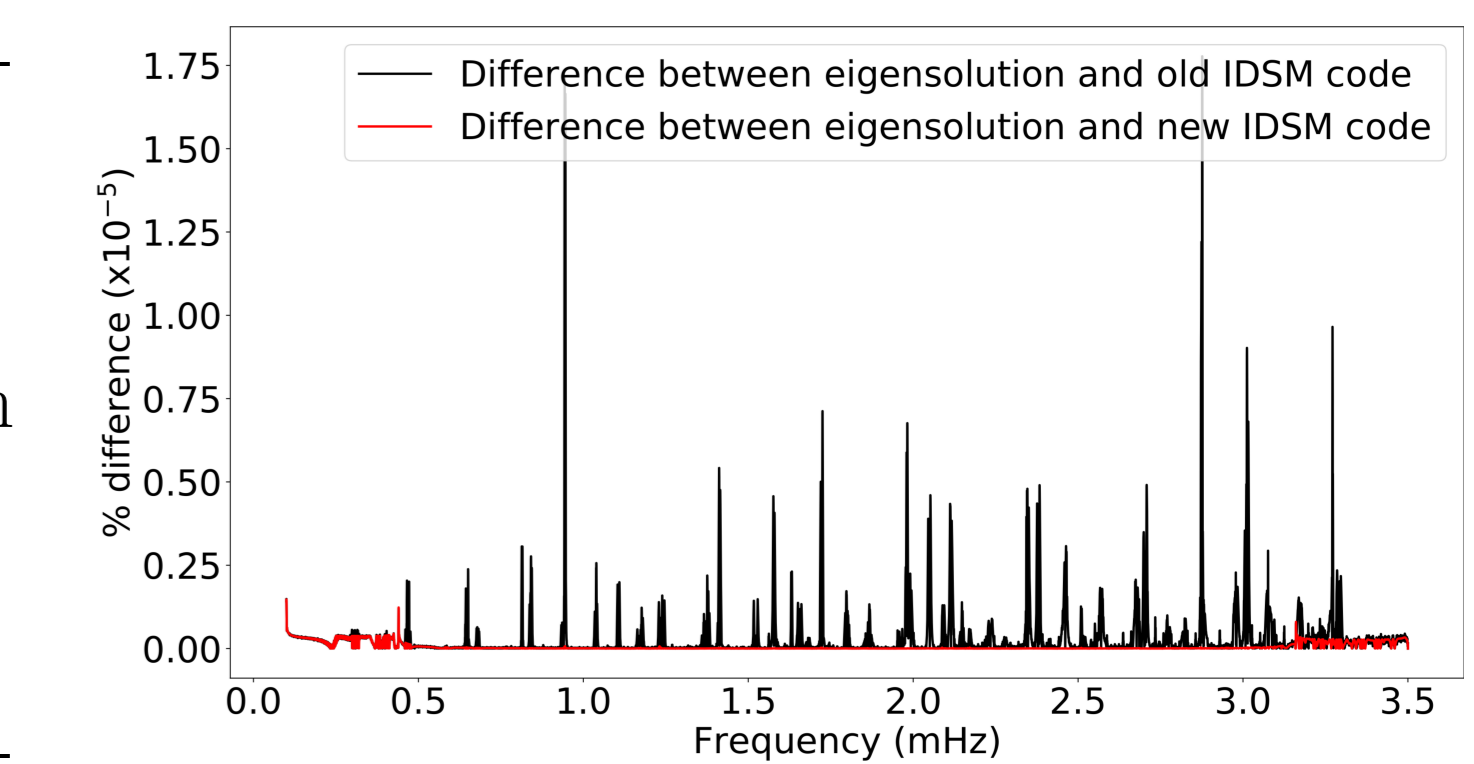


Fig. 5: Difference between IDSM and eigensolution

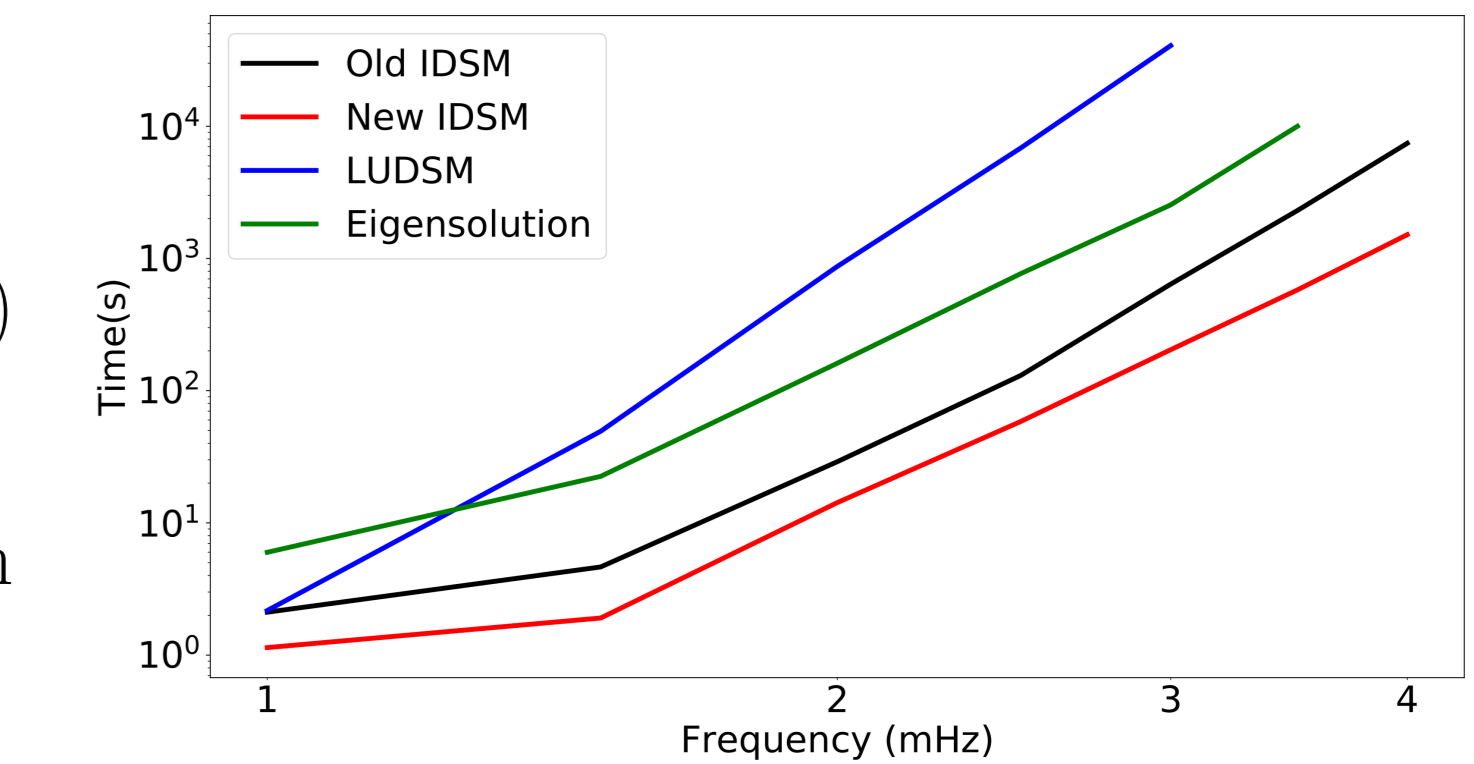


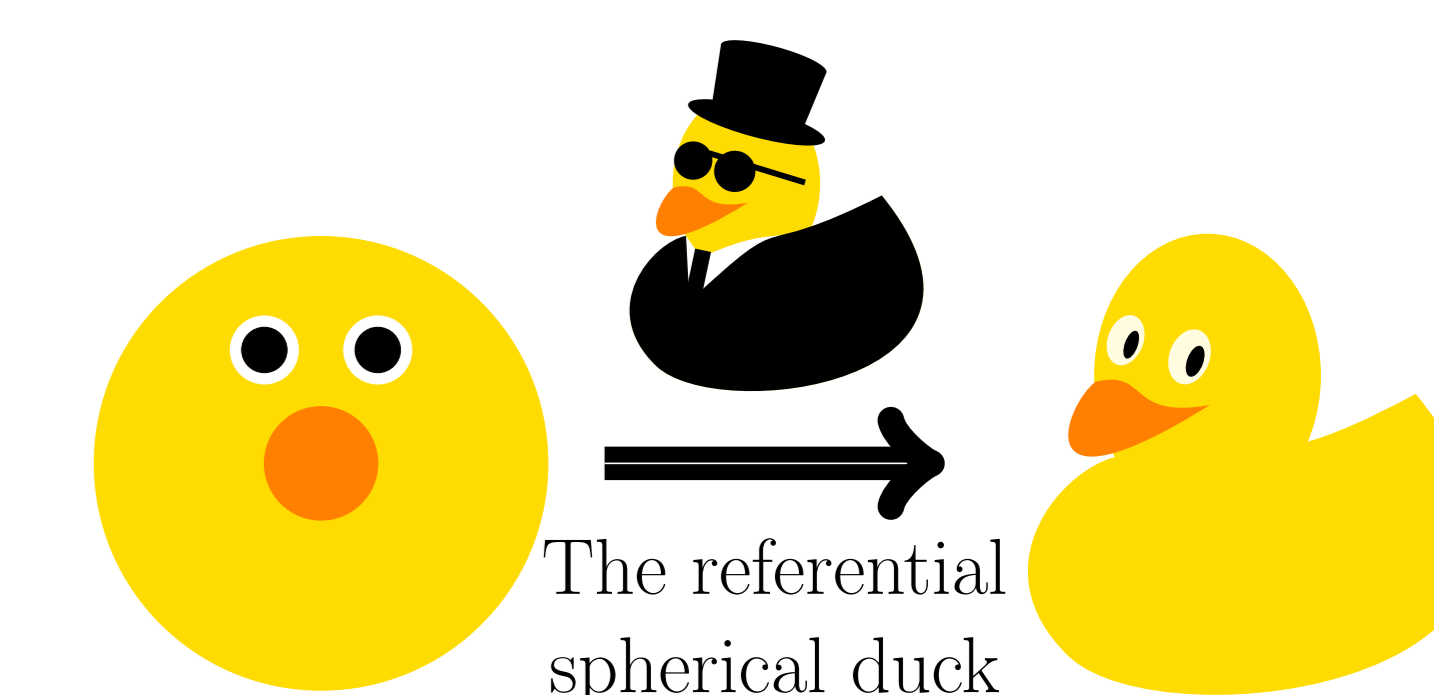
Fig. 6: Comparison of time scaling for various solvers

## 7. Future work

- Matrix elements for referential formulation to be calculated
- Set of basis functions cannot just be normal modes of 1D Earth
- Fully matrix-free implementation possible—efficiency and memory benefits for large coupling calculations
- Benchmarks:
  - (a) Traditional normal-mode codes with small density and topography variations
  - (b) Self-benchmarks with different  $\Phi$
- Implement adjoint calculations

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