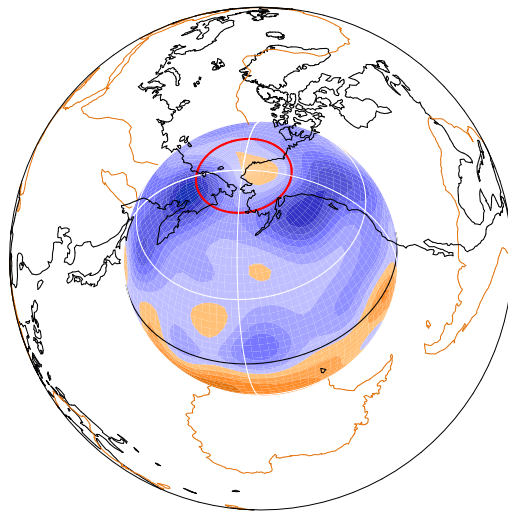


# LEOPACK



**itfvf**

Inhomogeneous **T**emperature **F**unction **V**ector **F**orm

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# 1 itfvf

## Inhomogeneous Temperature Function Vector Form

All of the time-stepping codes assume that the part of the temperature being solved for has uniform thermal boundaries. This means that at each of the inner and outer boundaries, the temperature ( $\Theta$ ) or the heat-flux ( $-\partial\Theta/\partial r$ ) is zero. Inner and outer boundaries may have either like or unlike boundary conditions. In order to impose a lateral variation in the temperature or heat-flux at either boundary, we therefore need to form a supplementary temperature function. This is referred to as  $T_a$ .

`itfvf` forms a function  $T_a(r, \theta, \phi)$  in the standard format (i.e. with a `.ints`, `.vecs` and `.xarr` file).  $T_a$  has an essentially arbitrary form apart from the conditions that it must satisfy either

$$T_a(r_i, \theta, \phi) = g_i(\theta, \phi) \quad (1)$$

or

$$\left. \frac{\partial T_a(r, \theta, \phi)}{\partial r} \right|_{r=r_i} = g_i(\theta, \phi) \quad (2)$$

at the inner boundary, and either

$$T_a(r_o, \theta, \phi) = g_o(\theta, \phi) \quad (3)$$

or

$$\left. \frac{\partial T_a(r, \theta, \phi)}{\partial r} \right|_{r=r_o} = g_o(\theta, \phi) \quad (4)$$

at the outer boundary: depending upon whether a constant temperature or constant heat-flux boundary condition is to be satisfied.

The functions  $g_i$  and  $g_o$  are both expanded as series of spherical harmonics:

$$g_i(\theta, \phi) = \frac{\sqrt{\varepsilon_i}}{N_i} \sum_{l=0}^{l=L} \sum_{m=0}^{m=l} \left[ g_{i,l}^{mc} \cos m\phi + g_{i,l}^{ms} \sin m\phi \right] P_l^m(\cos \theta) \quad (5)$$

and

$$g_o(\theta, \phi) = \frac{\sqrt{\varepsilon_o}}{N_o} \sum_{l=0}^{l=L} \sum_{m=0}^{m=l} \left[ g_{o,l}^{mc} \cos m\phi + g_{o,l}^{ms} \sin m\phi \right] P_l^m(\cos \theta), \quad (6)$$

where the associated Legendre function  $P_l^m(\cos \theta)$  satisfies the Schmidt quasi-normalisation condition

$$\int_0^\pi [P_l^m(\cos \theta)]^2 \sin \theta d\theta = \frac{2(2 - \delta_{m0})}{2l + 1}. \quad (7)$$

The user specifies the non-zero coefficients  $g_{i,l}^{mc}$ ,  $g_{i,l}^{ms}$ ,  $g_{o,l}^{mc}$  and  $g_{o,l}^{ms}$ , and **itfvf** calculates normalisation factors,  $N_i$  and  $N_o$ , such that

$$\int_0^{2\pi} \int_0^\pi [g_i(\theta, \phi)]^2 \sin \theta d\theta d\phi = \varepsilon_i \quad (8)$$

and

$$\int_0^{2\pi} \int_0^\pi [g_o(\theta, \phi)]^2 \sin \theta d\theta d\phi = \varepsilon_o. \quad (9)$$

Therefore, if a user specified that the only non-zero coefficients in the series (5) and (6) were to be  $g_{o,2}^{2c} = 3.0$  and  $g_{o,3}^{2s} = -3.0$  with  $\varepsilon_o = 1.0$ , the result would be exactly the same had the user specified the coefficients  $g_{o,2}^{2c} = 0.012$  and  $g_{o,3}^{2s} = -0.012$  with  $\varepsilon_o = 1.0$ . This gives the user a global control on the normalisation of the “strength” of the functions  $g_i$  and  $g_o$  without having to worry about the actual coefficients themselves, other than their relative sizes. I personally tend to set  $\varepsilon_o$  to unity to ensure a simple normalisation and then alter the strength of the boundary heating using the **SCAL** parameter in the inhomogeneous thermal boundary time-step programs. I have yet to impose a lateral thermal variation at the inner boundary, but designed **itfvf.f** with the possibility in mind.

The stand-alone source code version of the program is compiled by typing

```
make itfvf
```

within this directory. Once the executable is created, begin execution by typing

```
itfvf < inputfile
```

The inputs file must have the following format.

---

```
* input file for itfvf
root                               : Output filename stem
boundary_coeffs                    : File containing coeff.s
 40  2   4   0.53846154   1.5384615   1 : nr isp ifrmf ri ro ithebc
 1.0      1.0                               : epsi  epso
```

---

Any line in the input file beginning with an asterisk, **\***, is ignored by the program and can thus be used to enter comments and notes.

The arguments are as follows

- **filename stem**: First characters in output files to be generated by current run. Running **itfvf** with the above input file will create the files **root.ints**, **root.vecs** and **root.xarr**

- **coefficient file name.** This file contains the specifications for the coefficients  $g_{i,l}^{mc}$ ,  $g_{i,l}^{ms}$ ,  $g_{o,l}^{mc}$  and  $g_{o,l}^{ms}$ . Details regarding the form of this file are found below.
- **nr:** number of radial grid nodes to be used in the expression of our temperature function,  $T_a(r, \theta, \phi)$ .
- **isp:** flag for the spacing of these grid nodes. **isp** = 1 forces evenly spaced grid nodes from ESNAAS and **isp** = 2 forces Chebyshev zero spaced nodes from ZCPAAS (see Figure 2).
- **ifrmf:** flag for order of elements in the resulting solution vector. **ifrmf** may take the values 3 and 4. For explanation, see Figure (1).
- **ri:** Radius of the inner boundary,  $r_i$ .
- **ro:** Radius of the outer boundary,  $r_o$ .
- **ithebc:** Flag for thermal boundary conditions at inner and outer boundaries. **ithebc** = 1 means fixed temperature at both inner and outer boundaries. (In other words, we apply conditions 1 and 3.)  
**ithebc** = 2 means fixed temperature at the inner boundary and fixed flux at the outer. (In other words, we apply conditions 1 and 4.)  
**ithebc** = 3 means fixed flux at the inner boundary and fixed temperature at the outer. (In other words, we apply conditions 2 and 3.)
- **epsi:**  $\varepsilon_i$ .
- **epso:**  $\varepsilon_o$ .

Note that it is quite trivial to create a program which does not normalise the output, and uses the exact same coefficients in the input file.

## 1.1 Specifying coefficients for thermal heterogeneity

In the above input file, a filename is specified: **boundary\_coeffs**. This file specifies all coefficients  $g_{i,l}^{mc}$ ,  $g_{i,l}^{ms}$ ,  $g_{o,l}^{mc}$  and  $g_{o,l}^{ms}$ . Coefficients not listed are assumed to be zero.

This file contains one line for each coefficient. The first two characters of each line must contain either **IB** or **OB**: which specify either inner or outer boundary. These two characters must be exactly at the beginning of a line. The remaining part of the line contains, in any format, the numbers **L**, **M**, **ICS** and **COEF**.

**L** and **M** are clearly  $l$  and  $m$ : **ICS** is 1 to indicate a  $\cos m\phi$  harmonic (i.e.  $g_{i,l}^{mc}$  or  $g_{o,l}^{mc}$ ) and 2 to indicate a  $\sin m\phi$  harmonic (i.e.  $g_{i,l}^{ms}$  or  $g_{o,l}^{ms}$ ). **COEF** is ofcourse

just the value of the coefficient.

## Examples

The file

```
OB   2   2   1       1.0
```

sets all coefficients to zero except for  $g_{o,2}^{2c}$  which is set to unity.

The file

```
IB   2   2   1       1.0
OB   2   2   1       1.0
OB   3   2   2       1.0
```

sets  $g_{i,2}^{2c} = 1.0$ ,  $g_{o,2}^{2c} = 1.0$  and  $g_{o,3}^{2s} = 1.0$ . This fixes the contribution from  $g_{o,2}^{2c}$  relative to that from  $g_{o,3}^{2s}$ , whereas the contribution from  $g_{i,2}^{2c}$  is only fixed relative to the others by the parameters  $\varepsilon_i$  and  $\varepsilon_o$ .

## 1.2 Subprograms required for itfvf

### SUBS subroutines

```
vecop.f fopen.f fclose.f esnaas.f zcpaas.f shcanc.f
ithcar.f itfa.f hmfwt.f svfwt.f xarrwt.f fnamer.f
lmfind.f itfcf.f cntric.f
```

### SUBS integer function

```
indfun.f indshc.f
```

## 1.3 Run-time limitations

Several parameters are set at the outset which limit the physical size of the problem.

```
INTEGER NRMAX, NHMAX, ISVMAX, LHMAX, LHLH2M, NDCS, NITHMX
PARAMETER ( NRMAX = 100, LHMAX = 64, LHLH2M = LHMAX*(LHMAX+2),
1          NHMAX = LHLH2M+1, ISVMAX = NHMAX*NRMAX, NDCS = 1,
2          NITHMX = NHMAX )
```

If the values are insufficient, then change them and recompile. (Note that NDCS is not size dependent and should not be changed.)

- NRMAX is the maximum permitted number of radial grid nodes.
- LHMAX is the highest permitted spherical harmonic degree,  $l$ .

## 1.4 Sample runs of itfvf

The directory

`$LEOPACK_DIR/SAMPLERUNS/ITFVF`

contains example input files, coefficient files and model output. Do not under any circumstances edit these files, as these examples should serve as a control for the correct working of the code. After compiling the program, copy the `.input` and `.coeff` files to another directory, run the code and confirm that the output agrees with that in the directory.

### 1.4.1 Example a

We have a spherical shell with radius ratio  $r_o/r_i = 0.35$ . We want a fixed heat-flux with a  $Y_2^{2c}$  variation at the outer boundary only, 40 grid nodes with Chebyshev spacing and  $\varepsilon = 1$ .

```
itfvf < example_a.input
```

produces the three output files `example_aOUTPUT.ints`, `example_aOUTPUT.vecs` and `example_aOUTPUT.xarr`.

$$\begin{array}{cc}
(a) & \left[ \begin{array}{c} f_1(r_1) \\ f_2(r_1) \\ \vdots \\ f_j(r_1) \\ \vdots \\ f_{N_h}(r_1) \\ \dots \\ f_1(r_2) \\ f_2(r_2) \\ \vdots \\ f_j(r_2) \\ \vdots \\ f_{N_h}(r_2) \\ \dots \\ \vdots \\ \dots \\ f_1(r_i) \\ f_2(r_i) \\ \vdots \\ f_j(r_i) \\ \vdots \\ f_{N_h}(r_i) \\ \dots \\ \vdots \\ \dots \\ f_1(r_{N_r}) \\ f_2(r_{N_r}) \\ \vdots \\ f_j(r_{N_r}) \\ \vdots \\ f_{N_h}(r_{N_r}) \end{array} \right] \\
(b) & \left[ \begin{array}{c} f_1(r_1) \\ f_1(r_2) \\ \vdots \\ f_1(r_i) \\ \vdots \\ f_1(r_{N_r}) \\ \dots \\ f_2(r_1) \\ f_2(r_2) \\ \vdots \\ f_2(r_i) \\ \vdots \\ f_2(r_{N_r}) \\ \dots \\ \vdots \\ \dots \\ f_j(r_1) \\ f_j(r_2) \\ \vdots \\ f_j(r_i) \\ \vdots \\ f_j(r_{N_r}) \\ \dots \\ \vdots \\ \dots \\ f_{N_h}(r_1) \\ f_{N_h}(r_2) \\ \vdots \\ f_{N_h}(r_i) \\ \vdots \\ f_{N_h}(r_{N_r}) \end{array} \right]
\end{array}$$

Figure 1: Alternative arrangements of elements within the solution vector. (a) shows **IFORMF** = **3** and (b) shows **IFORMF** = **4**.

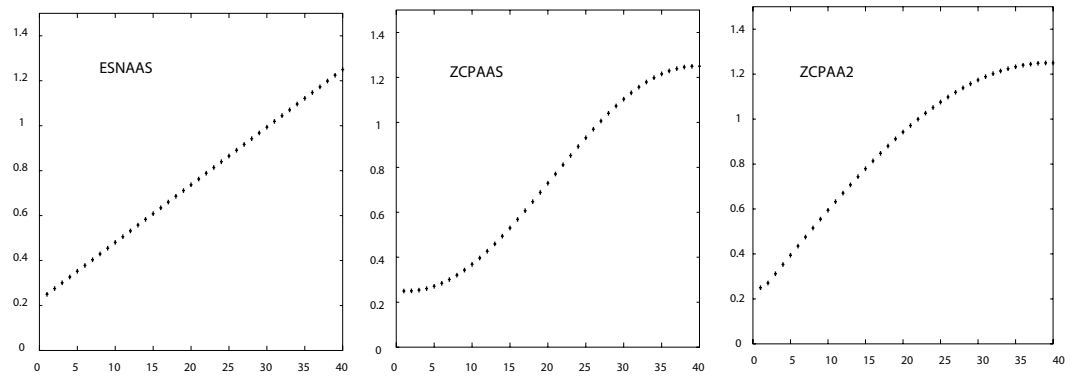


Figure 2: Distributions of radial grid nodes with  $r_i = 0.25$ ,  $r_o = 1.25$  and  $N_r = 40$ . Subroutine used to generate  $r$  values as shown.