In collaboration with:
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FreeFEM is a partial differential equation solver for non-linear multi-physics systems in 1D, 2D, 3D and 3D border domains (surface and curve).

Problems involving partial differential equations from several branches of physics, such as fluid-structure interactions, require interpolations of data on several meshes and their manipulation within one program. FreeFEM includes a fast interpolation algorithm and a language for the manipulation of data on multiple meshes.

FreeFEM is written in C++ and its language is a C++ idiom.

FreeFEM currently interfaces to the following libraries:

- ARPACK
- BLAS
- OpenBLAS
- FFTW 3.3.2
- Ipopt 3.12.4
- Gmm++ 4.2
- freeYams
- METIS
- ParMETIS
- Mmg
- mshmet
- MUMPS
- NLopt 2.2.4
- ScaLAPACK
- Scotch
- SuiteSparse
- SuperLU
- TetGen
- PETSc
- SLEPe
- HTool
- HPDDM
- BemTool
- ParMmg
1.1 Version 4.5: new features

1.1.1 Release, binaries packages

- Since the version 4.5, the FreeFEM binary packages provides with a compiled PETSc library.
- FreeFEM is now interfaced with ParMmg.

1.1.2 New meshes and FEM border

After Surface FEM, Line FEM is possible with a new mesh type meshL, P0 P1 P2 P1dc FE, basic FEM, mesh generation. This new development allows to treat a 1d problem, such as a problem described on a 3d curve.

Abstract about Line FEM in FreeFEM.

- new meshL type, refer to the section The type meshL in 3 dimension
  - new type of surface mesh: meshL
  - the functionalities on the meshL type, it is necessary to load the plugin "msh3".
  - generator of meshL segment, define multi border and buildmesh function.
  - basic transformation are available: movemesh, trunc, extract, checkmesh, change, AddLayers, glue of meshL.

It is possible to build the underlying meshL from a meshS with the function buildBdMesh: ThS=buildBdMesh(ThS) builds the boundary domain associated to the meshS ThS and extract it by the command meshL ThL=ThS. Gamma.

- new finite element space with curve finite element type
- FESpace P0 P1,P2,P1dc Lagrange finite elements and possible to add a custumed finite element with the classical method (like a plugin).
- as in the standard 2d, 3d, surface 3d case, the variational problem associated to surface PDE can be defined by using the keywords
  - problem
  - varf to access to matrix and RHS vector
  - available operators are int1d, on and the operator int0d to define a Neumann boundary condition
- visualisation tools
  - plot with plot of ffglut, medit meshes meshL and solutions
  - 2d or 3d view, with in 3d the option to visualize the element Normals at element (touch 'T') and the deformed domain according to it (touch '2').
  - loading, saving of meshes and solution at FreeFEM’s format
    - "mesh" mesh format file of Medit (P. Frey LJLL)
    - "msh" for mesh and "sol" data solution at freefem format
1.1.3 Boundary Element Method

Allows to define and solve a 2d/3d BEM formulation and rebuild the associated potential. The document is in construction.

1.2 Installation guide

To use FreeFEM, two installation methods are available: user access (binary package) and access developers (from the source code). Follow the section corresponding to your type of installation.

Note: Since the version 4.5, FreeFEM release provides with the last version of PETSc.

1.2.1 Using binary package

First, open the following web page download page and choose your platform: Linux, MacOS or Windows.

Note: Binary packages are available for Microsoft Windows, MacOS and some Linux distributions. Since the release 4.5, FreeFEM binaries provide with the current version of PETSc.

Install FreeFEM by double-clicking on the appropriate file. Under Linux and MacOS the install directory is one of the following /usr/local/bin, /usr/local/share/freefem++, /usr/local/lib/ff++

Windows installation

Note: The windows package is build for Window 7 64bits. The support ended for all releases under Windows 32 bits since the V4.

First download the windows installation executable, then double click to install FreeFEM. Install MSMPI for parallel version under window64 MS MPI V10.1.2, and install both msmpisdk.msi and MSMpiSetup.exe.

In most cases just answer yes (or type return) to all questions.

Otherwise in the Additional Task windows, check the box “Add application directory to your system path.” This is required otherwise the program ffglut.exe will not be found.

By now you should have two new icons on your desktop:

• FreeFem++ (VERSION).exe, the freefem++ application.
• FreeFem++ (VERSION) Examples, a link to the freefem++ examples folder.
where (VERSION) is the version of the files (for example 4.5).
By default, the installed files are in C:\Programs Files\FreeFem++. In this directory, you have all the .dll files and other applications: FreeFem++-nw.exe, ffglut.exe,... The syntax for the command-line tools are the same as those of FreeFem.exe.
To use FreeFEM binaries under Windows, two methods are possible:
  • Use the FreeFEM launcher (launchff++.exe)
  Warning: if you launch FreeFEM without filename script by double-clicking, your get a error due (it is bug of usage GetOpenFileName in win64).
  • In shell terminal (cmd, powershell, bash, ... ):
    – To launch sequential version:
      

```
C:\>"Program Files (x86)\FreeFem++\FreeFem++.exe" <mySequentialScript.edp
```

– To launch parallel version:

```
C:\>"Program Files\Microsoft MPI\Bin\mpiexec.exe" -n <nbProcs> C:\>"Program Files ...
˓
→(x86)\FreeFem++\FreeFem++-mpi.exe" <myParallelScript.edp
```

macOS X installation

Download the macOS X binary version file, extract all the files by double clicking on the icon of the file, go the the directory and put the FreeFem++.app application in the /Applications directory.
If you want terminal access to FreeFEM just copy the file FreeFem++ in a directory of your $PATH shell environment variable.

Ubuntu installation

**Note:** The Debian package is built for Ubuntu 16.04

Beforehand, install the following dependances libraries using the apt tool:

```
sudo apt-get install libgsl-dev libhdf5-dev
liblapack-dev libopenmpi-dev freeglut3-dev
```

Download the package FreeFEM .deb, install it by the command

```
dpkg -i FreeFEM_VERSION_Ubuntu_withPETSc_amd64.deb
```
FreeFEM is directly available in your terminal by the command “FreeFem++”.

Arch AUR package

An up-to-date package of FreeFEM for Arch is available on the Archlinux user repository.
To install it:
1.2.2 Compiling source code

Various versions of FreeFEM are possible:

- sequential and without plugins (contains in 3rdparty)
- parallel with plugins (and with PETSc).

**Note:** We advise you to use the package manager for macOS Homebrew to get the different packages required available here.

Compilation on OSX (>=10.13)

1. Install Xcode, Xcode Command Line tools and Xcode Additional Tools from the Apple website
2. Install gfortran from Homebrew
   ```bash
   brew cask install gfortran
   ```
   **Note:** If you have installed gcc via brew, gfortran comes with it and you do not need this line
3. To use FreeFEM parallel version, install openmpi or mpich
   ```bash
   # to install openmpi
   curl -L https://download.open-mpi.org/release/open-mpi/v4.0/openmpi-4.0.1.tar.gz --output openmpi-4.0.1.tar.gz
   tar xf openmpi-4.0.1.tar.gz
   cd openmpi-4.0.1/
   # to install mpich
   curl -L http://www.mpich.org/static/downloads/3.3.2/mpich-3.3.2.tar.gz --output mpich-3.3.2.tar.gz
   tar xf mpich-3.3.2.tar.gz
   cd mpich-3.3.2
   ```
   ```bash
   # with brew gcc gfortran compilers
   ./configure CC=clang CXX=clang++ FC=gfortran-9 F77=gfortran-9 --prefix=/where/you/want/to/have/files/installed
   # with LLVM gcc and brew gfortran compilers
   ./configure CC=gcc-9 CXX=g++-9 FC=gfortran-9 F77=gfortran-9 --prefix=/where/you/want/to/have/files/installed
   ```

```bash
make -j<nbProcs>
makewhattr -si
```
4. Install the minimal libraries for **FreeFEM**

   brew install m4 git flex bison

5. If you want build your own configure according your system, install autoconf and automake from Homebrew (optional, see note in step 10)

   brew install autoconf automake

6. To use **FreeFEM** with its plugins, install from Homebrew suitesparse, hdf5, cmake, wget

   brew install suitesparse hdf5 cmake wget

7. Install **gsl**

   curl -O http://mirror.cyberbits.eu/gnu/gsl/gsl-2.5.tar.gz
   tar zxvf gsl-2.5.tar.gz
   cd gsl-2.5
   ./configure
   make -j<nbProcs>
   make -j<nbProcs> install --prefix=/where/you/want/to/have/files/installed

8. Download the latest Git for Mac installer git and the **FreeFEM** source from the repository

   git clone https://github.com/FreeFem/FreeFem-sources.git

9. Configure your source code

   cd FreeFem-sources
   autoreconf -i

   **Note:** if your autoreconf version is too old, do tar zxvf AutoGeneratedFile.tar.gz

   • following your compilers

   // with brew gcc gfortran compilers
   ./configure --enable-download CC=clang CXX=clang++ F77=gfortran-9
   FC=gfortran-9 --prefix=/where/you/want/to/have/files/installed
   // with LLVM gcc and brew gfortran compilers
   ./configure --enable-download CC=clang CXX=clang++ F77=gfortran-9
   FC=gfortran-9 --prefix=/where/you/want/to/have/files/installed

10. Download the 3rd party packages to use FreeFEM plugins

    ./3rdparty/getall -a

    **Note:** All the third party packages have their own licence

11. If you want use PETSc/SLEPc and HPDDM (High Performance Domain Decomposition Methods)

    cd 3rdparty/ff-petsc
    make petsc-slepc // add SUDO=sudo if your installation directory is the default /usr/local
    (continues on next page)
12. Build your **FreeFEM** library and executable

```bash
make -j<nbProcs>
make -j<nbProcs> check
```

**Note:** `make check` is optionnally, but advise to check the validity of your **FreeFEM** building

13. **Install the FreeFEM application**  make install // add SUDO=sudo might be necessary

**Note:** it isn’t necessary to execute this last command, FreeFEM executable is available here your_installation/src/nw/FreeFem++ and mpi executable here your_installation/src/mpi/ff-mpirun.

### Compilation on Ubuntu

1. Install the following packages on your system

```bash
sudo apt-get update && sudo apt-get upgrade
sudo apt-get install cpp freeglut3-dev g++ gcc gfortran \
m4 make patch pkg-config wget python unzip \
liblapack-dev libhdf5-dev libgsl-dev \
autoconf automake autotools-dev bison flex gdb git cmake
# mpich is required for the FreeFEM parallel computing version
sudo apt-get install mpich
```

**Warning:** In the oldest distribution of Ubuntu, libgsl-dev does not exists, use libgsl2-dev instead

2. Download **FreeFEM** source from the repository

```bash
git clone https://github.com/FreeFem/FreeFem-sources.git
```

3. Autoconf

```bash
cd FreeFem-sources
autoreconf -i
```

**Note:** if your autoreconf version is too old, do `tar zxvf AutoGeneratedFile.tar.gz`

4. Configure

```bash
./configure --enable-download --enable-optim \
--prefix=/where/you/want/to/have/files/installed
```
Note: To see all the options, type ./configure --help

5. Download the 3rd party packages

   ./3rdparty/getall -a

Note: All the third party packages have their own licence

6. If you want use PETSc/SLEPC and HPDDM (High Performance Domain Decomposition Methods) for massively parallel computing

   cd 3rdparty/ff-petsc
   make petsc-slepc // add SUDO=sudo if your installation directory is the default /usr/local
   cd -
   ./reconfigure

7. Build your FreeFEM library and executable

   make -j<nbProcs>
   make -j<nbProcs> check

Note: make check is optionally, but advise to check the validity of your FreeFEM building

8. Install the executable

   make install

Note: it isn’t necessary to execute this last command, FreeFEM executable is available here your_installation/src/nw/FreeFem++ and mpi executable here your_installation/src/mpi/ff-mpirun

Compilation on Arch Linux

Warning: As Arch is in rolling release, the following information can be quickly outdated!

Warning: FreeFEM fails to compile using the newest version of gcc 8.1.0, use an older one instead.

1. Install the following dependencies:

   pacman -Syu
   pacman -S git openmpi gcc-fortran wget python
           freeglut m4 make patch gmm
           blas lapack hdf5 gsl fftw arpack suitesparse
           gnuplot autoconf automake bison flex gdb
           valgrind cmake texlive-most

1.2. Installation guide
2. Download the **FreeFEM** source from the repository

   ```bash
git clone https://github.com/FreeFem/FreeFem-sources.git
   ```

3. Autoconf

   ```bash
cd FreeFem-sources
autoreconf -i
   ```

4. Configure

   ```bash
./configure --enable-download --enable-optim
   ```

   **Note:** To see all the options, type `./configure --help`

5. Download the packages

   ```bash
./3rdparty/getall -a
   ```

   **Note:** All the third party packages have their own licence

6. If you want use HPDDM (High Performance Domain Decomposition Methods) for massively parallel computing, install PETSc/SLEPc

   ```bash
cd 3rdparty/ff-petsc
make petsc-slepc SUDO=sudo
cd -
./reconfigure
   ```

7. Compile the **FreeFEM** source

   ```bash
make
   ```

   **Note:** If your computer has many threads, you can run `make` in parallel using `make -j16` for 16 threads, for example.

   **Note:** Optionally, check the compilation with `make check`

8. Install the **FreeFEM** application

   ```bash
sudo make install
   ```

---

**Compilation on Linux with Intel software tools**

Follow the [guide](https://www.intel.com/content/www/us/en/intel-software-development-kits/)**
Compilation on Windows

**Warning:** The support ended for all releases under Windows 32 bits since the V4. We assume your development machine is 64-bit, and you want your compiler to target 64-bit windows by default.

1. Install the Microsoft MPI v10.1.2 (archived) (msmpisdk.msi and MSMpiSetup.exe)

2. Download msys2-x86_64-latest.exe (x86_64 version) and run it.

3. Install the version control system Git for Windows

4. In the MSYS2 shell, execute the following. Hint: if you right click the title bar, go to Options -> Keys and tick “Ctrl+Shift-letter shortcuts” you can use Ctrl+Shift+V to paste in the MSYS shell.

   ```sh
   pacman -Syuu
   ```

   Close the MSYS2 shell once you’re asked to. There are now 3 MSYS subsystems installed: MSYS2, MinGW32 and MinGW64. They can respectively be launched from C:devmsys64msys2.exe, C:devmsys64mingw32.exe and C:devmsys64mingw64.exe Reopen MSYS2 (doesn’t matter which version, since we’re merely installing packages). Repeatedly run the following command until it says there are no further updates. You might have to restart your shell again.

   ```sh
   pacman -Syuu
   ```

5. Now that MSYS2 is fully up-to-date, install the following dependancies

   - for 64 bit systems:

     ```sh
     pacman -S autoconf make automake-wrapper bison git \
     mingw-w64-x86_64-freeglut mingw-w64-x86_64-toolchain \
     mingw-w64-x86_64-openblas patch python perl pkg-config pkgfile \
     rebase tar time tzcode unzip which mingw-w64-x86_64-libmicrourtis \
     --ignore mingw-w64-x86_64-gcc-ada --ignore mingw-w64-x86_64-gcc-objc \
     --ignore mingw-w64-x86_64-gdb mingw-w64-x86_64-cmake --noconfirm
     ```

   - for 32 bit systems (**FreeFEM** lower than version 4):

     ```sh
     pacman -S autoconf automake-wrapper bash bash-completion \
     bison bsdtpio bsdtar bzip2 coreutils curl dash file filesystem \
     findutils flex gawk gcc gcc-fortran gcc-libs grep gzip inetutils \
     info less indir make man-db git mingw-w64-i686-freeglut \
     mingw-w64-i686-toolchain mingw-w64-i686-gsl mingw-w64-i686-hdf5 \
     mingw-w64-i686-openblas mintty msys2-keyring msys2-launcher-git \
     msys2-runtime ncurses pacman pacman-mirrors pactoys-git patch pax-git \
     perl pkg-config pkgfile rebase sed tar tftp-hpa time tzcode unzip \
     util-linux which
     ```

6. Open a MingW64 terminal (or MingW32 for old 32 bit **FreeFEM** version) and compile the **FreeFEM** source

   ```sh
   git clone https://github.com/FreeFem/FreeFem-sources
   cd FreeFem-sources
   autoreconf -i
   ./configure --enable-generic --enable-optim \
   --enable-download --enable-maintainer-mode \
   CXXFLAGS=-mtune=generic CFLAGS=-mtune=generic
   ```

   (continues on next page)
7. If you want use HPDDM (High Performance Domain Decomposition Methods) for massively parallel computing, install PETSc/SLEPc

```
    cd 3rdparty/ff-petsc
    make petsc-slepc SUDO=sudo
    cd -
    ./reconfigure
```

8. Download the 3rd party packages and build your FreeFEM library and executable

```
    ./3rdparty/getall -a
    make
    make check
    make install
```

**Note:** The FreeFEM executable (and some other like ffmedit,...) are in C:\msys64\mingw64\bin (or C:\msys32\mingw32\bin).

### 1.2.3 Environment variables and init file

FreeFEM reads a user's init file named freefem++.pref to initialize global variables: verbosity, includepath, loadpath.

**Note:** The variable verbosity changes the level of internal printing (0: nothing unless there are syntax errors, 1: few, 10: lots, etc. ...), the default value is 2.

The included files are found in the includepath list and the load files are found in the loadpath list.

The syntax of the file is:

```
    verbosity = 5
    loadpath += "~/Library/FreeFem++/lib"
    loadpath += "~/Users/hecht/Library/FreeFem++/lib"
    includepath += "~/Library/FreeFem++/edp"
    includepath += "~/Users/hecht/Library/FreeFem++/edp"
    # This is a comment
    load += "funcTemplate"
    load += "myfunction"
    load += "MUMPS_seq"
```

The possible paths for this file are

- under Unix and MacOs

```
    /etc/freefem++.pref
    ~/.freefem++.pref
    freefem++.pref
```

- under windows

```
    %FreeFEM%/freefem++.pref
    %FreeFEM%/freefem++.pref
    freefem++.pref
```

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We can also use shell environment variables to change verbosity and the search rule before the init files.

```bash
export FF_VERBOSITY=50
export FF_INCLUDEPATH="dir;;dir2"
export FF_LOADPATH="dir;;dir3"
```

**Note:** The separator between directories must be `;;` and not `;` because `;` is used under Windows.

**Note:** To show the list of init of FreeFEM, do

```bash
export FF_VERBOSITY=100;
./FreeFem++-nw
```

### 1.2.4 Coloring Syntax FreeFem++

**Atom**

In order to get the syntax highlighting in Atom, you have to install the FreeFEM language support.

You can do it directly in Atom: Edit -> Preferences -> Install, and search for `language-freefem-offical`.

To launch scripts directly from Atom, you have to install the `atom-runner` package. Once installed, modify the Atom configuration file (Edit -> Config...) to have something like that:

```json
"x":
  ...
  runner:
    extensions:
      edp: "FreeFem++"
    scopes:
      "Freefem++": "FreeFem++"
```

Reboot Atom, and use Alt+R to run a FreeFem++ script.

**Gedit**

In order to get the syntax highlighting in Gedit, you have to download the Gedit parser and copy it in `/usr/share/gtksourceview-3.0/language-specs/`.  

1.2. Installation guide
Textmate 2, an editor under macOS

To use the coloring FreeFEM syntax with the Textmate 2 editor on Mac 10.7 or better, download from macromates.com and download the textmate freefem++ syntax here (version June 2017). To install this parser, unzip Textmate2-ff++.zip and follow the explanation given in file How_To.rtf.

Notepad++, an editor under Windows

Read and follow the instruction, FREEFEM++ COLOR SYNTAX OF WINDOWS.

Emacs editor

For emacs editor you can download ff++-mode.el.

1.3 Download

1.3.1 Latest binary packages

FreeFEM v4.6 runs under macOS, Ubuntu, and 64-bit Windows.

<table>
<thead>
<tr>
<th>Operating System</th>
<th>FreeFEM Version</th>
<th>Size</th>
<th>Date</th>
</tr>
</thead>
<tbody>
<tr>
<td>macOS 10.10.5 or higher</td>
<td>4.5</td>
<td>412 MB</td>
<td>Feb 11, 2020</td>
</tr>
<tr>
<td>Ubuntu 16.04 or higher</td>
<td>4.6</td>
<td>212 MB</td>
<td>Mar 02, 2020</td>
</tr>
<tr>
<td>64-bit Windows</td>
<td>4.6</td>
<td>185 MB</td>
<td>Mar 02, 2020</td>
</tr>
<tr>
<td>Docker image</td>
<td>4.6</td>
<td>487 MB</td>
<td>Mar 02, 2020</td>
</tr>
<tr>
<td>Source 4.6</td>
<td>4.6</td>
<td>12.4 MB</td>
<td>Mar 02, 2020</td>
</tr>
<tr>
<td>previous releases</td>
<td>–</td>
<td>–</td>
<td>–</td>
</tr>
</tbody>
</table>

The source code is available on the FreeFEM GitHub Repository.

Note: The support ended for all releases under Windows 32 bits.

1.3.2 Syntax highlighters

<table>
<thead>
<tr>
<th>Lexer type</th>
<th>Version</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>Emacs</td>
<td>0.3</td>
<td>freefem-mode.el</td>
</tr>
<tr>
<td>Textmate 2</td>
<td>1.0</td>
<td>FreeFem.tmbundle</td>
</tr>
<tr>
<td>Gedit</td>
<td>1.0</td>
<td>fpp.lang</td>
</tr>
<tr>
<td>Atom</td>
<td>0.3</td>
<td>language-freefem or via the Atom package manager</td>
</tr>
<tr>
<td>Pygments</td>
<td>1.0</td>
<td>freefem.py</td>
</tr>
<tr>
<td>Vim</td>
<td>0.1</td>
<td>edp.vim</td>
</tr>
</tbody>
</table>
1.4 History

The project has evolved from MacFem, PCfem, written in Pascal. The first C version lead to freefem 3.4; it offered mesh adaptivity on a single mesh only.

A thorough rewriting in C++ led to freefem+ (freefem+ 1.2.10 was its last release), which included interpolation over multiple meshes (functions defined on one mesh can be used on any other mesh); this software is no longer maintained but is still in use because it handles a problem description using the strong form of the PDEs. Implementing the interpolation from one unstructured mesh to another was not easy because it had to be fast and non-diffusive; for each point, one had to find the containing triangle. This is one of the basic problems of computational geometry (see [PREPARATA1985] for example). Doing it in a minimum number of operations was the challenge. Our implementation is $O(n \log n)$ and based on a quadtree. This version also grew out of hand because of the evolution of the template syntax in C++.

We have been working for a few years now on FreeFEM, entirely re-written again in C++ with a thorough usage of template and generic programming for coupled systems of unknown size at compile time. Like all versions of freefem, it has a high level user friendly input language which is not too far from the mathematical writing of the problems.

The freefem language allows for a quick specification of any partial differential system of equations. The language syntax of FreeFEM is the result of a new design which makes use of the STL [STROUSTRUP2000], templates, and bison for its implementation; more details can be found in [HECHT2002]. The outcome is a versatile software in which any new finite elements can be included in a few hours; but a recompilation is then necessary. Therefore the library of finite elements available in FreeFEM will grow with the version number and with the number of users who program more new elements. So far we have discontinuous $P_0$ elements, linear $P_1$ and quadratic $P_2$ Lagrangian elements, discontinuous $P_1$ and Raviart-Thomas elements and a few others like bubble elements.

The development of FreeFEM through more than 30 years

1987
MacFem/PCFem the old ones (O. Pironneau in Pascal) no free.

1992
FreeFem rewrite in C++ (P1,P0 one mesh) O. Pironneau, D. Bernardi, F. Hecht (mesh adaptatation, bamg), C. Prudhomme.

1996
FreeFem+ rewrite in C++ (P1,P0 more mesh) O. Pironneau, D. Bernardi, F. Hecht (algebra of function).

1998
FreeFem++ rewrite with an other finite element kernel and an new language F. Hecht, O. Pironneau, K. Ohtsuka.
1999

FreeFem 3d (S. Del Pino), a first 3d version based on fictitious domain method.

2008

FreeFem++ v3 uses a new finite element kernel multidimensionnels: 1d, 2d, 3d.

2014

FreeFem++ v3.34 parallel version

2017

FreeFem++ v3.57 parallel version

2018

FreeFem++ v4: New matrix type, Surface element, New Parallel tools ...

1.5 Citation

1.5.1 If you use FreeFEM, please cite the following reference in your work:

BibTeX

```latex
@article{MR3043640,
    AUTHOR = {Hecht, F.},
    TITLE = {New development in FreeFem++},
    JOURNAL = {J. Numer. Math.},
    FJOURNAL = {Journal of Numerical Mathematics},
    VOLUME = {20},
    YEAR = {2012},
    NUMBER = {3-4},
    PAGES = {251--265},
    ISSN = {1570-2820},
    MRCLASS = {65Y15},
    MRNUMBER = {3043640},
    URL = {https://freefem.org/}
}
```

APA


ISO690


MLA

1.6 Authors

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</tbody>
</table>
And all the dedicated GitHub contributors

1.7 Contributing

1.7.1 Bug report

Concerning the FreeFEM documentation

Open an Issue on FreeFem-doc repository.

Concerning the FreeFEM compilation or usage

Open an Issue on FreeFem-sources repository.

1.7.2 Improve content

Ask one of the contributors for Collaborator Access or make a Pull Request.
The **FreeFEM** language is *typed*, polymorphic and reentrant with *macro generation*.

Every variable must be typed and declared in a statement, that is separated from the next by a semicolon ;.

The **FreeFEM** language is a C++ idiom with something that is more akin to LaTeX.

For the specialist, one key guideline is that **FreeFEM** rarely generates an internal finite element array, this was adopted for speed and consequently **FreeFEM** could be hard to beat in terms of execution speed, except for the time lost in the interpretation of the language (which can be reduced by a systematic usage of `varf` and `matrix` instead of `problem`).

The Development Cycle: Edit–Run/Visualize–Revise

Many examples and tutorials are given there after and in the *examples section*. It is better to study them and learn by example.

If you are a beginner in the finite element method, you may also have to read a book on variational formulations.

The development cycle includes the following steps:

**Modeling:** From strong forms of PDE to weak forms, one must know the variational formulation to use **FreeFEM**; one should also have an eye on the reusability of the variational formulation so as to keep the same internal matrices; a typical example is the time dependent heat equation with an implicit time scheme: the internal matrix can be factorized only once and **FreeFEM** can be taught to do so.

**Programming:** Write the code in **FreeFEM** language using a text editor such as the one provided in your integrated environment.

**Run:** Run the code (here written in file `mycode.edp`). That can also be done in terminal mode by :

```
FreeFem++ mycode.edp
```

**Visualization:** Use the keyword `plot` directly in `mycode.edp` to display functions while **FreeFEM** is running. Use the plot-parameter `wait=1` to stop the program at each plot.

**Debugging:** A global variable `debug` (for example) can help as in `wait=true` to `wait=false`.

```
bool debug = true;
border a(t=0, 2.*pi){x=cos(t); y=sin(t); label=1;};
border b(t=0, 2.*pi){x=0.8+0.3*cos(t); y=0.3*sin(t); label=2;};
```

(continues on next page)
Changing debug to false will make the plots flow continuously. Watching the flow of graphs on the screen (while drinking coffee) can then become a pleasant experience.

Error management

Error messages are displayed in the console window. They are not always very explicit because of the template structure of the C++ code (we did our best!). Nevertheless they are displayed at the right place. For example, if you forget parenthesis as in:

```cpp
bool debug = true;

mesh Th = square(10,10;
plot(Th);
```

then you will get the following message from FreeFEM:

```
2 : mesh Th = square(10,10;
Error line number 2, in file bb.edp, before token ;
parse error
current line = 2
syntax error
current line = 2
Compile error : syntax error
line number :2, ;
error Compile error : syntax error
line number :2, ;
code = 1 mpirank: 0
```

If you use the same symbol twice as in:

```cpp
real aaa = 1;
real aaa;
```

then you will get the message:

```
2 : real aaa; The identifier aaa exists
    the existing type is <Pd>
    the new type is <Pd>
```

If you find that the program isn’t doing what you want you may also use `cout` to display in text format on the console window the value of variables, just as you would do in C++.

The following example works:
Another trick is to *comment in and out* by using `//` as in C++. For example:

```plaintext
real aaa = 1;
// real aaa;
```

### 2.1 Getting started

For a given function $f(x, y)$, find a function $u(x, y)$ satisfying:

\[
-\Delta u(x, y) = f(x, y) \quad \text{for all } (x, y) \text{ in } \Omega
\]

\[
u(x, y) = 0 \quad \text{for all } (x, y) \text{ on } \partial \Omega
\]

(2.1)

Here $\partial \Omega$ is the boundary of the bounded open set $\Omega \subset \mathbb{R}^2$ and $\Delta u = \frac{\partial^2 u}{\partial x^2} + \frac{\partial^2 u}{\partial y^2}$.

We will compute $u$ with $f(x, y) = xy$ and $\Omega$ the unit disk. The boundary $C = \partial \Omega$ is defined as:

$$C = \{(x, y) | x = \cos(t), y = \sin(t), 0 \leq t \leq 2\pi\}$$

**Note:** In FreeFEM, the domain $\Omega$ is assumed to be described by the left side of its boundary.

The following is the **FreeFEM** program which computes $u$:

```plaintext
// Define mesh boundary
border C(t=0, 2*pi){
x = cos(t);
y = sin(t);
}

// The triangulated domain Th is on the left side of its boundary
mesh Th = buildmesh(C(50));

// The finite element space defined over Th is called here Vh
fespace Vh(Th, P1);
Vh u, v; // Define u and v as piecewise-P1 continuous functions

// Define a function f
func f = x*y;

// Get the clock in second
real cpu = clock();

// Define the PDE
solve Poisson(u, v, solver=LU)
    = int2d(Th) // The bilinear part
dx(u) * dx(v)
+ dy(u) * dy(v)
;
- int2d(Th) // The right hand side
```

(continues on next page)
Fig. 2.1: Poisson’s equation

As illustrated in Fig. 2.1b, we can see the isovalue of $u$ by using FreeFEM `plot` command (see line 29 above).

**Note:** The qualifier `solver=LU` (line 18) is not required and by default a multi-frontal LU is used.

The lines containing `clock` are equally not required.

**Tip:** Note how close to the mathematics FreeFEM language is.

Lines 19 to 24 correspond to the mathematical variational equation:

\[
\int_{T_h} \left( \frac{\partial u}{\partial x} \frac{\partial v}{\partial x} + \frac{\partial u}{\partial y} \frac{\partial v}{\partial y} \right) dx dy = \int_{T_h} f v dx dy
\]

for all $v$ which are in the finite element space $V_h$ and zero on the boundary $C$.

**Tip:** Change $P1$ into $P2$ and run the program.

This first example shows how FreeFEM executes with no effort all the usual steps required by the finite element method (FEM). Let’s go through them one by one.
On the line 2:

The boundary \( \Gamma \) is described analytically by a parametric equation for \( x \) and for \( y \). When \( \Gamma = \sum_{j=0}^{J} \Gamma_j \) then each curve \( \Gamma_j \) must be specified and crossings of \( \Gamma_j \) are not allowed except at end points.

The keyword \texttt{label} can be added to define a group of boundaries for later use (boundary conditions for instance). Hence the circle could also have been described as two half circle with the same label:

\begin{verbatim}
border Gamma1(t=0, pi){x=cos(t); y=sin(t); label=C};
border Gamma2(t=pi, 2.*pi){x=cos(t); y=sin(t); label=C};
\end{verbatim}

Boundaries can be referred to either by name (\texttt{Gamma1} for example) or by label (\texttt{C} here) or even by its internal number here 1 for the first half circle and 2 for the second (more examples are in \textit{Meshing Examples}).

On the line 5

The triangulation \( \mathcal{T}_h \) of \( \Omega \) is automatically generated by \texttt{buildmesh(C(50))} using 50 points on \( C \) as in Fig. 2.1a. The domain is assumed to be on the left side of the boundary which is implicitly oriented by the parametrization. So an elliptic hole can be added by typing:

\begin{verbatim}
border C(t=2.*pi, 0){x=0.1+0.3*cos(t); y=0.5*sin(t);};
\end{verbatim}

If by mistake one had written:

\begin{verbatim}
border C(t=0, 2.*pi){x=0.1+0.3*cos(t); y=0.5*sin(t);};
\end{verbatim}

then the inside of the ellipse would be triangulated as well as the outside.

\textbf{Note:} Automatic mesh generation is based on the Delaunay-V oronoi algorithm. Refinement of the mesh are done by increasing the number of points on \( \Gamma \), for example \texttt{buildmesh(C(100))}, because inner vertices are determined by the density of points on the boundary.

Mesh adaptation can be performed also against a given function \( f \) by calling \texttt{adaptmesh(Th,f)}.

Now the name \( \mathcal{T}_h \) (\texttt{Th} in \textit{FreeFEM}) refers to the family \( \{ T_k \}_{k=1, \ldots, n_t} \) of triangles shown in Fig. 2.1a.

Traditionally \( h \) refers to the mesh size, \( n_t \) to the number of triangles in \( \mathcal{T}_h \) and \( n_v \) to the number of vertices, but it is seldom that we will have to use them explicitly.

If \( \Omega \) is not a polygonal domain, a “skin” remains between the exact domain \( \Omega \) and its approximation \( \Omega_h = \bigcup_{k=1}^{n_t} T_k \). However, we notice that all corners of \( \Gamma_h = \partial \Omega_h \) are on \( \Gamma \).

On line 8:

A finite element space is, usually, a space of polynomial functions on elements, triangles here only, with certain matching properties at edges, vertices etc. Here \texttt{fespace Vh(Th, P1)} defines \( V_h \) to be the space of continuous functions which are affine in \( x, y \) on each triangle of \( T_h \).

As it is a linear vector space of finite dimension, basis can be found. The canonical basis is made of functions, called the \textit{hat function} \( \phi_k \), which are continuous piecewise affine and are equal to 1 on one vertex and 0 on all others. A typical hat function is shown on Fig. 2.2b.

\textbf{Note:} The easiest way to define \( \phi_k \) is by making use of the \textit{barycentric coordinates} \( \lambda_i(x, y), \ i = 1, 2, 3 \) of a point \( q = (x, y) \in T \), defined by \( \sum \lambda_i = 1 \), \( \sum \lambda_i q^i = q \) where \( q^i, \ i = 1, 2, 3 \) are the 3 vertices of \( T \). Then it is easy to see that the restriction of \( \phi_k \) on \( T \) is precisely \( \lambda_k \).
Then:

$$V_h(T_h, P_1) = \left\{ w(x,y) \middle| w(x,y) = \sum_{k=1}^{M} w_k \phi_k(x,y), w_k \text{ are real numbers} \right\}$$  \hspace{1cm} (2.2)

where $M$ is the dimension of $V_h$, i.e. the number of vertices. The $w_k$ are called the degrees of freedom of $w$ and $M$ the number of degree of freedom.

It is said also that the nodes of this finite element method are the vertices.

**Setting the problem**

On line 9, \verb|Vh u, v| declares that $u$ and $v$ are approximated as above, namely:

$$u(x,y) \simeq u_h(x,y) = \sum_{k=0}^{M-1} u_k \phi_k(x,y)$$  \hspace{1cm} (2.3)

On the line 12, the right hand side $f$ is defined analytically using the keyword \verb|func|.

Line 18 to 26 define the bilinear form of equation (2.1) and its Dirichlet boundary conditions.

This variational formulation is derived by multiplying (2.1) by $v(x,y)$ and integrating the result over $\Omega$:

$$- \int_{\Omega} \Delta u \, dx \, dy = \int_{\Omega} f \, v \, dx \, dy$$

Then, by Green’s formula, the problem is converted into finding $u$ such that

$$a(u,v) - \ell(f,v) = 0 \quad \forall v \text{ satisfying } v = 0 \text{ on } \partial \Omega.$$  \hspace{1cm} (2.4)

In **FreeFEM** the Poisson problem can be declared only as in:
\begin{verbatim}
Vh u,v; problem Poisson(u,v) = ...
and solved later as in:

Poisson; //the problem is solved here

or declared and solved at the same time as in:

Vh u,v; solve Poisson(u,v) = ...

and (2.4) is written with \( dx(u) = \partial u/\partial x \), \( dy(u) = \partial u/\partial y \) and:

\[
\int_\Omega \nabla u \cdot \nabla v \, dx \rightarrow \text{int2d}(Th) \left( \ dx(u) \cdot dx(v) + dy(u) \cdot dy(v) \right) \\
\int_\Omega f \, dv \rightarrow \text{int2d}(Th) ( f \cdot v ) \text{ (Notice here, u is unused)}
\]

**Warning:** In FreeFEM bilinear terms and linear terms should not be under the same integral indeed to construct the linear systems FreeFEM finds out which integral contributes to the bilinear form by checking if both terms, the unknown (here \( u \)) and test functions (here \( v \)) are present.

**Solution and visualization**

On line 15, the current time in seconds is stored into the real-valued variable \( cpu \).

Line 18, the problem is solved. Line 29, the visualization is done as illustrated in Fig. 2.1b.

(see Plot for zoom, postscript and other commands).

Line 32, the computing time (not counting graphics) is written on the console. Notice the C++-like syntax; the user needs not study C++ for using FreeFEM, but it helps to guess what is allowed in the language.

**Access to matrices and vectors**

Internally FreeFEM will solve a linear system of the type

\[
\sum_{j=0}^{M-1} A_{ij} u_j - F_i = 0, \quad i = 0, \cdots, M-1; \quad F_i = \int_\Omega f \phi_i \, dx \, dy
\]

which is found by using (2.3) and replacing \( v \) by \( \phi_i \) in (2.4). The Dirichlet conditions are implemented by penalty, namely by setting \( A_{ii} = 10^{30} \) and \( F_i = 10^{30} \times 0 \) if \( i \) is a boundary degree of freedom.

**Note:** The number \( 10^{30} \) is called \texttt{tgv} (très grande valeur or very high value in english) and it is generally possible to change this value, see the item \texttt{freefem’solve, tgv=}.

The matrix \( A = (A_{ij}) \) is called stiffness matrix. If the user wants to access \( A \) directly he can do so by using (see section Variational form, Sparse matrix, PDE data vector for details).

\begin{verbatim}
varf a(u,v)
  = int2d(Th) ( 
    dx(u) \cdot dx(v) 
    + dy(u) \cdot dy(v) 
  )
\end{verbatim}

(continues on next page)
The vector $F$ in (2.5) can also be constructed manually:

```plaintext
varf l(unused, v) = int2d(Th) (f*v) + on(C, unused=0);  
Vh F;

F[] = l(0, Vh); //F[] is the vector associated to the function F
```

The problem can then be solved by:

```plaintext
u[] = A^{-1}*F[]; //u[] is the vector associated to the function u
```

**Note:** Here $u$ and $F$ are finite element function, and $u[]$ and $F[]$ give the array of value associated ($u[] \equiv (u_i)_{i=0,...,M-1}$ and $F[] \equiv (F_i)_{i=0,...,M-1}$).

So we have:

$$u(x, y) = \sum_{i=0}^{M-1} u[][i] \phi_i(x, y), \quad F(x, y) = \sum_{i=0}^{M-1} F[][i] \phi_i(x, y)$$

where $\phi_i, i = 0,...,M-1$ are the basis functions of $Vh$ like in equation :eq: equation3, and $M = Vh.ndof$ is the number of degree of freedom (i.e. the dimension of the space $Vh$).

The linear system (2.5) is solved by UMFPACK unless another option is mentioned specifically as in:

```plaintext
Vh u, v;
problem Poisson(u, v, solver=CG) = int2d(...
```

meaning that `Poisson` is declared only here and when it is called (by simply writing `Poisson;`) then (2.5) will be solved by the Conjugate Gradient method.

### 2.2 Classification of partial differential equations

**Summary:** It is usually not easy to determine the type of a system. Yet the approximations and algorithms suited to the problem depend on its type:

- **Finite Elements compatible (LBB conditions) for elliptic systems**
- **Finite difference on the parabolic variable and a time loop on each elliptic subsystem of parabolic systems; better stability diagrams when the schemes are implicit in time.**
- **Upwinding, Petrov-Galerkin, Characteristics-Galerkin, Discontinuous-Galerkin, Finite Volumes for hyperbolic systems plus, possibly, a time loop.**

*When the system changes type, then expect difficulties (like shock discontinuities)!

**Elliptic, parabolic and hyperbolic equations**
A partial differential equation (PDE) is a relation between a function of several variables and its derivatives.

\[
F \left( \phi(x), \frac{\partial \phi}{\partial x_1}(x), \ldots, \frac{\partial \phi}{\partial x_d}(x), \frac{\partial^2 \phi}{\partial x_1^2}(x), \ldots, \frac{\partial^m \phi}{\partial x_d^m}(x) \right) = 0, \quad \forall x \in \Omega \subset \mathbb{R}^d
\]

The range of \(x\) over which the equation is taken, here \(\Omega\), is called the domain of the PDE. The highest derivation index, here \(m\), is called the order. If \(F\) and \(\phi\) are vector valued functions, then the PDE is actually a system of PDEs.

Unless indicated otherwise, here by convention one PDE corresponds to one scalar valued \(F\) and \(\phi\). If \(F\) is linear with respect to its arguments, then the PDE is said to be linear.

The general form of a second order, linear scalar PDE is

\[
\alpha \phi + A \cdot \nabla \phi + B : \nabla(\nabla \phi) = f \quad \text{in} \quad \Omega \subset \mathbb{R}^d,
\]

where \(\frac{\partial^2 \phi}{\partial x_i \partial x_j}\) and \(A : B\) means \(\sum_{i,j=1}^d a_{ij} b_{ij}\), \(f(x), \alpha(x) \in \mathbb{R}, \ a(x) \in \mathbb{R}^d, \ B(x) \in \mathbb{R}^{d \times d}\) are the PDE coefficients. If the coefficients are independent of \(x\), the PDE is said to have constant coefficients.

To a PDE we associate a quadratic form, by replacing \(\phi\) by \(1\), \(\partial \phi / \partial x_i\) by \(z_i\) and \(\partial^2 \phi / \partial x_i \partial x_j\) by \(z_i z_j\), where \(z\) is a vector in \(\mathbb{R}^d\):

\[
\alpha + A \cdot z + z^T B z = f.
\]

If it is the equation of an ellipse (ellipsoid if \(d \geq 2\)), the PDE is said to be elliptic; if it is the equation of a parabola or a hyperbola, the PDE is said to be parabolic or hyperbolic.

If \(B \equiv 0\), the degree is no longer 2 but 1, and for reasons that will appear more clearly later, the PDE is still said to be hyperbolic.

These concepts can be generalized to systems, by studying whether or not the polynomial system \(P(z)\) associated with the PDE system has branches at infinity (ellipsoids have no branches at infinity, paraboloids have one, and hyperboloids have several).

If the PDE is not linear, it is said to be non-linear. These are said to be locally elliptic, parabolic, or hyperbolic according to the type of the linearized equation.

For example, for the non-linear equation

\[
\frac{\partial^2 \phi}{\partial t^2} - \frac{\partial \phi}{\partial x} \frac{\partial^2 \phi}{\partial x^2} = 1
\]

we have \(d = 2, \ x_1 = t, \ x_2 = x\) and its linearized form is:

\[
\frac{\partial^2 u}{\partial t^2} - \frac{\partial u}{\partial x} \frac{\partial^2 u}{\partial x^2} - \frac{\partial \phi}{\partial x} \frac{\partial^2 u}{\partial x} = 0
\]

which for the unknown \(u\) is locally elliptic if \(\frac{\partial \phi}{\partial x} < 0\) and locally hyperbolic if \(\frac{\partial \phi}{\partial x} > 0\).

Tip: Laplace’s equation is elliptic:

\[
\Delta \phi \equiv \frac{\partial^2 \phi}{\partial x_1^2} + \frac{\partial^2 \phi}{\partial x_2^2} + \cdots + \frac{\partial^2 \phi}{\partial x_d^2} = f, \quad \forall x \in \Omega \subset \mathbb{R}^d
\]

Tip: The heat equation is parabolic in \(Q = \Omega \times ]0, T[ \subset \mathbb{R}^{d+1}\):

\[
\frac{\partial \phi}{\partial t} - \mu \Delta \phi = f \quad \forall x \in \Omega \subset \mathbb{R}^d, \quad \forall t \in ]0, T[
\]
Tip: If $\mu > 0$, the wave equation is hyperbolic:
\[
\frac{\partial^2 \varphi}{\partial t^2} - \mu \Delta \varphi = f \text{ in } Q.
\]

Tip: The convection diffusion equation is parabolic if $\mu \neq 0$ and hyperbolic otherwise:
\[
\frac{\partial \varphi}{\partial t} + a \nabla \varphi - \mu \Delta \varphi = f
\]

Tip: The biharmonic equation is elliptic:
\[
\Delta (\Delta \varphi) = f \text{ in } \Omega.
\]

Boundary conditions
A relation between a function and its derivatives is not sufficient to define the function. Additional information on the boundary $\Gamma = \partial \Omega$ of $\Omega$, or on part of $\Gamma$ is necessary. Such information is called a boundary condition.

For example:
\[
\varphi(x) \text{ given, } \forall x \in \Gamma,
\]
is called a Dirichlet boundary condition. The Neumann condition is
\[
\frac{\partial \varphi}{\partial n}(x) \text{ given on } \Gamma \text{ (or } n \cdot B \nabla \varphi, \text{ given on } \Gamma \text{ for a general second order PDE)}
\]
where $n$ is the normal at $x \in \Gamma$ directed towards the exterior of $\Omega$ (by definition $\frac{\partial \varphi}{\partial n} = \nabla \varphi \cdot n$).

Another classical condition, called a Robin (or Fourier) condition is written as:
\[
\varphi(x) + \beta(x) \frac{\partial \varphi}{\partial n}(x) \text{ given on } \Gamma.
\]

Finding a set of boundary conditions that defines a unique $\varphi$ is a difficult art.

In general, an elliptic equation is well posed (i.e. $\varphi$ is unique) with one Dirichlet, Neumann or Robin condition on the whole boundary.

Thus, Laplace’s equation is well posed with a Dirichlet or Neumann condition but also with:
\[
\varphi \text{ given on } \Gamma_1, \frac{\partial \varphi}{\partial n} \text{ given on } \Gamma_2, \Gamma_1 \cup \Gamma_2 = \Gamma, \Gamma_1 \cap \Gamma_2 = \emptyset.
\]

Parabolic and hyperbolic equations rarely require boundary conditions on all of $\Gamma \times [0, T]$. For instance, the heat equation is well posed with:
\[
\varphi \text{ given at } t = 0 \text{ and Dirichlet or Neumann or mixed conditions on } \partial \Omega.
\]

Here $t$ is time so the first condition is called an initial condition. The whole set of conditions is also called Cauchy condition.

The wave equation is well posed with:
\[
\varphi \text{ and } \frac{\partial \varphi}{\partial t} \text{ given at } t = 0 \text{ and Dirichlet or Neumann or mixed conditions on } \partial \Omega.
\]
2.3 Membrane

Summary: Here we shall learn how to solve a Dirichlet and/or mixed Dirichlet Neumann problem for the Laplace operator with application to the equilibrium of a membrane under load. We shall also check the accuracy of the method and interface with other graphics packages.

An elastic membrane $\Omega$ is attached to a planar rigid support $\Gamma$, and a force $f(x)dx$ is exerted on each surface element $dx = dx_1dx_2$. The vertical membrane displacement, $\varphi(x)$, is obtained by solving Laplace’s equation:

$$-\Delta \varphi = f \text{ in } \Omega$$

As the membrane is fixed to its planar support, one has:

$$\varphi|_{\Gamma} = 0$$

If the support wasn’t planar but had an elevation $z(x_1, x_2)$ then the boundary conditions would be of non-homogeneous Dirichlet type.

$$\varphi|_{\Gamma} = z$$

If a part $\Gamma_2$ of the membrane border $\Gamma$ is not fixed to the support but is left hanging, then due to the membrane’s rigidity the angle with the normal vector $n$ is zero; thus the boundary conditions are:

$$\varphi|_{\Gamma_1} = z, \quad \frac{\partial \varphi}{\partial n}|_{\Gamma_2} = 0$$

where $\Gamma_1 = \Gamma - \Gamma_2$; recall that $\frac{\partial \varphi}{\partial n} = \nabla \varphi \cdot n$. Let us recall also that the Laplace operator $\Delta$ is defined by:

$$\Delta \varphi = \frac{\partial^2 \varphi}{\partial x_1^2} + \frac{\partial^2 \varphi}{\partial x_2^2}$$

Todo: Check references

With such “mixed boundary conditions” the problem has a unique solution (see Dautray-Lions (1988), Strang (1986) and Raviart-Thomas (1983)). The easiest proof is to notice that $\varphi$ is the state of least energy, i.e.

$$E(\varphi) = \min_{\varphi \sim z \in V} E(v), \quad \text{with} \quad E(v) = \int_{\Omega} \left( \frac{1}{2} |\nabla v|^2 - fv \right)$$

and where $V$ is the subspace of the Sobolev space $H^1(\Omega)$ of functions which have zero trace on $\Gamma_1$. Recall that ($x \in \mathbb{R}^d, d = 2$ here):

$$H^1(\Omega) = \{ u \in L^2(\Omega) : \nabla u \in (L^2(\Omega))^d \}$$

Calculus of variation shows that the minimum must satisfy, what is known as the weak form of the PDE or its variational formulation (also known here as the theorem of virtual work)

$$\int_{\Omega} \nabla \varphi \cdot \nabla w = \int_{\Omega} f w \quad \forall w \in V$$

Next an integration by parts (Green’s formula) will show that this is equivalent to the PDE when second derivatives exist.
**Warning:** Unlike the previous version Freefem+ which had both weak and strong forms, FreeFEM implements only weak formulations. It is not possible to go further in using this software if you don’t know the weak form (i.e. variational formulation) of your problem: either you read a book, or ask help from a colleague or drop the matter. Now if you want to solve a system of PDE like \( A(u,v) = 0 \), \( B(u,v) = 0 \) don’t close this manual, because in weak form it is

\[
\int_{\Omega} (A(u,v)w_1 + B(u,v)w_2) = 0 \quad \forall w_1, w_2 ...
\]

**Example**

Let an ellipse have the length of the semimajor axis \( a = 2 \), and unitary the semiminor axis. Let the surface force be \( f = 1 \). Programming this case with FreeFEM gives:

```plaintext
// Parameters
real theta = 4.*pi/3.;
real a = 2.; //The length of the semimajor axis
real b = 1.; //The length of the semiminor axis
func z = x;

// Mesh
border Gamma1(t=0., theta){x=a*cos(t); y=b*sin(t);}
border Gamma2(t=theta, 2.*pi){x=a*cos(t); y=b*sin(t);}
mesh Th = buildmesh(Gamma1(100) + Gamma2(50));

// Fespace
fespace Vh(Th, P2); //P2 conforming triangular FEM
Vh phi, w, f=1;

// Solve
solve Laplace(phi, w)
   = int2d(Th)(
      dx(phi)*dx(w)
      + dy(phi)*dy(w)
   )
   - int2d(Th)(
      f*w
   )
   + on(Gamma1, phi=z)
;

// Plot
plot(phi, wait=true, ps="membrane.eps"); //Plot phi
plot(Th, wait=true, ps="membraneTh.eps"); //Plot Th

// Save mesh
savemesh(Th,"Th.msh");
```

A triangulation is built by the keyword `buildmesh`. This keyword calls a triangulation subroutine based on the Delaunay test, which first triangulates with only the boundary points, then adds internal points by subdividing the edges. How fine the triangulation becomes is controlled by the size of the closest boundary edges.

The PDE is then discretized using the triangular second order finite element method on the triangulation; as was briefly indicated in the previous chapter, a linear system is derived from the discrete formulation whose size is the number of vertices plus the number of mid-edges in the triangulation.

The system is solved by a multi-frontal Gauss LU factorization implemented in the package UMFPACK.
The keyword *plot* will display both $T_h$ and $\phi$ (remove $T_h$ if $\phi$ only is desired) and the qualifier *fill=true* replaces the default option (colored level lines) by a full color display.

```
plot(phi, wait=true, fill=true); //Plot phi with full color display
```

Results are on Fig. 2.3a and Fig. 2.3b.

Next we would like to check the results!

One simple way is to adjust the parameters so as to know the solutions. For instance on the unit circle $a=1$, $\varphi_e = \sin(x^2 + y^2 - 1)$ solves the problem when:

$$z = 0, f = -4(\cos(x^2 + y^2 - 1) - (x^2 + y^2) \sin(x^2 + y^2 - 1))$$

except that on $\Gamma_2 \partial_n \varphi = 2$ instead of zero. So we will consider a non-homogeneous Neumann condition and solve:

$$\int_{\Omega} \nabla \varphi \cdot \nabla w = \int_{\Omega} fw + \int_{\Gamma_2} 2w \quad \forall w \in V$$

We will do that with two triangulations, compute the $L^2$ error:

$$\epsilon = \int_{\Omega} |\varphi - \varphi_e|^2$$

and print the error in both cases as well as the log of their ratio an indication of the rate of convergence.
border Gamma1(t=0., theta){x=a*cos(t); y=b*sin(t);} 
border Gamma2(t=theta, 2.*pi){x=a*cos(t); y=b*sin(t);} 

// Error loop 
real[int] L2error(2); // an array of two values 
for(int n = 0; n < 2; n++) {
    // Mesh 
    mesh Th = buildmesh(Gamma1(20*(n+1)) + Gamma2(10*(n+1)));
    // Fespace 
    fespace Vh(Th, P2);
    Vh phi, w;
    // Solve 
    solve Laplace(phi, w) 
        = int2d(Th)( 
            dx(phi)*dx(w) 
            + dy(phi)*dy(w)
        ) 
        - int2d(Th) ( 
            f*w
        ) 
        - int1d(Th, Gamma2) ( 
            2*w
        ) 
        + on(Gamma1, phi=0) ;
    // Plot 
    plot(Th, phi, wait=true, ps="membrane.eps");
    // Error 
    L2error[n] = sqrt(int2d(Th)((phi-phiexact)^2));
}

// Display loop 
for(int n = 0; n < 2; n++)
    cout << "L2error " << n << " = " << L2error[n] << endl;
// Convergence rate 
cout << "convergence rate = " << log(L2error[0]/L2error[1])/log(2.) << endl;

The output is:

L2error 0 = 0.00462991 
L2error 1 = 0.00117128 
convergence rate = 1.9829 
times: compile 0.02s, execution 6.94s

We find a rate of 1.98, which is not close enough to the 3 predicted by the theory.

The Geometry is always a polygon so we lose one order due to the geometry approximation in $O(h^2)$.

Now if you are not satisfied with the .eps plot generated by FreeFEM and you want to use other graphic facilities, then you must store the solution in a file very much like in C++. It will be useless if you don’t save the triangulation as well, consequently you must do
For the triangulation the name is important: **the extension determines the format.**

For the triangulation the name is important: **the extension determines the format.**

![3D graph](image)

**Fig. 2.4:** The 3D version drawn by gnuplot from a file generated by **FreeFEM**

Still that may not take you where you want. Here is an interface with gnuplot (see: web site link) to produce the **Fig. 2.4.**

```
//to build a gnuplot data file
{
    ofstream ff("graph.txt");
    for (int i = 0; i < Th.nt; i++)
    {
        for (int j = 0; j < 3; j++)
            ff << Th[i][j].x << " " << Th[i][j].y << " " << phi[Vh(i,j)] << endl;
    }
}
```

We use the finite element numbering, where $Wh(i,j)$ is the global index of $j^{Th}$ degrees of freedom of triangle number $i$.

Then open gnuplot and do:

```
set palette rgbformulae 30,31,32
splot "graph.txt" w l pal
```

This works with $P2$ and $P1$, but not with $P1nc$ because the 3 first degrees of freedom of $P2$ or $P2$ are on vertices and not with $P1nc$.

### 2.3. Membrane
2.4 Heat Exchanger

Summary: Here we shall learn more about geometry input and triangulation files, as well as read and write operations.

The problem Let \( \{C_i\}_{1,2} \) be 2 thermal conductors within an enclosure \( C_0 \) (see Fig. 2.5).

The first one is held at a constant temperature \( u_1 \) the other one has a given thermal conductivity \( \kappa_2 \) 3 times larger than the one of \( C_0 \).

We assume that the border of enclosure \( C_0 \) is held at temperature \( 20^\circ C \) and that we have waited long enough for thermal equilibrium.

In order to know \( u(x) \) at any point \( x \) of the domain \( \Omega \), we must solve:

\[
\nabla \cdot (\kappa \nabla u) = 0 \quad \text{in} \quad \Omega, \quad u|_\Gamma = g
\]

where \( \Omega \) is the interior of \( C_0 \) minus the conductor \( C_1 \) and \( \Gamma \) is the boundary of \( \Omega \), that is \( C_0 \cup C_1 \).

Here \( g \) is any function of \( x \) equal to \( u_i \) on \( C_i \).

The second equation is a reduced form for:

\[
u = u_i \quad \text{on} \quad C_i, \quad i = 0, 1.
\]

The variational formulation for this problem is in the subspace \( H^1_0(\Omega) \subset H^1(\Omega) \) of functions which have zero traces on \( \Gamma \).

\[
u - g \in H^1_0(\Omega) : \int_\Omega \nabla u \nabla v = 0 \quad \forall v \in H^1_0(\Omega)
\]

Let us assume that \( C_0 \) is a circle of radius 5 centered at the origin, \( C_i \) are rectangles, \( C_1 \) being at the constant temperature \( u_1 = 60^\circ C \) (so we can only consider its boundary).
// Parameters
int C1=99;
int C2=98; //could be anything such that !=0 and C1!=C2

// Mesh
border C0(t=0., 2.*pi){x=5.*cos(t); y=5.*sin(t);}
border C11(t=0., 1.){x=1.+t; y=3.; label=C1;}
border C12(t=0., 1.){x=2.-t; y=3.; label=C1;}
border C13(t=0., 1.){x=2.-t; y=-3.; label=C1;}
border C14(t=0., 1.){x=1.; y=-3.+6.*t; label=C1;}
border C21(t=0., 1.){x=-2.+t; y=3.; label=C2;}
border C22(t=0., 1.){x=-1.; y=3.-6.*t; label=C2;}
border C23(t=0., 1.){x=-1.-t; y=-3.; label=C2;}
border C24(t=0., 1.){x=-2.; y=-3.+6.*t; label=C2;}
plot( C0(50) //to see the border of the domain
+ C11(5)+C12(20)+C13(5)+C14(20)
+ C21(-5)+C22(-20)+C23(-5)+C24(-20),
wait=true, ps="heatexb.eps");

mesh Th=buildmesh(C0(50)
+ C11(5)+C12(20)+C13(5)+C14(20)
+ C21(-5)+C22(-20)+C23(-5)+C24(-20));
plot(Th,wait=1);

// Fespace
fespace Vh(Th, P1);
Vh u, v;
Vh kappa=1 + 2*(x<-1)*(x>-2)*(y<3)*(y>-3);

// Solve
solve a(u, v)
= int2d(Th)(
kappa*
  dx(u)*dx(v)
  + dy(u)*dy(v)
)
+on(C0, u=20)
+on(C1, u=60)
;

// Plot
plot(u, wait=true, value=true, fill=true, ps="HeatExchanger.eps");

Note the following:

- C0 is oriented counterclockwise by t, while C1 is oriented clockwise and C2 is oriented counterclockwise. This is why C1 is viewed as a hole by buildmesh.

- C1 and C2 are built by joining pieces of straight lines. To group them in the same logical unit to input the boundary conditions in a readable way we assigned a label on the boundaries. As said earlier, borders have an internal number corresponding to their order in the program (check it by adding a cout << C22; above). This is essential to understand how a mesh can be output to a file and re-read (see below).

- As usual the mesh density is controlled by the number of vertices assigned to each boundary. It is not possible
Fig. 2.6: Heat exchanger

to change the (uniform) distribution of vertices but a piece of boundary can always be cut in two or more parts, for instance C12 could be replaced by C121+C122:

```c
  // border C12(t=0.,1.){x=2.; y=3.-6.*t; label=C1;}
  border C121(t=0.,0.7){x=2.; y=3.-6.*t; label=C1;}
  border C122(t=0.7,1.){x=2.; y=3.-6.*t; label=C1;}
  ...
  buildmesh(.../+/ C12(20) */ + C121(12) + C122(8) + ...);
```

Tip: Exercise:

Use the symmetry of the problem with respect to the x axes.

Triangulate only one half of the domain, and set homogeneous Neumann conditions on the horizontal axis.

Writing and reading triangulation files Suppose that at the end of the previous program we added the line

```c
  savemesh(Th, "condensor.msh");
```

and then later on we write a similar program but we wish to read the mesh from that file. Then this is how the condenser should be computed:

```c
  // Mesh
  mesh Sh = readmesh("condensor.msh");
  // Fespace
  fespace Wh(Sh, P1);
  Wh us, vs;
  // Solve
  solve b(us, vs)
    = int2d(Sh)(
```

(continues on next page)
Note that the names of the boundaries are lost but either their internal number (in the case of C0) or their label number (for C1 and C2) are kept.

### 2.5 Acoustics

**Summary:** Here we go to grip with ill posed problems and eigenvalue problems

Pressure variations in air at rest are governed by the wave equation:

\[
\frac{\partial^2 u}{\partial t^2} - c^2 \Delta u = 0
\]

When the solution wave is monochromatic (and that depends on the boundary and initial conditions), \( u \) is of the form

\[
u(x, t) = Re(v(x)e^{ikt})
\]

where \( v \) is a solution of Helmholtz’s equation:

\[
k^2 v + c^2 \Delta v = 0 \quad \text{in } \Omega
\]

\[
\frac{\partial v}{\partial n} |_{\Gamma} = g
\]

where \( g \) is the source.

Note the “+” sign in front of the Laplace operator and that \( k > 0 \) is real. This sign may make the problem ill posed for some values of \( \frac{c}{k} \), a phenomenon called “resonance”.

At resonance there are non-zero solutions even when \( g = 0 \). So the following program may or may not work:
// Solve
solve sound(u, v)
  = int2d(Th) (
    u*v * kc2
    - dx(u)*dx(v)
    - dy(u)*dy(v)
  )
  - int1d(Th, a4) ( g * v
  );

// Plot
plot(u, wait=1, ps="Sound.eps");

Results are on Fig. 2.7a. But when $k^2$ is an eigenvalue of the problem, then the solution is not unique:

- if $u_e \neq 0$ is an eigen state, then for any given solution $u + u_e$ is another solution.

To find all the $u_e$ one can do the following:

// Parameters
real sigma = 20; //value of the shift

// Problem
// OP = A - sigma B ; // The shifted matrix
varf op(u1, u2)
  = int2d(Th) (
    dx(u1)*dx(u2)
    + dy(u1)*dy(u2)
    - sigma * u1*u2
  )
  ;

varf b([u1], [u2])
  = int2d(Th) ( u1*u2
  ) ; // No Boundary condition see note \ref{note BC EV}

matrix OP = op(Vh, Vh, solver=Crout, factorize=1);
matrix B = b(Vh, Vh, solver=CG, eps=1e-20);

// Eigen values
int nev=2; // Number of requested eigenvalues near sigma
real[1int] ev(nev); // To store the nev eigenvalue
Vh[1int] eV(nev); // To store the nev eigenvector

int k=EigenValue(OP, B, sym=true, sigma=sigma, value=ev, vector=eV,
  tol=1e-10, maxit=0, ncv=0);

cout << ev(0) << " 2 eigen values " << ev(1) << endl;
v = eV[0];
plot(v, wait=true, ps="eigen.eps");
2.6 Thermal Conduction

**Summary**: Here we shall learn how to deal with a time dependent parabolic problem. We shall also show how to treat an axisymmetric problem and show also how to deal with a nonlinear problem.

**How air cools a plate**

We seek the temperature distribution in a plate $(0, L_x) \times (0, L_y) \times (0, L_z)$ of rectangular cross section $\Omega = (0, 6) \times (0, 1)$; the plate is surrounded by air at temperature $u_e$ and initially at temperature $u = u_0 + \frac{1}{2} u_1$. In the plane perpendicular to the plate at $z = L_z/2$, the temperature varies little with the coordinate $z$; as a first approximation the problem is 2D.

We must solve the temperature equation in $\Omega$ in a time interval $(0,T)$.

$$
\begin{align*}
\partial_t u - \nabla \cdot (\kappa \nabla u) &= 0 & \text{in } (0,T) \\
u(x, y, 0) &= u_0 + xu_1 & \text{on } \Gamma \\
\kappa \frac{\partial u}{\partial n} + \alpha (u - u_e) &= 0 & \text{on } \Gamma \times (0,T)
\end{align*}
$$

Here the diffusion $\kappa$ will take two values, one below the middle horizontal line and ten times less above, so as to simulate a thermostat.

The term $\alpha (u - u_e)$ accounts for the loss of temperature by convection in air. Mathematically this boundary condition is of Fourier (or Robin, or mixed) type.

The variational formulation is in $L^2(0,T; H^1(\Omega))$; in loose terms and after applying an implicit Euler finite difference approximation in time; we shall seek $u^n(x, y)$ satisfying for all $w \in H^1(\Omega)$:

$$
\int_\Omega \left( \frac{u^n - u^{n-1}}{\delta t} w + \kappa \nabla u^n \nabla w \right) + \int_{\Gamma} \alpha (u^n - u_n e) w = 0
$$

```plaintext
// Parameters
func u0 = 10. + 90.*x/6.;
func k = 1.8*(y<0.5) + 0.2;
real ue = 25.;
real alpha=0.25;
real T=5.;
real dt=0.1 ;

// Mesh
mesh Th = square(30, 5, [6.*x,y]);
```

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(continued from previous page)

```cpp
// Fespace
fespace Vh(Th, P1);
Vh u=u0, v, uold;

// Problem
problem thermic(u, v)
  = int2d(Th)(
    u*v/dt
    + k*(
      dx(u) * dx(v)
      + dy(u) * dy(v)
    )
  )
  + int1d(Th, 1, 3){
    alpha*u*v
  }
  - int1d(Th, 1, 3){
    alpha*ue*v
  }
  - int2d(Th){
    uold*v/dt
  }
  + on(2, 4, u=u0);

// Time iterations
ofstream ff("thermic.dat");
for(real t = 0; t < T; t += dt){
  uold = u; //equivalent to u^{n-1} = u^n
  thermic; //here the thermic problem is solved
  ff << u(3., 0.5) << endl;
  plot(u);
}
```

**Note:** We must separate by hand the bilinear part from the linear one.

**Note:** The way we store the temperature at point (3, 0.5) for all times in file thermic.dat. Should a one dimensional plot be required (you can use gnuplot tools), the same procedure can be used. For instance to print $x \mapsto \frac{\partial u}{\partial y}(x, 0.9)$ one would do:

```cpp
for(int i = 0; i < 20; i++)
  cout << dy(u)(6.0*i/20.0,0.9) << endl;
```

Results are shown on Fig. 2.8a and Fig. 2.8b.

### 2.6.1 Axisymmetry: 3D Rod with circular section

Let us now deal with a cylindrical rod instead of a flat plate. For simplicity we take $\kappa = 1$.

In cylindrical coordinates, the Laplace operator becomes ($r$ is the distance to the axis, $z$ is the distance along the axis,
(a) Temperature at $t = 4.9$.

**Fig. 2.8:** Thermal conduction

$\theta$ polar angle in a fixed plane perpendicular to the axis):

$$\Delta u = \frac{1}{r} \partial_r (r \partial_r u) + \frac{1}{r^2} \partial_{\theta \theta} u + \partial_{zz}^2.$$ Symmetry implies that we loose the dependence with respect to $\theta$; so the domain $\Omega$ is again a rectangle $[0, R] \times [0, \pi]$. We take the convention of numbering of the edges as in `square()` (1 for the bottom horizontal ...); the problem is now:

$$r \partial_t u - \partial_r (r \partial_r u) - \partial_z (r \partial_z u) = 0 \quad \text{in } \Omega$$

$$u|_{\Gamma_4} = u_0 + \frac{z}{L_z} (u_1 - u)$$

$$u|_{\Gamma_2} = u_1$$

$$\alpha (u - u_e) + \frac{\partial u}{\partial n}|_{\Gamma_1 \cup \Gamma_3} = 0$$

Note that the PDE has been multiplied by $r$.

After discretization in time with an implicit scheme, with time steps $dt$, in the FreeFEM syntax $r$ becomes $x$ and $z$ becomes $y$ and the problem is:

```cpp
problem thermaxi(u, v) = int2d(Th)(
(u*v/dt + dx(u)*dx(v) + dy(u)*dy(v)) * x
 ) + int1d(Th, 3) (
 alpha * x * u * v
 ) - int1d(Th, 3) (
 alpha * x * u_e * v
 ) - int2d(Th) (
 uold * v * x / dt
 ) + on(2, 4, u = u0);
```

2.6. Thermal Conduction
Note: The bilinear form degenerates at $x = 0$. Still one can prove existence and uniqueness for $u$ and because of this degeneracy no boundary conditions need to be imposed on $\Gamma_1$.

### 2.6.2 A Nonlinear Problem: Radiation

Heat loss through radiation is a loss proportional to the absolute temperature to the fourth power (Stefan’s Law). This adds to the loss by convection and gives the following boundary condition:

$$\kappa \frac{\partial u}{\partial n} + \alpha(u - u_e) + c[(u + 273)^4 - (u_e + 273)^4] = 0$$

The problem is nonlinear, and must be solved iteratively with fixed-point iteration where $m$ denotes the iteration index, a semi-linearization of the radiation condition gives

$$\frac{\partial u^{m+1}}{\partial n} + \alpha(u^{m+1} - u_e) + c(u^{m+1} + u_e + 546)((u^{m+1} + 273)^2 + (u_e + 273)^2) = 0,$$

because we have the identity $a^4 - b^4 = (a - b)(a + b)(a^2 + b^2)$.

The iterative process will work with $v = u - u_e$.

```plaintext
...  // Mesh
fespace Vh(Th, P1);
Vh vold, w, v=u0-ue, b,vp;

// Problem
problem thermradia(v, w)
  = int2d(Th)(
    v*w/dt
    + k*(dx(v) * dx(w) + dy(v) * dy(w))
  )
  + int1d(Th, 1, 3)(
    b*v*w
  )
  - int2d(Th)(
    vold*w/dt
  )
  - on(2, 4, v=u0-ue);

verbosity=0; // to remove spurious FREEfem print
for (real t=0; t<T; t+=dt) {
  vold[] = v[]; // just copy DoF's, faster than interpolation pv=v;
  for (int m = 0; m < 5; m++) {
    vp[] = v[]; // save previous state of commute error
    b = alpha + rad * (v + 2*uek) * ((v+uek)^2 + uek^2);
    thermradia;
    vp[] = v[];
    real err = vp[].linfty; // error value
    cout << " time " << t << " iter " << m << " err = " << vp[].linfty << endl;
    if (err < 1e-5) break; // if error is enough small break fixed-point loop
  }
  v[] += ue; // add a constant to all DoF's of v
}
plot(v);
```
2.7 Irrotational Fan Blade Flow and Thermal effects

**Summary:** Here we will learn how to deal with a multi-physics system of PDEs on a complex geometry, with multiple meshes within one problem. We also learn how to manipulate the region indicator and see how smooth is the projection operator from one mesh to another.

**Incompressible flow**

Without viscosity and vorticity incompressible flows have a velocity given by:

\[
\mathbf{u} = \left( \frac{\partial \psi}{\partial x}, \frac{\partial \psi}{\partial y} \right),
\]

where \( \psi \) is solution of \( \Delta \psi = 0 \)

This equation expresses both incompressibility (\( \nabla \cdot \mathbf{u} = 0 \)) and absence of vortex (\( \nabla \times \mathbf{u} = 0 \)).

As the fluid slips along the walls, normal velocity is zero, which means that \( \psi \) satisfies:

\( \psi \) constant on the walls.

One can also prescribe the normal velocity at an artificial boundary, and this translates into non constant Dirichlet data for \( \psi \).

**Airfoil**

Let us consider a wing profile \( S \) in a uniform flow. Infinity will be represented by a large circle \( C \) where the flow is assumed to be of uniform velocity; one way to model this problem is to write:

\[
\Delta \psi = 0 \text{ in } \Omega, \quad \psi|_S = -I, \quad \psi|_C = u_\infty x^\perp
\]

where \( \partial \Omega = C \cup S \) and \( I \) is the lift force.

**The NACA0012 Airfoil**

An equation for the upper surface of a NACA0012 (this is a classical wing profile in aerodynamics) is:

\[
y = 0.17735\sqrt{x} - 0.075597x - 0.212836x^2 + 0.17363x^3 - 0.06254x^4.
\]

(continues on next page)
(a) Zoom around the NACA0012 airfoil showing the streamlines (curve \( \psi = \text{constant} \)). To obtain such a plot use the interactive graphic command: “+” and p.

(b) Temperature distribution at time T=25 (now the maximum is at 90 instead of 120).

Fig. 2.9: The NACA0012 Airfoil

```
20 = int2d(Th) ( // scalar product
    grad(psi)'*grad(w)
    + on(C, psi = [uinfty1,uinfty2]'*[y,-x])
    + on(S, psi=-lift) // to get a correct value
); // Added to have a fine mesh at trail
21 plot(psi, wait=1);
```

A zoom of the streamlines are shown on Fig. 2.9a.

### 2.7.1 Heat Convection around the airfoil

Now let us assume that the airfoil is hot and that air is there to cool it. Much like in the previous section the heat equation for the temperature \( v \) is

\[
\frac{\partial v}{\partial t} - \nabla \cdot (\kappa \nabla v) + u \cdot \nabla v = 0, \quad v(t=0) = v_0, \quad \frac{\partial v}{\partial n}|_C = 0
\]

But now the domain is outside AND inside \( S \) and \( \kappa \) takes a different value in air and in steel. Furthermore there is convection of heat by the flow, hence the term \( u \cdot \nabla v \) above.

Consider the following, to be plugged at the end of the previous program:

```
// Parameters
real dt=0.05;
real nbT=50;
// Mesh
border D(t=0., 2.){x=1.+cos(theta)*t; y=+sin(theta)*t;} // Added to have a fine mesh
mesh Sh = buildmesh(C(25) + Splus(-90) + Sminus(-90) + D(200));
int steel=Sh(0.5,0).region, air=Sh(-1,0).region;
// Fespace
fespace Vh(Sh, P2);
```

(continues on next page)
fespace Wh(Sh, P1);
Wh v, vv;

fespace W0(Sh, P0);
W0 k=0.01*(region==air)+0.1*(region==steel);
W0 u1=dy(psi)*(region==air), u2=-dx(psi)*(region==air);
Wh vold = 120*(region==steel);

// Problem
int i;
problem thermic(v, vv, init=i, solver=LU) =
  int2d(Sh)(v*vv/dt + k*grad(v)'*grad(vv) + 10*(u1*dx(v)+u2*dy(v))*vv)
  - int2d(Sh)(vold*vv/dt);

for(i = 0; i < nbT; i++){
  v = vold;
  thermic;
  plot(v);
}

Note: How steel and air are identified by the mesh parameter region which is defined when buildmesh is called and takes an integer value corresponding to each connected component of $\Omega$;

How the convection terms are added without upwinding. Upwinding is necessary when the Pecley number $|u|L/\kappa$ is large (here is a typical length scale), The factor 10 in front of the convection terms is a quick way of multiplying the velocity by 10 (else it is too slow to see something).

The solver is Gauss’ LU factorization and when init $\neq 0$ the LU decomposition is reused so it is much faster after the first iteration.

2.8 Pure Convection : The Rotating Hill

Summary: Here we will present two methods for upwinding for the simplest convection problem. We will learn about Characteristics-Galerkin and Discontinuous-Galerkin Finite Element Methods.

Let $\Omega$ be the unit disk centered at $(0,0)$; consider the rotation vector field

$$u = [u_1, u_2], \quad u_1 = y, \quad u_2 = -x$$

Pure convection by $u$ is

$$\partial_t c + u \cdot \nabla c = 0 \quad \text{in } \Omega \times (0, T)$$
$$c(t=0) = c^0 \quad \text{in } \Omega.$$

The exact solution $c(x_t, t)$ at time $t$ en point $x_t$ is given by:

$$c(x_t, t) = c^0(x, 0)$$
where $x_t$ is the particle path in the flow starting at point $x$ at time $0$. So $x_t$ are solutions of

$$\dot{x}_t = u(x_t), \quad x_{t=0} = x, \quad \text{where} \quad \dot{x}_t = \frac{d(x_t)}{dt}$$

The ODE are reversible and we want the solution at point $x$ at time $t$ (not at point $x_t$) the initial point is $x_{-\tau}$, and we have

$$c(x,t) = c(0)(x_{-\tau},0)$$

The game consists in solving the equation until $T = 2\pi$, that is for a full revolution and to compare the final solution with the initial one; they should be equal.

### 2.8.1 Solution by a Characteristics-Galerkin Method

In FreeFEM there is an operator called `convect([u1,u2], dt, c)` which compute $c \circ X$ with $X$ is the convect field defined by $X(x) = x_{dt}$ where $x_{\tau}$ is particle path in the steady state velocity field $u = [u1,u2]$ starting at point $x$ at time $\tau = 0$, so $x_{\tau}$ is solution of the following ODE:

$$\dot{x}_{\tau} = u(x_{\tau}), x_{\tau=0} = x.$$  

When $u$ is piecewise constant; this is possible because $x_{\tau}$ is then a polygonal curve which can be computed exactly and the solution exists always when $u$ is divergence free; convect returns $c(x_{dt}) = C \circ X$.

```cpp
// Parameters
real dt = 0.17;

// Mesh
border C(t=0., 2.*pi) {x=cos(t); y=sin(t);};
mesh Th = buildmesh(C(100));

// Fespace
fespace Uh(Th, P1);
Uh cold, c = exp(-10*((x-0.3)^2 + (y-0.3)^2));
Uh u1 = y, u2 = -x;

// Time loop
real t = 0;
for (int m = 0; m < 2.*pi/dt; m++) {
    t += dt;
    cold = c;
    c = convect([u1, u2], -dt, cold);
    plot(c, cmm=" t="+t +", min="+c[].min", max="+c[].max");
}
```

**Note:** 3D plots can be done by adding the qualifier `dim=3` to the `plot` instruction.

The method is very powerful but has two limitations:

- it is not conservative
- it may diverge in rare cases when $|u|$ is too small due to quadrature error.
2.8.2 Solution by Discontinuous-Galerkin FEM

Discontinuous Galerkin methods take advantage of the discontinuities of $c$ at the edges to build upwinding. There are may formulations possible. We shall implement here the so-called dual-$P^1_{DC}$ formulation (see [ERN2006]):

$$
\int_{\Omega} \left( \frac{c^{n+1} - c^n}{\delta t} + u \cdot \nabla c \right) w + \int_E (c|n \cdot u| - \frac{1}{2} |n \cdot u| c) w = \int_{E_T^\Gamma} |n \cdot u| c w \quad \forall w
$$

where $E$ is the set of inner edges and $E_T^\Gamma$ is the set of boundary edges where $u \cdot n < 0$ (in our case there is no such edges). Finally $[c]$ is the jump of $c$ across an edge with the convention that $c^+$ refers to the value on the right of the oriented edge.

```cpp
// Parameters
real al=0.5;
real dt = 0.05;

// Mesh
border C(t=0., 2.*pi) {x=cos(t); y=sin(t);};
mesh Th = buildmesh(C(100));

// Fespace
defspace Vh(Th,P1dc);
Vh w, ccold, v1 = y, v2 = -x, cc = exp(-10*(x-0.3)^2 + (y-0.3)^2));

// Macro
macro n() (N.x*v1 + N.y*v2) // Macro without parameter

// Problem
problem Adual(cc, w) = int2d(Th)(
(c/dt+(v1*dx(cc)+v2*dy(cc)))*w
) + intalledges(Th)(
(1-nTonEdge)*w*(al*abs(n)-n/2)*jump(cc)
) - int2d(Th)(
ccold*w/dt
);

// Time iterations
for (real t = 0.; t < 2.*pi; t += dt){
  ccold = cc;
  Adual;
  plot(cc, fill=1, cmm="t="+t+", min="+cc[].min+", max="+ cc[].max);
}

// Plot
real [int] viso = [-0.2, -0.1, 0., 0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1., 1. -1];
plot(cc, wait=1, fill=1, ps="ConvectCG.eps", viso=viso);
plot(cc, wait=1, fill=1, ps="ConvectDG.eps", viso=viso);
```

Note: New keywords: intalledges to integrate on all edges of all triangles

$$
\int_{\partial T} \sum_{T \in \mathcal{Th}} \int_{\partial T}
$$
The rotating hill after one revolution with Characteristics-Galerkin

The rotating hill after one revolution with Discontinuous $P_1$ Galerkin

Fig. 2.10: Rotating hill

(so all internal edges are see two times), nTonEdge which is one if the triangle has a boundary edge and two otherwise, jump to implement $[c]$.

Results of both methods are shown on Fig. 2.10a nad Fig. 2.10b with identical levels for the level line; this is done with the plot-modifier viso.

Notice also the macro where the parameter $u$ is not used (but the syntax needs one) and which ends with a //; it simply replaces the name $n$ by $(N.x*v1+N.y*v2)$. As easily guessed $N.x, N.y$ is the normal to the edge.

Now if you think that DG is too slow try this:

```plaintext
// Mesh
border C(t=0., 2.*pi) {x=cos(t); y=sin(t);};
mesh Th = buildmesh(C(100));

fespace Vh(Th, P1); //P1,P2,P0,P1dc,P2dc, uncond stable
Vh vh,vo,u1 = y, u2 = -x, v = exp(-10*((x-0.3)^2+(y-0.3)^2));
real dt = 0.03, t=0, tmax=2*pi, al=0.5, alp=200;

problem A(v,vh) = int2d(Th)(v*vh/dt-v*(u1*dx(vh)+u2*dy(vh)))
  + intalledges(Th) (mean(v)*((N.x*u1+N.y*u2)
    + alp*jump(v)*abs((N.x*u1+N.y*u2))
  + int1d(Th, 1)((N.x*u1+N.y*u2)>0)*((N.x*u1+N.y*u2)*v*vh)
  - int2d(Th) (vo*vh/dt));

varf A dual(v,vh) = int2d(Th) ((v/dt+(u1*dx(v)+u2*dy(v)))*vh)
  + intalledges(Th) ((1-nTonEdge)*vh*(al*abs((N.x*u1+N.y*u2)
    -(N.x*u1+N.y*u2)/2)*jump(v));

varf rhs(vo,vh) = int2d(Th) (vo*vh/dt);
```

(continues on next page)
2.9 The System of elasticity

Elasticity

Solid objects deform under the action of applied forces:

a point in the solid, originally at \((x, y, z)\) will come to \((X, Y, Z)\) after some time; the vector \(u = (u_1, u_2, u_3) = (X - x, Y - y, Z - z)\) is called the displacement. When the displacement is small and the solid is elastic, Hooke’s law gives a relationship between the stress tensor \(\sigma(u) = (\sigma_{ij}(u))\) and the strain tensor \(\epsilon(u) = (\epsilon_{ij}(u))\)

\[
\sigma_{ij}(u) = \lambda \delta_{ij} \nabla \cdot u + 2\mu \epsilon_{ij}(u),
\]

where the Kronecker symbol \(\delta_{ij} = 1\) if \(i = j\), 0 otherwise, with

\[
\epsilon_{ij}(u) = \frac{1}{2} \left( \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right),
\]

and where \(\lambda, \mu\) are two constants that describe the mechanical properties of the solid, and are themselves related to the better known constants \(E\), Young’s modulus, and \(\nu\), Poisson’s ratio:

\[
\mu = \frac{E}{2(1 + \nu)}, \quad \lambda = \frac{E\nu}{(1 + \nu)(1 - 2\nu)}.
\]

Lamé’s system

Let us consider a beam with axis \(Oz\) and with perpendicular section \(\Omega\). The components along \(x\) and \(y\) of the strain \(u(x)\) in a section \(\Omega\) subject to forces \(f\) perpendicular to the axis are governed by:

\[-\mu \Delta u - (\mu + \lambda) \nabla (\nabla \cdot u) = f\text{ in } \Omega,
\]

where \(\lambda, \mu\) are the Lamé coefficients introduced above.

Remark, we do not use this equation because the associated variational form does not give the right boundary condition, we simply use:

\[-div(\sigma) = f\text{ in } \Omega,
\]

where the corresponding variational form is:

\[
\int_\Omega \sigma(u) : \epsilon(v) \, dx - \int_\Omega vf \, dx = 0;
\]

where \(:\) denotes the tensor scalar product, i.e. \(a : b = \sum_{i,j} a_{ij}b_{ij}\).
So the variational form can be written as:

\[ \int_{\Omega} \lambda \nabla u \cdot \nabla v + 2\mu \epsilon(u) : \epsilon(v) \, dx - \int_{\Omega} vf \, dx = 0; \]

**Tip:** Consider an elastic plate with the undeformed rectangle shape \([0, 20] \times [-1, 1]\).

The body force is the gravity force \(f\) and the boundary force \(g\) is zero on lower, upper and right sides. The left vertical side of the beam is fixed. The boundary conditions are:

\[
\sigma \cdot n = g = 0 \quad \text{on } \Gamma_1, \Gamma_4, \Gamma_3, \quad u = 0 \quad \text{on } \Gamma_2
\]

Here \(u = (u, v)\) has two components.

The above two equations are strongly coupled by their mixed derivatives, and thus any iterative solution on each of the components is risky. One should rather use FreeFEM's system approach and write:

```cpp
// Parameters
real E = 21e5;
real nu = 0.28;
real f = -1;

// Mesh
mesh Th = square(10, 10, [20*x, 2*y-1]);

// Fespace
fespace Vh(Th, P2);
Vh u, v;
Vh uu, vv;

// Macro
real sqrt2=sqrt(2.);
macro epsilon(u1,u2) [dx(u1),dy(u2),(dy(u1)+dx(u2))/sqrt2] //
    The sqrt2 is because we want: epsilon(u1,u2)'* epsilon(v1,v2) = epsilon(u):epsilon(v)
macro div(u,v) ( dx(u)+dy(v) ) //

// Problem
real mu= E/(2*(1+nu));
real lambda = E*nu/(1+nu)*((1-2*nu));
solve lame([u, v], [uu, vv])
  = int2d(Th) ( 
    lambda * div(u, v) * div(uu, vv)
    + 2.*mu * ( epsilon(u,v)' * epsilon(uu,vv) )
  )
  - int2d(Th) ( 
    f*vv
  )
+ on(4, u=0, v=0);

// Plot
real coef=100;
```

(continues on next page)
2.10 The System of Stokes for Fluids

In the case of a flow invariant with respect to the third coordinate (two-dimensional flow), flows at low Reynolds number (for instance micro-organisms) satisfy,

\[-\Delta \mathbf{u} + \nabla p = 0 \]
\[\nabla \cdot \mathbf{u} = 0\]

where \(\mathbf{u} = (u_1, u_2)\) is the fluid velocity and \(p\) its pressure.

The driven cavity is a standard test. It is a box full of liquid with its lid moving horizontally at speed one. The pressure and the velocity must be discretized in compatible finite element spaces for the LBB conditions to be satisfied:

\[
\sup_{p \in P_h} \frac{(\mathbf{u}, \nabla p)}{|p|} \geq \beta |\mathbf{u}| \quad \forall \mathbf{u} \in U_h
\]
// Parameters
int nn = 30;

// Mesh
def Th = square(nn, nn);

// Fespace
fespace Uh(Th, P1);
Uh u, v;
Uh uu, vv;

fespace Ph(Th, P1);
Ph p, pp;

// Problem
solve stokes ([u, v, p], [uu, vv, pp])
  = int2d(Th)(
    dx(u) * dx(uu)
    + dy(u) * dy(uu)
    + dx(v) * dx(vv)
    + dy(v) * dy(vv)
    + dx(p) * uu
    + dy(p) * vv
    + pp * (dx(u) + dy(v))
    - 1e-10 * p * pp
  )
  + on(1, 2, 4, u=0, v=0)
  + on(3, u=1, v=0)
;

// Plot
plot([u, v], p, wait=1);

---

**Note:** We add a stabilization term \(-10e-10 * p * pp\) to fix the constant part of the pressure.

Results are shown on Fig. 2.12.

### 2.11 A projection algorithm for the Navier-Stokes equations

**Summary:** Fluid flows require good algorithms and good triangulations. We show here an example of a complex algorithm and an first example of mesh adaptation.

An incompressible viscous fluid satisfies:

\[
\begin{align*}
\partial_t u + u \cdot \nabla u + \nabla p - \nu \Delta u &= 0 & \text{in } \Omega \times [0, T] \\
\nabla \cdot u &= 0 & \text{in } \Omega \times [0, T] \\
\left. u \right|_{t=0} &= u^0 \\
\left. u \right|_{\Gamma} &= u_{\Gamma}
\end{align*}
\]

A possible algorithm, proposed by Chorin, is:

\[
\begin{align*}
\frac{1}{\Delta t} [u^{m+1} - u^m - \nu \Delta u^m] + \nabla p^m - \nu \Delta u^m &= 0 \\
\left. u \right|_{\Gamma} &= u_{\Gamma} \\
\nu \partial_n \left. u \right|_{\Gamma_{out}} &= 0
\end{align*}
\]
Fig. 2.12: Solution of Stokes’ equations for the driven cavity problem, showing the velocity field and the pressure level lines.
\[ -\Delta p^{m+1} = -\nabla \cdot u^m o X^m \]
\[ \partial_n p^{m+1} = 0 \quad \text{on } \Gamma \]
\[ p^{m+1} = 0 \quad \text{on } \Gamma_{out} \]

where \( u_0 X(x) = u(x - u(x) \delta t) \) since \( \partial_t u + u \cdot \nabla u \) is approximated by the method of characteristics, as in the previous section.

We use the Chorin’s algorithm with free boundary condition at outlet (i.e. \( p = 0, \nu \partial_n u = 0 \)), to compute a correction, \( q \), to the pressure.

\[ -\Delta q = \nabla \cdot u \]
\[ q = 0 \quad \text{on } \Gamma_{out} \]

and define

\[ u^{m+1} = \tilde{u} + P\nabla q \delta t \]
\[ p^{m+1} = p^m - q \]

where \( \tilde{u} \) is the \((u^{m+1}, v^{m+1})\) of Chorin’s algorithm, and where \( P \) is the \( L^2 \) projection with mass lumping (a sparse matrix).

The backward facing step

The geometry is that of a channel with a backward facing step so that the inflow section is smaller than the outflow section. This geometry produces a fluid recirculation zone that must be captured correctly.

This can only be done if the triangulation is sufficiently fine, or well adapted to the flow.

**Note:** There is a technical difficulty in the example: the output B.C. Here we put \( p = 0 \) and \( \nu \partial_n u = 0 \).

```c
// Parameters
verbosity = 0;
int nn = 1;
real nu = 0.0025;
dt = 0.2;
real epsv = 1e-6;
real epsu = 1e-6;
real epsp = 1e-6;

// Mesh
border a0(t=1, 0){x=-2; y=t; label=1;}
border a1(t=-2, 0){x=t; y=0; label=2;}
border a2(t=0, -0.5){x=0; y=t; label=2;}
border a3(t=0, 1){x=18*t^1.2; y=-0.5; label=2;}
border a4(t=-0.5, 1){x=18; y=t; label=3;}
border a5(t=1, 0){x=-2+20*t; y=1; label=4;}

mesh Th = buildmesh(a0(3*nn) + a1(20*nn) + a2(10*nn) + a3(150*nn) + a4(5*nn) +
                    a5(100*nn));
plot(Th);

// Fespace
fespace Vh(Th, P1);
Vh w;
Vh u = 0, v = 0;
Vh p = 0;
Vh q = 0;
```

(continues on next page)
// Definition of Matrix dtMx and dtMMy
matrix dtM1x, dtM1y;

// Macro
macro BuildMat()
{
   /* for memory management */
   varf vM(unused, v) = int2d(Th)(v);
   varf vdx(u, v) = int2d(Th)(v*dx(u)*dt);
   varf vdy(u, v) = int2d(Th)(v*dy(u)*dt);

   real[int] Mlump = vM(0, Vh);
   real[int] one(Vh.n dof); one = 1;
   real[int] M1 = one ./ Mlump;
   matrix dM1 = M1;
   matrix Mdx = vdx(Vh, Vh);
   matrix Mdy = vdy(Vh, Vh);
   dtM1x = dM1*Mdx;
   dtM1y = dM1*Mdy;
}

// Build matrices
BuildMat

// Time iterations
real err = 1.;
real outflux = 1.;
for(int n = 0; n < 300; n++){
   // Update
   Vh uold=u, vold=v, pold=p;
   // Solve
   solve pb4u (u, w, init=n, solver=CG, eps=epsu)
      = int2d(Th)(
            u*w/dt
         + nu*(dx(u)*dx(w) + dy(u)*dy(w))
        )
   -int2d(Th)(
              convect([uold, vold], -dt, uold)/dt*w
         - dx(p)*w
        )
   + on(1, u=4*y*(1-y))
   + on(2, 4, u=0)
   ;
   plot(u);

   solve pb4v (v, w, init=n, solver=CG, eps=epsv)
      = int2d(Th)(
            v*w/dt
         + nu*(dx(v)*dx(w) + dy(v)*dy(w))
        )
   -int2d(Th)(
              convect([uold,vold],-dt,vold)/dt*w
         - dy(p)*w
        )
   +on(1, 2, 3, 4, v=0)
   ;
}
solve pb4p (q, w, solver=CG, init=n, eps=epsp)
   = int2d(Th) (dx(q)*dx(w)+dy(q)*dy(w))
   - int2d(Th) (dx(u)+dy(v))*w/dt
   + on(3, q=0);

//to have absolute epsilon in CG algorithm.
epsv = -abs(epsv);
epsu = -abs(epsu);
epsp = -abs(epsp);
p = pold-q;
u[] += dtM1x*q[];
v[] += dtM1y*q[];

// Mesh adaptation
if (n%50 == 49) {
   Th = adaptmesh(Th, [u, v], q, err=0.04, nbvx=100000);
   plot(Th, wait=true);
   BuildMat // Rebuild mat.
}

// Error & Outflux
err = sqrt(int2d(Th) (square(u-uold)+square(v-vold))/Th.area);
outflux = int1d(Th) ([u,v]'*[N.x,N.y]);
cout << " iter " << n << " Err L2 = " << err << " outflux = " << outflux << endl;
if(err < 1e-3) break;

// Verification
assert(abs(outflux)< 2e-3);

// Plot
plot(p, wait=1, ps="NSprojP.eps");
plot(u, wait=1, ps="NSprojU.eps");

Rannacher’s projection algorithm: result on an adapted mesh, Fig. 2.13a, showing the pressure, Fig. 2.13b, and the horizontal velocity Fig. 2.13c for a Reynolds number of 400 where mesh adaptation is done after 50 iterations on the first mesh.

### 2.12 Newton Method for the Steady Navier-Stokes equations

The problem is find the velocity field \( \mathbf{u} = (u_i)_{i=1}^d \) and the pressure \( p \) of a Flow satisfying in the domain \( \Omega \subset \mathbb{R}^d (d = 2, 3) \):

\[
(u \cdot \nabla)u - \nu \Delta u + \nabla p = 0 \\
\nabla \cdot u = 0
\]

where \( \nu \) is the viscosity of the fluid, \( \nabla = (\partial_i)_{i=1}^d \), the dot product is \( \cdot \), and \( \Delta = \nabla \cdot \nabla \) with the same boundary conditions (\( u \) is given on \( \Gamma \)).
The weak form is find \( u, p \) such that for \( \forall v \) (zero on \( \Gamma \)), and \( \forall q \):

\[
\int_{\Omega} ((u \cdot \nabla)u) \cdot v + \nu \nabla u : \nabla v - p \nabla \cdot v - q \nabla \cdot u = 0
\]

The Newton Algorithm to solve nonlinear problem is:

Find \( u \in V \) such that \( F(u) = 0 \) where \( F : V \mapsto V \).

1. choose \( u_0 \in \mathbb{R}^n \),
2. for ( \( i = 0; i < \text{niter}; i = i + 1 \))
   1. solve \( DF(u_i)w_i = F(u_i) \);
   2. \( u_{i+1} = u_i - w_i \);
break \( ||w_i|| < \varepsilon \).

Where \( DF(u) \) is the differential of \( F \) at point \( u \), this is a linear application such that:

\[
F(u + \delta) = F(u) + DF(u)\delta + o(\delta)
\]

For Navier Stokes, \( F \) and \( DF \) are:

\[
F(u, p) = \int_{\Omega} ((u \cdot \nabla)u) \cdot v + \nu \nabla u : \nabla v - p \nabla \cdot v - q \nabla \cdot u
\]

\[
DF(u, p)(\delta u, \delta p) = \int_{\Omega} ((\delta u \cdot \nabla)u) \cdot v + ((u \cdot \nabla)\delta u) \cdot v + \nu \nabla \delta u : \nabla v - \delta p \nabla \cdot v - q \nabla \cdot \delta u
\]

So the Newton algorithm become:

```plaintext
// Parameters
real R = 5.;
real L = 15.;
real nu = 1./50.;
real nufinal = 1/200.;
real cnu = 0.5;
real eps = 1e-6;
verbosity = 0;

// Mesh
```

(continues on next page)
border cc(t=0, 2*pi){x=cos(t)/2.; y=sin(t)/2.; label=1;}
border ce(t=pi/2, 3*pi/2){x=cos(t)+R; y=sin(t)+R; label=1;}
border beb(tt=0, 1){real t=tt^1.2; x=t*L; y=-R; label=1;}
border beo(t=-R, R){x=L; y=t; label=0;}
border bei(t=-R/4, R/4){x=L/2; y=t; label=0;}

mesh Th = buildmesh(cc(-50) + ce(30) + beb(20) + beu(20) + beo(10) + bei(10));

//bounding box for the plot
func bb = [[-1,-2],[4,2]];

// Fespace
fespace Xh(Th, P2);
Xh u1, u2;
Xh v1, v2;
Xh du1, du2;
Xh ulp, u2p;

fespace Mh(Th, P1);
Mh p;
Mh q;
Mh dp;
Mh pp;

// Macro
macro Grad(u1,u2) [dx(u1), dy(u1), dx(u2),dy(u2)] //
macro UgradV(u1,u2,v1,v2) [u1,u2]'*[dx(v1),dy(v1)],
[1,1]'*[dx(v2),dy(v2)];
macro div(u1,u2) (dx(u1) + dy(u2)) //

// Initialization
u1 = (x^2+y^2) > 2;
u2 = 0;

// Viscosity loop
while(1){
    int n;
    real err=0;
    // Newton loop
    for (n = 0; n < 15; n++){
        // Newton
        solve Oseen ([du1, du2, dp], [v1, v2, q])
           = int2d(Th) {
               nu * (Grad(du1,du2)' + Grad(v1,v2))
                + UgradV(du1,du2, u1, u2)'*[v1,v2]
                + UgradV( u1, u2,du1,du2)'*[v1,v2]
                - div(du1,du2) * q
                - div(v1,v2) * dp
                - 1e-8*dp*q //stabilization term
            } - int2d(Th) {
               nu * (Grad(u1,u2)' + Grad(v1,v2))
                + UgradV(u1,u2, u1, u2)'*[v1,v2]
                - div(u1,u2) * q
                - div(v1,v2) * p
            }
        }
    }
}

(continues on next page)
\begin{verbatim}
+ on(1, du1=0, du2=0)

; u1[] -= du1[];

u2[] -= du2[];

p[] -= dp[];

real Lu1=u1[].linfty, Lu2=u2[].linfty, Lp=p[].linfty;

err = du1[].linfty/Lu1 + du2[].linfty/Lu2 + dp[].linfty/Lp;

cout << n << " err = " << err << " " << eps " rey = " << 1./nu << endl;
if(err < eps) break; //converge
if( n>3 && err > 10.) break; //blowup
}

if(err < eps){ //converge: decrease \$\nu\$ (more difficult)

// Plot
plot([u1, u2], p, wait=1, cmm=" rey = " + 1./nu , coef=0.3, bb=bb);

// Change nu
if( nu == nufinal) break;
if( n < 4) cnu = cnu^1.5; //fast converge => change faster

nu = max(nufinal, nu* cnu); //new viscosity

// Update
u1p = u1;
u2p = u2;
p = pp;
}
else{ //blowup: increase \$\nu\$ (more simple)

assert(cnu< 0.95); //the method finally blowup

// Recover nu
nu = nu/cnu;
cnu= cnu^(1./1.5); //no conv. => change lower

nu = nu* cnu; //new viscosity

cout << " restart nu = " << nu << " " <<'Rey = " << 1./nu " " (cnu = " << cnu << " " ) "n";

// Recover a correct solution
u1 = u1p;
u2 = u2p;
p = pp;
}

Note: We use a trick to make continuation on the viscosity $\nu$, because the Newton method blowup owe start with the final viscosity $\nu$.
$\nu$ is gradually increased to the desired value.
\end{verbatim}
2.13 A Large Fluid Problem

A friend of one of us in Auroville-India was building a ramp to access an air conditioned room. As I was visiting the
construction site he told me that he expected to cool air escaping by the door to the room to slide down the ramp and
refrigerate the feet of the coming visitors. I told him “no way” and decided to check numerically.

The fluid velocity and pressure are solution of the Navier-Stokes equations with varying density function of the tem-
perature.

The geometry is trapezoidal with prescribed inflow made of cool air at the bottom and warm air above and so are the
initial conditions; there is free outflow, slip velocity at the top (artificial) boundary and no-slip at the bottom. However
the Navier-Stokes cum temperature equations have a RANS $k - \epsilon$ model and a Boussinesq approximation for the
buoyancy. This comes to:

$$
\partial_t \theta + u \nabla \theta - \nabla \cdot (k_T \nabla \theta) = 0 \\
\partial_t u + u \nabla u - \nabla \cdot (\mu_T \nabla u) + \nabla p + \rho(\theta - \theta_0) \epsilon_2 = 0 \\
\nabla \cdot u = 0 \\
\mu_T = c_\mu \frac{\epsilon^2}{\kappa_T} \\
\kappa_T = \kappa \mu_T \\
\partial_t k + u \nabla k + \epsilon - \nabla \cdot (\mu_T \nabla k) = \frac{\mu_T}{2} |\nabla u + \nabla u^T|^2 \\
\partial_t \epsilon + u \nabla \epsilon + c_2 \frac{\epsilon^2}{\kappa_T} - \frac{\epsilon}{c_T} \nabla \cdot (\mu_T \nabla \epsilon) = \frac{\nu_T}{2} |\nabla u + \nabla u^T|^2
$$

We use a time discretization which preserves positivity and uses the method of characteristics $(X^m(x) \approx x - u^m(x) \delta t)$

\[ \frac{1}{\delta t} (u^{m+1} - u^m \circ X^m) - \nabla \cdot (\mu_T \nabla u^{m+1}) + \nabla p^{m+1} + \rho(\theta^{m+1} - \theta_0) \epsilon_2 = 0 \]

\[ \nabla \cdot u^{m+1} = 0 \]

\[ \frac{1}{\delta t} (k^{m+1} - k^m \circ X^m) + \frac{1}{k_m} \frac{e^m}{k_m^2} - \nabla \cdot (\mu_T \nabla k^{m+1}) = \frac{2}{\mu_T} |\nabla u^m + \nabla u^m|^2 \]

\[ \frac{1}{\delta t} (\epsilon^{m+1} - \epsilon^m \circ X^m) + c_2 \frac{\epsilon^m}{k_m^2} - \frac{\epsilon}{c_T} \nabla \cdot (\mu_T \nabla \epsilon^{m+1}) = \frac{\nu_T}{2} |\nabla u^m + \nabla u^m|^2 \]

\[ \mu_T^{m+1} = c_\mu \frac{k^{m+1}}{\epsilon^{m+1}} \]

\[ \kappa_T^{m+1} = \kappa \mu_T^{m+1} \]

In variational form and with appropriated boundary conditions the problem is:

\begin{verbatim}
load "iovtk"
verbosity=0;
\end{verbatim}

(continues on next page)
5 // Parameters
6 int nn = 15;
7 int nnPlus = 5;
8 real l = 1.;
9 real L = 15.;
10 real hSlope = 0.1;
11 real H = 6.;
12 real h = 0.5;
13
14 real reylnods = 500;
15 real beta = 0.01;
16
17 real eps = 9.81 / 303.;
18 real nu = 1;
19 real numu = nu / sqrt(0.09);
20 real nuep = pow(nu, 1.5) / 4.1;
21 real dt = 0.;
22
23 real Penalty = 1.e-6;
24
25 // Mesh
26 border b1(t=0, l){x=t; y=0;}
27 border b2(t=0, L-l){x=l+t; y=-hSlope*t;}
28 border b3(t=-hSlope*(L-l), H){x=L; y=t;}
29 border b4(t=L, 0){x-t; y=H;}
30 border b5(t=H, h){x=0; y-t;}
31 border b6(t=h, 0){x=0; y-t;}
32
33 mesh Th = buildmesh(b1(nnPlus*nn*l) + b2(nn*sqrt((L-l)^2 + (hSlope*(L-l))^2)) + b3(nn*(H-\rightarrow hSlope*(L-l))) + b4(nn*L) + b5(nn*(H-h)) + b6(nnPlus*nn*h));
34 plot(Th);
35
36 // Fespaces
37 fespace Vh2(Th, P1b);
38 Vh2 Ux, Uy;
39 Vh2 Vx, Vy;
40 Vh2 Upx, Upy;
41
42 fespace Vh(Th, P1);
43 Vh p=0, q;
44 Vh Tp, T=35;
45 Vh k=0.0001, kp=k;
46 Vh ep=0.0001, epp=ep;
47
48 fespace V0h(Th, P0);
49 V0h muT=1;
50 V0h prodk, prode;
51 Vh kappa=0.25e-4, stress;
52
53 // Macro
54 macro grad(u) [dx(u), dy(u)];
55 macro Grad(U) [grad(U#x), grad(U#y)];
56 macro Div(U) (dx(U#x) + dy(U#y));
57
58 // Functions
59 func g = (x) * (1-x) * 4;
60
(continues on next page)
// Problem
real alpha = 0.;

problem Temperature(T, q)
    = int2d(Th)(
        alpha * T * q
        + kappa * grad(T)' * grad(q)
    )
    + int2d(Th)(
        - alpha * convect([Upx, Upy], -dt, Tp) * q
    )
    + on(b6, T=25)
    + on(b1, b2, T=30)
;

problem KineticTurbulence(k, q)
    = int2d(Th)(
        (epp/kp + alpha) * k * q
        + muT * grad(k)' * grad(q)
    )
    + int2d(Th)(
        prodk * q
        - alpha * convect([Upx, Upy], -dt, kp) * q
    )
    + on(b5, b6, k=0.00001)
    + on(b1, b2, k=beta*numu*stress)
;

problem ViscosityTurbulence(ep, q)
    = int2d(Th)(
        (1.92*epp/kp + alpha) * ep * q
        + muT * grad(ep)' * grad(q)
    )
    + int1d(Th, b1, b2)(
        T * q * 0.001
    )
    + int2d(Th)(
        prodX * q
        - alpha * convect([Upx, Upy], -dt, epp) * q
    )
    + on(b5, b6, ep=0.00001)
    + on(b1, b2, ep=beta*nuep*pow(stress,1.5))
;

// Initialization with stationary solution
solve NavierStokes ([Ux, Uy, p], [Vx, Vy, q])
    = int2d(Th)(
        alpha * [Ux, Uy]' * [Vx, Vy]
        + muT * (Grad(U) : Grad(V))
        + p * q * Penalty
        - p * Div(V)
        - Div(U) * q
    )
    + int1d(Th, b1, b2, b4)(
        Ux * Vx * 0.1
    )
    + int2d(Th)(
(continues on next page)
eps *(T-35) * Vx
- alpha*convect([Upx, Upy], -dt, Upx)*Vx
- alpha*convect([Upx, Upy], -dt, Upy)*Vy
)
+ on(b6, Ux=3, Uy=0)
+ on(b5, Ux=0, Uy=0)
+ on(b1, b4, Uy=0)
+ on(b2, Uy=-Upx*N.x/N.y)
+ on(b3, Uy=0)
;
plot([Ux, Uy], p, value=true, coef=0.2, cmm="[Ux, Uy] - p");
{
    real[int] xx(21), yy(21), pp(21);
    for (int i = 0 ; i < 21; i++){
        yy[i] = i/20.;
        xx[i] = Ux(0.5,i/20.);
        pp[i] = p(i/20.,0.999);
    }
    cout << " " << yy << endl;
    plot([xx, yy], wait=true, cmm="Ux x=0.5 cup");
    plot([yy, pp], wait=true, cmm="p y=0.999 cup");
}
// Initialization
dt = 0.1; //probably too big
int nbiter = 3;
real coefdt = 0.25^(1./nbiter);
real coefcut = 0.25^(1./nbiter);
real cut = 0.01;
real tol = 0.5;
real coeftol = 0.5^(1./nbiter);
nu = 1./reylnods;
T = T - 10*((x<1)*(y<0.5) + (x>=1)*(y+0.1*(x-1)<0.5));
// Convergence loop
real T0 = clock();
for (int iter = 1; iter <= nbiter; iter++){
    cout << "Iteration " << iter << " - dt = " << dt << endl;
    alpha = 1/dt;
    // Time loop
    real t = 0.;
    for (int i = 0; i <= 500; i++){
        t += dt;
        cout << "Time step " << i << " - t = " << t << endl;
        // Update
        Upx = Ux;
        Upy = Uy;
        kp = k;
        epp = ep;
        Tp = max(T, 25); //for beauty only should be removed
        Tp = min(Tp, 35); //for security only should be removed
        kp = max(k, 0.0001); epp = max(ep, 0.0001); //to be secure: should not be active
muT = 0.09*kp*kp/epp;

// Solve NS
NavierStokes;

// Update
prode = -0.128*kp*(pow(2*dx(Ux),2)+pow(2*dy(Uy),2)+2*pow(dx(Uy)+dy(Ux),2))/2;
prodk = -prode*kp/epp*0.09/0.128;
kappa = muT/0.41;
stress = abs(dy(Ux));

// Solve k-eps-T
KineticTurbulence;
ViscosityTurbulence;
Temperature;

// Plot
plot(T, value=true, fill=true);
plot([Ux, Uy], p, coef=0.2, cmm=" [Ux, Uy] - p", WindowIndex=1);

// Time
cout << "Time = " << clock()-T0 << endl;

// Check
if (iter >= nbiter) break;

// Adaptmesh
Th = adaptmesh(Th, [dx(Ux), dy(Ux), dx(Ux), dy(Uy)], splitpbedge=1, abserro=0, 
cutoff=cut, err=tol, inquire=0, ratio=1.5, hmin=1./1000);
plot(Th);

// Update
dt = dt * coefdt;
tol = tol * coeftol;
cut = cut * coeffcut;

cout << "Total Time = " << clock()-T0 << endl;

2.14 An Example with Complex Numbers

In a microwave oven heat comes from molecular excitation by an electromagnetic field. For a plane monochromatic wave, amplitude is given by Helmholtz's equation:

\[ \beta v + \Delta v = 0. \]

We consider a rectangular oven where the wave is emitted by part of the upper wall. So the boundary of the domain is made up of a part \( \Gamma_1 \) where \( v = 0 \) and of another part \( \Gamma_2 = [c, d] \) where for instance \( v = \sin \left( \frac{\pi y - c}{c - d} \right) \).

Within an object to be cooked, denoted by \( B \), the heat source is proportional to \( v^2 \). At equilibrium, one has:

\[ -\Delta \theta = v^2 I_B \]

where \( I_B \) is 1 in the object and 0 elsewhere.
(a) Temperature at time step 100

(b) Velocity at time step 100

(c) Temperature at time step 200

(d) Velocity at time step 200

(e) Temperature at time step 300

(f) Velocity at time step 300

(g) Temperature at time step 400

(h) Velocity at time step 400

(i) Temperature at time step 500

(j) Velocity at time step 500

Fig. 2.15: A large fluid problem
In the program below $\beta = 1/(1 - i/2)$ in the air and $2/(1 - i/2)$ in the object $(i = \sqrt{-1})$:

```cpp
// Parameters
int nn = 2;
real a = 20.;
real b = 20.;
real c = 15.;
real d = 8.;
real e = 2.;
real l = 12.;
real f = 2.;
real g = 2.;

// Meshorder a0(t=0, 1){x=a*t; y=0; label=1;}
border a1(t=1, 2){x=a; y=b*(t-1); label=1;}
border a2(t=2, 3){x=a*(3-t); y=b; label=1;}
border a3(t=3, 4){x=0; y=b-(b-c)*(t-3); label=1;}
border a4(t=4, 5){x=0; y=c-(c-d)*(t-4); label=2;}
border a5(t=5, 6){x=0; y=d*(6-t); label=1;}
border b0(t=0, 1){x=a-f+e*(t-1); y=g; label=3;}
border b1(t=1, 4){x=a-f; y=g+l*(t-1)/3; label=3;}
border b2(t=4, 5){x=a-f-e*(t-4); y=l+g; label=3;}
border b3(t=5, 8){x=a-e-f; y=l+g-l*(t-5)/3; label=3;}
mesh Th = buildmesh(a0(10*nn) + a1(10*nn) + a2(10*nn) + a3(10*nn) + a4(10*nn) + a5(10*nn) + b0(5*nn) + b1(10*nn) + b2(5*nn) + b3(10*nn));
real meat = Th(a-f-e/2, g+l/2).region;
real air = Th(0.01, 0.01).region;
plot(Th, wait=1);

// Fespace
fespace Vh(Th, P1);
Vh R=(region-air)/(meat-air);
Vh<complex> v, w;
Vh vr, vi;

fespace Uh(Th, P1);
Uh u, uu, ff;

// Problem
solve muwave(v, w) = int2d(Th)({
v*w*(1+R)
  - (dx(v)*dx(w) + dy(v)*dy(w))*(1 - 0.5i)
}) + on(1, v=0) + on(2, v=sin(pi*(y-c))/(c-d)))
;
vr = real(v);
vi = imag(v);

// Plot
plot(vr, wait=1, ps="rmuonde.ps", fill=true);
plot(vi, wait=1, ps="imuonde.ps", fill=true);
```

(continues on next page)
Results are shown on Fig. 2.16a, Fig. 2.16b and Fig. 2.16c.

2.15 Optimal Control

Thanks to the function BFGS it is possible to solve complex nonlinear optimization problem within FreeFEM. For example consider the following inverse problem

\[ J(b, c, d) = \int_E (u - u_d)^2 \]

\[ -\nabla(\kappa(b, c, d) \cdot \nabla u) = 0 \]

\[ u|_\Gamma = u_\Gamma \]

where the desired state \( u_d \), the boundary data \( u_\Gamma \) and the observation set \( E \subseteq \Omega \) are all given. Furthermore let us assume that:

\[ \kappa(x) = 1 + bI_B(x) + cI_C(x) + dI_D(x) \quad \forall x \in \Omega \]

where \( B, C, D \) are separated subsets of \( \Omega \).

To solve this problem by the quasi-Newton BFGS method we need the derivatives of \( J \) with respect to \( b, c, d \). We self explanatory notations, if \( \delta b, \delta c, \delta d \) are variations of \( b, c, d \) we have:

\[ \delta J \approx 2 \int_E (u - u_d) \delta u \]

\[ -\nabla(\kappa \cdot \nabla \delta u) \approx \nabla(\delta \kappa \cdot \nabla u) \]

\[ \delta u|_\Gamma = 0 \]

Obviously \( J'_b \) is equal to \( \delta J \) when \( \delta b = 1, \delta c = 0, \delta d = 0 \), and so on for \( J'_c \) and \( J'_d \).

All this is implemented in the following program:

```cpp
// Mesh
border aa(t=0, 2*pi){x=5*cos(t); y=5*sin(t);};
border bb(t=0, 2*pi){x=cos(t); y=sin(t);};
border cc(t=0, 2*pi){x=-3+cos(t); y=sin(t);};
border dd(t=0, 2*pi){x=cos(t); y=-3+sin(t);};

mesh th = buildmesh(aa(70) + bb(35) + cc(35) + dd(35));

// Fespace
```
Fig. 2.16: Microwave
fespace Vh(th, P1);
Vh Ib=((x^2+y^2)<1.0001),
Ic=((x+3)^2+y^2)<1.0001),
Id=((x^2+(y+3)^2)<1.0001),
Ie=((x-1)^2+y^2)<=4),
ud, u, uh, du;

// Problem
real[int] z(3);
problem A(u, uh)
  = int2d(th)(
    (1+z[0]*Ib+z[1]*Ic+z[2]*Id)*(dx(u)*dx(uh) + dy(u)*dy(uh))
  )
  + on(aa, u=x^3-y^3)
;

// Solve
z[0]=2; z[1]=3; z[2]=4;
A;
ud = u;

ofstream f("J.txt");
func real J(real[int] & Z){
  for(int i = 0; i < z.n; i++)
    z[i] = Z[i];
  A;
  real s = int2d(th)(Ie*(u-ud)^2);
  f << s << " ";
  return s;
}

// Problem BFGS
real[int] dz(3), dJdz(3);
problem B (du, uh)
  = int2d(th)(
    (1+z[0]*Ib+z[1]*Ic+z[2]*Id)*(dx(du)*dx(uh) + dy(du)*dy(uh))
  )
  + int2d(th)(
    (dz[0]*Ib+dz[1]*Ic+dz[2]*Id)*(dx(u)*dx(uh) + dy(u)*dy(uh))
  )
  +on(aa, du=0)
;

func real[int] DJ(real[int] & Z){
  for(int i = 0; i < z.n; i++){
    for(int j = 0; j < dz.n; j++)
      dz[j] = 0;
    dz[i] = 1;
    B;
    dJdz[i] = 2*int2d(th)(Ie*(u-ud)*du);
  }
  return dJdz;
}

real[int] Z(3);
for(int j = 0; j < z.n; j++)
  Z[j]=1;

(continues on next page)
In this example the sets $B, C, D, E$ are circles of boundaries $bb, cc, dd, ee$ and the domain $\Omega$ is the circle of boundary $aa$.

The desired state $u_d$ is the solution of the PDE for $b = 2, c = 3, d = 4$. The unknowns are packed into array $z$.

**Note:** It is necessary to recopy $Z$ into $z$ because one is a local variable while the other one is global.

The program found $b = 2.00125, c = 3.00109, d = 4.00551$.

Fig. 2.17a and Fig. 2.17b show $u$ at convergence and the successive function evaluations of $J$.

Note that an *adjoint state* could have been used. Define $p$ by:

\[
-\nabla \cdot (\kappa \nabla p) = 2I_E(u - u_d)
\]

\[p|_\Gamma = 0\]

Consequently:

\[
\delta J = - \int_\Omega (\nabla \cdot (\kappa \nabla p)) \delta u
\]

\[
= \int_\Omega (\kappa \nabla p \cdot \nabla \delta u)
\]

\[
= - \int_\Omega (\delta \kappa \nabla p \cdot \nabla u)
\]
Then the derivatives are found by setting \( \delta b = 1, \delta c = \delta d = 0 \) and so on:

\[
J_b' = - \int_B \nabla p \cdot \nabla u \\
J_c' = - \int_C \nabla p \cdot \nabla u \\
J_d' = - \int_D \nabla p \cdot \nabla u 
\]

**Note:** As BFGS stores an \( M \times M \) matrix where \( M \) is the number of unknowns, it is dangerously expensive to use this method when the unknown \( x \) is a Finite Element Function. One should use another optimizer such as the NonLinear Conjugate Gradient NLCG (also a key word of FreeFEM).

### 2.16 A Flow with Shocks

Compressible Euler equations should be discretized with Finite Volumes or FEM with flux up-winding scheme but these are not implemented in FreeFEM. Nevertheless acceptable results can be obtained with the method of characteristics provided that the mean values \( \bar{f} = \frac{1}{2} \left( f^+ + f^- \right) \) are used at shocks in the scheme, and finally mesh adaptation.

\[
\partial_t \rho + \bar{u} \nabla \rho + \bar{p} \nabla \cdot u = 0 \\
\bar{p} (\partial_t u + \frac{\gamma - 1}{\rho} \nabla \cdot u + \nabla p) = 0 \\
\partial_t p + \bar{u} \nabla p + (\gamma - 1) \bar{p} \nabla \cdot u = 0
\]

One possibility is to couple \( u, p \) and then update \( \rho \), i.e.:

\[
\frac{1}{\rho_m} (p_m^{m+1} - p_m^m \circ X^m) + \nabla \cdot u^{m+1} = 0 \\
\frac{\gamma - 1}{\rho_m} (u^{m+1} - u^m \circ X^m) + \nabla p^{m+1} = 0 \\
\rho^{m+1} = \rho^m \circ X^m + \frac{\gamma - 1}{\rho_m} (p^{m+1} - p^m \circ X^m)
\]

A numerical result is given on Fig. 2.18 and the FreeFEM script is

```cpp
// Parameters
verbosity = 1;
int anew = 1;
int m = 5;
real x0 = 0.5;
real y0 = 0.2;
real rr = 0.2;
real dt = 0.01;
real u0 = 2.;
real err0 = 0.00625;
real pena = 2.;

// Mesh
border ccc(t=0, 2){x=2-t; y=1;};
border ddd(t=0, 1){x=0; y=1-t;};
border aaal(t=0, x0-rr){x=t; y=0;};
border cercle(t=pi, 0){x=x0+rr*cos(t); y=y0+rr*sin(t);}
border aa2(t=x0+rr, 2){x=t; y=0;};
border bbb(t=0, 1){x=2; y=t;};

mesh Th;
if (anew)
    Th = buildmesh (ccc(5*m) + ddd(3*m) + aaal(2*m) + cercle(5*m) + aa2(5*m) +
                    -> bbb(2*m));
```

(continues on next page)
else
    Th = readmesh("Th_circle.mesh"); plot(Th);

// fespace
fespace Wh(Th, P1);
Wh u, v;
Wh u1, v1;
Wh uh, vh;

fespace Vh(Th, P1);
Vh r, rh, r1;

// Macro
macro dn(u) (
    N.x*dx(u)+N.y*dy(u))
//

// Initialization
if(anew){
    u1 = u0;
    v1 = 0;
    r1 = 1;
}
else{
    ifstream g("u.txt"); g >> u1[];
    ifstream gg("v.txt"); gg >> v1[];
    ifstream ggg("r.txt"); ggg >> r1[];
    plot(u1, ps="eta.eps", value=1, wait=1);
    err0 = err0/10;
    dt = dt/10;
}

// Problem
problem euler(u, v, r, uh, vh, rh)
    = int2d(Th)(
        (u*uh + v*vh + r*rh)/dt
        + ((dx(r)*uh + dy(r)*vh) - (dx(rh)*u + dy(rh)*v))
    )
    + int2d(Th)(
        rh*convect([u1,v1],-dt,r1)
        + uh*convect([u1,v1],-dt,u1)
        + vh*convect([u1,v1],-dt,v1)
    )/dt
    + int1d(Th, 6)(
        rh*u
    )
    + on(2, r=0)
    + on(2, u=u0)
    + on(2, v=0);

// Iterations
int j = 80;
for(int k = 0; k < 3; k++){
    if(k==20){
        err0 = err0/10;
        dt = dt/10;
    }
2.17 Time dependent schema optimization for heat equations

First, it is possible to define variational forms, and use this forms to build matrix and vector to make very fast script (4 times faster here).

For example solve the *ThermalConduction* problem, we must solve the temperature equation in $\Omega$ in a time interval $(0,T)$.

\[
\begin{align*}
\frac{\partial u}{\partial t} - \nabla \cdot (\kappa \nabla u) &= 0 & \text{in } \Omega \times (0, T) \\
u(x, y, 0) &= u_0 + xu_1 & \text{on } \Gamma_{24} \times (0, T) \\
\kappa \frac{\partial u}{\partial n} + \alpha (u - u_e) &= 0 & \text{on } \Gamma \times (0, T)
\end{align*}
\]

The variational formulation is in $L^2(0, T; H^1(\Omega))$; we shall seek $u^n$ satisfying:

\[
\forall w \in V_0 : \int_{\Omega} \frac{u^n - u_{n-1}}{\delta t} w + \kappa \nabla u^n \nabla w + \int_{\Gamma} \alpha (u^n - u_{ue}) w = 0
\]

where $V_0 = \{ w \in H^1(\Omega)/w|_{\Gamma_{24}} = 0 \}$.

So, to code the method with the matrices $A = (A_{ij})$, $M = (M_{ij})$, and the vectors $u^n, b^n, b', b'', b_{cd}$ (notation if $w$ is a vector then $w_i$ is a component of the vector).

\[
u^n = A^{-1}b^n, \quad b' = b_0 + Mu^{n-1}, \quad b'' = \frac{1}{\varepsilon} b_{cd}, \quad b_i' = \begin{cases} b_i' & \text{if } i \in \Gamma_{24} \\ b_i'' & \text{else} \end{cases}
\]
Fig. 2.18: Pressure for a Euler flow around a disk at Mach 2 computed by (2.6)
Where with $\frac{1}{\varepsilon} = \text{tgv} = 10^{30}$:

\[
A_{ij} = \begin{cases} 
\frac{1}{\varepsilon} & \text{if } i \in \Gamma_{24}, \text{ and } j = i \\
\int_\Omega w_j w_i / dt + k(\nabla w_j, \nabla w_i) + \int_{\Gamma_{13}} \alpha w_j w_i & \text{else} 
\end{cases}
\]

\[
M_{ij} = \begin{cases} 
\frac{1}{\varepsilon} & \text{if } i \in \Gamma_{24}, \text{ and } j = i \\
\int_\Omega w_j w_i / dt & \text{else} 
\end{cases}
\]

\[
b_{0,i} = \int_\Omega \alpha u_{ue} w_i \\
b_{cl} = u_0 
\]

the initial data

The Fast version script:

```c
... 
Vh u0 = fu0, u = u0;
```

Create three variational formulation, and build the matrices $A,M$.

```c
varf vthermic (u, v) = int2d(Th)(
    u*v/dt
    + k*(dx(u)*dx(v) + dy(u)*dy(v))
) + int1d(Th, 1, 3){
    alpha*u*v
} + on(2, 4, u=1);

varf vthermic0 (u, v) = int1d(Th, 1, 3){
    alpha*ue*v
};

varf vMass (u,v) = int2d(Th)(
    u*v/dt
) + on(2, 4, u=1);

real tgv = 1e30;
matrix A = vthermic(Vh, Vh, tgv=tgv, solver=CG);
matrix M = vMass(Vh, Vh);
```

Now, to build the right hand size; we need 4 vectors.

```c
real[int] b0 = vthermic0(0,Vh); //constant part of RHS
real[int] bcn = vthermic(0,Vh); //tgv on Dirichlet part
real[int] bcl = tgv*u0[]; //the Dirichlet B.C. part
```

// The fast loop
```c
for(real t = 0; t < T; t += dt){
    real[int] b = b0; //the RHS
    b += M*u[]; //add the the time dependent part
    b = bcn ? bcl : b; //do $\forall i$: b[i] = bcn[i] ? bcl[i] : b[i];
    u[] = A^-1*b; //solve linear problem
}
```

(continues on next page)
2.18 Tutorial to write a transient Stokes solver in matrix form

Consider the following script to solve a time dependent Stokes problem in a cavity:

```cpp
// Parameters
real nu = 0.1;
real T=1.;
real dt = 0.1;

// Mesh
mesh Th = square(10, 10);

// Fespace
fespace Vh(Th, P2);
Vh u, v, uu, vv;
Vh uold=0, vold=0;

fespace Qh(Th, P1);
Qh p, pp;

// Problem
problem stokes (u, v, p, uu, vv, pp) = int2d(Th) (u*uu+v*vv)/dt + nu*(dx(u)*dx(uu) + dy(u)*dy(uu) + dx(v)*dx(vv) + dy(v)*dy(vv)) - p*pp*1.e-6 + p*(dx(uu) + dy(vv)) + pp*(dx(u) + dy(v))
   - int2d(Th) (uold*uu+vold*vv)/dt + on(1, 2, 4, u=0, v=0) + on(3, u=1, v=0);

// Time loop
int m, M = T/dt;
for(m = 0; m < M; m++)
{
    stokes;
    uold = u;
    vold = v;
}

// Plot
plot(p, [u, v], value=true, wait=true, cmm="t=#{m*dt}");
```

Every iteration is in fact of the form $A[u, v, p] = B[uold, vold, pold] + b$ where $A, B$ are matrices and $b$ is a vector containing the boundary conditions. $A, B, b$ are constructed by:
FreeFEM Documentation, Release 4.6

```plaintext
fespace Xh(Th, [P2, P2, P1]);

varf aa ([u, v, p], [uu, vv, pp])
  = int2d(Th) 
    (u*uu+v*vv)/dt 
    + nu*(dx(u)*dx(uu) + dy(u)*dy(uu) + dx(v)*dx(vv) + dy(v)*dy(vv)) 
    - p*pp*1.e-6 
    - p*(dx(uu) + dy(vv)) 
    - pp*(dx(u) + dy(v)) 
  + on(1, 2, 4, u=0, v=0) 
  + on(3, u=1, v=0);

varf bb ([uold, vold, pold], [uu, vv, pp])
  = int2d(Th) 
    (uold*uu+vold*vv)/dt 
  //+ on(1, 2, 4, uold=0, vold=0) 
  //+ on(3, uold=1, vold=0);

varf bcl ([uold, vold, pold], [uu, vv, pp])
  = on(1, 2, 4, uold=0, vold=0) 
  + on(3, uold=1, vold=0);

matrix A = aa(Xh, Xh, solver=UMFPACK);
matrix B = bb(Xh, Xh);
real[int] b = bcl(0, Xh);

Note that the boundary conditions are not specified in bb. Removing the comment // would cause the compiler to multiply the diagonal terms corresponding to a Dirichlet degree of freedom by a very large term (tgv); if so b would not be needed, on the condition that uold = 1 on boundary 3 initially. Note also that b has a tgv on the Dirichlet nodes, by construction, and so does A.

The loop will then be:

```plaintext
real[int] sol(Xh.ndof), aux(Xh.ndof);
for (m = 0; m < M; m++)
  aux = B*sol; aux += b;
  sol = A^-1 * aux;
```

There is yet a difficulty with the initialization of sol and with the solution from sol. For this we need a temporary vector in $X_h$ and here is a solution:

```plaintext
Xh [w1, w2, wp] = [uold, vold, pp];
sol = w1[]; //cause also the copy of w2 and wp
for (m = 0; m < M; m++)
  aux = B*sol; aux += b;
  sol = A^-1 * aux;

w1[] = sol; u = w1; v = w2; p = wp;
plot(p, [u, v], value=true, wait=true, cmm="t"=m*dt);
```

The freefem team agrees that the line sol=w1[]; is mysterious as it copies also w2 and wp into sol. Structured data such as vectors of $X_h$ here cannot be written component by component. Hence w1=u is not allowed.

2.18. Tutorial to write a transient Stokes solver in matrix form
2.19 Wifi Propagation

2.19.1 Summary

In this tutorial, we will study the wifi signal power in a flat. An awesome flat is especially designed for the experiment, with 2 walls:

![Flat Diagram](image)

Fig. 2.19: Flat

Even if the flat seems small enough to be covered by wifi everywhere, it is still interesting to study where the signal’s power is the lowest. We will study where to put the hotspot to get the best coverage, and as we’re a bit lazy we will only put it next to the left wall.

2.19.2 Physics

In a nutshell, the Wifi is a electromagnetic wave that contains a signal: Internet data. Electromagnetic waves are well known by physicists and are ruled by the 4 Maxwell equations which give you the solution for $E$, the electrical field, and $B$, the magnetic field, in space but also in time.

We don’t care about the time here, because the signal period is really short so our internet quality will not change with time. Without time, we’re looking for stationaries solutions, and the Maxwell equations can be simplified to one equation, the Helmholtz one:

$$\nabla^2 E + \frac{k^2}{n^2} E = 0$$

Where $k$ is the angular wavenumber of the wifi signal, and $n$ the refractive index of the material the wave is in.

Indeed, the main point of this study is the impact of walls on the signal’s power, where the $n$ is different from air (where it is 1). In walls, the refractive index is a complex number in which the two parts have a physic interpretation:

- The real part defines the reflection of the wall (the amount of signal that doesn’t pass).
- The imaginary part defines the absorption of the wall (the amount that disappears).

The wifi hotspot (simulated by a simple circle) will be the boundary condition, with a non null value for our electrical field.
2.19.3 Coding

The domain

In order to create the domain of experimentation, we need to create \texttt{border} objects, like this:

```plaintext
real a = 40, b = 40, c = 0.5;
border a00(t=0, 1) { x=a*t; y=0; label=1; }
border a10(t=0, 1) { x=a; y=b*t; label=1; }
border a20(t=1, 0) { x=a*t; y=b; label=1; }
border a30(t=1, 0) { x=0; y=b*t; label=1; }
border a01(t=0, 1) { x=c+(a-c*2)*t; y=c; label=1; }
border a11(t=0, 1) { x=a-c; y=c+(b-c*2)*t; label=1; }
border a21(t=1, 0) { x=c+(a-c*2)*t; y=b-c; label=1; }
border a31(t=1, 0) { x=c; y=c+(b-c*2)*t; label=1; }

real p = 5, q = 20, d = 34, e = 1;
border b00(t=0, 1) { x=p+d*t; y=q; label=3; }
border b10(t=0, 1) { x=p+d; y=q+e*t; label=3; }
border b20(t=1, 0) { x=p+d*t; y=q+e; label=3; }
border b30(t=1, 0) { x=p; y=q+e*t; label=3; }

real r = 30, s =1 , j = 1, u = 15;
border c00(t=0, 1) { x=r+j*t; y=s; label=3; }
border c10(t=0, 1) { x=r+j; y=s+u*t; label=3; }
border c20(t=1, 0) { x=r+j*t; y=s+u; label=3; }
border c30(t=1, 0) { x=r; y=s+u*t; label=3; }
```

Let's create a mesh

```plaintext
int n=13;
mesh Sh = buildmesh(a00(10*n) + a10(10*n) + a20(10*n) + a30(10*n) + a01(10*n) + a11(10*n) + a21(10*n) + a31(10*n) + b00(5*n) + b10(5*n) + b20(5*n) + b30(5*n) + c00(5*n) + c10(5*n) + c20(5*n) + c30(5*n));
plot(Sh, wait=1);
```

So we are creating a mesh, and plotting it:

There is currently no wifi hotspot, and as we want to resolve the equation for a multiple number of position next to the left wall, let's do a for loop:

```plaintext
int bx;
for (bx = 1; bx <= 7; bx++){
    border C(t=0, 2*pi) { x=2+cos(t); y=bx*5+sin(t); label=2; }
    mesh Th = buildmesh(a00(10*n) + a10(10*n) + a20(10*n) + a30(10*n) + a01(10*n) + a11(10*n) + a21(10*n) + a31(10*n) + b00(5*n) + b10(5*n) + b20(5*n) + b30(5*n) + c00(5*n) + c10(5*n) + c20(5*n) + c30(5*n));
```

The border \texttt{C} is our hotspot and as you can see a simple circle. \texttt{Th} is our final mesh, with all borders and the hotspot. Let's resolve this equation!

```plaintext
fespace Vh(Th, P1);
func real wall() {
    (continues on next page)

2.19. Wifi Propagation
```
Fig. 2.20: Mesh

(continued from previous page)

```cpp
if (Th(x,y).region == Th(0.5,0.5).region || Th(x,y).region == Th(7,20.5).region ||
    Th(x,y).region == Th(30.5,2).region) { return 1; }
else { return 0; }
```

```cpp
Vh<complex> v,w;
randinit(900);
Vh wallreflexion = randreal1();
Vh<complex> wallabsorption = randreal1()*0.5i;
Vh k = 6;
```

```cpp
cout << "Reflection of walls : " << wallreflexion << "\n";
cout << "Absorption of walls : " << wallabsorption << "\n";
```

```cpp
problem muwave(v,w) =
int2d(Th)(
    (v*w*k^2)/(1+(wallreflexion+wallabsorption)*wall())^2
    - (dx(v)*dx(w)+dy(v)*dy(w))
) + on(2, v=1)
```

```cpp
muwave;
Vh vm = log(real(v)^2 + imag(v)^2);
plot(vm, wait=1, fill=true, value=0, nbiso=65);
```

A bit of understanding here:

- The `fespace` keyword defines a finite elements space, no need to know more here.
- The function `wall` return 0 if in air and 1 if in a wall (x and y are global variables).
• For this example, random numbers are used for the reflection and the absorption.
• The problem is defined with `problem` and we solve it by calling it.

Finally, I plotted the log of the module of the solution $v$ to see the signal’s power, and here we are:

![Fig. 2.21: Solution](image)

Beautiful isn’t it? This is the first position for the hotspot, but there are 6 others, and the electrical field is evolving depending on the position. You can see the other positions here:

### 2.20 Plotting in Matlab and Octave

#### 2.20.1 Overview

In order to create a plot of a FreeFEM simulation in Matlab© or Octave two steps are necessary:

- The mesh, the finite element space connectivity and the simulation data must be exported into files
- The files must be imported into the Matlab / Octave workspace. Then the data can be visualized with the `ffmatlib` library

The steps are explained in more detail below using the example of a stripline capacitor.

**Note:** Finite element variables must be in P1 or P2. The simulation data can be 2D or 3D.

#### 2.20.2 2D Problem

Consider a stripline capacitor problem which is also shown in Fig. 2.23. On the two boundaries (the electrodes) $C_A$, $C_K$ a Dirichlet condition and on the enclosure $C_B$ a Neumann condition is set. The electrostatic potential $u$ between
Fig. 2.22: Wifi propagation
the two electrodes is given by the Laplace equation

$$\Delta u(x, y) = 0$$

and the electrostatic field $E$ is calculated by

$$E = -\nabla u$$

```plaintext
int CA=3, CK=4, CB=5;
real w2=1.0, h=0.4, d2=0.5;

border bottomA(t=-w2,w2) { x=t; y=d2; label=CA; };
border rightA(t=d2,d2+h) { x=w2; y=t; label=CA; };
border topA(t=w2,-w2) { x=t; y=d2+h; label=CA; };
border leftA(t=d2+h,d2) { x=-w2; y=t; label=CA; };
border bottomK(t=-w2,w2) { x=t; y=-d2-h; label=CK; };
border rightK(t=-d2-h,-d2) { x=w2; y=t; label=CK; };
border topK(t=w2,-w2) { x=t; y=-d2; label=CK; };
border leftK(t=-d2,-d2-h) { x=-w2; y=t; label=CK; };
border enclosure(t=0,2*pi) { x=5*cos(t); y=5*sin(t); label=CB; }

int n=15;

mesh Th = buildmesh(enclosure(3*n) +
    bottomA(-w2*n)+topA(-w2*n)+rightA(-h*n)+leftA(-h*n)+
    bottomK(-w2*n)+topK(-w2*n)+rightK(-h*n)+leftK(-h*n));

fespace Vh(Th,P1);
Vh u,v;
real u0=2.0;

problem Laplace(u,v,solver=LU) =
    int2d(Th)(dx(u)*dx(v) + dy(u)*dy(v))
    + on(CA,u=u0)+on(CK,u=0);
real error=0.01;
for (int i=0;i<1;i++)
    Laplace;
    Th=adaptmesh(Th,u,err=error);
    error=error/2.0;
}
Laplace;
Vh Ex, Ey;
Ex = -dx(u);
Ey = -dy(u);
plot(u,[Ex,Ey],wait=true);
```

### 2.20.3 Exporting Data

The mesh is stored with the FreeFEM command `savemesh()`, while the connectivity of the finite element space and the simulation data are stored with the macro commands `ffSaveVh()` and `ffSaveData()`. These two commands are located in the `ffmatlib.idp` file which is included in the `ffmatlib`. Therefore, to export the stripline capacitor data the following statement sequence must be added to the FreeFEM code:

#### 2.20. Plotting in Matlab and Octave
```csharp
#include "ffmatlib.idp"

//Save mesh
savemesh(Th,"capacitor.msh");
//Save finite element space connectivity
ffSaveVh(Th,Vh,"capacitor_vh.txt");
//Save some scalar data
ffSaveData(u,"capacitor_potential.txt");
//Save a 2D vector field
ffSaveData2(Ex,Ey,"capacitor_field.txt");
```

### 2.20.4 Importing Data

The mesh file can be loaded into the Matlab / Octave workspace using the `ffreadmesh()` command. A mesh file consists of three main sections:

1. The mesh points as nodal coordinates
2. A list of boundary edges including boundary labels
3. List of triangles defining the mesh in terms of connectivity

The three data sections mentioned are returned in the variables `p`, `b` and `t`. The finite element space connectivity and the simulation data can be loaded using the `ffreaddata()` command. Therefore, to load the example data the following statement sequence must be executed in Matlab / Octave:

```csharp
%Add ffmatlib to the search path
addpath('add here the link to the ffmatlib');
%Load the mesh
[p,b,t,nv,nbe,nt,labels]=ffreadmesh('capacitor.msh');
%Load the finite element space connectivity
vh=ffreaddata('capacitor_vh.txt');
%Load scalar data
u=ffreaddata('capacitor_potential.txt');
%Load 2D vector field data
[Ex,Ey]=ffreaddata('capacitor_field.txt');
```

### 2.20.5 2D Plot Examples

`ffpdeplot()` is a plot solution for creating patch, contour, quiver, mesh, border, and region plots of 2D geometries. The basic syntax is:

```csharp
[handles,varargout] = ffpdeplot(p,b,t,varargin)
```

`varargin` specifies parameter name / value pairs to control the plot behaviour. A table showing all options can be found in the `ffmatlib` documentation. A small selection of possible plot commands is given as follows:

- Plot of the boundary and the mesh:

```csharp
ffpdeplot(p,b,t,'Mesh','on','Boundary','on');
```

- Patch plot (2D map or density plot) including mesh and boundary:

```csharp
ffpdeplot(p,b,t,'VhSeq',vh,'XYData',u,'Mesh','on','Boundary','on', ... 
  'XLim',[-2 2], 'YLim',[-2 2]);
```
Fig. 2.23: Boundary and Mesh

Fig. 2.24: Patch Plot with Mesh
• 3D surf plot:

```matlab
ffpdepplot(p,b,t,'VhSeq',vh,'XYData',u,'ZStyle','continuous', ...
    'Mesh','off');
lighting gouraud;
view([-47,24]);
camlight('headlight');
```

![3D Surf Plot](image)

**Fig. 2.25:** 3D Surf Plot

• Contour (isovalue) and quiver (vector field) plot:

```matlab
ffpdepplot(p,b,t,'VhSeq',vh,'XYData',u,'Mesh','off','Boundary','on', ...
    'XGrid',[-2 2],'YGrid',[-2 2],'Contour','on','CColor','r', ...
    'XYStyle','off','CGridParam',[150,150],'ColorBar','off', ...
    'FlowData',[Ex,Ey],'FGridParam',[24,24]);
```

![Contour and Quiver Plot](image)

**Fig. 2.26:** Contour and Quiver Plot

Download run through example:

Matlab / Octave file
FreeFEM script

2.20.6 3D Plot Examples

3D problems are handled by the `ffpdeplot3D()` command, which works similarly to the `ffpdeplot()` command. In particular in three-dimensions cross sections of the solution can be created. The following example shows a cross-sectional problem of a three-dimensional parallel plate capacitor.

![3D Plot Example](image)

*Fig. 2.27: Slice on a 3D Parallel Plate Capacitor*

Download run through example:
Matlab / Octave file
FreeFEM script

2.20.7 References

- Octave
- Matlab
- ffmatlib
CHAPTER
THREE

DOCUMENTATION

The fruit of a long maturing process, freefem, in its last avatar, FreeFEM, is a high level integrated development environment (IDE) for numerically solving partial differential equations (PDE) in dimension 1, 2, 3 and surface and line 3D. It is the ideal tool for teaching the finite element method but it is also perfect for research to quickly test new ideas or multi-physics and complex applications.

FreeFEM has an advanced automatic mesh generator, capable of a posteriori mesh adaptation; it has a general purpose elliptic solver interfaced with fast algorithms, such as the multi-frontal method UMFPACK, SuperLU, MUMPS. Hyperbolic and parabolic problems are solved by iterative algorithms prescribed by the user with the high level language of FreeFEM. It has several triangular finite elements, including discontinuous elements. Everything is there in FreeFEM to prepare research quality reports with online color display, zooming and other features as well as postscript printouts.

This manual is meant for students at a Masters level, for researchers at any level, and for engineers (including financial engineering) with some understanding of variational methods for partial differential equations.

Introduction

A partial differential equation is a relation between a function of several variables and its (partial) derivatives. Many problems in physics, engineering, mathematics and even banking are modeled by one or several partial differential equations.

FreeFEM is a software to solve these equations numerically. As its name implies, it is a free software (see the copyrights for full detail) based on the Finite Element Method; it is not a package, it is an integrated product with its own high level programming language. This software runs on all UNIX OS (with g++ 3.3 or later, and OpenGL), on Window XP, Vista and 7, 8, 10 and on MacOS 10 intel.

Moreover FreeFEM is highly adaptive. Many phenomena involve several coupled systems. Fluid-structure interactions, Lorentz forces for aluminum casting and ocean-atmosphere problems are three such systems. These require different finite element approximations and polynomial degrees, possibly on different meshes. Some algorithms like the Schwarz’ domain decomposition method also requires data interpolation on multiple meshes within one program. FreeFEM can handle these difficulties, i.e. arbitrary finite element spaces on arbitrary unstructured and adapted bi-dimensional meshes.

The characteristics of FreeFEM are:

- Problem description (real or complex valued) by their variational formulations, with access to the internal vectors and matrices if needed.
- Multi-variables, multi-equations, bi-dimensional and three-dimensional static or time dependent, linear or non-linear coupled systems; however the user is required to describe the iterative procedures which reduce the problem to a set of linear problems.
- Easy geometric input by analytic description of boundaries by pieces; however this part is not a CAD system; for instance when two boundaries intersect, the user must specify the intersection points.
• Automatic mesh generator, based on the Delaunay-Voronoi algorithm; the inner point density is proportional to
the density of points on the boundaries [GEORGE1996].
• Metric-based anisotropic mesh adaptation. The metric can be computed automatically from the Hessian of any
FreeFEM function [HECHT1998].
• High level user friendly typed input language with an algebra of analytic and finite element functions.
• Multiple finite element meshes within one application with automatic interpolation of data on different meshes
and possible storage of the interpolation matrices.
• A large variety of triangular finite elements: linear, quadratic Lagrangian elements and more, discontinuous P1
and Raviart-Thomas elements, elements of a non-scalar type, the mini-element,... (but no quadrangles).
• Tools to define discontinuous Galerkin finite element formulations P0, P1dc, P2dc and keywords: jump, mean,
intalledges.
• A large variety of linear direct and iterative solvers (LU, Cholesky, Crout, CG, GMRES, UMFPACK, MUMPS,
SuperLU, ... ) and eigenvalue and eigenvector solvers (ARPACK).
• Near optimal execution speed (compared with compiled C++ implementations programmed directly).
• Online graphics, generation of .txt,.eps,.gnu, mesh files for further manipulations of input and output data.
• Many examples and tutorials: elliptic, parabolic and hyperbolic problems, Navier-Stokes flows, elasticity, fluid
structure interactions, Schwarz’s domain decomposition method, eigenvalue problem, residual error indicator,
... 
• A parallel version using MPI

3.1 Notations

Here mathematical expressions and corresponding FreeFEM commands are explained.

3.1.1 Generalities

• \([\delta_{ij}]\) Kronecker delta (0 if \(i \neq j\), 1 if \(i = j\) for integers \(i, j\))
• \([\forall]\) for all
• \([\exists]\) there exists
• [i.e.] that is
• [PDE] partial differential equation (with boundary conditions)
• [0] the empty set
• [\(\mathbb{N}\)] the set of integers \((a \in \mathbb{N} \Leftrightarrow \text{int } a)\), int means long int inside FreeFEM
• [\(\mathbb{R}\)] the set of real numbers \((a \in \mathbb{R} \Leftrightarrow \text{real } a)\), double inside FreeFEM
• [\(\mathbb{C}\)] the set of complex numbers \((a \in \mathbb{C} \Leftrightarrow \text{complex } a)\), complex<double>
• [\(\mathbb{R}^d\)] \(d\)-dimensional Euclidean space
### 3.1.2 Sets, Mappings, Matrices, Vectors

Let $E, F, G$ be three sets and $A$ the subset of $E$.

- $\{x \in E | P\}$ the subset of $E$ consisting of the elements possessing the property $P$
- $[E \cup F]$ the set of elements belonging to $E$ or $F$
- $[E \cap F]$ the set of elements belonging to $E$ and $F$
- $[E \setminus A]$ the set $\{x \in E | x \notin A\}$
- $[E + F]$ $E \cup F$ with $E \cap F = \emptyset$
- $[E \times F]$ the Cartesian product of $E$ and $F$
- $[E^n]$ the $n$-th power of $E (E^2 = E \times E, E^n = E \times E^{n-1})$
- $[f : E \to F]$ the mapping form $E$ into $F$, i.e., $E \ni x \mapsto f(x) \in F$
- $[I_E \text{ or } I]$ the identity mapping in $E$, i.e., $I(x) = x \ \forall x \in E$
- $[f \circ g]$ for $f : F \to G$ and $g : E \to F$, $E \ni x \mapsto (f \circ g)(x) = f(g(x)) \in G$ (see Elementary function)
- $[f|A]$ the restriction of $f : E \to F$ to the subset $A$ of $E$
- $[\{a_k\}]$ column vector with components $a_k$
- $[\langle a_k \rangle]$ row vector with components $a_k$
- $[\langle a_k \rangle^T]$ denotes the transpose of a matrix $(a_k)$, and is $\{a_k\}$
- $[\{a_{ij}\}]$ matrix with components $a_{ij}$, and $(a_{ij})^T = (a_{ji})$

### 3.1.3 Numbers

For two real numbers $a, b$

- $[a, b]$ is the interval $\{x \in \mathbb{R} | a \leq x \leq b\}$
- $]a, b]$ is the interval $\{x \in \mathbb{R} | a < x \leq b\}$
- $[a, b[$ is the interval $\{x \in \mathbb{R} | a \leq x < b\}$
- $]a, b[$ is the interval $\{x \in \mathbb{R} | a < x < b\}$

### 3.1.4 Differential Calculus

- $[\partial f / \partial x]$ the partial derivative of $f : \mathbb{R}^d \to \mathbb{R}$ with respect to $x$ ($dx(ξ)$)
- $[\nabla f]$ the gradient of $f : \Omega \to \mathbb{R}$, i.e., $\nabla f = (\partial f / \partial x, \partial f / \partial y)$
- $[\text{div}(f) \text{ or } \nabla \cdot f]$ the divergence of $f : \Omega \to \mathbb{R}^d$, i.e., $\text{div}(f) = \partial f_1 / \partial x + \partial f_2 / \partial y$
- $[\Delta f]$ the Laplacian of $f : \Omega \to \mathbb{R}$, i.e., $\Delta f = \partial^2 f / \partial x^2 + \partial^2 f / \partial y^2$

### 3.1.5 Meshes

- $[\Omega]$ usually denotes a domain on which PDE is defined
- $[\Gamma]$ denotes the boundary of $\Omega$, i.e., $\Gamma = \partial \Omega$ (keyword border, see Border)
• \([\mathcal{T}_h]\) the triangulation of \(\Omega\), i.e., the set of triangles \(T_k\), where \(h\) stands for mesh size (keyword \texttt{mesh}, \texttt{buildmesh}, see \textit{Mesh Generation})
• \([n_t]\) the number of triangles in \(\mathcal{T}_h\) (get by \texttt{Th.nt})
• \([\Omega_h]\) denotes the approximated domain \(\Omega_h = \bigcup_{k=1}^{n_t} T_k\) of \(\Omega\). If \(\Omega\) is polygonal domain, then it will be \(\Omega = \Omega_h\)
• \([\Gamma_h]\) the boundary of \(\Omega_h\)
• \([n_v]\) the number of vertices in \(\mathcal{T}_h\) (get by \texttt{Th.nv})
• \([n_b]\) the number of boundary element in \(\mathcal{T}_h\) (get by \texttt{Th.nbe})
• \([|\Omega_h|]\) the measure (area or volume) in \(\mathcal{T}_h\) (get by \texttt{Th.measure})
• \([|\partial\Omega_h|]\) the measure of the border (length or area) in \(\mathcal{T}_h\) (get by \texttt{Th.bordermeasure})
• \([h_{\text{min}}]\) the minimum edge size of \(\mathcal{T}_h\) (get by \texttt{Th.hmin})
• \([h_{\text{max}}]\) the maximum edge size of \(\mathcal{T}_h\) (get by \texttt{Th.hmax})
• \([q_i q_j]\) the segment connecting \(q_i\) and \(q_j\)
• \([q_{k1}, q_{k2}, q_{k3}]\) the vertices of a triangle \(T_k\) with anti-clock direction (get the coordinate of \(q_{kj}\) by \texttt{(Th[k-1][j-1].x, Th[k-1][j-1].y)})
• \([I_{\Omega}]\) the set \(\{i \in \mathbb{N} | q_i \notin \Gamma_h\}\)

### 3.1.6 Functional Spaces

• \([L^2(\Omega)]\) the set \(\left\{w(x,y) \left| \int_{\Omega} |w(x,y)|^2 \, dx \, dy < \infty \right\} \)

\[
\text{norm: } ||w||_{0, \Omega} = \left(\int_{\Omega} |w(x,y)|^2 \, dx \, dy \right)^{1/2}
\]

\[
\text{scalar product: } (v, w) = \int_{\Omega} vw
\]

• \([H^1(\Omega)]\) the set \(\left\{w \in L^2(\Omega) \left| \int_{\Omega} (|\partial w/\partial x|^2 + |\partial w/\partial y|^2) \, dx \, dy < \infty \right\} \)

\[
\text{norm: } ||w||_{1, \Omega} = (||w||_{0, \Omega}^2 + ||\nabla w||_{0, \Omega}^2)^{1/2}
\]

• \([H^m(\Omega)]\) the set \(\left\{w \in L^2(\Omega) \left| \int_{\Omega} \frac{\partial^{|\alpha|} w}{\partial x^{\alpha_1} \partial y^{\alpha_2}} \in L^2(\Omega), \forall \alpha = (\alpha_1, \alpha_2) \in \mathbb{N}^2, |\alpha| = \alpha_1 + \alpha_2 \right\} \)

\[
\text{scalar product: } (v, w)_{1, \Omega} = \sum_{|\alpha| \leq m} \int_{\Omega} D^\alpha v D^\alpha w
\]

• \([H_0^1(\Omega)]\) the set \(\{w \in H^1(\Omega) \mid u = 0 \text{ on } \Gamma\}\)

\([L^2(\Omega)^2]\) denotes \(L^2(\Omega) \times L^2(\Omega)\), and also \(H^1(\Omega)^2 = H^1(\Omega) \times H^1(\Omega)\)
3.1.7 Finite Element Spaces

- \([V_h]\) denotes the finite element space created by \texttt{fespace Vh(Th, *)} in \texttt{FreeFEM} (see \textit{Finite Elements} for \(*\))

- \([\Pi_h f]\) the projection of the function \(f\) into \(V_h\) means \(v = P_{ih}(f) \cdot \{v\}\) for FE-function \(v\) in \(V_h\) means the column vector \((v_1, \ldots, v_M)^T\) if \(v = v_1 \phi_1 + \cdots + v_M \phi_M\), which is shown by \texttt{fespace Vh(Th, P2); Vh v; cout << v[] << endl;}

3.2 Mesh Generation

In this section, operators and tools on meshes are presented.

FreeFEM type for mesh variable:

- 1D mesh: \texttt{meshL}
- 2D mesh: \texttt{mesh}
- 3D volume mesh: \texttt{mesh3}
- 3D border meshes
  - 3D surface \texttt{meshS}
  - 3D curve \texttt{meshL}

Through this presentation, the principal commands for the mesh generation and links between \texttt{mesh} - \texttt{mesh3} - \texttt{meshS} - \texttt{meshL} are described.

3.2.1 The type mesh in 2 dimension

Commands for 2d mesh Generation

The \texttt{FreeFEM} type to define a 2d mesh object is \texttt{mesh}.

The command \texttt{square}

The command \texttt{square} triangulates the unit square.

The following generates a \(4 \times 5\) grid in the unit square \([0, 1]^2\). The labels of the boundaries are shown in Fig. 3.1.

```
mesh Th = square(4, 5);
```

To construct a \(n \times m\) grid in the rectangle \([x_0, x_1] \times [y_0, y_1]\), proceed as follows:

```
real x0 = 1.2;
real x1 = 1.8;
real y0 = 0;
real y1 = 1;
int n = 5;
real m = 20;
mesh Th = square(n, m, [x0+(x1-x0)*x, y0+(y1-y0)*y]);
```

\textbf{Note:} Adding the named parameter \texttt{flags=icase} with \texttt{icase}:
Fig. 3.1: Boundary labels of the mesh by square(10,10)

0. will produce a mesh where all quads are split with diagonal \( x - y = \text{constant} \)
1. will produce a Union Jack flag type of mesh
2. will produce a mesh where all quads are split with diagonal \( x + y = \text{constant} \)
3. same as in case 0, except two corners where the triangles are the same as case 2, to avoid having 3 vertices on the boundary
4. same as in case 2, except two corners where the triangles are the same as case 0, to avoid having 3 vertices on the boundary

```c
mesh Th = square(n, m, [x0+(x1-x0)*x0, y0+(y1-y0)*y1], flags=icase);
```

Note: Adding the named parameter `label=\text{labs}` will change the 4 default label numbers to `\text{labs[i-1]}`, for example `\text{int[int] labs=[11, 12, 13, 14]}`, and adding the named parameter `region=10` will change the region number to 10, for instance (v 3.8).

To see all of these flags at work, check Square mesh example:

```c
for (int i = 0; i < 5; ++i){
  int[int] labs = [11, 12, 13, 14];
  mesh Th = square(3, 3, \text{flags=i}, label=\text{labs}, region=10);
  plot(Th, \text{wait=1, cmm="square flags = "+i });
}
```

The command buildmesh

mesh building with border

Boundaries are defined piecewise by parametrized curves. The pieces can only intersect at their endpoints, but it is possible to join more than two endpoints. This can be used to structure the mesh if an area touches a border and create new regions by dividing larger ones:
int upper = 1;
int others = 2;
int inner = 3;

border C01(t=0, 1){x=0; y=-1+t; label=upper;}
border C02(t=0, 1){x=1.5-1.5*t; y=-1; label=upper;}
border C03(t=0, 1){x=1.5; y=-t; label=upper;}
border C04(t=0, 1){x=1+0.5*t; y=0; label=others;}
border C05(t=0, 1){x=0.5+0.5*t; y=0; label=others;}
border C06(t=0, 1){x=0.5*t; y=-0.5; label=inner;}
border C11(t=0, 1){x=0.5; y=-0.5*t; label=inner;}
border C12(t=0, 1){x=0.5+0.5*t; y=-0.5; label=inner;}
border C13(t=0, 1){x=1; y=-0.5+0.5*t; label=inner;}

int n = 10;
plot(C01(-n) + C02(-n) + C03(-n) + C04(-n) + C05(-n) + C06(-n) + C11(n) + C12(n) + C13(n), wait=true);

mesh Th = buildmesh(C01(-n) + C02(-n) + C03(-n) + C04(-n) + C05(-n) + C06(-n) + C11(n) + C12(n) + C13(n), wait=true);
plot(Th, wait=true);

cout << "Part 1 has region number " << Th(0.75, -0.25).region << endl;
cout << "Part 2 has region number " << Th(0.25, -0.25).region << endl;

Borders and mesh are respectively shown in Fig. 3.2a and Fig. 3.2b.

Triangulation keywords assume that the domain is defined as being on the left (resp right) of its oriented parameterized boundary

\[ \Gamma_j = \{(x,y) \mid x = \varphi_x(t), y = \varphi_y(t), a_j \leq t \leq b_j\} \]

To check the orientation plot \( t \mapsto (\varphi_x(t), \varphi_y(t)), t_0 \leq t \leq t_1 \). If it is as in Fig. 3.3, then the domain lies on the shaded area, otherwise it lies on the opposite side.

The general expression to define a triangulation with \texttt{buildmesh} is

```
mesh Mesh_Name = buildmesh(Gamma1(m1)+...+GammaJ(mj), OptionalParameter);
```
where \( m_j \) are positive or negative numbers to indicate how many vertices should be on \( \Gamma_j \), \( \Gamma = \bigcup_{j=1}^{J} \Gamma_j \), and the optional parameter (see also References), separated with a comma, can be:

- \( \text{nbvtx} = \text{int} \), to set the maximum number of vertices in the mesh.
- \( \text{fixedborder} = \text{bool} \), to say if the mesh generator can change the boundary mesh or not (by default the boundary mesh can change; beware that with periodic boundary conditions (see. Finite Element), it can be dangerous.

The orientation of boundaries can be changed by changing the sign of \( m_j \).

The following example shows how to change the orientation. The example generates the unit disk with a small circular hole, and assigns “1” to the unit disk (“2” to the circle inside). The boundary label **must be non-zero**, but it can also be omitted.

```plaintext
border a(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
border b(t=0, 2*pi){x=0.3+0.3*cos(t); y=0.3*sin(t); label=2;}
plot(a(50) + b(30)); // to see a plot of the border mesh
mesh Thwithouthole = buildmesh(a(50) + b(30));
mesh Thwithhole = buildmesh(a(50) + b(-30));
plot(Thwithouthole, ps="Thwithouthole.eps");
plot(Thwithhole, ps="Thwithhole.eps");
```

**Note:** Notice that the orientation is changed by \( b(-30) \) in the 5th line. In the 7th line, \( \text{ps}="\text{fileName}" \) is used to generate a postscript file with identification shown on the figure.

**Note:** Borders are evaluated only at the time `plot` or `buildmesh` is called so the global variables are defined at this time. In this case, since \( r \) is changed between the two border calls, the following code will not work because the first border will be computed with \( r=0.3 \):

```plaintext
real r=1;
border a(t=0, 2*pi){x=r*cos(t); y=r*sin(t); label=1;}
r=0.3;
border b(t=0, 2*pi){x=r*cos(t); y=r*sin(t); label=1;}
mesh Thwithhole = buildmesh(a(50) + b(-30)); // bug (a trap) because
// the two circles have the same radius = :math:`0.3`
```

**mesh building with array of border**
Sometimes it can be useful to make an array of the border, but unfortunately it is incompatible with the FreeFEM syntax. To bypass this problem, if the number of segments of the discretization \( n \) is an array, we make an implicit loop on all of the values of the array, and the index variable \( i \) of the loop is defined after the parameter definition, like in

\[
\text{border } a(t=0, 2*\pi; i)\ldots
\]

A first very small example:

```plaintext
1 border a(t=0, 2*pi; i){
2 \hspace{1cm} x=(i+1)*cos(t);
3 \hspace{1cm} y=(i+1)*sin(t);
4 \hspace{1cm} label=1;}
5 int[int] nn = [10, 20, 30];
6 plot(a(nn)); //plot 3 circles with 10, 20, 30 points
```

And a more complex example to define a square with small circles:

```plaintext
1 real[int] xx = [0, 1, 1, 0],
2 yy = [0, 0, 1, 1];
3 //radius, center of the 4 circles
4 real[int] RC = [0.1, 0.05, 0.05, 0.1],
5 XC = [0.2, 0.8, 0.2, 0.8],
6 YC = [0.2, 0.8, 0.8, 0.2];
7 int[int] NC = [-10,-11,-12,13]; //list number of \( \pm \) segments of the 4 circles borders
8
9 border bb(t=0, 1; i) {
10 \hspace{1cm} // i is the index variable of the multi border loop
11 \hspace{1cm} \hspace{1cm} int ii = (i+1)%4;
12 \hspace{1cm} \hspace{1cm} real t1 = 1-t;
13 \hspace{1cm} \hspace{1cm} x = xx[i]*t1 + xx[ii]*t;
14 \hspace{1cm} \hspace{1cm} y = yy[i]*t1 + yy[ii]*t;
15 \hspace{1cm} \hspace{1cm} label = 0;
16 \hspace{1cm} }
17
18 border cc(t=0, 2*pi; i) {
19 \hspace{1cm} (continues on next page)
```

Fig. 3.4: Mesh with a hole
\[ x = RC[i] \cdot \cos(t) + XC[i]; \]
\[ y = RC[i] \cdot \sin(t) + YC[i]; \]
\[ \text{label} = i + 1; \]

`int[int] nn = [4, 4, 5, 7]; // 4 border, with 4, 4, 5, 7 segment respectively`
`plot(bb(nn), cc(NC), \text{wait}=1);`
`mesh \text{th} = \text{buildmesh}(bb(nn) + cc(NC));`
`plot(th, \text{wait}=1);`

### Mesh Connectivity and data

The following example explains methods to obtain mesh information.

```cpp
// Mesh
mesh Th = square(2, 2);

cout << "// Get data of the mesh" << endl;
{
    int NbTriangles = Th.nt;
    real MeshArea = Th.measure;
    real BorderLength = Th.bordermeasure;

    cout << "Number of triangle(s) = " << NbTriangles << endl;
    cout << "Mesh area = " << MeshArea << endl;
    cout << "Border length = " << BorderLength << endl;
}

// Th(i) return the vertex i of Th
// Th[k] return the triangle k of Th
// Th[k][i] return the vertex i of the triangle k of Th
for (int i = 0; i < NbTriangles; i++)
    for (int j = 0; j < 3; j++)
        cout << i << " " << j << " - Th[i][j] = " << Th[i][j]
            << " x = " << Th[i][j].x
            << " y = " << Th[i][j].y
            << " label=" << Th[i][j].label << endl;
}

// Hack to get vertex coordinates" << endl;
{
    fespace fem1(Th, P1);
    fem1 Thx=x,Thy=y;

    int NbVertices = Th.nv;
    cout << "Number of vertices = " << NbVertices << endl;

    for (int i = 0; i < NbVertices; i++)
        cout << "Th(" << i << ") : " << Th[i].x << " Th(i).y << " " << Th(i).
            ->label
            << endl << \"old method: " << Thx[i] << " Thy[i]" << endl;
}

// Method to find information of point (0.55,0.6)" << endl;
{
    int TNumber = Th(0.55, 0.6).nuTriangle; //the triangle number
    int RLabel = Th(0.55, 0.6).region; //the region label

    (continues on next page)"
cout << "Triangle number in point (0.55, 0.6): " << TNumber << endl;
cout << "Region label in point (0.55, 0.6): " << RLabel << endl;
}

cout << "// Information of triangle" << endl;
{
    int TNumber = Th(0.55, 0.6).nuTriangle;
    real TArea = Th[TNumber].area; //triangle area
    real TRegion = Th[TNumber].region; //triangle region
    real TLabel = Th[TNumber].label; //triangle label, same as region for triangles

    cout << "Area of triangle " << TNumber <<": " << TArea << endl;
    cout << "Region of triangle " << TNumber <<": " << TRegion << endl;
    cout << "Label of triangle " << TNumber <<": " << TLabel << endl;
}

cout << "// Hack to get a triangle containing point x, y or region number (old method) " << endl;
{
    fespace fem0(Th, P0);
    fem0 TNumbers; //a P0 function to get triangle numbering
    for (int i = 0; i < Th.nt; i++)
        TNumbers[i][i] = i;
    fem0 RNumbers = region; //a P0 function to get the region number

    int TNumber = TNumbers(0.55, 0.6); // Number of the triangle containing (0.55, 0, 6)
    int RNumber = RNumbers(0.55, 0.6); // Number of the region containing (0.55, 0, 6)

cout << "Point (0.55,0.6) :": " << endl;
    cout << "Triangle number = " << TNumber << endl;
    cout << "Region number = " << RNumber << endl;
}

cout << "// New method to get boundary information and mesh adjacent" << endl;
{
    int k = 0;
    int l=1;
    int e=1;

    // Number of boundary elements
    int NbBoundaryElements = Th.nbe;
    cout << "Number of boundary element = " << NbBoundaryElements << endl;
    // Boundary element k in {0, ..., Th.nbe}
    int BoundaryElement = Th.be(k);
    cout << "Boundary element " << k << " = " << BoundaryElement << endl;
    // Vertice l in {0, 1} of boundary element k
    int Vertex = Th.be(k)[l];
    cout << "Vertex " << l << " of boundary element " << k << " = " << Vertex << endl;
    // Triangle containing the boundary element k
    int Triangle = Th.be(k).Element;
    cout << "Triangle containing the boundary element " << k << " = " << Triangle << endl;
    // Triangle edge number containing the boundary element k
    int Edge = Th.be(k).whoInElement;
    cout << "Triangle edge number containing the boundary element " << k << " = " << Edge << endl;
}
// Adjacent triangle of the triangle k by edge e
int Adjacent = Th[k].adj(e); // The value of e is changed to the corresponding
   // edge in the adjacent triangle
   cout << "Adjacent triangle of the triangle " << k << " by edge " << e << " = " << Adjacent
   // The value of e is changed to the corresponding
   Adjacent << endl;
   cout << "\tCorresponding edge = " << e << endl;
   // If there is no adjacent triangle by edge e, the same triangle is returned
   // Th[k] == Th[k].adj(e)
   // Else a different triangle is returned
   // Th[k] != Th[k].adj(e)
}

// Print mesh connectivity
{
    int NbTriangles = Th.nt;
    for (int k = 0; k < NbTriangles; k++)
        cout << k << " : "
             << int(Th[k][0]) << " "
             << int(Th[k][1])
             << " , label "
             << Th[k].label << endl;

    for (int k = 0; k < NbTriangles; k++)
        for (int e = 0, ee; e < 3; e++)
            // set ee to e, and ee is change by method adj,
            cout << k << " "
                 << ee << " => "
                 << int(Th[k].adj((ee=e))) << " "
                 << ee
                 << ", adj: "
                 << (Th[k].adj((ee=e)) != Th[k]) << endl;

    int NbBoundaryElements = Th.nbe;
    for (int k = 0; k < NbBoundaryElements; k++)
        cout << k << " : "
             << Th.be(k)[0] << " "
             << Th.be(k)[1]
             << " , label "
             << Th.be(k).label
             << " , triangle "
             << int(Th.be(k).Element)
             << " , "
             << Th.be(k).whoinElement << endl;

    real[int] bb(4);
    boundingbox(Th, bb);
    cout << "boundingbox:" << endl;
    cout << "xmin = " << bb[0]
         << ", xmax = " << bb[1]
         << ", ymin = " << bb[2]
         << ", ymax = " << bb[3] << endl;
}

The output is:

// Get data of the mesh
Number of triangle = 8
Mesh area = 1
Border length = 4
0 0 - Th[i][j] = 0, x = 0, y= 0, label=4
0 1 - Th[i][j] = 1, x = 0.5, y= 0, label=1
0 2 - Th[i][j] = 4, x = 0.5, y= 0.5, label=0
1 0 - Th[i][j] = 0, x = 0, y= 0, label=4
1 1 - Th[i][j] = 4, x = 0.5, y= 0.5, label=0
1 2 - Th[i][j] = 3, x = 0, y= 0.5, label=4
2 0 - Th[i][j] = 1, x = 0.5, y= 0, label=1
2 1 - Th[i][j] = 2, x = 1, y= 0, label=2
FreeFEM Documentation, Release 4.6

3.2. Mesh Generation

// Hack to get vertex coordinates
Number of vertices = 9

// Old method:

// Old method:

// New method to get boundary information and mesh adjacent
Number of boundary element = 8

// Print mesh connectivity
0 : 0 1 4, label 0

// Method to find the information of point (0.55, 0.6)
Point (0.55, 0.6):
  Triangle number = 7
  Region number = 0

// New method to get boundary information and mesh adjacent
Number of boundary element = 8

// Print mesh connectivity
0 : 0 1 4, label 0
The real characteristic function of a mesh is \( \chi(Th) \) in 2D and 3D where:

\[ \chi(Th)(P) = \begin{cases} 1 & \text{if } P \in Th \\ 0 & \text{if } P \notin Th \end{cases} \]

**The keyword “triangulate”**

**FreeFEM** is able to build a triangulation from a set of points. This triangulation is a Delaunay mesh of the convex hull of the set of points. It can be useful to build a mesh from a table function.

The coordinates of the points and the value of the table function are defined separately with rows of the form: \( x \ y \ f(x, y) \) in a file such as:
The third column of each line is left untouched by the `triangulate` command. But you can use this third value to define a table function with rows of the form: \( x \ y \ f(x,y) \).

The following example shows how to make a mesh from the file `xyf` with the format stated just above. The command `triangulate` only uses the 1st and 2nd columns.

```plaintext
// Build the Delaunay mesh of the convex hull
mesh Thxy=triangulate("xyf"); //points are defined by the first 2 columns of file

// Plot the created mesh
plot(Thxy);

// Fespace
fespace Vhxy(Thxy, P1);
Vhxy fxy;

// Reading the 3rd column to define the function fxy
{
    ifstream file("xyf");
    real xx, yy;
    for(int i = 0; i < fxy.n; i++)
        file >> xx >> yy >> fxy[i][i]; //to read third row only.
        //xx and yy are just skipped
}
```

(continues on next page)
// Plot
plot(fxy);

One new way to build a mesh is to have two arrays: one for the \( x \) values and the other for the \( y \) values.

```plaintext
//set two arrays for the x's and y's
Vhxy xx = x, yy = y;
//build the mesh
mesh Th = triangulate(xx[], yy[]);
```

### 2d Finite Element space on a boundary

To define a Finite Element space on a boundary, we came up with the idea of a mesh with no internal points (called empty mesh). It can be useful to handle Lagrange multipliers in mixed and mortar methods.

So the function `emptymesh` removes all the internal points of a mesh except points on internal boundaries.

```plaintext
{| border a(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
| mesh Th = buildmesh(a(20));
| Th = emptymesh(Th);
| plot(Th);
|

It is also possible to build an empty mesh of a pseudo subregion with `emptymesh(Th, ssd)` using the set of edges from the mesh `Th`; an edge \( e \) is in this set when, with the two adjacent triangles \( e = t1 \cap t2 \) and \( ssd[T1] \neq ssd[T2] \) where `ssd` refers to the pseudo region numbering of triangles, they are stored in the `int[int]` array of size "the number of triangles".

```plaintext
{| mesh Th = square(10, 10);
| int[int] ssd(Th.nt);
| //build the pseudo region numbering
| for(int i = 0; i < ssd.n; i++){
|     int iq = i/2; //because 2 triangles per quad
|     int ix = iq%10;
|     int iy = iq/10;
|     ssd[i] = 1 + (ix>=5) + (iy>=5)*2;
| }
| //build empty with all edges $e=T1 \cap T2$ and $ssd[T1] \neq ssd[T2]$
| Th = emptymesh(Th, ssd);
| //plot
| plot(Th);
| savemesh(Th, "emptymesh.msh");
|
```

### Remeshing

#### The command `movemesh`

Meshes can be translated, rotated, and deformed by `movemesh`; this is useful for elasticity to watch the deformation due to the displacement \( \Phi(x, y) = (\Phi_1(x, y), \Phi_2(x, y)) \) of shape.
It is also useful to handle free boundary problems or optimal shape problems. If \( \Omega \) is triangulated as \( T_h(\Omega) \), and \( \Phi \) is a displacement vector then \( \Phi(T_h) \) is obtained by:

```plaintext
mesh Th = movemesh(Th, [Phi1, Phi2]);
```

Sometimes the transformed mesh is invalid because some triangles have flipped over (meaning it now has a negative area). To spot such problems, one may check the minimum triangle area in the transformed mesh with `checkmovemesh` before any real transformation.

For example:

\[
\begin{align*}
\Phi_1(x, y) &= x + k \times \sin(y \times \pi) / 10 \\
\Phi_2(x, y) &= y + k \times \cos(y \pi) / 10
\end{align*}
\]

for a big number \( k > 1 \).

```plaintext
verbosity = 4;

// Parameters
real coef = 1;

// Mesh
border a(t=0, 1) {x=t; y=0; label=1;};
border b(t=0, 0.5) {x=1-t; y=t; label=1;};
border c(t=0, 0.5) {x=1-t; y=0.5; label=1;};
border d(t=0.5, 1) {x=0.5; y=t; label=1;};
border e(t=0.5, 1) {x=1-t; y=1; label=1;};
border f(t=0, 1) {x=0; y=1-t; label=1;};
mesh Th = buildmesh(a(6) + b(4) + c(4) + d(4) + e(4) + f(6));
plot(Th, wait=true, fill=true, ps="Lshape.eps");

// Function
func uu = sin(y*pi)/10;
func vv = cos(x*pi)/10;

// Checkmovemesh
real minT0 = checkmovemesh(Th, [x, y]); //return the min triangle area
while(1){ // find a correct move mesh
    real minT = checkmovemesh(Th, [x*coef=uu, y*coef=vv]);
    if (minT > minT0/5) break; //if big enough
}```
Note: Consider a function \( u \) defined on a mesh \( \text{Th} \). A statement like \( \text{Th} = \text{movemesh}(\text{Th}...) \) does not change \( u \) and so the old mesh still exists. It will be destroyed when no function uses it. A statement like \( u = u \) redefines \( u \) on the new mesh \( \text{Th} \) with interpolation and therefore destroys the old \( \text{Th} \), if \( u \) was the only function using it.

Now, we give an example of moving a mesh with a Lagrangian function \( u \) defined on the moving mesh.
### The command `hTriangle`

This section presents the way to obtain a regular triangulation with FreeFEM.

For a set $S$, we define the diameter of $S$ by

$$\text{diam}(S) = \sup\{|x - y|; x, y \in S\}$$

The sequence $\{\mathcal{T}_h\}_{h \to 0}$ of $\Omega$ is called regular if they satisfy the following:

1. $\lim_{h \to 0} \max\{\text{diam}(T_k)| T_k \in \mathcal{T}_h\} = 0$
2. There is a number $\sigma > 0$ independent of $h$ such that $\frac{\rho(T_k)}{\text{diam}(T_k)} \geq \sigma$ for all $T_k \in \mathcal{T}_h$ where $\rho(T_k)$ are the diameter of the inscribed circle of $T_k$.

We put $h(\mathcal{T}_h) = \max\{\text{diam}(T_k)| T_k \in \mathcal{T}_h\}$, which is obtained by

```cpp
mesh Th = .........;
fespace Ph(Th, P0);
Ph h = hTriangle;
cout << "size of mesh = " << h[].max << endl;
```

### The command `adaptmesh`

The function:

$$f(x, y) = 10.0x^3 + y^3 + \tan^{-1}[\frac{\varepsilon}{(\sin(5.0y) - 2.0x)}], \varepsilon = 0.0001$$

sharply varies in value and the initial mesh given by one of the commands in the Mesh Generation part cannot reflect its sharp variations.

```cpp
// Parameters
real eps = 0.0001;
real h = 1;
 realtime hmin = 0.05;
func f = 10.0*x^3 + y^3 + h*atan2(eps, sin(5.0*y)-2.0*x);

// Mesh
mesh Th = square(5, 5, [-1+2*x, -1+2*y]);
```

### 3.2. Mesh Generation

(continued on next page)
FreeFEM uses a variable metric/Delaunay automatic meshing algorithm.

The command:

```plaintext
mesh ATh = adaptmesh(Th, f);
```

create the new mesh ATh adapted to the Hessian

\[
D^2 f = \left( \frac{\partial^2 f}{\partial x^2}, \frac{\partial^2 f}{\partial x \partial y}, \frac{\partial^2 f}{\partial y^2} \right)
\]

of a function (formula or FE-function).

Mesh adaptation is a very powerful tool when the solution of a problem varies locally and sharply.

Here we solve the Poisson's problem, when \( f = 1 \) and \( \Omega \) is an L-shape domain.

Tip: The solution has the singularity \( r^{3/2} \), \( r = |x - \gamma| \) at the point \( \gamma \) of the intersection of two lines \( bc \) and \( bd \) (see Fig. 3.9a).
FreeFEM Documentation, Release 4.6

3.2. Mesh Generation

```
// Parameters
real error = 0.1;

// Mesh
border ba(t=0, 1){x=t; y=0; label=1;}
border bb(t=0, 0.5){x=1; y=t; label=1;}
border bc(t=0, 0.5){x=1-t; y=0.5; label=1;}
border bd(t=0.5, 1){x=0.5; y=t; label=1;}
border be(t=0.5, 1){x=1-t; y=1; label=1;}
border bf(t=0, 1){x=0; y=1-t; label=1;}
mesh Th = buildmesh(ba(6) + bb(4) + bc(4) + bd(4) + be(4) + bf(6));

// Fespace
fespace Vh(Th, P1);
Vh u, v;

// Function
func f = 1;

// Problem
problem Poisson(u, v, solver=CG, eps=1.e-6)
  = int2d(Th)(
    dx(u)*dx(v)
    + dy(u)*dy(v)
  )
  - int2d(Th)(
    f*v
  )
  + on(1, u=0);

// Adaptmesh loop
for (int i = 0; i < 4; i++){
  Poisson;
}
```

Fig. 3.9: Mesh adaptation
To speed up the adaptation, the default parameter \texttt{err} of \texttt{adaptmesh} is changed by hand; it specifies the required precision, so as to make the new mesh finer or coarser.

The problem is coercive and symmetric, so the linear system can be solved with the conjugate gradient method (parameter \texttt{solver=CG}) with the stopping criteria on the residual, here \texttt{eps=1.e-6}).

By \texttt{adaptmesh}, the slope of the final solution is correctly computed near the point of intersection of \texttt{bc} and \texttt{bd} as in Fig. 3.9b.

This method is described in detail in [HECHT1998]. It has a number of default parameters which can be modified. If \texttt{f1, f2} are functions and \texttt{thold, Thnew} are meshes:

\begin{verbatim}
1 Thnew = adaptmesh(Thold, f1 ... );
2 Thnew = adaptmesh(Thold, f1,f2 ... );
3 Thnew = adaptmesh(Thold, [f1,f2] ... );
\end{verbatim}

The additional parameters of \texttt{adaptmesh} are:

See \textit{Reference part} for more informations

- \texttt{hmin}=\textbf{Minimum edge size}. Its default is related to the size of the domain to be meshed and the precision of the mesh generator.
- \texttt{hmax}=\textbf{Maximum edge size}. It defaults to the diameter of the domain to be meshed.
- \texttt{err}=\textbf{P}_1 interpolation error level (0.01 is the default).
- \texttt{errg}=\textbf{Relative geometrical error}. By default this error is 0.01, and in any case it must be lower than \(1/\sqrt{2}\). Meshes created with this option may have some edges smaller than the \(-\texttt{hmin}\) due to geometrical constraints.
- \texttt{nbvx}=\textbf{Maximum number of vertices generated by the mesh generator} (9000 is the default).
- \texttt{nbsmooth}=\textbf{number of iterations of the smoothing procedure} (5 is the default).
- \texttt{nbjacy}=\textbf{number of iterations in a smoothing procedure} during the metric construction, 0 means no smoothing, 6 is the default.
- \texttt{ratio}=\textbf{ratio for a prescribed smoothing on the metric}. If the value is 0 or less than 1.1 no smoothing is done on the metric. 1.8 is the default. If \texttt{ratio > 1.1}, the speed of mesh size variations is bounded by \(\log(\texttt{ratio})\).

\textbf{Note:} As \texttt{ratio} gets closer to 1, the number of generated vertices increases. This may be useful to control the thickness of refined regions near shocks or boundary layers.

- \texttt{omega}=\textbf{relaxation parameter for the smoothing procedure}. 1.0 is the default.
- \texttt{iso}=\textbf{If true, forces the metric to be isotropic}. \texttt{false} is the default.
• **abserror**= If false, the metric is evaluated using the criteria of equi-repartition of relative error. false is the default. In this case the metric is defined by:

\[ \mathcal{M} = \left( \frac{1}{\text{err \ coef}^2} \frac{|\mathcal{H}|}{\max(CutOff, |\eta|)} \right)^p \]

Otherwise, the metric is evaluated using the criteria of equi-distribution of errors. In this case the metric is defined by:

\[ \mathcal{M} = \left( \frac{1}{\text{err \ coef}^2} \frac{|\mathcal{H}|}{\sup(\eta) - \inf(\eta)} \right)^p . \]

• **cutoff**= lower limit for the relative error evaluation. 1.0e-6 is the default.

• **verbosity**= informational messages level (can be chosen between 0 and \( \infty \)). Also changes the value of the global variable verbosity (obsolete).

• **inquire**= To inquire graphically about the mesh. false is the default.

• **splitpbedge**= If true, splits all internal edges in half with two boundary vertices. true is the default.

• **maxsubdiv**= Changes the metric such that the maximum subdivision of a background edge is bound by \( \text{val} \).

  Always limited by 10, and 10 is also the default.

• **rescaling**= if true, the function, with respect to which the mesh is adapted, is rescaled to be between 0 and 1. true is the default.

• **keepbackvertices**= if true, tries to keep as many vertices from the original mesh as possible. true is the default.

• **IsMetric**= if true, the metric is defined explicitly. false is the default. If the 3 functions \( m_{11}, m_{12}, m_{22} \) are given, they directly define a symmetric matrix field whose Hessian is computed to define a metric. If only one function is given, then it represents the isotropic mesh size at every point.

  For example, if the partial derivatives \( f_{xx} = \frac{\partial^2 f}{\partial x^2}, f_{xy} = \frac{\partial^2 f}{\partial x \partial y}, f_{yy} = \frac{\partial^2 f}{\partial y^2} \) are given, we can set \( \text{Th} = \text{adaptmesh} \)(\text{Th}, f_{xx}, f_{xy}, f_{yy}, \text{IsMetric}=1, \text{nbvx}=10000, \text{hmin}=\text{hmin}) ; \)

• **power**= exponent power of the Hessian used to compute the metric. 1 is the default.

• **thetamax**= minimum corner angle in degrees. Default is 10° where the corner is \( ABC \) and the angle is the angle of the two vectors \( AB, BC \), (0 imply no corner, 90 imply perpendicular corner, ...).

• **splitin2**= boolean value. If true, splits all triangles of the final mesh into 4 sub-triangles.

• **metric**= an array of 3 real arrays to set or get metric data information. The size of these three arrays must be the number of vertices. So if \( m_{11}, m_{12}, m_{22} \) are three P1 finite elements related to the mesh to adapt, you can write: \( \text{metric}=[m_{11}[],m_{12}[],m_{22}[]] \) (see file convecapt.edp for a full example)

• **nomeshgeneration**= If true, no adapted mesh is generated (useful to compute only a metric).

• **periodic**= Writing periodic=[\([4, y]\), \([2, y]\), \([1, x]\), \([3, x]\)]; builds an adapted periodic mesh.

  The sample builds a biperiodic mesh of a square. (see periodic finite element spaces, and see the Sphere example for a full example)

We can use the command adaptmesh to build a uniform mesh with a constant mesh size. To build a mesh with a constant mesh size equal to \( \frac{1}{30} \) try:
mesh Th=square(2, 2); //the initial mesh
plot(Th, wait=true, ps="square-0.eps");

Th = adaptmesh(Th, 1./30., IsMetric=1, nbvx=10000);
plot(Th, wait=true, ps="square-1.eps");

Th = adaptmesh(Th, 1./30., IsMetric=1, nbvx=10000); //More the one time du to
Th = adaptmesh(Th, 1./30., IsMetric=1, nbvx=10000); //Adaptation bound 'maxsubdiv='
plot(Th, wait=true, ps="square-2.eps");

The command trunc

Two operators have been introduced to remove triangles from a mesh or to divide them. Operator trunc has the following parameters:

- boolean function to keep or remove elements
- label= sets the label number of new boundary item, one by default.
- split= sets the level of triangle splitting. Each triangle is split in $n \times n$, one by default.

To create the mesh Th3 where all triangles of a mesh Th are split in $3 \times 3$, just write:

mesh Th3 = trunc(Th, 1, split=3);

The following example construct all “truncated” meshes to the support of the basic function of the space $Vh$ (cf. $\text{abs}(u)>0$), split all the triangles in $5 \times 5$, and put a label number to 2 on a new boundary.

// Mesh
mesh Th = square(3, 3);

// Fespace
fespace Vh(Th, P1);
Vh u=0;

// Loop on all degrees of freedom
int n=u.n;
for (int i = 0; i < n; i++){
    u[i][i] = 1; // The basis function i
    plot(u, wait=true);
    mesh Sh1 = trunc(Th, abs(u)>1.e-10, split=5, label=2);
    plot(Th, Sh1, wait=true, ps="trunc"+i".eps");
    u[i][i] = 0; // reset
}

The command change

This command changes the label of elements and border elements of a mesh.

Changing the label of elements and border elements will be done using the keyword change. The parameters for this command line are for two dimensional and three dimensional cases:

- refe= is an array of integers to change the references on edges
- reft= is an array of integers to change the references on triangles
- label= is an array of integers to change the 4 default label numbers
Fig. 3.10: Mesh adaptation

(a) Initial mesh  
(b) First iteration  
(c) Last iteration

Fig. 3.11: Trunc

(a) Mesh of support the function P1 number 0, split in $5 \times 5$  
(b) Mesh of support the function P1 number 6, split in $5 \times 5$
• region= is an array of integers to change the default region numbers
• renumv= is an array of integers, which explicitly gives the new numbering of vertices in the new mesh. By default, this numbering is that of the original mesh
• renumt= is an array of integers, which explicitly gives the new numbering of elements in the new mesh, according the new vertices numbering given by renumv=. By default, this numbering is that of the original mesh
• flabel= is an integer function given the new value of the label
• fregion= is an integer function given the new value of the region
• rmledges= is an integer to remove edges in the new mesh, following a label
• rmInternalEdges= is a boolean, if equal to true to remove the internal edges. By default, the internal edges are stored

These vectors are composed of $n_l$ successive pairs of numbers $O, N$ where $n_l$ is the number (label or region) that we want to change. For example, we have:

\[
label = [O_1, N_1, ..., O_{n_l}, N_{n_l}]
\]
\[
region = [O_1, N_1, ..., O_{n_l}, N_{n_l}]
\]

An application example is given here:

```plaintext
1 // Mesh
2 mesh Th1 = square(10, 10);
3 mesh Th2 = square(20, 10, [x+1, y]);
4 int[int] r1=[2,0];
5 plot(Th1, wait=true);
6 Th1 = change(Th1, label=r1); //change the label of Edges 2 in 0.
7 plot(Th1, wait=true);
8 // boundary label: 1 -> 1 bottom, 2 -> 1 right, 3->1 top, 4->1 left boundary label is 1
9 int[int] re=[1,1, 2,1, 3,1, 4,1]
10 Th2=change(Th2, re);
11 plot(Th2, wait=1); 
```

The command `splitmesh`

Another way to split mesh triangles is to use `splitmesh`, for example:

```plaintext
1 // Mesh
2 border a(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
3 mesh Th = buildmesh(a(20));
4 plot(Th, wait=true, ps="NotSplittedMesh.eps");
5 // Splitmesh
6 Th = splitmesh(Th, 1 + 5*(square(x-0.5) + y^2));
7 plot(Th, wait=true, ps="SplittedMesh.eps");
```

Meshing Examples
Fig. 3.12: Split mesh

Tip: Two rectangles touching by a side

```cpp
border a(t=0, 1){x=t; y=0;};
border b(t=0, 1){x=1; y=t;};
border c(t=1, 0){x=t; y=1;};
border d(t=1, 0){x=0; y=t;};
border cl(t=0, 1){x=t; y=1;};
border e(t=0, 0.2){x=1; y=1+t;};
border f(t=1, 0){x=t; y=1.2;};
border g(t=0.2, 0){x=0; y=1+t;};
int n=1;

mesh th = buildmesh(a(10*n) + b(10*n) + c(10*n) + d(10*n));
mesh TH = buildmesh(cl(10*n) + e(5*n) + f(10*n) + g(5*n));
plot(th, TH, ps="TouchSide.esp");
```

Tip: NACA0012 Airfoil

```cpp
border upper(t=0, 1){x=t; y=0.17735*sqrt(t) - 0.075597*t - 0.212836*(t^2) + 0.17363*(t^3) + 0.06254*(t^4);}
border lower(t=1, 0){x = t; y = -(0.17735*sqrt(t) - 0.075597*t - 0.212836*(t^2) + 0.17363*(t^3) + 0.06254*(t^4));}
border c(t=0, 2*pi){x=0.8*cos(t) + 0.5; y=0.8*sin(t);}
mesh Th = buildmesh(c(30) + upper(35) + lower(35));
plot(Th, ps="NACA0012.eps", bw=true);
```

Tip: Cardioid

3.2. Mesh Generation
Fig. 3.13: Two rectangles touching by a side

Fig. 3.14: NACA0012 Airfoil
real b = 1, a = b;
border C(t=0, 2*pi){x=(a+b)*cos(t)-b*cos((a+b)*t/b); y=(a+b)*sin(t)-b*sin((a+b)*t/b);}
mesh Th = buildmesh(C(50));
plot(Th, ps="Cardioid.eps", bw=true);

Fig. 3.15: Domain with Cardioid curve boundary

Tip: Cassini Egg

border C(t=0, 2*pi) {x=(2*cos(2*t)+3)*cos(t); y=(2*cos(2*t)+3)*sin(t);}
mesh Th = buildmesh(C(50));
plot(Th, ps="Cassini.eps", bw=true);

Fig. 3.16: Domain with Cassini egg curve boundary

Tip: By cubic Bezier curve

// A cubic Bezier curve connecting two points with two control points
func real bzi(real p0, real p1, real q1, real q2, real t)
return p0*(1-t)^3 + q1*3*(1-t)^2*t + q2*3*(1-t)*t^2 + p1*t^3;

real[int] p00 = [0, 1], p01 = [0, -1], q00 = [-2, 0.1], q01 = [-2, -0.5];
real[int] p11 = [1,-0.9], q10 = [0.1, -0.95], q11= [0.5, -1];
real[int] p21 = [2, 0.7], q20 = [3, -0.4], q21 = [4, 0.5];
real[int] q30 = [0.5, 1.1], q31 = [1.5, 1.2];

border G1(t=0, 1){
  x=bzi(p00[0], p01[0], q00[0], q01[0], t);
  y=bzi(p00[1], p01[1], q00[1], q01[1], t);
}

border G2(t=0, 1){
  x=bzi(p01[0], p11[0], q10[0], q11[0], t);
  y=bzi(p01[1], p11[1], q10[1], q11[1], t);
}

border G3(t=0, 1){
  x=bzi(p11[0], p21[0], q20[0], q21[0], t);
  y=bzi(p11[1], p21[1], q20[1], q21[1], t);
}

border G4(t=0, 1){
  x=bzi(p21[0], p00[0], q30[0], q31[0], t);
  y=bzi(p21[1], p00[1], q30[1], q31[1], t);
}

int m = 5;

mesh Th = buildmesh(G1(2*m) + G2(m) + G3(3*m) + G4(m));
plot(Th, ps="Bezier.eps", bw=true);

---

**Tip:** Section of Engine

real a = 6., b = 1., c = 0.5;

border L1(t=0, 1){x=-a; y=1+b-2*(1+b)*t;}
border L2(t=0, 1){x=a+2*a*t; y=-1-b*(x/a)*(x/a)*(3-2*abs(x)/a ) ;}
border L3(t=0, 1){x=a; y=-1-b*(1+b)*t; }
border L4(t=0, 1){x=-a-t; y=0;}
border L5(t=0, pi){x=c+sin(t)/2; y=c/2-c*cos(t)/2;}
border L6(t=0, 1){x=a+t; y=c;}
border L7(t=0, 1){x=a; y=c+(1+b-c)*t;}
border L8(t=0, 1){x=a-2*a*t; y=1+b*(x/a)*(x/a)*(3-2*abs(x)/a );}

---

Fig. 3.17: Boundary drawn by Bezier curves
FreeFEM Documentation, Release 4.6

(continued from previous page)

```
12 plot(Th, ps="Engine.eps", bw=true);
```

**Fig. 3.18:** Section of Engine

---

**Tip:** Domain with U-shape channel

```
1 real d = 0.1; //width of U-shape
2 border L1(t=0, 1-d){x=-1; y=-d-t;}
3 border L2(t=0, 1-d){x=-1; y=1-t;}
4 border B(t=0, 2){x=-1-t; y=-1;}
5 border C1(t=0, 1){x=t-1; y=d;}
6 border C2(t=0, 2+d){x=0; y=d-t;}
7 border C3(t=0, 1){x=-t; y=-d;}
8 border R(t=0, 2){x=1; y=-1+t;}
9 border T(t=0, 2){x=1-t; y=1;}
10 int n = 5;
11 mesh Th = buildmesh(L1(n/2) + L2(n/2) + B(n) + C1(n) + C2(3) + C3(n) + R(n) + T(n));
12 plot(Th, ps="U-shape.eps", bw=true);
```

**Fig. 3.19:** Domain with U-shape channel changed by $d$

---

**Tip:** Domain with V-shape cut

3.2. Mesh Generation 121
real dAg = 0.02; // angle of V-shape
border C(t=dAg, 2*pi-dAg){x=cos(t); y=sin(t);} real[int] pa(2), pb(2), pc(2);
pa[0] = cos(dAg);
pa[1] = sin(dAg);
pb[0] = cos(2*pi-dAg);
pb[1] = sin(2*pi-dAg);
pc[0] = 0;
pc[1] = 0;
border seg1(t=0, 1){x=(1-t)*pb[0]+t*pc[0]; y=(1-t)*pb[1]+t*pc[1];}
border seg2(t=0, 1){x=(1-t)*pc[0]+t*pa[0]; y=(1-t)*pc[1]+t*pa[1];}
mesh Th = buildmesh(seg1(20) + C(40) + seg2(20));
plot(Th, ps="V-shape.eps", bw=true);

Fig. 3.20: Domain with V-shape cut changed by dAg

Tip: Smiling face

real d=0.1; int m = 5; real a = 1.5, b = 2, c = 0.7, e = 0.01;
border F(t=0, 2*pi){x=a*cos(t); y=b*sin(t);}
border E1(t=0, 2*pi){x=0.2*cos(t)-0.5; y=0.2*sin(t)+0.5;}
border E2(t=0, 2*pi){x=0.2*cos(t)+0.5; y=0.2*sin(t)+0.5;}
func real st(real t){
  return sin(pi*t) - pi/2;
}
border C1(t=0.5, 0.5){x=(1-t)*c*cos(st(t)); y=(1-t)*c*sin(st(t));}
border C2(t=0, 1){x=((1-d)+t)*c*cos(st(0.5)); y=((1-d)+t)*c*sin(st(0.5));}
border C3(t=0.5, -0.5){x=c*cos(st(t)); y=c*sin(st(t));}
border C4(t=0, 1){x=(1-t)*c*cos(st(-0.5)); y=(1-t)*c*sin(st(-0.5));}
border C0(t=0, 2*pi){x=0.1*cos(t); y=0.1*sin(t);}
mesh Th = buildmesh(F(10*m) + C1(2*m) + C2(3) + C3(2*m) + C4(3) + C0(m) + E1(-2*m) + E2(-2*m));
plot(Th, ps="SmileFace.eps", bw=true);

Fig. 3.21: Smiling face (Mouth is changeable)

Tip: 3 points bending

```
// Square for Three-Point Bend Specimens fixed on Fix1, Fix2
// It will be loaded on Load.

real a = 1, b = 5, c = 0.1;
int n = 5, m = b*n;

border Left(t=0, 2*a){x=-b; y=a-t;}
border Bot1(t=0, b/2-c){x=-b+t; y=-a;}
border Fix1(t=0, 2*c){x=-b/2-c+t; y=-a;}
border Bot2(t=0, b-2*c){x=-b/2+c+t; y=-a;}
border Fix2(t=0, 2*c){x=b/2-c+t; y=-a;}
border Bot3(t=0, b/2-c){x=b/2+c-t; y=-a;}
border Right(t=0, 2*a){x=b; y=-a+t;}
border Top1(t=0, b-c){x=b-t; y=a;}
border Load(t=0, 2*c){x=c-t; y=a;}
border Top2(t=0, b-c){x=-c-t; y=a;}

mesh Th = buildmesh(Left(n) + Bot1(m/4) + Fix1(5) + Bot2(m/2) + Fix2(5) + Bot3(m/4) + Right(n) + Top1(m/2) + Load(10) + Top2(m/2));
plot(Th, ps="ThreePoint.eps", bw=true);
```
3.2.2 The type mesh3 in 3 dimension

Note: Up to the version 3, FreeFEM allowed to consider a surface problem such as the PDE is treated like boundary conditions on the boundary domain (on triangles describing the boundary domain). With the version 4, in particular 4.2.1, a completed model for surface problem is possible, with the definition of a surface mesh and a surface problem with a variational form on domain (with triangle elements) and application of boundary conditions on border domain (describing by edges). The keywords to define a surface mesh is meshS.

3d mesh generation

Note: For 3D mesh tools, put load "msh3" at the top of the .edp script.

The command cube

The function cube like its 2d function square is a simple way to build cubic objects, it is contained in plugin msh3 (import with load "msh3").

The following code generates a $3 \times 4 \times 5$ grid in the unit cube $[0, 1]^3$.

```plaintext
mesh3 Th = cube(3, 4, 5);
```

By default the labels are:
1. face $y = 0$,
2. face $x = 1$,
3. face $y = 1$,
4. face $x = 0$,
5. face $z = 0$,
6. face $z = 1$

and the region number is 0.

A full example of this function to build a mesh of cube $[-1, 1]^3$ with face label given by $(ix + 4*(iy + 1) + 16*(iz + 1))$ where $(ix, iy, iz)$ are the coordinates of the barycenter of the current face, is given below.
```cpp
load "msh3"

int[int] l6 = [37, 42, 45, 40, 25, 57];
int r11 = 11;
mesh3 Th = cube(4, 5, 6, [x*2-1, y*2-1, z*2-1], label=16, flags =3, region=r11);
cout << "Volume = " << Th.measure <<", border area = " << Th.bordermeasure << endl;

int err = 0;
for(int i = 0; i < 100; ++i){
    real s = int2d(Th,i)(1.);
    real sx = int2d(Th,i)(x);
    real sy = int2d(Th,i)(y);
    real sz = int2d(Th,i)(z);
    if(s){
        int ix = (sx/s+1.5);
        int iy = (sy/s+1.5);
        int iz = (sz/s+1.5);
        int ii = (ix + 4*(iy+1) + 16*(iz+1) );
        //value of ix,iy,iz => face min 0, face max 2, no face 1
        cout << "Label = " << i <<", s = " << s <<" \ s " << ix << iy << iz << " : " ;<< iii <<"\n"
    }
}
real volr11 = int3d(Th,r11)(1.);
cout << "Volume region = " << 11 << " : " << volr11 << endl;
if((volr11 - Th.measure )>1e-8) err++;
plot(Th, fill=false);
cout << "Nb err = " << err << endl;
assert(err==0);
```

The output of this script is:

```plaintext
Enter: BuildCube
kind = 3 n tet Cube = 6 / n slip 6 19
Cube  nv=210 nt=720 nbe=296
Out: BuildCube
Volume = 8, border area = 24
Label = 25, s = 4 110 : 25
Label = 37, s = 4 101 : 37
Label = 40, s = 4 011 : 40
Label = 42, s = 4 211 : 42
Label = 45, s = 4 121 : 45
Label = 57, s = 4 112 : 57
Volume region = 11: 8
Nb err = 0
```

The command **buildlayers**

This mesh is obtained by extending a two dimensional mesh in the z-axis.

The domain $\Omega_{3d}$ defined by the layer mesh is equal to $\Omega_{3d} = \Omega_{2d} \times [z_{min}, z_{max}]$ where $\Omega_{2d}$ is the domain defined by the two dimensional meshes. $z_{min}$ and $z_{max}$ are functions of $\Omega_{2d}$ in $\mathbb{R}$ that defines respectively the lower surface and upper surface of $\Omega_{3d}$.

### 3.2. Mesh Generation
Fig. 3.23: The 3D mesh of function `cube(4, 5, 6, flags=3)`

Fig. 3.24: Example of Layer mesh in three dimensions.
For a vertex of a two dimensional mesh \( V_i^{2d} = (x_i, y_i) \), we introduce the number of associated vertices in the \( z \)-axis \( M_i + 1 \).

We denote by \( M \) the maximum of \( M_i \) over the vertices of the two dimensional mesh. This value is called the number of layers (if \( \forall i, M_i = M \) then there are \( M \) layers in the mesh of \( \Omega_{3d} \)). \( V_i^{2d} \) generated \( M + 1 \) vertices which are defined by:

\[
\forall j = 0, \ldots, M, \quad V_{i,j}^{3d} = (x_i, y_i, \theta_i(z_{i,j}))
\]

where \( (z_{i,j})_{j=0,\ldots,M} \) are the \( M + 1 \) equidistant points on the interval \([z\min(V_i^{2d}), z\max(V_i^{2d})]\):

\[
z_{i,j} = j \delta \alpha + z\min(V_i^{2d}), \quad \delta \alpha = \frac{z\max(V_i^{2d}) - z\min(V_i^{2d})}{M}.
\]

The function \( \theta_i \), defined on \([z\min(V_i^{2d}), z\max(V_i^{2d})]\), is given by:

\[
\theta_i(z) = \begin{cases} 
\theta_i,0 & \text{if } z = z\min(V_i^{2d}), \\
\theta_i,j & \text{if } z \in [\theta_i,j-1, \theta_i,j],
\end{cases}
\]

with \( (\theta_i,j)_{j=0,\ldots,M} \) are the \( M_i + 1 \) equidistant points on the interval \([z\min(V_i^{2d}), z\max(V_i^{2d})]\).

Set a triangle \( K = (V_{i,j}^{3d}, V_{i,j+1}^{3d}, V_{i+1,j+1}^{3d}) \) of the two dimensional mesh. \( K \) is associated with a triangle on the upper surface (resp. on the lower surface) of layer mesh:

\((V_{i,j}^{3d}, V_{i+1,j}^{3d}, V_{i+1,j+1}^{3d})\) (resp. \((V_{i+1,j}^{3d}, V_{i+1,j+1}^{3d}, V_{i,j+1}^{3d})\)).

Also \( K \) is associated with \( M \) volume prismatic elements which are defined by:

\[
\forall j = 0, \ldots, M, \quad H_j = (V_{i,j}^{3d}, V_{i,j+1}^{3d}, V_{i+1,j+1}^{3d}, V_{i+1,j}^{3d}, V_{i,j+1}^{3d}).
\]

These volume elements can have some merged point:

- 0 merged point : prism
- 1 merged points : pyramid
- 2 merged points : tetrahedra
- 3 merged points : no elements

The elements with merged points are called degenerate elements. To obtain a mesh with tetrahedra, we decompose the pyramid into two tetrahedra and the prism into three tetrahedra. These tetrahedra are obtained by cutting the quadrilateral face of pyramid and prism with the diagonal which have the vertex with the maximum index (see [HECHT1992] for the reason of this choice).

The triangles on the middle surface obtained with the decomposition of the volume prismatic elements are the triangles generated by the edges on the border of the two dimensional mesh. The label of triangles on the border elements and tetrahedra are defined with the label of these associated elements.

The arguments of \texttt{buildlayers} is a two dimensional mesh and the number of layers \( M \).

The parameters of this command are:

- \texttt{zbound = \texttt{[zmin, zmax]}} where \texttt{zmin} and \texttt{zmax} are functions expression. These functions define the lower surface mesh and upper mesh of surface mesh.

- \texttt{coef = A function expression between [0,1]}. This parameter is used to introduce degenerate element in mesh.

The number of associated points or vertex \( V_i^{2d} \) is the integer part of \( \texttt{coef}(V_i^{2d})M \).

- \texttt{region=} This vector is used to initialize the region of tetrahedra.

This vector contains successive pairs of the 2d region number at index \( 2i \) and the corresponding 3d region number at index \( 2i + 1 \), like \texttt{change}.
• **labelmid**= This vector is used to initialize the 3d labels number of the vertical face or mid face from the 2d label number.

This vector contains successive pairs of the 2d label number at index \(2i\) and the corresponding 3d label number at index \(2i + 1\), like *change*.

• **labelup**= This vector is used to initialize the 3d label numbers of the upper/top face from the 2d region number.

This vector contains successive pairs of the 2d region number at index \(2i\) and the corresponding 3d label number at index \(2i + 1\), like *change*.

• **labeldown**= Same as the previous case but for the lower/down face label.

Moreover, we also add post processing parameters that allow to moving the mesh. These parameters correspond to parameters *transfo*, *facemerge* and *ptmerge* of the command line *movemesh*.

The vector *region*, *labelmid*, *labelup* and *labeldown* These vectors are composed of \(nl\) successive pairs of number \(O_i, N_i\) where \(nl\) is the number (label or region) that we want to get.

An example of this command is given in the *Build layer mesh example*.

**Tip:** Cube

```plaintext
//Cube.idp
load "medit"
load "msh3"

func mesh3 Cube (int[int] &NN, real[int, int] &BB, int[int, int] &L){
  real x0 = BB(0,0), x1 = BB(0,1);
  real y0 = BB(1,0), y1 = BB(1,1);
  real z0 = BB(2,0), z1 = BB(2,1);

  int nx = NN[0], ny = NN[1], nz = NN[2];

  // 2D mesh
  mesh Thx = square(nx, ny, [x0+(x1-x0)*x, y0+(y1-y0)*y]);

  // 3D mesh
  int[int] rup = [0, L(2,1)], rdown=[0, L(2,0)];
  int[int] rmid=[1, L(1,0), 2, L(0,1), 3, L(1,1), 4, L(0,0)];
  mesh3 Th = buildlayers(Thx, nz, zbound=[z0,z1],
    labelmid=rmid, labelup = rup, labeldown = rdown);
  return Th;
}
```

**Tip:** Unit cube

```plaintext
#include "Cube.idp"

int[int] NN = [10,10,10]; //the number of step in each direction
real [int, int] BB = [[0,1],[0,1],[0,1]]; //the bounding box
int [int, int] L = [[1,2],[3,4],[5,6]]; //the label of the 6 face left,right, front,→back, down, right
mesh3 Th = Cube(NN, BB, L);
medit("Th", Th);
```
Fig. 3.25: The mesh of a cube made with cube.epd

Tip: Cone
An axisymtric mesh on a triangle with degenerateness

```plaintext
load "msh3"
load "medit"

// Parameters
real RR = 1;
real HH = 1;
int nn = 10;

// 2D mesh
border Taxe(t=0, HH){x=t; y=0; label=0;}
border Hypo(t=1, 0){x=HH*t; y=RR*t; label=1;}
border Vert(t=0, RR){x=HH; y=t; label=2;}
mesh Th2 = buildmesh(Taxe(HH*nn) + Hypo(sqrt(HH*HH+RR*RR)*nn) + Vert(RR*nn));
plot(Th2, wait=true);

// 3D mesh
real h = 1./nn;
int MaxLayersT = (int(2*pi*RR/h))/4; //number of layers
real zminT = 0;
real zmaxT = 2*pi; //height 2*pi
func fx = y*cos(z);
func fy = y*sin(z);
func fz = x;
int[int] r1T = [0,0], r2T = [0,0,2,2], r4T = [0,2];
//trick function:
```
(continues on next page)
27 //The function defined the proportion
class func deg = max(.01, y/max(x/HH, 0.4)/RR);
29
28 mesh3 Th3T = buildlayers(Th2, coef=deg, MaxLayersT,
29    zbound=[zminT, zmaxT], transfo=[fx, fy, fz],
30    facemerge=0, region=r1T, labelmid=r2T);
31
32 medit("cone", Th3T);

Fig. 3.26: The mesh of a cone made with cone.edp

Tip: Buildlayer mesh

load "msh3"
load "TetGen"
load "medit"

// Parameters
int C1 = 99;
int C2 = 98;

// 2D mesh
border C01(t=0, pi){x=t; y=0; label=1;}
border C02(t=0, 2*pi){x=pi; y=t; label=1;}
border C03(t=0, pi){x=pi-t; y=2*pi; label=1;}
border C04(t=0, 2*pi){x=0; y=2*pi-t; label=1;}

border C11(t=0, 0.7){x=0.5+t; y=2.5; label=C1;}
border C12(t=0, 2){x=1.2; y=2.5+t; label=C1;}
border C13(t=0, 0.7){x=1.2-t; y=4.5; label=C1;}
border C14(t=0, 2){x=0.5; y=4.5-t; label=C1;}

border C21(t=0, 0.7){x=2.3+t; y=2.5; label=C2;}
border C22(t=0, 2){x=3; y=2.5+t; label=C2;}
border C23(t=0, 0.7){x=3-t; y=4.5; label=C2;}
border C24(t=0, 2){x=2.3; y=4.5-t; label=C2;}

(continues on previous page)
```cpp
mesh Th = buildmesh(C01(10) + C02(10) + C03(10) + C04(10)
    + C21(-5) + C22(-5) + C23(-5) + C24(-5));

mesh Ths = buildmesh(C01(10) + C02(10) + C03(10) + C04(10)
    + C11(5) + C12(5) + C13(5) + C14(5));

// Construction of a box with one hole and two regions
func zmin = 0.;
func zmax = 1.;
int MaxLayer = 10;

func XX = x*cos(y);
func YY = x*sin(y);
func ZZ = z;

int[int] r1 = [0, 41], r2 = [98, 98, 99, 99, 1, 56];
int[int] r3 = [4, 12]; // the triangles of upper surface mesh
    // generated by the triangle in the 2D region
    // of mesh Th of label 4 as label 12
int[int] r4 = [4, 45]; // the triangles of lower surface mesh
    // generated by the triangle in the 2D region
    // of mesh Th of label 4 as label 45.

mesh3 Th3 = buildlayers(Th, MaxLayer, zbound=[zmin, zmax], region=r1,
    labelmid=r2, labelup=r3, labeldown=r4);
medit("box 2 regions 1 hole", Th3);

// Construction of a sphere with TetGen
func XX1 = cos(y)*sin(x);
func YY1 = sin(y)*sin(x);
func ZZ1 = cos(x);

real[int] domain = [0., 0., 0., 0., 0.001];
string test = "paACQ";
cout << "test = " "test << endl;

mesh3 Th3sp = tetgtransfo(Ths, transfo=[XX1, YY1, ZZ1],
    switch=test, nbofregions=1, regionlist=domain);
medit("sphere 2 regions", Th3sp);
```

### Remeshing

**Note:** if an operation on a mesh3 is performed then the same operation is applied on its surface part (its meshS associated)

### The command change

This command changes the label of elements and border elements of a mesh. It’s the equivalent command in 2d mesh case.
Changing the label of elements and border elements will be done using the keyword `change`. The parameters for this command line are for two dimensional and three dimensional cases:

- `reftet` is a vector of integer that contains successive pairs of the old label number to the new label number.
- `refface` is a vector of integer that contains successive pairs of the old region number to new region number.
- `flabel` is an integer function given the new value of the label.
- `fregion` is an integer function given the new value of the region.
- `rmInternalFaces` is a boolean, equal true to remove the internal faces.
- `rmlfaces` is a vector of integer, where triangle’s label given are remove of the mesh.

These vectors are composed of \( n_l \) successive pairs of numbers \( O, N \) where \( n_l \) is the number (label or region) that we want to change. For example, we have:

- `label = [O_1, N_1, ..., O_{n_l}, N_{n_l}]`
- `region = [O_1, N_1, ..., O_{n_l}, N_{n_l}]`

An example of use:

```plaintext
1 // Mesh
2 mesh3 Th1 = cube(10, 10);
3 mesh3 Th2 = cube(20, 10, [x+1, y, z]);
4
5 int[int] rl=[2,0];
6 plot(Th1, wait=true);
7 Th1 = change(Th1, label=rl); //change the label of Edges 2 in 0.
8 plot(Th1, wait=true);
9
10 // boundary label: 1 -> 1 bottom, 2 -> 1 right, 3->1 top, 4->1 left boundary label is...
11 int[int] re=[1,1, 2,1, 3,1, 4,1]
12 Th2=change(Th2, re);
13 plot(Th2, wait=1);
```

**The command `trunc`**

This operator have been introduce to remove a piece of mesh or/and split all element or for a particular label element.

The three named parameter - boolean function to keep or remove elements - `split` sets the level \( n \) of triangle splitting, each triangle is splitted in \( n \times n \) (one by default) - freefem:`label` sets the label number of new boundary item (1 by default).

An example of use:

```plaintext
1 load "msh3"
2 load "medit"
3 int nn=8;
4 mesh3 Th=cube(nn,nn,nn);
5 // remove the small cube $\{x<0.5 \mid (y<0.5) \mid (z<0.5)\}$
6 Th= trunc(Th, (x<0.5) \& (y<0.5) \& (z<0.5)), split=3, label=3);
7 medit("cube", Th);
```
The command **movemesh**

3D meshes can be translated, rotated, and deformed using the command line `movemesh` as in the 2D case (see section `movemesh`). If $\Omega$ is tetrahedrized as $T_h(\Omega)$, and $\Phi(x, y) = (\Phi_1(x, y, z), \Phi_2(x, y, z), \Phi_3(x, y, z))$ is the transformation vector then $\Phi(T_h)$ is obtained by:

```plaintext
mesh3 Th = movemesh(Th, [Phi1, Phi2, Phi3], ...);
```

```plaintext
mesh3 Th = movemesh3(Th, transfo=[Phi1, Phi2, Phi3], ...);  // syntax with transfo=
```

The parameters of `movemesh` in three dimensions are:

- **transfo=** sets the geometric transformation $\Phi(x, y) = (\Phi_1(x, y, z), \Phi_2(x, y, z), \Phi_3(x, y, z))$
- **region=** sets the integer labels of the tetrahedra. 0 by default.
- **label=** sets the labels of the border faces. This parameter is initialized as the label for the keyword `change`.
- **facemerge=** An integer expression. When you transform a mesh, some faces can be merged. This parameter equals to one if the merges’ faces is considered. Otherwise it equals to zero. By default, this parameter is equal to 1.
- **ptmerge =** A real expression. When you transform a mesh, some points can be merged. This parameter is the criteria to define two merging points. By default, we use

$$ptmerge = 1e-7 \ Vol(B),$$

where $B$ is the smallest axis parallel boxes containing the discretion domain of $\Omega$ and $Vol(B)$ is the volume of this box.

- **orientation =** An integer expression equal 1, give the orientation of the triangulation, elements must be in the reference orientation (counter clock wise) equal -1 reverse the orientation of the tetrahedra

**Note:** The orientation of tetrahedra are checked by the positivity of its area and automatically corrected during the building of the adjacency.

An example of this command can be found in the *Poisson’s equation 3D example*.

```plaintext
load "medit"
include "cube.idp"
int[int] Nxyz=[20,5,5];
real [int,int] Bxyz=[[0.,5.],[0.,1.],[0.,1.]];
int [int,int] Lxyz=[[1,2],[2,2],[2,2]];
real E = 21.5e4;
real sigma = 0.29;
real mu = E/(2*(1+sigma));
real lambda = E*sigma/((1+sigma)*(1-2*sigma));
real gravity = -0.05;
real sqrt2=sqrt(2.);
mesh3 Th=Cube(Nxyz,Bxyz,Lxyz);
fespace Vh(Th,[P1,P1,P1]);
Vh [u1,u2,u3], [v1,v2,v3];
macro epsilon(u1,u2,u3) [dx(u1),dy(u2),dz(u3),(dz(u2)+dy(u3))/sqrt2,(dz(u1)+dx(u3))/sqrt2,dy(u1)+dx(u2)] // EOM
macro div(u1,u2,u3) ( dx(u1)+dy(u2)+dz(u3) ) // EOM
```

(continues on next page)
solve  Lame([u1,u2,u3],[v1,v2,v3]) =
    int3d(Th) ( lambda*div(u1,u2,u3)*div(v1,v2,v3) 
    + 2.*mu*( epsilon(u1,u2,u3)'*epsilon(v1,v2,v3) ) 
    ) 
    - int3d(Th) (gravity*v3) 
    + on(1,u1=0,u2=0,u3=0);

real  dmax = u1[].max;
real  coef  = 0.1/dmax;

int[int] ref2 = [1,0,2,0]; // array
mesh3 Thm = movemesh(Th, [x+u1*coef,y+u2*coef,z+u3*coef], label=ref2);
// mesh3 Thm = movemesh3(Th, transfo=[x+u1*coef,y+u2*coef,z+u3*coef], label=ref2); older syntax
Thm = change(Thm, label=ref2);
plot(Th, Thm, wait=1, cmm="coef amplification = "+coef");

movemesh doesn’t use the prefix transfo= [...], the geometric transformation is directly given by [...]
the arguments list

**The command extract**

This command offers the possibility to extract a boundary part of a mesh3

- refface, is a vector of integer that contains a list of triangle face references, where the extract function must be apply.

- label, is a vector of integer that contains a list of tetrahedra label

```cpp
load "msh3"
int  nn = 30;
int[int] labs = [1, 2, 2, 1, 1, 2]; // Label numbering
mesh3 Th = cube(nn, nn, nn, label=labs);
// extract the surface (boundary) of the cube
int[int] llabs = [1, 2];
mesh3 ThS = extract(Th, label=llabs);
```

**The command buildSurface**

This new function allows to build the surface mesh of a volume mesh, under the condition the surface is the boundary of the volume. By definition, a mesh3 is defined by a list of vertices, tetrahedron elements and triangle border elements. buildSurface function create the meshS corresponding, given the list vertices which are on the border domain, the triangle elements and build the list of edges. Remark, for a closed surface mesh, the edges list is empty.

**The command movemesh23**

A simple method to tranform a 2D mesh in 3D Surface mesh. The principe is to project a two dimensional domain in a three dimensional space, 2d surface in the (x,y,z)-space to create a surface mesh 3D, meshS.
**Warning:** Since the release 4.2.1, the FreeFEM function `movemesh23` returns a meshS type.

This corresponds to translate, rotate or deform the domain by a displacement vector of this form $\Phi(x, y) = (\Phi_1(x, y), \Phi_2(x, y), \Phi_3(x, y))$.

The result of moving a two dimensional mesh $Th_2$ by this three dimensional displacement is obtained using:

```cpp
**meshS** Th3 = movemesh23(Th2, transfo=[\Phi(1), \Phi(2), \Phi(3)]);```

The parameters of this command line are:

- `transfo=[\Phi_1, \Phi_2, \Phi_3]` sets the displacement vector of transformation $\Phi(x, y) = [\Phi_1(x, y), \Phi_2(x, y), \Phi_3(x, y)]$.
- `label` sets an integer label of triangles.
- `orientation` sets an integer orientation to give the global orientation of the surface of mesh. Equal 1, give a triangulation in the reference orientation (counter clock wise) equal -1 reverse the orientation of the triangles
- `ptmerge` = A real expression. When you transform a mesh, some points can be merged. This parameter is the criteria to define two merging points. By default, we use

$$ptmerge = 1e-7 \ Vol(B),$$

where $B$ is the smallest axis, parallel boxes containing the discretized domain of $\Omega$ and $\Vol(B)$ is the volume of this box.

We can do a “gluing” of surface meshes using the process given in Change section. An example to obtain a three dimensional mesh using the command line `tetg` and `movemesh23` is given below.

```cpp
load "msh3"
load "tetgen"

// Parameters
real x10 = 1.;
real x11 = 2.;
real y10 = 0.;
real y11 = 2.*pi;
func ZZ1min = 0;
func ZZ1max = 1.5;
func XX1 = x;
func YY1 = y;
real x20 = 1.;
real x21 = 2.;
real y20=0.;
real y21=1.5;
func ZZ2 = y;
func XX2 = x;
func YY2min = 0.;
func YY2max = 2.*pi;
real x30=0.;
real x31=2*pi;
real y30=0.;
```

(continues on next page)
real y31=1.5;
func XX3min = 1.;
func XX3max = 2.;
func YY3 = x;
func ZZ3 = y;

// Mesh
mesh Thsq1 = square(5, 35, [x10+(x11-x10)*x, y10+(y11-y10)*y]);
mesh Thsq2 = square(5, 8, [x20+(x21-x20)*x, y20+(y21-y20)*y]);
mesh Thsq3 = square(35, 8, [x30+(x31-x30)*x, y30+(y31-y30)*y]);

// Mesh 2D to 3D surface
meshS Th31h = movemesh23(Thsq1, transfo=[XX1, YY1, ZZ1max], orientation=1);
meshS Th31b = movemesh23(Thsq1, transfo=[XX1, YY1, ZZ1min], orientation=-1);
meshS Th32h = movemesh23(Thsq2, transfo=[XX2, YY2max, ZZ2], orientation=-1);
meshS Th32b = movemesh23(Thsq2, transfo=[XX2, YY2min, ZZ2], orientation=1);
meshS Th33h = movemesh23(Thsq3, transfo=[XX3max, YY3, ZZ3], orientation=1);
meshS Th33b = movemesh23(Thsq3, transfo=[XX3min, YY3, ZZ3], orientation=-1);

// Gluing surfaces
meshS Th33 = Th31h + Th31b + Th32h + Th32b + Th33h + Th33b;
plot(Th33, cmm="Th33");

// Tetrahelize the interior of the cube with TetGen
real[int] domain = [1.5, pi, 0.75, 145, 0.0025];
meshS Thfinal = tetg(Th33, switch="paAAQY", regionlist=domain);
plot(Thfinal, cmm="Thfinal");

// Build a mesh of a half cylindrical shell of interior radius 1, and exterior radius 2 and a height of 1.5
func mv2x = x*cos(y);
func mv2y = x*sin(y);
func mv2z = z;
meshS Thmv2 = movemesh(Thfinal, transfo=[mv2x, mv2y, mv2z], facemerge=0);
plot(Thmv2, cmm="Thmv2");

3d Meshing examples

Tip: Lake
load "msh3"
load "medit"

// Parameters
int nn = 5;

// 2D mesh
border cc(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
mesh Th2 = buildmesh(cc(100));

// 3D mesh
Tip: Hole region

```
load "msh3"
load "TetGen"
load "medit"

// 2D mesh
mesh Th = square(10, 20, [x*pi/2, 2*y*pi]); // ]-pi/2, pi/2[X]0,2pi[

// 3D mesh
//parametrization of a sphere
func f1 = cos(x)*cos(y);
func f2 = cos(x)*sin(y);
func f3 = sin(x);
//partial derivative of the parametrization
func f1x = sin(x)*cos(y);
func f1y = -cos(x)*sin(y);
func f2x = -sin(x)*sin(y);
func f2y = cos(x)*cos(y);
func f3x = cos(x);
func f3y = 0;
//M = DF^t DF
func m11 = f1x^2 + f2x^2 + f3x^2;
func m21 = f1x*f1y + f2x*f2y + f3x*f3y;
func m22 = f1y^2 + f2y^2 + f3y^2;
func perio = [[4, y], [2, y], [1, x], [3, x]];
real hh = 0.1;
real vv = 1/square(hh);
verbosity = 2;
Th = adaptmesh(Th, m11*vv, m21*vv, IsMetric=1, periodic=perio);
Th = adaptmesh(Th, m11*vv, m21*vv, m22*vv, IsMetric=1, periodic=perio);
plot(Th, wait=true);

//construction of the surface of spheres
real Rmin = 1.;
func f1min = Rmin*f1;
func f2min = Rmin*f2;
func f3min = Rmin*f3;
```

(continues on next page)
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```cpp
meshS ThSsph = movemesh23(Th, transfo=[f1min, f2min, f3min]);
real Rmax = 2.0;
func f1max = Rmax*f1;
func f2max = Rmax*f2;
func f3max = Rmax*f3;
meshS ThSsph2 = movemesh23(Th, transfo=[f1max, f2max, f3max]);

//gluing meshes
meshS ThS = ThSsph + ThSsph2;

Tip: Build a 3d mesh of a cube with a balloon
```

```cpp
load "msh3"
load "TetGen"
load "medit"
include "MeshSurface.idp"

// Parameters
real hs = 0.1; //mesh size on sphere
int[int] N = [20, 20, 20];
real [int,int] B = [[-1, 1], [-1, 1], [-1, 1]];
int [int,int] L = [[1, 2], [3, 4], [5, 6]];

// Meshes
meshS ThH = SurfaceHex(N, B, L, 1);
meshS ThS = Sphere(0.5, hs, 7, 1);
meshS ThHS = ThH + ThS;
medit("Hex-Sphere", ThHS);
real voltet = (hs^3)/6.0;

cout << "voltet = " << voltet << endl;
real[int] domain = [0, 0, 0, 1, voltet, 0, 0, 0.7, 2, voltet];

Tip: Build a 3d mesh of a cube with a balloon
```
3.2.3 The type meshS in 3 dimension

**Warning:** Since the release 4.2.1, the surface mesh3 object (list of vertices and border elements, without tetrahedra elements) is remplaced by meshS type.

### Commands for 3d surface mesh generation

#### The command `square3`

The function `square3` like the function `square` in 2d is the simple way to a build the unit square plan in the space $\mathbb{R}^p$. To use this command, it is necessary to load the pluging `msh3` (need load "msh3"). A square in 3d consists in building a 2d square which is projected from $\mathbb{R}^p$ to $\mathbb{R}^q$. The parameters of this command line are:

- `n,m` generates a `n×m` grid in the unit square
- `[.,.,.]` is $[\Phi_1, \Phi_2, \Phi_3]$ is the geometric transformation from $\mathbb{R}^p$ to $\mathbb{R}^q$. By default, $[\Phi_1, \Phi_2, \Phi_3] = [x, y, 0]$
- `orientation= equal 1`, gives the orientation of the triangulation, elements are in the reference orientation (counter clock wise) equal -1 reverse the orientation of the triangles it’s the global orientation of the surface 1 extern (-1 intern)

```c
real R = 3, r=1;
real h = 0.2; //
int nx = R*2*\pi/h;
int ny = r*2*\pi/h;
func torex= (R+r*cos(y*\pi/2))*cos(x*\pi/2);
func torey= (R+r*cos(y*\pi/2))*sin(x*\pi/2);
func torez= r*sin(y*\pi/2);

meshS ThS=square3(nx,ny,[torex,torey,torez],orientation=-1) ;
```

![Image](https://via.placeholder.com/150)

(a) The surface mesh of the hex with internal sphere  
(b) The tetrahedral mesh of the cube with internal ball

Fig. 3.27: Cube sphere
The following code generates a $3 \times 4 \times 5$ grid in the unit cube $[0,1]^3$ with a clockwise triangulation.

### Surface mesh builders

Adding at the top of a FreeFEM script include "MeshSurface.idp", constructors of sphere, ellipsoid, surface mesh of a 3d box are available.

- **SurfaceHex(N, B, L, orient)**
  - this operator allows to build the surface mesh of a 3d box
  - int[int] N=[nx,ny,nz]; // the number of seg in the 3 direction
  - real [int,int] B=[[xmin,xmax],[ymin,ymax],[zmin,zmax]]; // bounding box
  - int [int,int] L=[[1,2],[3,4],[5,6]]; // the label of the 6 face left, right, front, back, down, right
  - orient the global orientation of the surface 1 extern (-1 intern),
  - returns a meshS type

- **Ellipsoide (RX, RY, RZ, h, L, OX, OY, OZ, orient)**
  - h is the mesh size
  - L is the label
  - orient the global orientation of the surface 1 extern (-1 intern)
  - OX, OY, OZ are real numbers to give the Ellipsoide center (optimal, by default is (0,0,0))
  - where RX, RY, RZ are real numbers such as the parametric equations of the ellipsoid is:
  - returns a meshS type

\[
\forall u \in \left[\frac{-\pi}{2}, \frac{\pi}{2}\right] \text{ and } v \in [0, 2\pi], \begin{cases}\ x = Rx \cos(u) \cos(v) + Ox \\
y = Ry \cos(u) \sin(v) + Oy \\
z = Rz \sin(v) + OZ\end{cases}
\]

- **Sphere(R, h, L, OX, OY, OZ, orient)**
  - where R is the radius of the sphere,
  - OX, OY, OZ are real numbers to give the Ellipsoide center (optimal, by default is (0,0,0))
  - h is the mesh size of the sphere
  - L is the label the sphere
  - orient the global orientation of the surface 1 extern (-1 intern)
  - returns a meshS type

```c
func meshS SurfaceHex(int[int] & N, real[int,int] & B, int[int,int] & L, int orientation) {...
    real x0=B(0,0), x1=B(0,1);
    real y0=B(1,0), y1=B(1,1);
    real z0=B(2,0), z1=B(2,1);
    int nx=N[0], ny=N[1], nz=N[2];
    mesh Thx = square(ny,nz, [y0+(y1-y0)*x, z0+(z1-z0)*y]);
    mesh Thy = square(nx,nz, [x0+(x1-x0)*x, z0+(z1-z0)*y]);
}
```

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(continued from previous page)

```c
mesh Thz = square(nx, ny, [x0*(1-x0)*x, y0*(1-y0)*y]);

int[int] refx=[0,L(0,0)], refX=[0,L(0,1)];  // Xmin, Ymax faces labels
int[int] refy=[0,L(1,0)], refY=[0,L(1,1)];  // Ymin, Ymax faces labels
int[int] refz=[0,L(2,0)], refZ=[0,L(2,1)];  // Zmin, Zmax faces labels

meshS Thx0 = movemesh23(Thx, transfo=[x0, x, y], orientation=-orientation, label=refx);
meshS Thx1 = movemesh23(Thx, transfo=[x1, x, y], orientation=+orientation, label=refX);
meshS Thy0 = movemesh23(Thy, transfo=[x, y0, y], orientation=-orientation, label=refy);
meshS Thy1 = movemesh23(Thy, transfo=[x, y1, y], orientation=+orientation, label=refY);
meshS Thz0 = movemesh23(Thz, transfo=[x, y, z0], orientation=-orientation, label=refz);
meshS Thz1 = movemesh23(Thz, transfo=[x, y, z1], orientation=+orientation, label=refZ);

meshS Th = Thx0+Thx1+Thy0+Thy1+Thz0+Thz1;
return Th;
```

```c
}

func meshS Ellipsoide (real RX, real RY, real RZ, real h, int L, real Ox, real Oy, real Oz, int orientation)
{
  mesh Th=square(10,20, [x*pi/2, 2*y*pi]);  // $\pi/2, \pi$
  // a parametrization of a sphere
  func f1 = RX*cos(x)*cos(y);
  func f2 = RY*cos(x)*sin(y);
  func f3 = RZ*sin(x);
  // partial derivative
  func f1x= -RX*sin(x)*cos(y);
  func f1y= -RX*cos(x)*sin(y);
  func f2x= -RY*sin(x)*sin(y);
  func f2y= +RY*cos(x)*cos(y);
  func f3x=-RZ*cos(x);
  func f3y=0;
  // the metric on the sphere $ M = DF^t \cdot DF $
  func m11=f1x^2+f2x^2+f3x^2;
  func m21=f1x*f1y+f2x*f2y+f3x*f3y;
  func m22=f1y^2+f2y^2+f3y^2;
  func perio=[[4,y],[2,y],[1,x],[3,x]];  // to store the periodic condition
  real hh=h;  // hh mesh size on unite sphere
  real vv= 1/square(hh);
  Th=adaptmesh(Th,m11*vv,m21*vv,m22*vv,
              IsMetric=1,periodic=perio);
  Th=adaptmesh(Th,m11+vv,m21+vv,m22+vv,
              IsMetric=1,periodic=perio);
  Th=adaptmesh(Th,m11+vv,m21+vv,m22+vv,
              IsMetric=1,periodic=perio);
  Th=adaptmesh(Th,m11+vv,m21+vv,m22+vv,
              IsMetric=1,periodic=perio);
  int[int] ref=[0,L];
  meshS ThS=movemesh23(Th, transfo=[f1,f2,f3], orientation=orientation, refface=ref);
  ThS=mmgs(ThS, hhmin=h, hhmax=h, hgrad=2.);
  return ThS;
}

func meshS Ellipsoide (real RX, real RY, real RZ, real h, int L, int orientation)
{
  return Ellipsoide (RX, RY, RZ, h, L, 0., 0., 0., orientation);
}

func meshS Sphere(real R, real h, int L, int orientation)
{
  return Ellipsoide(R,R,h,L,orientation);
}
```

(continues on next page)

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func meshS Sphere(real R, real h, int L, real Ox, real Oy, real Oz, int orientation) {
    return Ellipsoide(R, R, h, L, Ox, Oy, Oz, orientation);
}

2D mesh generators combined with movemesh23

FreeFEM’s meshes can be built by the composition of the movemesh23 command from a 2d mesh generation. The operation is a projection of a 2d plane in \( \mathbb{R}^3 \) following the geometric transformation \([ \Phi_1, \Phi_2, \Phi_3 ]\).

```plaintext
load "msh3"
real l = 3;
border a(t=-l,l){x=t; y=-l; label=1;};
border b(t=-l,l){x=l; y=t; label=1;};
border c(t=l,-l){x=t; y=l; label=1;};
border d(t=l,-l){x=-l; y=t; label=1;};
int n = 100;
border i(t=0,2*pi)(x=1.1*cos(t); y=1.1*sin(t); label=5);;
mesh th= buildmesh(a(n)+b(n)+c(n)+d(n)+i(-n));
```

Remeshing

The command trunc

This operator allows to define a meshS by truncating another one, i.e. by removing triangles, and/or by splitting each triangle by a given positive integer s. In a FreeFEM script, this function must be called as follows:

```plaintext
meshS TS2=trunc(TS1, boolean function to keep or remove elements, split = s, label = ...)
```

The command has the following arguments:

- boolean function to keep or remove elements
- split= sets the level n of triangle splitting. each triangle is splitted in n × n ( one by default)
- label= sets the label number of new boundary item (1 by default)

An example of how to call the function

```plaintext
real R = 3, r=1;
real h = 0.2; //
int nx = R*2*pi/h;
int ny = r*2*pi/h;
func torex= (R+r*cos(y*pi*2))*cos(x*pi*2);
func torey= (R+r*cos(y*pi*2))*sin(x*pi*2);
func torez= r*sin(y*pi*2);
// build a tore
meshS ThS=square3(nx,ny,[torex,torey,torez]) ;
ThS=trunc(ThS, (x < 0.5) | (y < 0.5) | (z > 1.), split=4);
```
The command **movemesh**

Like 2d and 3d type meshes in **FreeFEM**, meshS can be translated, rotated or deformed by an application \([\Phi_1, \Phi_2, \Phi_3]\). The image \(T_h(\Omega)\) is obtained by the command **movemeshS**.

The parameters of movemeshS are:

- **transfo=** sets the geometric transformation \(\Phi(x, y) = (\Phi_1(x, y, z), \Phi_2(x, y, z), \Phi_3(x, y, z))\)
- **region=** sets the integer labels of the triangles. 0 by default.
- **label=** sets the labels of the border edges. This parameter is initialized as the label for the keyword `change`.
- **edgemerge=** An integer expression. When you transform a mesh, some triangles can be merged and fix the parameter to 1, else 0. By default, this parameter is equal to 1.
- **ptmerge =** A real expression. When you transform a mesh, some points can be merged. This parameter is the criteria to define two merging points. By default, we use

\[
ptmerge = 1e - 7 \ Vol(B),
\]

where \(B\) is the smallest axis parallel boxes containing the discretion domain of \(\Omega\) and \(Vol(B)\) is the volume of this box.
- **orientation =** An integer expression equal 1, give the orientation of the triangulation, elements must be in the reference orientation (counter clock wise) equal -1 reverse the orientation of the triangles. It’s the global orientation of the normals at the surface 1 extern (-1 intern)

Example of using

```plaintext
meshS Th1 = square3(n, n, [2*x, y, 1], orientation=-1);
meshS Th2=movemeshS(Th1, transf=[x, y, z]);
meshS Th3=movemesh(Th1, [x, y, z]);
```

The command **change**

Equivalent for a 2d or 3d mesh, the command **change** changes the label of elements and border elements of a meshS. The parameters for this command line are:

- **reftri=** is a vector of integer that contains successive pairs of the old label number to the new label number for elements.
- **refedge=** is a vector of integer that contains successive pairs of the old region number to new region number for boundary elements.
- **fillabel=** is an integer function given the new value of the label.
- **fregion=** is an integer function given the new value of the region.
- **rmledges=** is a vector of integer, where edge’s label given are remove of the mesh

These vectors are composed of \(n_l\) successive pairs of numbers \(O, N\) where \(n_l\) is the number (label or region) that we want to change. For example, we have:

\[
\text{label} = [O_1, N_1, \ldots, O_{n_l}, N_{n_l}]
\]
\[
\text{region} = [O_1, N_1, \ldots, O_{n_l}, N_{n_l}]
\]
Link with a mesh3

In topology and mathematics, the boundary of a subset S of a topological space X is the set of points which can be approached both from S and from the outside of S. The general definitions to the boundary of a subset S of a topological space X are:

- the closure of S without the interior of S $\partial S = S \setminus S^o$.
- the intersection of the closure of S with the closure of its complement $\partial S = S \cap (X \setminus S)$.
- the set of points p of X such that every neighborhood of p contains at least one point of S and at least one point not of S.

More concretely in FreeFEM, the gestion of a 3D mesh is as follows. Let be $\Omega$ a subset of $\mathbb{R}^3$ and $\partial \Omega$ is boundary, the finite element discretization $\Omega_h$ of this domain gives:

- a mesh3 type, denotes Th3, meshing the volume domain. It contains all the nodes, the tetrahedrons $\Omega_i$ such as $\Omega_h = \bigcup_i \Omega_i$ and the list of triangles describing the boundary domain
- a meshS type, denotes ThS, meshing the boundary of the volume domain. Typically, containing the nodes belonging to the boundary of Th3 and, if it exists the boundary triangles and the edges.

Remark: Condition of meshS existence | In FreeFEM, a meshS can be defined in 2 cases such as:

- Th3 $\subset$ ThS where it exactly describes the bounder of Th3.
- a meshS is an explicite surface mesh given by a list of vertices, triangle finite elements and boundary edge elements (can be optional follows the geometry domain)

Note: Hence, if an input mesh (.msh freefem or .mesh format) contains a list of vertices, tetrahedra, triangles and edges, FreeFEM builds a mesh3 whitch contains explicitly a surface mesh type meshS.

The command Gamma

The command Gamma allows to build and manipulate the border mesh independly of a volume mesh such as the surface is described by triangle elements and edges border elements in 3d. Use this function, suppose that the mesh3 object even contains the geometric description of its surface. That means, the input mesh explicitly contains the list of vertices, tetrahedra, triangles and edges. In case where the surface mesh doesn’t exist, before calling Gamma, must build it by calling the buildSurface function (see the next function description).

```c
1  load "msh3"
2  int n = 10;
3  int nvb = (n+1)^3 - (n-1)^3; // Nb boundary vertices
4  int ntb = n*n*12; // Nb of Boundary triangle
5  mesh3 Th=cube(n,n,n);
6  Th = buildBdMesh(Th); // build the surface mesh
7  // build Th1, the surface of Th, defined by triangles elements and edges border_elements list
8  meshS Th1 = Th.Gamma;
```

The command buildBdMesh

Let Th3 a volume mesh (mesh3 type) ; such as the geometry description is a list of vertices, tetrahedra elements and triangle border elements. FreeFEM can generate the surface mesh associated to Th3. The intern mechanism
of **FreeFEM** created directly the meshS associated to Th3 and accessible by the command `meshS ThS = Th3;`.

**The command **savesurfacemesh**

Available for 3d meshes, the command `savesurfacemesh` saves the entire surface of a 3d volume `mesh3` at the format `.mesh`. Two possibilities about the `mesh3` surface:

- the geometric surface isn’t explicite, that means the `mesh3` doesn’t contain surface elements (triangles) and border surface elements (edge). The surface is defined by the border of the volume. Hence, `savesurfacemesh` returns the list of vertices and faces of the volume mesh, according to a local numbring at the border mesh.

- the geometric surface is explicite and known by the `mesh3` type. This may be due to the nature of the data mesh (list of vertices, tetrahedra, triangles, edges) or a surface building by `FreeFEM` with the calling of `buildSurface` operator. In this case, `savesurfacemesh` allows to save the real geometry of the surface 3d mesh (list of vertices, triangles, edges)

**Example of use**

```plaintext
1 load "msh3"
2 mesh3 Th3=cube(10,15,5);
3 savemesh(Th3, "surf.mesh");
4 savesurfacemesh(Th3, "surfreal.mesh");
5 mesh3 ThS3 = trunc(Th3, 1, split=3);
6 meshS ThSS = ThS3.Gamma;
7 savesurfacemesh(ThS3, "surfacesplit.mesh");
8 savemesh(ThSS,"GammaSplit.mesh" );
```

**volume mesh and meshS=NULL**

`savesurfmesh(Th,filename_mesh)` write in the file the vertices list and the triangle list (face of the volum mesh) according to a numbering in local surface

`savesurfmesh(Th,filename_points,filename_faces)` The operation does the same thing that the first exept to

**Glue of meshS meshes**

A surface 3d mesh can be the result of the generation of several assembled meshes, with caution of the right orientation at the merged interfaces.

```plaintext
1 meshS Th1 = square3(n,n, [2*x,y,1],orientation=-1);
2 meshS Th2 = square3(n,n, [2*x,y,0],orientation=1);
3 meshS Th11 = square3(n,n, [2*x,1,y],orientation=1);
4 meshS Th22 = square3(n,n, [2*x,0,y],orientation=-1);
5 meshS Th5 = square3(n,n, [1,y,x]);
6 meshS Th6 = square3(n,n, [2,y,x],orientation=1);
7 meshS Th = Th1+Th2+Th11+Th22+Th5+Th6;
8 assert(Th.nbnomanifold==40);
```

**Warning:** For the moment, the case of no manifold mesh are not considered in FreeFEM. To check if the meshS contains no manifold elements, the command `nbnomanifold`. 

---

### 3.2. Mesh Generation

145
3.2.4 The type meshL in 3 dimension

Commands for 3d curve mesh generation

The command segment

The function segment is a basic command to define a curve in 3D space.

The parameters of this command line are:

- n generates a n subsegments from the unit line
- \([.,.,.]\) is \([\Phi_1, \Phi_2, \Phi_3]\) is the geometric transformation from \(\mathbb{R}^3\) to \(\mathbb{R}^3\). By default, \([\Phi_1, \Phi_2, \Phi_3]\) = \([x,0,0]\)
- orientation= equal 1, gives the orientation of the triangulation, elements are in the reference orientation (counter clock wise) equal -1 reverse the orientation of the triangles it’s the global orientation of the surface 1 extern (-1 intern)
- cleanmesh= is a boolean, allowing remove the duplicated nodes
- removeduplicate= is a boolean, allowing remove the duplicated elements and border elements
- precismesh this parameter is the criteria to define two merging points. By default, it value is 1e-7 and define the smallest axis parallel boxes containing the discretion domain of \(\Omega\)

By default, the border points are marked by label 1 and 2.

```
real R = 3, r=1;
real h = 0.1; //
int nx = R*2*pi/h;
func torex = (R+r*cos(y*pi*2))*cos(x*pi*2);
func torey = (R+r*cos(y*pi*2))*sin(x*pi*2);
func torez = r*sin(y*pi*2);
meshL Th=segment(nx,[torex,torey,torez],removeduplicate=true) ;
```

The following code generates a 10 subsegments from the unit line with a clock wise triangulation, according to the geometric transformation \([torex,torey,torez]\) and removing the duplicated points/elements

The command buildmesh

This operator allows to define a curve mesh from multi-borders. The domain can be defined by a parametrized curve (keyword border), such as Th1 in the following example or piecewise by parametrized curves, such as the construction of the mesh Th2.

The pieces can only intersect at their endpoints, but it is possible to join more than two endpoints.

```
load "msh3"

// conical helix
border E1(t=0, 10.*pi){x=(1.*t)*cos(t); y=-(1.*t)*sin(t); z=t;}
meshL Th1=buildmeshL(E1(1000));

int upper = 1, others = 2, inner = 3, n = 10;
border D01(t=0, 1) {x=0; y=-1+t; }
border D02(t=0, 1){x=1.5-1.5*t; y=-1; z=3;label=upper;}
border D03(t=0, 1){x=1.5; y=-t; z=3;label=upper;}
border D04(t=0, 1){x=1+0.5*t; y=0; z=3;label=others;}
```
Remeshing

The command **trunc**

This operator allows to define a mesh by truncating another one, i.e. by removing segments, and/or by splitting each element by a given positive integer \( s \). Here, an example to use this function:

```plaintext
meshL ThL2 = trunc(ThL1, boolean function to keep or remove elements, split = s, label = ...)  
```

The command has the following arguments:

- **boolean function to keep or remove elements**
- **split** = sets the level \( n \) of edge splitting, each edge is splitted in \( n \) subpart (one by default)
- **label** = sets the label number of new boundary item (1 by default)
- **new2old**
- **old2new**
- **renum**
- **orientation** = equal 1, gives the orientation of the triangulation, elements are in the reference orientation (counter clock wise) equal -1 reverse the orientation of the triangles it's the global orientation of the surface 1 extern (-1 intern)
- **cleanmesh** = is a boolean, allowing remove the duplicated nodes
- **removeduplicate** = is a boolean, allowing remove the duplicated elements and border elements
- **precismesh** = this parameter is the criteria to define two merging points. By default, it value is 1e-7 and define the smallest axis parallel boxes containing the discretion domain of \( \Omega \)

An example of how to call this function:

```plaintext
int nx=10;
meshL Th = segment(nx, [5.*x, cos(pi*x), sin(pi*x)]);
Th = trunc(Th, (x < 0.5) | (y < 0.5) | (z > 1.), split=4);
```

The command **movemesh**

This is the classical mesh transformation \texttt{FreeFEM} function, meshL can be deformed by an application \([ \Phi_1, \Phi_2, \Phi_3 \]). The image \( T_h(\Omega) \) is obtained by the command \texttt{movemeshL}.

The parameters of movemesh are:

- **transfo** = sets the geometric transformation \( \Phi(x, y) = (\Phi_1(x, y, z), \Phi_2(x, y, z), \Phi_3(x, y, z)) \)
- **refedge** = sets the integer labels of the triangles. 0 by default.
FreeFEM Documentation, Release 4.6

- \textbf{refpoint=} sets the labels of the border points. This parameter is initialized as the label for the keyword \texttt{change}.

- \textbf{precismesh} this parameter is the criteria to define two merging points. By default, it value is 1e-7 and define the smallest axis parallel boxes containing the discretion domain of $\Omega$.

- \textbf{orientation = An integer expression} equal 1, give the orientation of the triangulation, elements must be in the reference orientation (counter clock wise) equal -1 reverse the orientation of the triangles. It’s the global orientation of the normals at the surface 1 extern (-1 intern).

- \textbf{cleanmesh=} is a boolean, allowing remove the duplicated nodes

- \textbf{removeduplicate=} is a boolean, allowing remove the duplicated elements and border elements

\textbf{Note:} The definition of the geometric transformation depends on the space dimension of the studied problem. It means that, with curve FEM, it’s possible to treat a real 1D problem (space coordinate is $x$) then the transformation is given by $x: \rightarrow F(x)$, that means $\left[F_x\right]$ and $F_y=F_z=0$ in FreeFEM function.

Example of using

```
1  int nx=100;
2  meshL Th=Sline(nx);
3  meshL Th31=movemesh(Th, [x]);
4  meshL Th32=movemesh(Th, [x,-x*(x-1)]);
5  meshL Th3=Th31+Th32;
```

\textbf{The command change}

The command change changes the label of elements and border elements of a \texttt{meshL}.

The parameters for this command line are:

- \textbf{refedge=} is a vector of integer that contains successive pairs of the old label number to the new label number for elements.

- \textbf{refpoint=} is a vector of integer that contains successive pairs of the old region number to new region number for boundary elements.

- \textbf{flabel=} is an integer function given the new value of the label.

- \textbf{fregion=} is an integer function given the new value of the region.

- \textbf{rmlpoint=} is a vector of integer, where edge’s label given are remove of the mesh

These vectors are composed of $n_l$ successive pairs of numbers $O, N$ where $n_l$ is the number (label or region) that we want to change. For example, we have:

\[
\text{label} = [O_1, N_1, ..., O_{n_l}, N_{n_l}]
\]
\[
\text{region} = [O_1, N_1, ..., O_{n_l}, N_{n_l}]
\]

\textbf{The commands buildBdMesh and Gamma}

The command Gamma allows to extract the border mesh independently of a surface mesh. With this function, the constructed border mesh contains the full geometric description of the boundary surface. In case where the border mesh doesn’t exist, before calling Gamma, must build it by calling the buildBdMesh function (see the next function description).
Glue of meshL meshes

An assembling of meshL is possible thanks to the operator +. The result returns a meshL, with caution of the right orientation at the merged interfaces. Here, the function checkmesh can be called.

```plaintext
int n = 10;
meshL ThL = extract(Th, [x+2, y*5], refedge=ll);
```

**Warning:** For the moment, the case of no manifold mesh are not considered in FreeFEM. To check if the meshL contains no manifold elements, the command nbnomanifold.

The command extract

This operator allows to extract a labeled piece or the entire border of a 2D mesh and project it in 3D. Optionally, a geometric transformation can be applied.

```plaintext
mesh Th = square(10, 10);
int ll = [4];
meshL ThL = extract(Th, [x+2, y*5], refedge=ll);
```

The commands rebuildBorder

This operator, used in the last example, allows to reconstruted the border elements following a special criteria ridgeangledetection. By default, it value is \( \frac{\pi}{5} \cdot arctan(1) \approx 40 \), the diedral angle for a decahedron.

The commands checkmesh

This function is available for all 3D meshes. It checkes and validates the a given mesh, allows to remove duplicate vertices and/or elements and border elements. The possible arguments are

- `precismesh` = this parameter is the criteria to define two merging points. By default, it value is 1e-7 and define the smallest axis parallel boxes containing the discretion domain of \( \Omega \)
- `removeduplicate` = is a boolean, allowing remove the duplicated elements and border elements
- `rebuildboundary` = is a boolean, allowing rebuild the border elements (in case of incomplete list given by the mesh)
Example:

```c
mesh3 Th = checkmesh(Th);
```

**TetGen: A tetrahedral mesh generator**

TetGen is a software developed by Dr. Hang Si of Weierstrass Institute for Applied Analysis and Stochastics in Berlin, Germany [HANG2006]. TetGen is free for research and non-commercial use. For any commercial license utilization, a commercial license is available upon request to Hang Si.

This software is a tetrahedral mesh generator of a three dimensional domain defined by its boundary (a surface). The input domain takes into account a polyhedral or a piecewise linear complex. This tetrahedralization is a constrained Delaunay tetrahedralization.

The method used in TetGen to control the quality of the mesh is a Delaunay refinement due to Shewchuk [SHEWCHUK1998]. The quality measure of this algorithm is the Radius-Edge Ratio (see Section 1.3.1 [HANG2006] for more details). A theoretical bound of this ratio of the Shewchuk algorithm is obtained for a given complex of vertices, constrained segments and facets of surface mesh, with no input angle less than 90 degrees. This theoretical bound is 2.0.

The launch of TetGen is done with the keyword `tetg`. The parameters of this command line is:

- `reftet` = sets the label of tetrahedra.
- `label` = is a vector of integers that contains the old labels number at index $2i$ and the new labels number at index $2i + 1$ of Triangles. This parameter is initialized as a label for the keyword `change`.
- `switch` = A string expression. This string corresponds to the command line switch of TetGen see Section 3.2 of [HANG2006].
- `nbofholes` = Number of holes (default value: “size of holelist / 3”).
- `holelist` = This array corresponds to `holelist` of TetGenio data structure [HANG2006]. A real vector of size $3 * nbofholes$. In TetGen, each hole is associated with a point inside this domain. This vector is $x^h_i, y^h_i, z^h_i, x^h_{i+1}, y^h_{i+1}, z^h_{i+1}, \cdots$, where $x^h_i, y^h_i, z^h_i$ is the associated point with the $i^{th}$ hole.
- `nbofregions` = Number of regions (default value: “size of regionlist / 5”).
- `regionlist` = This array corresponds to `regionlist` of TetGenio data structure [HANG2006]. The attribute and the volume constraint of region are given in this real vector of size $5 * nbofregions$. The $i^{th}$ region is described by five elements: $x-$coordinate, $y-$coordinate and $z-$coordinate of a point inside this domain ($x_i, y_i, z_i$); the attribute ($at_i$) and the maximum volume for tetrahedra ($mvol_i$) for this region. The `regionlist` vector is: $x_1, y_1, z_1, at_1, mvol_1, x_2, y_2, z_2, at_2, mvol_2, \cdots$.
- `noffacetcl` = Number of facets constraints “size of facetcl / 2”).
- `facetcl` = This array corresponds to `facetconstraintlist` of TetGenio data structure [HANG2006]. The $i^{th}$ facet constraint is defined by the facet marker $Ref_i^{fc}$ and the maximum area for faces $marea_i^{fc}$. The `facetcl` array is: $Ref_1^{fc}, marea_1^{fc}, Ref_2^{fc}, marea_2^{fc}, \cdots$.

This parameters has no effect if switch q is not selected.

Principal switch parameters in TetGen:

- `p` Tetrahedralization of boundary.
- `q` Quality mesh generation. The bound of Radius-Edge Ratio will be given after the option q. By default, this value is 2.0.
• **a Constructs with the volume constraints on tetrahedra.** These volumes constraints are defined with the bound of the previous switch \( q \) or in the parameter \( \text{regionlist} \).

• **A Attributes reference to region given in the regionlist.** The other regions have label 0.
  The option \( AA \) gives a different label at each region. This switch works with the option \( p \). If option \( r \) is used, this switch has no effect.

• **r Reconstructions and Refines a previously generated mesh.** This character is only used with the command line \text{tetgreconstruction}.

• **Y This switch preserves the mesh on the exterior boundary.**
  This switch must be used to ensure a conformal mesh between two adjacent meshes.

• **YY This switch preserves the mesh on the exterior and interior boundary.**

• **C** The consistency of the result’s mesh is testing by TetGen.

• **CC** The consistency of the result’s mesh is testing by TetGen and also constrained checks of Delaunay mesh (if \( p \) switch is selected) or the consistency of Conformal Delaunay (if \( q \) switch is selected).

• **V Give information of the work of TetGen.** More information can be obtained in specified \( \text{VV} \) or \( \text{VVV} \).

• **Q Quiet:** No terminal output except errors

• **M** The coplanar facets are not merging.

• **T Sets a tolerance for coplanar test.** The default value is \( 1e - 8 \).

• **d** Intersections of facets are detected.

To obtain a tetrahedral mesh with TetGen, we need the surface mesh of a three dimensional domain. We now give the command line in \text{FreeFEM} to construct these meshes.

**The keyword tetgtransfo**

This keyword corresponds to a composition of command line \text{tetg} and \text{movemesh23}.

\[
\text{tetgtransfo}(\text{Th2, transfo}=[\Phi(1), \Phi(2), \Phi(3)], ...) = \text{tetg}(\text{Th3surf, ...}),
\]

where \( \text{Th3surf} = \text{movemesh23}(\text{Th2, transfo}=[\Phi(1), \Phi(2), \Phi(3)]) \) and \( \text{Th2} \) is the input two dimensional mesh of \text{tetgtransfo}.

The parameters of this command line are, on one hand, the parameters \text{label, switch, regionlist, nboffacetcl, facetcl} of keyword \text{tetg} and on the other hand, the parameter \text{ptmerge} of keyword \text{movemesh23}.

**Note:** To use \text{tetgtransfo}, the result's mesh of \text{movemesh23} must be a closed surface and define one region only. Therefore, the parameter \text{regionlist} is defined for one region.

An example of this keyword can be found in line 61 of the \text{Build layer mesh example}.

**The keyword tetgconvexhull**

\text{FreeFEM}, using TetGen, is able to build a tetrahedralization from a set of points. This tetrahedralization is a Delaunay mesh of the convex hull of the set of points.

The coordinates of the points can be initialized in two ways. The first is a file that contains the coordinate of points...
\( X_i = (x_i, y_i, z_i) \). This file is organized as follows:

\[
\begin{array}{ccc}
  n_v & x_1 & y_1 & z_1 \\
  & x_2 & y_2 & z_2 \\
  & \vdots & \vdots & \vdots \\
  & x_{n_v} & y_{n_v} & z_{n_v}
\end{array}
\]

The second way is to give three arrays that correspond respectively to the \( x \)--coordinates, \( y \)--coordinates and \( z \)--coordinates.

The parameters of this command line are:

- **switch**= A string expression. This string corresponds to the command line `switch` of TetGen see Section 3.2 of [HANG2006].
- **reftet**= An integer expression. Set the label of tetrahedra.
- **label**= An integer expression. Set the label of triangles.

In the string `switch`, we can’t used the option `p` and `q` of TetGen.

### Reconstruct/Refine a 3d mesh with TetGen

Meshes in three dimension can be refined using TetGen with the command line `tetgreconstruction`.

The parameter of this keyword are

- **region**= an integer array that changes the region number of tetrahedra. This array is defined as the parameter `reftet` in the keyword change.
- **label**= an integer array that changes the label of boundary triangles. This array is defined as the parameter `label` in the keyword change.
- **sizeofvolume**= a reel function. This function constraints the volume size of the tetrahedra in the domain (see Isotrope mesh adaption section to build a 3d adapted mesh).

The parameters `switch`, `nbofregions`, `regionlist`, `nboffacetcl` and `facetcl` of the command line which call TetGen (`tetg`) is used for `tetgrefine`.

In the parameter `switch=`, the character `r` should be used without the character `p`.

For instance, see the manual of TetGen [HANG2006] for effect of `r` to other character.

The parameter `regionlist` defines a new volume constraint in the region. The label in the `regionlist` will be the previous label of region.

This parameter and `nbofregions` can’t be used with the parameter `sizeofvolume`.

**Example refinesphere.edp**

```plaintext
load "msh3"
load "tetgen"
load "medit"

mesh Th=square(10,20,[x*pi-pi/2,2*y*pi]); // $\frac{-\pi}{2},\frac{-\pi}{2}\times0,\rightarrow2\pi[$

// a parametrization of a sphere
func f1 =cos(x)*cos(y);
func f2 =cos(x)*sin(y);
func f3 = sin(x);
```

(continues on next page)
// partial derivative of the parametrization DF
func f1x=sin(x)*cos(y);
func f1y=-cos(x)*sin(y);
func f2x=-sin(x)*sin(y);
func f2y=cos(x)*cos(y);
func f3x=cos(x);
func f3y=0;

// $ M = D^t F D $
func m11=f1x^2+f2x^2+f3x^2;
func m21=f1x*f1y+f2x*f2y+f3x*f3y;
func m22=f1y^2+f2y^2+f3y^2;

func perio=[[4,y],[2,y],[1,x],[3,x]];
real hh=0.1;
real vv= 1/square(hh);
verbosity=2;
Th=adaptmesh(Th,m11*vv,m21*vv,m22*vv,
IsMetric=1,periodic=perio);
plot(Th,wait=1);
verbosity=2;

// construction of the surface of spheres
real Rmin = 1.;
func f1min = Rmin*f1;
func f2min = Rmin*f2;
func f3min = Rmin*f3;

meshS ThS=movemesh23(Th,transfo=[f1min,f2min,f3min]);
real[int] domain = [0.,0.,0.,145,0.01];
mesh3 Th3sph=tetg(ThS,switch="paAAQYY",nbofregions=1,regionlist=domain);

int[int] newlabel = [145,18];
real[int] domainrefine = [0.,0.,0.,145,0.0001];
mesh3 Th3sphrefine=tetgreconstruction(Th3sph,switch="raAQ",region=newlabel,
˓nbofregions=1,regionlist=domainrefine,sizeofvolume=0.0001);

int[int] newlabel2 = [145,53];
func fsize = 0.01/(( 1 + 5*sqrt( (x-0.5)^2+(y-0.5)^2+(z-0.5)^2 ) )^3);
mesh3 Th3sphrefine2=tetgreconstruction(Th3sph,switch="raAQ",region=newlabel2,
˓sizeofvolume=fsize);

medit("sphere",Th3sph,wait=1);
medit("spheredefinedomain",wait=1,Th3sphrefine);
medit("sphererefinelocal",wait=1,Th3sphrefine2);

// FFCS: testing 3d plots
plot(Th3sph);
plot(Th3sphrefine);
plot(Th3sphrefine2);
3.2.5 Read/Write Statements for meshes

2d case

format of mesh data

Users who want to read a triangulation made elsewhere should see the structure of the file generated below:

```plaintext
border C(t=0, 2*pi){x=cos(t); y=sin(t);} 
mesh Th1 = buildmesh(C(10)); 
savemesh(Th1, "mesh.msh"); 
mesh Th2=readmesh("mesh.msh");
```

The mesh is shown on Fig. 3.28.

The information about Th are saved in the file mesh.msh whose structure is shown on Table 3.1. An external file contains a mesh at format .mesh can be read by the command `readmesh(file_name)`.

There, $n_v$ denotes the number of vertices, $n_t$ the number of triangles and $n_s$ the number of edges on boundary.

For each vertex $q_i$, $i = 1, \cdots, n_v$, denoted by $(q_{ix}, q_{iy})$ the $x$-coordinate and $y$-coordinate.

Each triangle $T_k$, $k = 1, \cdots, n_t$ has three vertices $q_{k1}, q_{k2}, q_{k3}$ that are oriented counter-clockwise.

The boundary consists of 10 lines $L_i$, $i = 1, \cdots, 10$ whose end points are $q^{i1}, q^{i2}$.

In the Fig. 3.28, we have the following.

$n_v = 14, n_t = 16, n_s = 10$

$q^1 = (-0.309016994375, 0.951056516295)$

$\cdots$

$q^{14} = (-0.309016994375, -0.951056516295)$

The vertices of $T_1$ are $q^9, q^{12}, q^{10}$. 

Fig. 3.28: Mesh by `buildmesh(C(10))`
The vertices of $T_{16}$ are $q^9, q^{10}, q^6$.
The edge of the 1st side $L_1$ are $q^6, q^5$.

The edge of the 10th side $L_{10}$ are $q^{10}, q^6$.

### Table 3.1: The structure of mesh_sample.msh

<table>
<thead>
<tr>
<th>Content of the file</th>
<th>Explanation</th>
</tr>
</thead>
<tbody>
<tr>
<td>14 16 10</td>
<td>$n_v$, $n_t$, $n_e$</td>
</tr>
<tr>
<td>-0.309016994375 0.951056516295 1</td>
<td>$q_x^1 q_y^1$ boundary label = 1</td>
</tr>
<tr>
<td>0.309016994375 0.951056516295 1</td>
<td>$q_x^2 q_y^2$ boundary label = 1</td>
</tr>
<tr>
<td>...</td>
<td>...</td>
</tr>
<tr>
<td>-0.309016994375 -0.951056516295 1</td>
<td>$q_x^{14} q_y^{14}$ boundary label = 1</td>
</tr>
<tr>
<td>9 12 10 0</td>
<td>$1_1$ $1_2$ $1_3$ region label = 0</td>
</tr>
<tr>
<td>5 9 6 0</td>
<td>$2_1$ $2_2$ $2_3$ region label = 0</td>
</tr>
<tr>
<td>...</td>
<td>...</td>
</tr>
<tr>
<td>9 10 6 0</td>
<td>$10_1$ $16_2$ $16_3$ region label = 0</td>
</tr>
<tr>
<td>6 5 1</td>
<td>$1_1$ $1_2$ boundary label = 1</td>
</tr>
<tr>
<td>5 2 1</td>
<td>$2_1$ $2_2$ boundary label = 1</td>
</tr>
<tr>
<td>...</td>
<td>...</td>
</tr>
<tr>
<td>10 6 1</td>
<td>$10_1$ $10_2$ boundary label = 1</td>
</tr>
</tbody>
</table>

In FreeFEM there are many mesh file formats available for communication with other tools such as emc2, modulef, ... (see Mesh format chapter).

The extension of a file implies its format. More details can be found on the file format .msh in the article by F. Hecht “bamg: a bidimensional anisotropic mesh generator” [HECHT1998_2].

A mesh file can be read into FreeFEM except that the names of the borders are lost and only their reference numbers are kept. So these borders have to be referenced by the number which corresponds to their order of appearance in the program, unless this number is overwritten by the keyword label. Here are some examples:

```plaintext
// Parameters
int n = 10;

// Mesh
border floor(t=0, 1){x= t; y=0; label=1;};
border right(t=0, 1){x=1; y=t; label=5;};
border ceiling(t=1, 0){x=t; y=1; label=5;};
border left(t=1, 0){x=0; y=t; label=5;};

mesh th = buildmesh(floor(n) + right(n) + ceiling(n) + left(n));

// save mesh in different formats
savemesh(th, "toto.am_fmt"); // format "formatted Marrocco"
savemesh(th, "toto.Th"); // format database db mesh "bamg"
savemesh(th, "toto.msh"); // format freefem
savemesh(th, "toto.nopo"); // modulef format

// Fespace
fespace fempl(th, P1);
fempl f = sin(x)*cos(y);
fempl g;
```

(continues on next page)
//save the fespace function in a file
{
    ofstream file("f.txt");
    file << f[] << endl;
} //the file is automatically closed at the end of the block
//read a file and put it in a fespace function
{
    ifstream file("f.txt");
    file >> g[] ;
} //the file is equally automatically closed

// Plot
plot(g);

// Mesh 2
//read the mesh for freefem format saved mesh
mesh th2 = readmesh("toto.msh");

// Fespace 2
fespace Vh2(th2, P1);
Vh2 u, v;

// Problem
//solve:
// $u + \Delta u = g$ in $\Omega$
// $u=0$ on $\Gamma_1$
// $\frac{\partial u }{\partial n} = g$ on $\Gamma_2$
solve Problem(u, v)
    = int2d(th2) (u*v
        - dx(u)*dx(v)
        - dy(u)*dy(v)
    )
    + int2d(th2) (-g*v
    )
    + int1d(th2, 5) (g*v
    )
    + on(1, u=0)
;
// Plot
plot(th2, u);

Input/output for a mesh

- the command readmesh

The function readmesh allows to build a mesh from a data file

```cpp
mesh Th=readmeshS("Th.mesh");
mesh Thff = readmesh("Thff.msh"); // FreeFEM format
```

- the command savemesh
The function `savemesh` allows to export a mesh

```plaintext
savemesh(Th, "Th.mesh")
savemesh(Thff, "Thff.mesh") // FreeFEM format
savemesh(th, "toto.msh"); // format freefem
savemesh(th, "toto.am_fmt"); // format "formated Marrocco"
savemesh(th, "toto.Th"); // format database db mesh "bamg"
savemesh(th, "toto.nopo"); // modulef format
savemesh(Th, "mm", [x, y, u]); // save surface mesh for medit, see for example minimal-surf.edp
exec("medit mm;rm mm.bb mm.faces mm.points");
```

- the command `vtkloadS`

The function `vtkload` allows to build a mesh from a data mesh at vtk format mesh

```plaintext
load "iovtk"
mesh Th=vtkloadS("mymesh.vtk");
```

- the command `savevtk`

The function `savevtk` allows to export a mesh to a data mesh at vtk format mesh

```plaintext
load "iovtk"
savevtk("Th.vtk",Th);
```

- the command `gmshload`

The function `gmshloadS` allows to build a mesh from a data mesh file at formatmsh (GMSH)

```plaintext
load "gmsh"
mesh Th=gmshload("mymesh.msh");
```

- the command `savegmsh`

The function `savegmsh` allows to export a mesh to a data mesh msh (GMSH)

```plaintext
load "gmsh"
savegmsh(Th, "Th");
```

3d case

format of mesh data

In three dimensions, the file mesh format supported for input and output files by FreeFEM are the extension .msh and .mesh. These formats are described in the Mesh Format section.
Extension file .msh

The structure of the files with extension .msh in 3D is given by:

\[
\begin{array}{ccc}
q^1_x & q^1_y & q^1_z \\
q^2_x & q^2_y & q^2_z \\
\vdots & \vdots & \vdots \\
q^n_x & q^n_y & q^n_z \\
\end{array}
\]

Vertex label

\[
\begin{array}{cccc}
1_1 & 1_2 & 1_3 & 1_4 \\
2_1 & 2_2 & 2_3 & 2_4 \\
\vdots & \vdots & \vdots & \vdots \\
(n_{tet})_1 & (n_{tet})_2 & (n_{tet})_3 & (n_{tet})_4 \\
\end{array}
\]

region label

\[
\begin{array}{cccc}
1_1 & 1_2 & 1_3 & 1_4 \\
2_1 & 2_2 & 2_3 & 2_4 \\
\vdots & \vdots & \vdots & \vdots \\
(n_{tri})_1 & (n_{tri})_2 & (n_{tri})_3 & (n_{tri})_4 \\
\end{array}
\]

boundary label

In this structure, \( n_v \) denotes the number of vertices, \( n_{tet} \) the number of tetrahedra and \( n_{tri} \) the number of triangles.

For each vertex \( q^i \), \( i = 1, \ldots, n_v \), we denote by \((q^i_x, q^i_y, q^i_z)\) the \( x \)-coordinate, the \( y \)-coordinate and the \( z \)-coordinate.

Each tetrahedra \( T_k, k = 1, \ldots, n_{tet} \) has four vertices \((q^k_1, q^k_2, q^k_3, q^k_4)\).

The boundary consists of a union of triangles. Each triangle \( tri_j, j = 1, \ldots, n_{tri} \) has three vertices \((q^{j1}, q^{j2}, q^{j3})\).

extension file .mesh

The data structure for a three dimensional mesh is composed of the data structure presented in Mesh Format section and a data structure for the tetrahedra. The tetrahedra of a three dimensional mesh are referred using the following field:

```
1 Tetrahedra
2 NbTetrahedra
3 Vertex1 Vertex2 Vertex3 Vertex4 Label
4 ...
5 Vertex1 Vertex2 Vertex3 Vertex4 Label
6 Triangles
7 NbTriangles
8 Vertex1 Vertex2 Vertex3 Label
9 ...
10 Vertex1 Vertex2 Vertex3 Label
```

This field is express with the notation of Mesh Format section.

Input/output for a mesh3

- the command `readmesh3`

The function `readmesh3` allows to build a mesh3 from a data file

```
1 mesh3 Th3=readmesh3("Th3.mesh");
```

- the command `savemesh`

The function `savemesh` allows to export a mesh3

```
1 savemesh(Th3,"Th3.mesh")
2 savemesh(Th3ff,"Th3ff.msh") // FreeFEM format
```
• the command \texttt{vtkload3}

The function \texttt{vtkload3} allows to build a mesh3 from a data mesh at vtk format mesh

\begin{verbatim}
load "iovtk"
mesh3 Th3=vtkload3("mymesh.vtk");
\end{verbatim}

• the command \texttt{savevtk}

The function \texttt{savevtk} allows to export a mesh3 to a data mesh at vtk format mesh

\begin{verbatim}
load "iovtk"
savevtk("Th3.vtk",Th3);
\end{verbatim}

• the command \texttt{gmshload3}

The function \texttt{gmshload3} allows to build a mesh3 from a data mesh file at format msh (GMSH)

\begin{verbatim}
load "gmsh"
mesh3 Th3=gmshload3("mymesh.msh");
\end{verbatim}

• the command \texttt{savegmsh}

The function \texttt{savegmsh} allows to export a mesh3 to a data mesh msh (GMSH)

\begin{verbatim}
load "gmsh"
savegmsh(Th3, "Th3");
\end{verbatim}

\section*{Surface 3d case}

\subsection*{format of mesh data}

Like 2d and 3d, the input and output format files supported by FreeFEM are the extension .msh and .mesh. These formats are described in the \textit{Mesh Format section.}

\textbf{Extension file .msh} The structure of the files with extension .msh in surface 3D is given by:

\begin{verbatim}
n_v n_tri n_edges q_1x q_1y q_1z Vertexlabel
q_2x q_2y q_2z Vertexlabel
... ... ...
q_nvx q_nvy q_nvz Vertexlabel
r_1 1_2 1_3 regionlabel
2_1 2_2 2_3 regionlabel
... ... ...
(n_tri)_1 (n_tri)_2 (n_tri)_3 regionlabel
1_1 1_2 boundarylabel
2_1 2_2 boundarylabel
... ... ...
(n_edge)_1 (n_edge)_2 boundarylabel
\end{verbatim}

In this structure, \(n_v\) denotes the number of vertices, \(n_{tet}\) the number of tetrahedra and \(n_{tri}\) the number of triangles.

For each vertex \(q^i, i = 1, \cdots, n_v\), we denote by \((q^i_x, q^i_y, q^i_z)\) the \(x\)-coordinate, the \(y\)-coordinate and the \(z\)-coordinate.

Each tetrahedra \(T_k, k = 1, \cdots, n_{tet}\) has four vertices \(q^{k_1}, q^{k_2}, q^{k_3}, q^{k_4}\).
The boundary consists of a union of triangles. Each triangle $bc_j, j = 1, \cdots, n_{tri}$ has three vertices $q^{j_1}, q^{j_2}, q^{j_3}$.

**extension file .mesh** The data structure for a three dimensional mesh is composed of the data structure presented in *Mesh Format section* and a data structure for the tetrahedra. The tetrahedra of a three dimensional mesh are referred using the following field:

```
MeshVersionFormatted 2
Dimension 3
Vertices
NbVertices
(v0)x (v0)y (v0)z
...
(vn)x (vn)y (vn)z

Triangles
NbTriangles
Vertex1 Vertex2 Vertex3 Label
...
Vertex1 Vertex2 Vertex3 Label

Edges
NbEdges
Vertex1 Vertex2 Label
...
Vertex1 Vertex2 Label

End
```

This field is express with the notation of *Mesh Format section*.

**Input/output for a meshS**

- the command `readmesh3`

  The function `readmesh3` allows to build a meshS from a data file

```
1 meshS ThS=readmesh3("ThS.mesh");
2 meshS Th3ff = readmesh3("ThSff.msh"); // FreeFEM format
```

- the command `savemesh`

  The function `savemesh` allows to export a meshS

```
1 savemesh(ThS,"ThS.mesh")
2 savemesh(ThSff,"ThSff.msh") // FreeFEM format
```

- the command `vtkloadS`

  The function `vtkloadS` allows to build a meshS from a data mesh at vtk format mesh

```
1 load "iovtk"
2 meshS ThS=vtkloadS("mymesh.vtk");
```

- the command `savevtk`

  The function `savevtk` allows to export a meshS to a data mesh at vtk format mesh
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```plaintext
load "iovtk"
savvtk("ThS.vtk",ThS);
```

- the command `gmshloadS`

The function `gmshloadS` allows to build a `meshS` from a data mesh file at format `msh` (GMSH)

```plaintext
load "gmsh"
meshS ThS=gmshloadS("mymesh.msh");
```

- the command `savegmsh`

The function `savegmsh` allows to export a `meshS` to a data mesh `msh` (GMSH)

```plaintext
load "gmsh"
savegmsh(ThS, "ThS");
```

### 3.2.6 Medit

The keyword `medit` allows to display a mesh alone or a mesh and one or several functions defined on the mesh using the Pascal Frey’s freeware `medit`. `medit` opens its own window and uses OpenGL extensively. Naturally to use this command `medit` must be installed.

A visualization with `medit` of scalar solutions $f_1$ and $f_2$ continuous, piecewise linear and known at the vertices of the mesh $Th$ is obtained using:

```plaintext
medit("sol1 sol2", Th, f1, f2, order=1);
```

The first plot named `sol1` display $f_1$. The second plot names `sol2` display $f_2$.

The arguments of the function `medit` are the name of the different scenes (separated by a space) of `medit`, a mesh and solutions.

Each solution is associated with one scene. The scalar, vector and symmetric tensor solutions are specified in the format described in the section dealing with the keyword `savesol`.

The parameters of this command line are:

- `order=0` if the solution is given at the center of gravity of elements. 1 is the solution is given at the vertices of elements.
- `meditff=` set the name of execute command of `medit`. By default, this string is `medit`.
- `save=` set the name of a file `.sol` or `.solb` to save solutions.

This command line allows also to represent two different meshes and solutions on them in the same windows. The nature of solutions must be the same. Hence, we can visualize in the same window the different domains in a domain decomposition method for instance. A visualization with `medit` of scalar solutions $h_1$ and $h_2$ at vertices of the mesh $Th_1$ and $Th_2$ respectively are obtained using:

```plaintext
medit("sol2domain", Th1, h1, Th2, h2, order=1);
```

Tip: `Medit`

```plaintext
load "medit"
```

// Initial Problem:

(continues on next page)
4 // Resolution of the following EDP:
5 // -Delta u_s = f on \Omega = \{(x,y) | 1 <= sqrt(x^2+y^2) <= 2\}
6 // -Delta u_1 = f_1 on \Omega_1 = \{(x,y) | 0.5 <= sqrt(x^2+y^2) <= 1.\}
7 // u = 1 on Gamma
8 // Null Neumann condition on Gamma_1 and on Gamma_2
9 // We find the solution u by solving two EDP defined on domain Omega and Omega_1
10 // This solution is visualize with medit
11 verbosity=3;
12
13 // Mesh
14 border Gamma(t=0, 2*pi){x=cos(t); y=sin(t); label=1;};
15 border Gamma1(t=0, 2*pi){x=2*cos(t); y=2*sin(t); label=2;};
16 border Gamma2(t=0, 2*pi){x=0.5*cos(t); y=0.5*sin(t); label=3;};
17
18 mesh Th = buildmesh(Gamma(40) + Gamma(-40));  //Omega
19 mesh Th1 = buildmesh(Gamma1(40) + Gamma2(-40));  //Omega_1
20
21 // Fespace
22 fespace Vh(Th, P2);
23 func f = sqrt(x*x + y*y);
24 Vh us, v;
25
26 fespace Vh1(Th1, P2);
27 func f1 = 10*sqrt(x*x+y*y);
28 Vh1 u1, v1;
29
30 // Macro
31 macro Grad2(us) [dx(us), dy(us)] // EOM
32
33 // Problem
34 problem Lap2dOmega (us, v, init=false)
35 = int2d(Th)(
36    Grad2(v)' * Grad2(us)
37 ) - int2d(Th)(
38    f*v
39 ) + on(1, us=1)
40;
41
42 problem Lap2dOmega1 (u1, v1, init=false)
43 = int2d(Th1)(
44    Grad2(v1)' * Grad2(u1)
45 ) - int2d(Th1)(
46    f1*v1
47 ) + on(1, u1=1)
48;
49
50 // Solve
51 Lap2dOmega;
52 Lap2dOmega1;
53
54 // Plot with medit
55 medit("solution", Th, us, Th1, ul, order=1, save="testsavemedit.solb");
3.2.7 Mshmet

Mshmet is a software developed by P. Frey that allows to compute an anisotropic metric based on solutions (i.e. Hessian-based). This software can return also an isotropic metric. Moreover, mshmet can also construct a metric suitable for levelset interface capturing. The solution can be defined on 2D or 3D structured/unstructured meshes. For example, the solution can be an error estimate of a FE solution.

Solutions for mshmet are given as an argument. The solution can be a func, a vector func, a symmetric tensor, a fespace function, a fespace vector function and a fespace symmetric tensor. The symmetric tensor argument is defined as this type of data for datasol argument. This software accepts more than one solution.

For example, the metric $M$ computed with mshmet for the solution $u$ defined on the mesh $Th$ is obtained by writing:

```plaintext
fespace Vh(Th, P1);
Vh u; // a scalar fespace function
real[int] M = mshmet(Th, u);
```

The parameters of the keyword mshmet are:

- **normalization** = (b) do a normalization of all solution in $[0, 1]$.
- **aniso** = (b) build anisotropic metric if 1 (default 0: isotropic)
- **levelset** = (b) build metric for levelset method (default: false)
- **verbosity** = (l) level of verbosity
- **nbregul** = (l) number of regularization’s iteration of solutions given (default 0).
- **hmin** = (d)
- **hmax** = (d)
- **err** = (d) level of error.
- **width** = (d) the width
- **metric** = a vector of double. This vector contains an initial metric given to mshmet. The structure of the metric vector is described in the next paragraph.
- **loptions** = a vector of integer of size 7. This vector contains the integer parameters of mshmet (for expert only).
  - loptions(0): normalization (default 1).
  - loptions(1): isotropic parameters (default 0). 1 for isotropic metric results otherwise 0.
  - loptions(2): level set parameters (default 0). 1 for building level set metric otherwise 0.
  - loptions(3): debug parameters (default 0). 1 for turning on debug mode otherwise 0.
  - loptions(4): level of verbosity (default 10).
  - loptions(5): number of regularization’s iteration of solutions given (default 0).
  - loptions(6): previously metric parameter (default 0). 1 for using previous metric otherwise 0.
- **doptions** = a vector of double of size 4. This vector contains the real parameters of mshmet (for expert only).
  - doptions(0): hmin : min size parameters (default 0.01).
  - doptions(1): hmax : max size parameters (default 1.0).
The result of the keyword `mshmet` is a `real[int]` which contains the metric computed by `mshmet` at the different vertices $V_i$ of the mesh.

With $n_v$ the number of vertices, the structure of this vector is:

$$M_{iso} = (m(V_0), m(V_1), \ldots, m(V_{n_v}))^t$$

for an isotropic metric $m$. For a symmetric tensor metric $h = \begin{pmatrix} m_{11} & m_{12} & m_{13} \\ m_{21} & m_{22} & m_{23} \\ m_{31} & m_{32} & m_{33} \end{pmatrix}$, the parameters metric is:

$$M_{aniso} = (H(V_0), \ldots, H(V_{n_v}))^t$$

where $H(V_i)$ is the vector of size 6 defined by $[m_{11}, m_{21}, m_{22}, m_{31}, m_{32}, m_{33}]$

Tip: `mshmet`

```plaintext
load "mshmet"
load "medit"
load "msh3"

// Parameters
real error = 0.01;
func zmin = 0;
func zmax = 1;
int MaxLayer = 10;

// Mesh
border a(t=0, 1.0){x=t; y=0; label=1;};
border b(t=0, 0.5){x=1; y=t; label=2;};
border c(t=0, 0.5){x=1-t; y=0.5; label=3;};
border d(t=0.5, 1){x=0.5; y=t; label=4;};
border e(t=0.5, 1){x=1-t; y=1; label=5;};
border f(t=0.0, 1){x=0; y=1-t; label=6;};
mesh Th = buildmesh(a(6) + b(4) + c(4) + d(4) + e(4) + f(6));
mesh3 Th3 = buildlayers(Th, MaxLayer, zbound=[zmin, zmax]);

// Fespace
fespace Vh3(Th3, P2);
Vh3 u3, v3;
fespace Vh3P1(Th3, P1);
Vh3P1 usol;

// Problem
problem Problem2(u3, v3, solver=sparesolver)
  = int3d(Th3)(
    u3*v3+1.0e-10
    + dx(u3)*dx(v3)
    + dy(u3)*dy(v3)
    + dz(u3)*dz(v3)
  )
  - int3d(Th3)(
    v3
)
```

(continues on next page)
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(continued from previous page)

```cpp
38
39  )
40  on(0, 1, 2, 3, 4, 5, 6, u3=0)
41  
42  // Solve
43  Problem2;
44  cout << u3[].min << " " << u3[].max << endl;
45  medit("Sol", Th3, u3);
46
47  real[int] bb = mshmet(Th3,u3);
48  cout << "Metric:" << bb << endl;
49  for (int ii = 0; ii < Th3.nv; iii++)
50    usol[][ii] = bb[ii];
51  medit("Metric", Th3, usol);

3.2.8 FreeYams

FreeYams is a surface mesh adaptation software which is developed by P. Frey. This software is a new version of yams. The adapted surface mesh is constructed with a geometric metric tensor field. This field is based on the intrinsic properties of the discrete surface.

Also, this software allows to construct a simplification of a mesh. This decimation is based on the Hausdorff distance between the initial and the current triangulation. Compared to the software yams, FreeYams can be used also to produce anisotropic triangulations adapted to levelset simulations. A technical report on freeYams documentation is available here.

To call FreeYams in FreeFEM, we used the keyword freeyams. The arguments of this function are the initial mesh and/or metric. The metric with freeyams are a func, a fespace function, a symmetric tensor function, a symmetric tensor fespace function or a vector of double (real[int]). If the metric is a vector of double, this data must be given in metric parameter. Otherwise, the metric is given in the argument.

For example, the adapted mesh of Thinit defined by the metric u defined as fespace function is obtained by writing:

```cpp
1  fespace Vh(Thinit, P1);
2  Vh u;
3  mesh3 Th = freeyams(Thinit, u);
```

The symmetric tensor argument for freeyams keyword is defined as this type of data for datasol argument.

- aniso= (b) aniso or iso metric (default 0, iso)
- mem= (l) memory of for freeyams in Mb (default -1, freeyams choose)
- hmin= (d)
- hmax= (d)
- gradation= (d)
- option= (l)
  - 0 : mesh optimization (smoothing+swapping)
  - 1 : decimation+enrichment adapted to a metric map. (default)

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- -1: decimation adapted to a metric map.
- 2: decimation+enrichment with a Hausdorff-like method
- -2: decimation with a Hausdorff-like method
- 4: split triangles recursively.
- 9: No-Shrinkage Vertex Smoothing

• `ridgeangle` = (d)
• `absolute` = (b)
• `verbosity` = (i)

• `metric` = vector expression. This parameter contains the metric at the different vertices on the initial mesh. With \( n_v \) being the number of vertices, this vector is:

\[
M_{iso} = (m(V_0), m(V_1), \ldots, m(V_{n_v}))^t
\]

for a scalar metric \( m \). For a symmetric tensor metric \( h = \begin{pmatrix} m_{11} & m_{12} & m_{13} \\ m_{21} & m_{22} & m_{23} \\ m_{31} & m_{32} & m_{33} \end{pmatrix} \), the parameter `metric` is:

\[
M_{aniso} = (H(V_0), \ldots, H(V_{n_v}))^t
\]

where \( H(V_i) \) is the vector of size 6 defined by \([m_{11}, m_{21}, m_{22}, m_{31}, m_{32}, m_{33}]\)

• `loptions` = a vector of integer of size 13. This vector contains the integer options of FreeYams. (just for the expert)

  - `loptions(0)`: anisotropic parameter (default 0). If you give an anisotropic metric 1 otherwise 0.
  - `loptions(1)`: Finite Element correction parameter (default 0). 1 for no Finite Element correction otherwise 0.
  - `loptions(2)`: Split multiple connected points parameter (default 1). 1 for splitting multiple connected points otherwise 0.
  - `loptions(3)`: maximum value of memory size in Mbytes (default -1: the size is given by freeyams).
  - `loptions(4)`: set the value of the connected component which we want to obtain. (Remark: freeyams give an automatic value at each connected component).
  - `loptions(5)`: level of verbosity
  - `loptions(6)`: Create point on straight edge (no mapping) parameter (default 0). 1 for creating point on straight edge otherwise 0.
  - `loptions(7)`: validity check during smoothing parameter. This parameter is only used with No-Shrinkage Vertex Smoothing optimization (optimization option parameter 9). 1 for No validity checking during smoothing otherwise 0.
  - `loptions(8)`: number of desired’s vertices (default -1).
  - `loptions(9)`: number of iteration of optimizations (default 30).
  - `loptions(10)`: no detection parameter (default 0). 1 for detecting the ridge on the mesh otherwise 0. The ridge definition is given in the parameter `doptions(12)`.
  - `loptions(11)`: no vertex smoothing parameter (default 0). 1 for smoothing the vertices otherwise 0.
- loptions(12): Optimization level parameter (default 0).
- 0: mesh optimization (smoothing+swapping)
- 1: decimation+enrichment adapted to a metric map.
- -1: decimation adapted to a metric map.
- 2: decimation+enrichment with a Hausdorff-like method
- -2: decimation with a Hausdorff-like method
- 4: split triangles recursively.
- 9: No-Shrinkage Vertex Smoothing

• doptions= a vector of double of size 11. This vectors contains the real options of freeyams.
  - doptions(0): Set the geometric approximation (Tangent plane deviation) (default 0.01).
  - doptions(1): Set the lambda parameter (default -1).
  - doptions(2): Set the mu parameter (default -1).
  - doptions(3): Set the gradation value (Mesh density control) (default 1.3).
  - doptions(4): Set the minimal size(hmin) (default -2.0: the size is automatically computed).
  - doptions(5): Set the maximal size(hmax) (default -2.0: the size is automatically computed).
  - doptions(6): Set the tolerance of the control of Chordal deviation (default -2.0).
  - doptions(7): Set the quality of degradation (default 0.599).
  - doptions(8): Set the declic parameter (default 2.0).
  - doptions(9): Set the angular walton limitation parameter (default 45 degree).
  - doptions(10): Set the angular ridge detection (default 45 degree).

Tip: freeyams

```c
load "msh3"
load "medit"
load "freeyams"

// Parameters
int nn = 20;
real zmin = 0;
real zmax = 1;

// Mesh
mesh Th2 = square(nn, nn);
int[int] rup = [0, 2], rdown = [0, 1];
int[int] rmid = [1, 1, 2, 1, 3, 1, 4, 1];
mesh3 Th = buildlayers(Th2, nn, zbound=[zmin, zmax], reffacemid=rmid, reffaceup=rup,
                        rreffacelow=rdown);
mesh3 Th3 = freeyams(Th);
medit("SurfaceMesh", Th3);
```

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3.2.9 mmg3d

Todo: mmg3d-v4.0

Mmg3d is a 3D remeshing software developed by C. Dobrzynski and P. Frey. This software allows to remesh an initial mesh made of tetrahedra. This initial mesh is adapted to a geometric metric tensor field or to a displacement vector (moving rigid body). The metric can be obtained with mshmet.

Note:
- If no metric is given, an isotropic metric is computed by analyzing the size of the edges in the initial mesh.
- If a displacement is given, the vertices of the surface triangles are moved without verifying the geometrical structure of the new surface mesh.

The parameters of mmg3d are:

- **options= vector expression.** This vector contains the option parameters of mmg3d. It is a vector of 6 values, with the following meaning:
  - Optimization parameters: (default 1)
    - 0: mesh optimization.
    - 1: adaptation with metric (deletion and insertion vertices) and optimization.
    - -1: adaptation with metric (deletion and insertion vertices) without optimization.
    - 4: split tetrahedra (be careful modify the surface).
    - 9: moving mesh with optimization.
    - -9: moving mesh without optimization.
  - Debug mode: (default 0)
    - 1: turn on debug mode.
    - 0: otherwise.
  - Specify the size of bucket per dimension (default 64)
  - Swapping mode: (default 0)
    - 1: no edge or face flipping.
    - 0: otherwise.
  - Insert points mode: (default 0)
    - 1: no edge splitting or collapsing and no insert points.
    - 0: otherwise.
  - 5. Verbosity level (default 3)

- **memory= integer expression.** Set the maximum memory size of new mesh in Mbytes. By default the number of maximum vertices, tetrahedra and triangles are respectively 500 000, 3000 000, 100000 which represent approximately a memory of 100 Mo.
• **metric= vector expression.** This vector contains the metric given at mmg3d. It is a vector of size $nv$ or 6 $nv$ respectively for an isotropic and anisotropic metric where $nv$ is the number of vertices in the initial mesh. The structure of metric vector is described in the *mshmet*.

• **displacement= $[\Phi_1, \Phi_2, \Phi_3]$** set the displacement vector of the initial mesh $\Phi(x, y) = [\Phi_1(x, y), \Phi_2(x, y), \Phi_3(x, y)]$.

• **displVect=** sets the vector displacement in a vector expression. This vector contains the displacement at each point of the initial mesh. It is a vector of size 3 $nv$.

**Tip:** mmg3d

```plaintext
load "msh3"
load "medit"
load "mmg3d"
include "Cube.idp"

// Parameters
int n = 6;
int[int] Nxyz = [12, 12, 12];
real [int, int] Bxyz = [[0., 1.], [0., 1.], [0., 1.]];
int [int, int] Lxyz = [[1, 1], [2, 2], [2, 2]];

// Mesh
mesh3 Th = Cube(Nxyz, Bxyz, Lxyz);
real[int] isometric(Th.nv);
for (int ii = 0; ii < Th.nv; ii++)
    isometric[ii] = 0.17;

mesh3 Th3 = mmg3d(Th, memory=100, metric=isometric);
// Plot
medit("Initial", Th);
medit("Isometric", Th3);
```

**Tip:** Falling spheres

```plaintext
load "msh3"
load "TetGen"
load "medit"
load "mmg3d"
include "MeshSurface.idp"

// Parameters
real hs = 0.8;
int[int] N = [4/hs, 8/hs, 11.5/hs];
real [int, int] B = [[-2, 2], [-2, 6], [-10, 1.5]];
int [int, int] L = [[311, 311], [311, 311], [311, 311]];
int[int] opt = [9, 0, 64, 0, 0, 3];
real[int] vit=[0, 0, -0.3];
func zero = 0.0;
func dep = vit[2];
```

(continues on next page)
// Meshes
meshS ThH = SurfaceHex(N, B, L, 1);
meshS ThSg = Sphere(1, hs, 300, -1);
meshS ThSd = Sphere(1, hs, 310, -1);
ThSd = movemesh(ThSd, [x, 4+y, z]);
meshS ThHS = ThH + ThSg + ThSd;
medit("ThHS", ThHS);

real voltet = (hs^3)/6.;
real[int] domain = [0, 0, -4, 1, voltet];
real [int] holes = [0, 0, 0, 0, 4, 0];

mesh3 Th = tetg(ThHS, switch="pqaAAYQ", nboregions=1, regionlist=domain,
   nbofholes=2, holelist=holes);
medit("Box-With-two-Ball", Th);

// Fespace
fespace Vh(Th, P1);
Vh uh,vh;

// Macro
macro Grad(u) [dx(u),dy(u),dz(u)]

// Problem
problem Lap (uh, vh, solver=CG)
   = int3d(Th) (Grad(uh)' * Grad(vh))
   + on(310, 300, uh=dep)
   + on(311, uh=0.)
;

// Falling loop
for(int it = 0; it < 29; it++){
   cout << " ITERATION " << it << endl;
   // Solve
   Lap;
   // Plot
   plot(Th, uh);
   // Sphere falling
   Th = mmg3d(Th, options=opt, displacement=[zero, zero, uh], memory=1000);
}

3.2.10 A first 3d isotropic mesh adaptation process

Tip: Adaptation 3D
load "msh3"
load "TetGen"
load "mshmet"
load "medit"
// Parameters
int nn = 6;
int[int] 11111 = [1, 1, 1, 1]; // labels
int[int] 101 = [0, 1];
int[int] 111 = [1, 1];
real errm = 1e-2; // level of error

// Mesh
mesh3 Th3 = buildlayers(square(nn, nn, region=0, label=11111),
nn, zbound=[0, 1], labelmid=111, labelup=101, labeldown=101);
Th3 = trunc(Th3, (x<0.5) | (y < 0.5) | (z < 0.5), label=1); // remove the ]0.5,1[^3 cube

// Fespace
fespace Vh(Th3, P1);
Vh u, v, usol, h;

// Macro
macro Grad(u) [dx(u), dy(u), dz(u)] // EOM

// Problem
problem Poisson (u, v, solver=CG)
= int3d(Th3)(
    Grad(u)' * Grad(v)
) - int3d(Th3)(
    l*v
) + on(1, u=0)
;

// Loop
for (int ii = 0; ii < 5; ii++){
    // Solve
    Poisson;
    cout << "u min, max = " << u[].min << " " << u[].max << endl;
    h=0.; // for resizing h[] because the mesh change
    h[] = mshmet(Th3, u, normalization=1, aniso=0, nbregul=1, hmin=1e-3, hmax=0.3, err=errm);
    cout << "h min, max = " << h[].min << " " << h[].max << " " << h[].n << " " << Th3.nv << endl;
    plot(u, wait=true);
    errm *= 0.8; // change the level of error
    cout << "Th3 " << Th3.nv << " " << Th3.nt << endl;
    Th3 = tetgreconstruction(Th3, switch="raAQ", sizeofvolume=h*h*h/6.); // rebuild mesh
    medit("U-adap-iso-"+ii, Th3, u, wait=true);
}
### 3.2.11 Build a 2d mesh from an isoline

The idea is to get the discretization of an isoline of fluid meshes, this tool can be useful to construct meshes from image. First, we give an example of the isovalue meshes 0.2 of analytical function $\sqrt{(x - 1/2)^2 + (y - 1/2)^2}$, on unit square.

```plaintext
load "isoline"

real[int,int] xy(3,1); //to store the isoline points
int[int] be(1); //to store the begin, end couple of lines

mesh Th = square(10,10);
fespace Vh(Th, P1);
Vh u = sqrt(square(x-0.5) + square(y-0.5));
real iso = 0.2;
real[int] viso = [iso];
plot(u, viso-viso, Th); //to see the iso line

int nbc = isoline(Th, u, xy, close=1, iso=iso, beginend=be, smoothing=0.1);
```

The `isoline` parameters are `Th` the mesh, the expression $u$, the bidimensional array `xy` to store the list coordinate of the points. The list of named parameter are:

- `iso=` value of the isoline to compute (0 is the default value)
- `close=` close the isoline with the border (default true), we add the part of the mesh border such the value is less than the isovalue
- `smoothing=` number of smoothing process is the $l''s$ where $l$ is the length of the current line component, $r$ the ratio, $s$ is smoothing value. The smoothing default value is 0.
- `ratio=` the ratio (1 by default).
- `eps=` relative $\epsilon$ (default 1e-10)
- `beginend=` array to get begin, end couple of each of sub line (resize automatically)
- `file=` to save the data curve in data file for gnuplot

In the array `xy` you get the list of vertices of the isoline, each connex line go from $i = i_0^c, \ldots, i_1^c - 1$ with $i_0^c = be(2*c)$ $i_1^c = be(2*c+1)$, and where $x_i = xy(0,i), y_i = xy(1,i), l_i = xy(2,i)$.

Here $l_i$ is the length of the line (the origin of the line is point $i_0^c$).

The sense of the isoline is such that the upper part is at the left size of the isoline. So here : the minimum is a point 0.5, 0.5 so the curve 1 turn in the clockwise sense, the order of each component are sort such that the number of point by component is decreasing.

```plaintext
cout << "Number of the line component = " << nbc << endl;
cout << "Number of points = " << xy.m << endl;
cout << "be = " << be << endl;

// shows the lines component
for (int c = 0; c < nbc; ++c){
  int i0 = be[2*c], i1 = be[2*c+1]-1;
  cout << "Curve " << c << endl;
  for(int i = i0; i <= i1; ++i)
    cout << "x= " << xy(0,i) << " y= " << xy(1,i) << " s= " << xy(2,i) << ",", " curve "+c);
plot([xy(0, i0:i1), xy(1, i0:i1)], wait=true, viso-viso, cmm=" curve "+c);
}
```

(continues on next page)
We also have a new function to easily parametrize a discrete curve defined by the couple \( b_e, x_y \).

```plaintext
border Curve0(t=0, 1){
    int c=0; // component 0
    int i0=be[2*c], i1=be[2*c+1]-1;
    P=Curve(xy, i0, il, t); // Curve 0
    label=1;
}

border Curvel(t=0, 1){
    int c=1; // component 1
    int i0=be[2*c], i1=be[2*c+1]-1;
    P=Curve(xy, i0, il, t); // Curve 1
    label=1;
}
plot(Curvel(100)); // show curve
mesh Th = buildmesh(Curvel(-100));
plot(Th, wait=true);
```

Secondly, we use this idea to build meshes from an image, we use the plugins `ppm2rnm` to read `pgm` a gray scale image and then we extract the gray contour at level 0.25.

**Tip:** Leman lake

```plaintext
load "ppm2rnm"
load "isoline"

// Parameters
string leman = "LemanLake.pgm";
real AreaLac = 580.03; // in km^2
real hsize = 5;
real[int, int] Curves(3, 1);
int[int] be(1);
int nc; // nb of curve
{
    real[int, int] ff1(lemman); // read image
    // and set it in a rect. array
    int nx = ff1.n, ny = ff1.m;
    // build a Cartesian mesh such that the origin is in the right place.
    mesh Th = square(nx-1, ny-1, [(nx-1)*x, (ny-1)*(1-y)]);
    // warning the numbering of the vertices (x,y) is
    // given by $i = x/nx + nx*y/ny$
    fespace Vh(Th, P1);
    Vh f1;
    f1[] = ff1; // transform array in finite element functions.
    nc = isoline(Th, f1, iso=0.25, close=1, Curves, beginend=be, smoothing=0.1, 
    →ratio=0.5);
}

// The longest isoline: the lake
int ic0 = be(0), ic1 = be(1)-1;
```

(continues on next page)
3.3 Finite element

As stated in tutorials, FEM approximates all functions \( w \) as:

\[
w(x, y) \simeq w_0 \phi_0(x, y) + w_1 \phi_1(x, y) + \cdots + w_{M-1} \phi_{M-1}(x, y)
\]

with finite element basis functions \( \phi_k(x, y) \) and numbers \( w_k \) \( (k = 0, \cdots, M - 1) \). The functions \( \phi_k(x, y) \) are constructed from the triangle \( T_{ik} \), and called shape functions.

In FreeFEM, the finite element space:

\[
V_h = \{ w \mid w_0 \phi_0 + w_1 \phi_1 + \cdots + w_{M-1} \phi_{M-1}, w_i \in \mathbb{R} \}
\]

is easily created by

- in 2d
  
  ```c
  fespace IDspace(IDmesh,IDFE);
  ```

or with \( \ell \) pairs of periodic boundary conditions in 2D:
fespace IDspace(IDmesh,<IDFE>,
    periodic=[[la1, sa1], [lb1, sbl],
    ... [lak, sak], [lbk, sbl]]);

• in 3D:

fespace IDspace(IDmesh3,<IDFE>,
    periodic=[[la1, sa1, ta1], [lb1, sb1, tbl],
    ... [lak, sak, tak], [lbk, sbl, tbl]]);

• in surface 3D:

fespace IDspace(IDmeshS,<IDFE>,
    periodic=[[la1, sa1, ta1], [lb1, sb1, tbl],
    ... [lak, sak, tak], [lbk, sbl, tbl]]);

where IDspace is the name of the space (e.g. \(V_h\)), IDmesh IDmesh3 IDmeshS `is respectively the name of the associated :freefem:`mesh, mesh3, meshS` and <IDFE> is an identifier of finite element type.

In 2D we have a pair of periodic boundary conditions, if \([la_i, sa_i], [lb_i, sb_i]\) is a pair of int, and the 2 labels \(la_i\) and \(lb_i\) refer to 2 pieces of boundary to be in equivalence.

If \([la_i, sa_i], [lb_i, sb_i]\) is a pair of real, then \(sa_i\) and \(sb_i\) give two common abscissa on the two boundary curves, and two points are identified as one if the two abscissa are equal.

In 2D, we have a pair of periodic boundary conditions, if \([la_i, sa_i, ta_i], [lb_i, sb_i, tb_i]\) is a pair of int, the 2 labels \(la_i\) and \(lb_i\) define the 2 pieces of boundary to be in equivalence.

If \([la_i, sa_i, ta_i], [lb_i, sb_i, tb_i]\) is a pair of real, then \(sa_i\), \(ta_i\) and \(sb_i\), \(tb_i\) give two common parameters on the two boundary surfaces, and two points are identified as one if the two parameters are equal.

Note: The 2D mesh of the two identified borders must be the same, so to be sure, use the parameter fixedborder=true in buildmesh command (see fixedborder).

### 3.3.1 List of the types of finite elements

As of today, the known types of finite elements are:

- \([P0]\) piecewise constant discontinuous finite element (2d, 3d, surface 3d), the degrees of freedom are the barycenter element value.

\[
P_0^h = \{ v \in L^2(\Omega) \mid \text{for all } K \in \mathcal{T}_h \text{ there is } \alpha_K \in \mathbb{R} : v|_K = \alpha_K \} \tag{3.2}
\]

- \([P1]\) piecewise linear continuous finite element (2d, 3d, surface 3d), the degrees of freedom are the vertices values.

\[
P_1^h = \{ v \in H^1(\Omega) \mid \forall K \in \mathcal{T}_h, v|_K \in P_1 \} \tag{3.3}
\]

- \([P1dc]\) piecewise linear discontinuous finite element (2d, 3d with load"Element_P1dc1")
\[ P_{1,h}^1 = \{ v \in L^2(\Omega) \mid \forall K \in T_h, v|_K \in P_1 \} \quad (3.4) \]

**Warning:** Due to an interpolation problem, the degree of freedom is not the vertices but three vertices which move inside \( T(X) = G + .99(X - G) \) where \( G \) is the barycenter.

- \([P1b]\) piecewise linear continuous finite element plus bubble (2d, 3d)

  **The 2D Case:**

  \[ P_{1,b,h}^1 = \{ v \in H^1(\Omega) \mid \forall K \in T_h, v|_K \in P_1 \oplus \text{Span}\{\lambda^K_0, \lambda^K_1, \lambda^K_2\} \} \quad (3.5) \]

  **The 3D Case:**

  \[ P_{1,b,h}^1 = \{ v \in H^1(\Omega) \mid \forall K \in T_h, v|_K \in P_1 \oplus \text{Span}\{\lambda^K_0, \lambda^K_1, \lambda^K_2, \lambda^K_3\} \} \quad (3.6) \]

  where \( \lambda^K_i, i = 0, \ldots, d \) are the \( d + 1 \) barycentric coordinate functions of the element \( K \) (triangle or tetrahedron).

- \([P1bl,P1bl3d]\) piecewise linear continuous finite element plus linear bubble (with load "Element_P1bl" 2d, 3d).

  The bubble is built by splitting the \( K \), a barycenter in \( d + 1 \) sub element. (need load "Element_P1bl")

- \([P2, P2]\) piecewise \( P_2 \) continuous finite element (2d, 3d, surface 3d)

  \[ P_{2,h}^2 = \{ v \in H^1(\Omega) \mid \forall K \in T_h, v|_K \in P_2 \} \]

  where \( P_2 \) is the set of polynomials of \( \mathbb{R}^2 \) of degrees \( \leq 2 \).

- \([P2b, P2b3d]\) piecewise \( P_2 \) continuous finite element plus bubble (2d, 3d with load "Element_P2bule3")

  **The 2D Case:**

  \[ P_{2,b,h}^2 = \{ v \in H^1(\Omega) \mid \forall K \in T_h, v|_K \in P_2 \oplus \text{Span}\{\lambda^K_0, \lambda^K_1, \lambda^K_2\} \} \]

  **The 3D Case:**

  \[ P_{2,b,h}^2 = \{ v \in H^1(\Omega) \mid \forall K \in T_h, v|_K \in P_2 \oplus \text{Span}\{\lambda^K_0, \lambda^K_1, \lambda^K_2, \lambda^K_3\} \} \]

- \([P2dc]\) piecewise \( P_2 \) discontinuous finite element (2d)

  \[ P_{2,dc,h}^2 = \{ v \in L^2(\Omega) \mid \forall K \in T_h, v|_K \in P_2 \} \]
Warning: Due to an interpolation problem, the degree of freedom is not the six P2 nodes but six nodes which move inside $T(X) = G + .99(X - G)$ where $G$ is the barycenter.

- \([P2h]\) quadratic homogeneous continuous (without \(P1\)).
- \([P3]\) piecewise $P_3$ continuous finite element (2d) (needs load "Element_P3")

$$\mathcal{P}_h^3 = \{ v \in H^1(\Omega) \mid \forall K \in \mathcal{T}_h, v|_K \in P_3 \}$$

where $P_3$ is the set of polynomials of $\mathbb{R}^2$ of degrees $\leq 3$.
- \([P3dc]\) piecewise $P_3$ discontinuous finite element (2d) (needs load "Element_P3dc")

$$\mathcal{P}_{dc}^3 = \{ v \in L^2(\Omega) \mid \forall K \in \mathcal{T}_h, v|_K \in P_3 \}$$

where $P_3$ is the set of polynomials of $\mathbb{R}^2$ of degrees $\leq 3$.
- \([P4]\) piecewise $P_4$ continuous finite element (2d) (needs load "Element_P4")

$$\mathcal{P}_h^4 = \{ v \in H^1(\Omega) \mid \forall K \in \mathcal{T}_h, v|_K \in P_4 \}$$

where $P_4$ is the set of polynomials of $\mathbb{R}^2$ of degrees $\leq 4$.
- \([P4dc]\) piecewise $P_4$ discontinuous finite element (2d) (needs load "Element_P4dc")

$$\mathcal{P}_{dc}^4 = \{ v \in L^2(\Omega) \mid \forall K \in \mathcal{T}_h, v|_K \in P_3 \}$$

where $P_3$ is the set of polynomials of $\mathbb{R}^2$ of degrees $\leq 3$.
- \([P0Edge]\) piecewise $P_0$ discontinuous finite element (2d) contained on each edge of the mesh.
- \([P1Edge]\) piecewise $P_1$ discontinuous finite element (2d) (needs load "Element_PkEdge") $P_1$ on each edge of the mesh.
- \([P2Edge]\) piecewise $P_2$ discontinuous finite element (2d) (needs load "Element_PkEdge") $P_2$ on each edge of the mesh.
- \([P3Edge]\) piecewise $P_3$ discontinuous finite element (2d) (needs load "Element_PkEdge") $P_3$ on each edge of the mesh.
- \([P4Edge]\) piecewise $P_4$ discontinuous finite element (2d) (needs load "Element_PkEdge") $P_4$ on each edge of the mesh.
- \([P5Edge]\) piecewise $P_5$ discontinuous finite element (2d) (needs load "Element_PkEdge") $P_5$ on each edge of the mesh.
- \([P2Morley]\) piecewise $P_2$ non conform finite element (2d) (needs load "Morley")

$$\mathcal{P}_h^2 = \left\{ v \in L^2(\Omega) \mid \forall K \in \mathcal{T}_h, v|_K \in P_3; \begin{cases} v \text{ continuous at vertices,} \\ \partial_n v \text{ continuous at middle of edge,} \end{cases} \right\}$$

where $P_3$ is the set of polynomials of $\mathbb{R}^2$ of degrees $\leq 2$.  

3.3. Finite element
**Warning:** To build the interplant of a function $u$ (scalar) for this finite element, we need the function and 2 partial derivatives $(u, u_x, u_y)$, creating this vectorial finite element with 3 components $(u, u_x, u_y)$.

See our example for solving the BiLaplacien problem:

```plaintext
load "Morley"

// Parameters
int nn = 10;
real h = 0.01;
real f = 1;

// Mesh
mesh Th = square(nn, nn);
Th = adaptmesh(Th, h, IsMetric=1);

// Fespace
fespace Vh(Th, P2Morley); //The Morley finite element space
Vh [u, ux, uy], [v, vx, vy];

// Macro
macro bilaplacien(u, v) (dxx(u)*dxx(v) + dyy(u)*dyy(v) + 2.*dxy(u)*dxy(v)) //

// Problem
solve bilap ([u, ux, uy], [v, vx, vy])
   = int2d(Th)(
       bilaplacien(u, v)
   )
   - int2d(Th)(
       f*v
   )
   + on(1, 2, 3, 4, u=0, ux=0, uy=0)
;

// Plot
plot(u, cmm="u");
```

- $[\text{HCT}]\ P_3 C^1$ conforms finite element (2d) (needs load "Element_HCT") one 3 sub triangles.

  Lets call $\mathcal{T}_h^\Delta$ the sub mesh of $\mathcal{T}_h$ where all triangles are split in 3 at the barycenter.

  $$\mathbb{P}_h^{HCT} = \left\{ v \in C^1(\Omega) \mid \forall K \in \mathcal{T}_h^\Delta, v|_K \in P_3 \right\}$$

  where $P_3$ is the set of polynomials of $\mathbb{R}^2$ of degrees $\leq 3$.

  The degrees of freedom are the values of the normal derivative at the mid-point of each edge $[\text{BERNADOU1980}]$.

**Warning:** To build the interplant of a function $u$ (scalar) for this finite element, we need the function and 2 partial derivatives $(u, u_x, u_y)$, creating this vectorial finite element with 3 components $(u, u_x, u_y)$ like in previous finite element.
• [P2BR] (needs load "BernadiRaugel") the Bernadi Raugel Finite Element is a Vectorial element (2d) with 2 components, see [BERNARDI1985].

It is a 2D coupled Finite Element, where the Polynomial space is \( P^2_2 \) with 3 normal bubble edge functions (\( P_3 \)). There are 9 degrees of freedom:

- 2 components at each of the 3 vertices and
- the 3 flux on the 3 edges.

• [RT0, RT03d] Raviart-Thomas finite element of degree 0.

The 2D Case:

\[
RT0_h = \left\{ v \in H(\text{div}) \mid \forall K \in T_h, v|_K(x, y) = \left[ \frac{\alpha}{\alpha_K} + \beta K \right] \frac{x}{y} \right\} \tag{3.7}
\]

The 3D Case:

\[
RT0_h = \left\{ v \in H(\text{div}) \mid \forall K \in T_h, v|_K(x, y, z) = \left[ \frac{\alpha}{\alpha_K} + \beta K \right] \frac{x}{y} \right\} \tag{3.8}
\]

where by writing \( \text{div} w = \sum_{i=1}^d \partial w_i / \partial x_i \) with \( w = (w_i)_{i=1}^d \):

\[
H(\text{div}) = \{ w \in L^2(\Omega)^d \mid \text{div} w \in L^2(\Omega) \}
\]

and where \( \alpha_K, \beta_K \) are real numbers.

• [RT0Ortho] Raviart-Thomas Orthogonal, or Nedelec finite element type I of degree 0 in dimension 2

\[
RT0Ortho_h = \left\{ v \in H(\text{curl}) \mid \forall K \in T_h, v|_K(x, y) = \left[ \frac{\alpha}{\alpha_K} + \beta K \right] \frac{-y}{x} \right\} \tag{3.9}
\]

• [Edge03d] 3d Nedelec finite element or Edge Element of degree 0.

\[
Edge0_h = \left\{ v \in H(\text{Curl}) \mid \forall K \in T_h, v|_K(x, y, z) = \left[ \frac{\alpha}{\alpha_K} + \beta K \right] \times \frac{x}{y} \right\} : \text{label} : \text{eq} : \text{Edge03d}
\]

where by writing \( \text{curl} w = \left[ \partial w_2 / \partial x_3 - \partial w_3 / \partial x_2, \partial w_3 / \partial x_1 - \partial w_1 / \partial x_3, \partial w_1 / \partial x_2 - \partial w_2 / \partial x_1 \right] \) with \( w = (w_i)_{i=1}^d \):

\[
H(\text{curl}) = \{ w \in L^2(\Omega)^d \mid \text{curl} w \in L^2(\Omega)^d \}
\]

and \( \alpha_1, \alpha_2, \alpha_3, \beta_1, \beta_2, \beta_3 \) are real numbers.

• [Edge13d] (needs load "Element_Mixte3d") 3d Nedelec finite element or Edge Element of degree 1.

• [Edge23d] (needs load "Element_Mixte3d") 3d Nedelec finite element or Edge Element of degree 2.

• [P1nc] piecewise linear element continuous at the mid-point of the edge only in 2D (Crouzeix-Raviart Finite Element 2D).

• [P2nc] piecewise quadratic plus a P3 bubble element with the continuity of the 2 moments on each edge (needs load "Element_P2nc")

• [RT1] (needs load "Element_Mixte")

\[
RT1_h = \left\{ v \in H(\text{div}) \mid \forall K \in T_h, \alpha_1, \alpha_2, \beta_K \in P^2_1, P_0, v|_K(x, y) = \left[ \frac{\alpha}{\alpha_K} + \beta K \right] \frac{x}{y} \right\} \tag{3.10}
\]

3.3. Finite element
• \([\text{RT1Ortho}]\) (needs load "Element_Mixte")

\[
RT_{1h} = \left\{ v \in H(\text{curl}) \mid \forall K \in \mathcal{T}_h, \alpha_K^1, \alpha_K^2, \beta_K \in P^2, \, v|_K(x,y) = \frac{\alpha_K^1}{\alpha_K^2} + \beta_K \right\}
\] (3.11)

• \([\text{RT2}]\) (needs load "Element_Mixte")

\[
RT_{2h} = \left\{ v \in H(\text{div}) \mid \forall K \in \mathcal{T}_h, \alpha_K^1, \alpha_K^2, \beta_K \in P^2, \, v|_K(x,y) = \frac{\alpha_K^1}{\alpha_K^2} + \beta_K \right\}
\] (3.12)

• \([\text{RT2Ortho}]\) (needs load "Element_Mixte")

\[
RT_{2h} = \left\{ v \in H(\text{curl}) \mid \forall K \in \mathcal{T}_h, \alpha_K^1, \alpha_K^2, \beta_K \in P^2, \, v|_K(x,y) = \frac{\alpha_K^1}{\alpha_K^2} + \beta_K \right\}
\] (3.13)

• \([\text{BDM1}]\) (needs load "Element_Mixte") the Brezzi-Douglas-Marini finite element:

\[
BDM_{1h} = \left\{ v \in H(\text{div}) \mid \forall K \in \mathcal{T}_h, \, v|_K \in P^2 \right\}
\] (3.14)

• \([\text{BDM1Ortho}]\) (needs load "Element_Mixte") the Brezzi-Douglas-Marini Orthogonal also call Nedelec of type II, finite element

\[
BDM_{1Ortho} = \left\{ v \in H(\text{curl}) \mid \forall K \in \mathcal{T}_h, \, v|_K \in P^2 \right\}
\] (3.15)

• \([\text{FEQF}]\) (needs load "Element_QF") the finite element to store functions at default quadrature points (so the quadrature is \(qf5pT\) in 2D and is \(qfV5\) in 3d).

For over quadrature you have the following corresponding finite element’s quadrature formula.

- FEQF1 \(\rightarrow\) \(qf1pT\),
- FEQF2 \(\rightarrow\) \(qf2pT\),
- FEQF5 \(\rightarrow\) \(qf5pT\),
- FEQF7 \(\rightarrow\) \(qf7pT\),
- FEQF9 \(\rightarrow\) \(qf9pT\),
- FEQF13d \(\rightarrow\) \(qfV1\),
- FEQF23d \(\rightarrow\) \(qfV2\),
- FEQF53d \(\rightarrow\) \(qfV5\),

You can use this element to optimize the storage and reuse of functions with a long formula inside an integral for non linear processes.
3.3.2 Use of fespace in 2D

With the 2D finite element spaces

\[ X_h = \{ v \in H^1([0,1]^2) \mid \forall K \in T_h \, v|_K \in P_1 \} \]

\[ X_{ph} = \{ v \in X_h \mid v(0) = v(1), v(1) = v(0) \} \]

\[ M_h = \{ v \in H^1([0,1]^2) \mid \forall K \in T_h \, v|_K \in P_2 \} \]

\[ R_h = \{ v \in H^1([0,1]^2)^2 \mid \forall K \in T_h \, v|_K(x,y) = \frac{\partial v}{\partial x} + \gamma \frac{\partial v}{\partial y} \} \]

when \( T_h \) is a mesh 10 \times 10 of the unit square \([0,1]^2\), we only write in FreeFEM:

```freefem
mesh Th = square(10, 10);
spaced Xh(Th, P1); //scalar FE
spaced Xph(Th, P1, 4, periodic=[[2, y], [4, y], [1, x], [3, x]]); //bi-periodic FE
spaced Mh(Th, P2); //scalar FE
spaced Rh(Th, RT0); //vectorial FE
```

where Xh, Mh, Rh expresses finite element spaces (called FE spaces) \( X_h \), \( M_h \), \( R_h \), respectively.

To use FE-functions \( u_h, v_h \in X_h, p_h, q_h \in M_h \) and \( U_h, V_h \in R_h \), we write:

```freefem
Xh uh, vh;
Xph uph, vph;
Mh ph, qh;
Rh [Uxh, Uyh], [Vxh, Vyh];
Xh[int] Uh(10); //array of 10 functions in Xh
Rh[int] [Wxh, Wyh](10); //array of 10 functions in Rh
Wxh[5](0.5,0.5); //the 6th function at point (0.5, 0.5)
Wxh[5][]; //the array of the degree of freedom of the 6th function
```

The functions \( U_h, V_h \) have two components so we have

\[ U_h = \begin{bmatrix} U_{xh} \\ U_{yh} \end{bmatrix} \quad \text{and} \quad V_h = \begin{bmatrix} V_{xh} \\ V_{yh} \end{bmatrix} \]

3.3.3 Use of fespace in 3D

With the 3D finite element spaces

\[ X_h = \{ v \in H^1([0,1]^3) \mid \forall K \in T_h \, v|_K \in P_1 \} \]

\[ X_{ph} = \{ v \in X_h \mid v(0) = v(1), v(1) = v(0) \} \]

\[ M_h = \{ v \in H^1([0,1]^3) \mid \forall K \in T_h \, v|_K \in P_2 \} \]

\[ R_h = \{ v \in H^1([0,1]^3)^2 \mid \forall K \in T_h \, v|_K(x,y,z) = \frac{\partial v}{\partial x} + \gamma \frac{\partial v}{\partial y} + \delta \frac{\partial v}{\partial z} \} \]

when \( T_h \) is a mesh 10 \times 10 \times 10 of the unit cubic \([0,1]^3\), we write in FreeFEM:
//label: 0 up, 1 down, 2 front, 3 left, 4 back, 5 right
int nn=10;

mesh3 Th=buildlayers(square(nn,nn,region=0),nn,
    zbound=[zmin,zmax], labelmid=rmid, reffaceup = rup,
    reffacelow = rdown);

fespace Xh(Th, P1); //scalar FE
// a FE space with full periodic condition in 3 axes
fespace Xph(Th, P1, periodic=[[1,y,z],[2,y,z],
    [3,x,z],[4,x,z],[5,x,y],[6,x,y]]);

fespace Mh(Th, P2); //scalar FE
fespace Rh(Th, RT03d); //vectorial FE

where Xh, Mh, Rh expresses finite element spaces (called FE spaces) \( X_h \), \( M_h \), \( R_h \), respectively.

The functions \( U_h, V_h \) have two components so we have

### 3.3.4 Use of fespace in surface 3D

With the 3D finite element spaces

\[
X_h = \{ v \in H^1([0,1]^3) | \forall K \in T_h \quad v|_K \in P_1 \}
\]

```plaintext
mesh3 Th = square3(10, 10);
fespace Xh(Th, P1); //scalar FE
```

where \( X_h \) expresses finite element spaces (called FE spaces) \( X_h \), respectively.

To use FE-functions \( u_h, v_h \in X_h, p_h, q_h \in M_h \) and \( U_h, V_h \in R_h \), we write:

```plaintext
Xh uh, vh;
Xh[int] Uh(10);  //array of 10 functions in Xh
```

### 3.3.5 Finite Element functions

To define and use FE-functions \( u_h, v_h \in X_h, p_h, q_h \in M_h \) and \( U_h, V_h \in R_h \), we write:

```plaintext
Xh uh, vh;
Xph uph, vph;
Mh ph, qh;
Rh [Uxh, Uyh, Uyzh], [Vxh, Vyh, Vyzh];
Xh[int] Uh(10);  //array of 10 functions in Xh
Rh[int] [Wxh, Wyh, Wzh](10);  //array of 10 functions in Rh
Wxh[5](0.5,0.5,0.5);  //the 6th function at point (0.5, 0.5, 0.5)
Wxh[5][];  //the array of the degree of freedom of the 6th function
```

The functions \( U_h, V_h \) have three components, so we have:

\[
U_h = \begin{bmatrix} (U_h)_x \\ (U_h)_y \\ (U_h)_z \end{bmatrix} \quad \text{and} \quad V_h = \begin{bmatrix} (V_h)_x \\ (V_h)_y \\ (V_h)_z \end{bmatrix}
\]

**Note:** One challenge of the periodic boundary condition is that the mesh must have equivalent faces.
The buildlayers mesh generator splits each quadrilateral face with the diagonal passing through the vertex with maximum number, so to be sure to have the same mesh one both face periodic the 2D numbering in corresponding edges must be compatible (for example the same variation).

By Default, the numbering of square vertex is correct.

To change the mesh numbering you can use the change function like:

```plaintext
1 { 2 3 int[int] old2new(0:Th.nv-1); //array set on 0, 1, .., nv-1 4 fespace Vh2(Th, P1); 5 Vh2 sorder = x+y; //choose an order increasing on 4 square borders with x or y 6 sort(sorder[], old2new); //build the inverse permutation 7 int[int] new2old = old2new^-1; //inverse the permutation 8 Th = change(Th, renumv=new2old); 9 }
```

The full example is in examples.

### 3.3.6 Lagrangian Finite Elements

**P0-element**

For each triangle \((d=2)\) or tetrahedron \((d=3)\) \(T_k\), the basis function \(\phi_k\) in \(Vh(Th, \ P0)\) is given by:

\[
\phi_k(x) = \begin{cases} 
1 & \text{if } (x) \in T_k \\
0 & \text{if } (x) \notin T_k
\end{cases}
\]

If we write:

```plaintext
1 Vh(Th, P0); 2 Vh fh = f(x, y);
```

then for vertices \(q^k, i = 1, 2, ..d + 1\) in Fig. 3.30, \(f_h\) is built as \(f_h = f_h(x, y) = \sum_k f \left( \frac{\sum_i q^{k_i}}{d+1} \right) \phi_k\)

See Fig. 3.31b for the projection of \(f(x, y) = \sin(\pi x) \cos(\pi y)\) on \(Vh(Th, \ P0)\) when the mesh \(Th\) is a \(4 \times 4\)-grid of \([-1,1]^2\) as in Fig. 3.31a.

**P1-element**

For each vertex \(q^i\), the basis function \(\phi_i\) in \(Vh(Th, \ P1)\) is given by:

\[
\phi_i(x, y) = a_i^k + b_i^k x + c_i^k y \text{ for } (x, y) \in T_k, \\
\phi_i(q^j) = 1, \quad \phi_i(q^j) = 0 \text{ if } i \neq j
\]

The basis function \(\phi_{k_i}(x, y)\) with the vertex \(q^{k_i}\) in Fig. 3.30 at point \(p = (x, y)\) in triangle \(T_k\) simply coincide with the barycentric coordinates \(\lambda_i^k\) (area coordinates):

\[
\phi_{k_i}(x, y) = \lambda_i^k(x, y) = \frac{\text{area of triangle}(p, q^{k_2}, q^{k_3})}{\text{area of triangle}(q^{k_2}, q^{k_2}, q^{k_3})}
\]

If we write:

```plaintext
3.3. Finite element
```
\[ \varphi_{k_1}(x, y) = \lambda_{k_1}(x, y)(2\lambda_{k_1}(x, y) - 1) \]

and for the mid-point \( q^{k_2} \):

\[ \varphi_{k_2}(x, y) = 4\lambda_{k_1}(x, y)\lambda_{k_1}(x, y) \]

If we write:

1. \( Vh(Th, P_2) \);
2. \( Vh\ f_h = f(x, y) \);

then:

\[ f_h = f_h(x, y) = \sum_{i=1}^{M} f(q^i)\varphi_i(x, y) \quad \text{(summation over all vertex or mid-point)} \]

See \textit{Projection to Vh(Th, P2)} for the projection of \( f(x, y) = \sin(\pi x)\cos(\pi y) \) into \( Vh(Th, P_2) \).
3.3. Finite element

**Fig. 3.31:** Finite element $\mathcal{P}_0$

**Fig. 3.32:** Finite elements $\mathcal{P}_1$, $\mathcal{P}_2$
3.3.7 Surface Lagrangian Finite Elements

Definition of the surface P1 Lagragian element

To build the surface Pk-Lagrange, the main idea is to consider the usual 2d Lagrangian Finite Elements; and its writing in barycentric coordinates; apply a space transformation and barycentric properties. The FreeFEM finite elements for surface problem are: \( P_0 \rightarrow P_1 \rightarrow P_2 \rightarrow P_3 \).

0) Notation

- Let \( \hat{\mathcal{K}} \) be the shape triangle in the space \( \mathbb{R}^2 \) of vertice \((i_0, i_1, i_2)\)
- Let \( K \) be a triangle of the space \( \mathbb{R}^3 \) of vertice \((A_0, A_1, A_2)\)
- \( x_q \) a quadrature point on \( K \)
- \( X_q \) a quadrature point on \( A \)
- \( P_{1d} \) designates 2d P1 Lagrangian Finite Elements
- \( P_{1s} \) designates surface 3d P1 Lagrangian Finite Elements
- \((\lambda_i)_{i=0}^{2} \) shape fonction of \( \hat{\mathcal{K}} \) (P1d)
- \((\psi_i)_{i=0}^{2} \) shape fonction of of \( K \) (P1s)

1) Geometric transformation: from the current FE to the reference FE

Let be \( \hat{x} = \begin{pmatrix} \hat{x} \\ \hat{y} \end{pmatrix} \) a point of the triangle \( \hat{\mathcal{K}} \subset \mathbb{R}^2 \) and \( X = \begin{pmatrix} x \\ y \\ z \end{pmatrix} \) a point of the triangle \( K \subset \mathbb{R}^3 \), where \( \hat{x} \) and \( X \) are expressed in barycentric coordinates.

The motivation here is to parameterize the 3d surface mesh to the reference 2d triangle, thus to be reduced to a finite element 2d P1. Let’s define a geometric transformation \( F \), such as \( F: \mathbb{R}^2 \rightarrow \mathbb{R}^3 \).

However, thus defines transformation \( F \) as not bijective.

So, consider the following approximation

\[
\hat{F}: \mathbb{R}^2 \rightarrow \mathbb{R}^3 \\
\hat{x} \rightarrow X \\
\begin{pmatrix} x \\ y \\ 0 \end{pmatrix} \rightarrow \begin{pmatrix} \frac{A_0 \cdot A_1}{A_0 \cdot A_2} \\ \frac{A_0 \cdot A_2}{A_0 \cdot A_1} \end{pmatrix} (\hat{x} - A_0)
\]

where \( \wedge \) denote the usual vector product.

Note: \( \frac{A_0 \cdot A_1}{A_0 \cdot A_2} = \begin{pmatrix} n_x \\ n_y \end{pmatrix} \) defines the normal to the tangent plane generated by \( (A_0, A_0 \cdot A_1, A_0 \cdot A_2) \).

The affine transformation \( \hat{F} \) allows you to pass from the 2d reference triangle, which we project in \( \mathbb{R}^3 \) to the 3d current triangle, discretizing the surface we note \( \Gamma \).

Then \( \hat{F}^{-1} \) is well defined and allows to return to the reference triangle \( \hat{\mathcal{K}} \), to the usual coordinates of \( \mathbb{R}^2 \) completed by the coordinate \( z = 0 \).

2) Interpolation element fini
Remember that the reference geometric element for the finite element $P_1$ is that we are building is the reference triangle $\hat{K}$ in the vertex plane $(i_0, i_1, i_2)$, which we project into space by posing $z = 0$ by the membrane hypothesis.

Hence $i_0 = \begin{pmatrix} 0 \\ 0 \\ 0 \end{pmatrix}$, $i_1 = \begin{pmatrix} 1 \\ 0 \\ 0 \end{pmatrix}$, $i_2 = \begin{pmatrix} 0 \\ 1 \\ 0 \end{pmatrix}$.

Let $X$ be a point of the current triangle $K$, we have $X = \tilde{F}(\hat{x})$. The barycentric coordinates of $X$ in $K$ are given by:

$$X = \sum_{i=0}^{2} A_i \lambda_i(\hat{x})$$

- $A_i$ the points of the current triangle $K$
- $\lambda_i$ basic functions $P_1$ in 2d
- $\lambda_0(x, y) = 1 - x - y$
- $\lambda_1(x, y) = x$
- $\lambda_2(x, y) = y$

We need to define a quadrature formula for the finite element approximation. The usual formulation for a 2d triangle will be used by redefining the quadrature points $X_q = x_q = \begin{pmatrix} \hat{x}_q \\ \hat{y}_q \\ 0 \end{pmatrix}$.

3) The Lagrangian $P_1$ functions and theirs 1st order derivatives

The finite element interpolation gives us the following relationship: $\psi(X) = F^{-1}(\psi)(F^{-1}(X))$. To find the expression of the basic functions $\psi$ on the current triangle $K$, it is sufficient to use the inverse of the transformation $\tilde{F}$ to get back to the reference triangle $\hat{K}$. However in FreeFEM, the definition of the reference finite element, the current geometry is based on barycentric coordinates in order not to use geometric transformation. $\tilde{F}$. The method used here is geometric and based on the properties of the vector product and the area of the current triangle $K$.

i) The shape functions

Let be the triangle $K$ of vertices $i_0, i_1, i_2 \subset \mathbb{R}^3$ and $(\lambda_i)_{i=0}^{2}$ the local barycentric coordinates at $K$. The normal is defined as the tangent plane generated by $(A_0, A_0A_1, A_0A_2)$, $\vec{n} = A_0A_1 \wedge A_0A_2$ avec $||\vec{n}|| = 2 \text{mes} (\hat{K})$.

Le denotes the operator $V$, defines the usual vector product of $\mathbb{R}^3$ such as $V(A, B, C) = (B - A) \wedge (C - A)$

The mixed product of three vectors $u, v, w$, noté $[u, v, w]$, is the determinant of these three vectors in any direct orthonormal basis, thus $(A \wedge V, C) = \det (A, B, C)$ with $(.,.)$ is the usual scalar product of $\mathbb{R}^3$. Let Ph $\mathbb{R}^3$ and P his projected in the triangle $K$ such as:

Let's lay the sub-triangles as follows:

3.3. Finite element
\[ K_0 = (P, A_1, A_2) \]
\[ K_1 = (A_0, P, A_2) \]
\[ K_2 = (A_0, A_1, P) \]

with \( K = K_0 \cup K_1 \cup K_2 \).

**Note:**

**Properties in \( \mathbb{R}^3 \)**

- Let \( \vec{n} \) be the normal to the tangent plane generated by \( (A_0, \overrightarrow{A_0A_1}, \overrightarrow{A_0A_2}) \)
- \( \vec{n} = \overrightarrow{A_0A_1} \land \overrightarrow{A_0A_2} \)
- By definition, \( A = \frac{1}{2} \| < \vec{n}, \vec{n} > \| \) and the vectorial area by \( A^S = \frac{1}{2} < \vec{n}, \vec{n} > \) hence \( A^S(PBC) = \frac{1}{2} < \vec{n}_0, \vec{n} > \), with \( \vec{n}_0 \) the normal vector to the plane (PBC)

Let’s define the respective vector areas

- \( \vec{N}_0(P) = V(P, A_1, A_2) \) the vectorial area of K0
- \( \vec{N}_1(P) = V(A_0, P, A_2) \) the vectorial area of K1
- \( \vec{N}_2(P) = V(A_0, A_1, P) \) the vectorial area of K2

By definition, in 3d, the barycentric coordinates are given by algebraic area ratio: 

\[ \lambda_i(P) = \frac{\langle \vec{N}_i(P), \vec{N} \rangle}{\langle \vec{N}, \vec{N} \rangle} \]

Note that \( \langle \vec{N}_i(P), \vec{N} \rangle = 2 \text{ sign } \text{mes}(K_i) \| \vec{N} \| \) and \( \langle \vec{N}, \vec{N} \rangle = 2 \text{ sign mes}(K) \| \vec{N} \| \), with sign the orientation of the current triangle compared to the reference triangle.

We find the finite element interpolation, \( P = \sum_{i=0}^{2} \lambda_i(P)A_i \).

**ii) 1st order derivatives of Lagrangian P1 FE**

Let \( \vec{Y} \) be any vector of \( \in \mathbb{R}^3 \).

\[ (\vec{N}_2(P), \vec{Y}) = ((A_1 - A_0) \land (P - A_0), Y) \]
\[ = \det(A_1 - A_0, P - A_0, Y) \]
\[ = \det(A_1 - A_0, P, Y) - \det(A_1 - A_0, A_0, Y) \]
Let’s calculate the differential of \((\vec{N}_2(P), Y), \forall Y\)

\[
D_P(\vec{N}_2(P), \vec{Y}) = \det(A_1 - A_0, P', Y) dP
\]

\[
\nabla_P(\vec{N}_2(P), \vec{Y}) = \det (A_1 - A_0, P', \vec{Y})
\]

\[
= -\det(A_1 - A_0, \vec{Y}, P')
\]

\[
= - (A_1 - A_0) \wedge Y, P'
\]

\[
= \vec{Y} \wedge (A_1 - A_0)
\]

Consider in particular \(Y = \vec{N}\), then

\[
\nabla_P(\vec{N}_2(P), \vec{N}) = \vec{N} \wedge (A_1 - A_0)
\]

\[
= \vec{N} \wedge E_2
\]

\[
= -\det(A_1 - A_0, \vec{N}, P')
\]

This leads to \(\nabla_P \lambda_2(P) = \frac{\vec{N} \wedge E_2}{(\vec{N}, \vec{N})}\) .

By similar calculations for \(\vec{N}_0(P)\) et \(\vec{N}_1(P)\)

\[
\nabla_P \lambda_i(P) = \frac{(\vec{N} \wedge E_i)}{(\vec{N}, \vec{N})}
\]

**Note:** With the definition of the surface gradient and the 2d Pk-Lagrange FE used barycentric coordinates, surface Pk-Lagrange FE are trivial.

### 3.3.8 P1 Nonconforming Element

Refer to [THOMASSET2012] for details; briefly, we now consider non-continuous approximations so we will lose the property:

\[
w_h \in V_h \subset H^1(\Omega)
\]

If we write:

```plaintext
forall (Th, P1nc);
forall fh = f(x,y);
```

then:

\[
f_h = f_h(x,y) = \sum_{i=1}^{n_v} f(m^i) \phi_i(x,y) \text{ (summation over all midpoint)}
\]

Here the basis function \(\phi_i\) associated with the mid-point \(m^i = (q^{ki} + q^{ki+1})/2\) where \(q^{ki}\) is the \(i\)-th point in \(T_k\), and we assume that \(j+1 = 0\) if \(j = 3\):

\[
\phi_i(x,y) = a_i^k + b_i^k x + c_i^k y \text{ for } (x,y) \in T_k,
\]

\[
\phi_i(m^i) = 1, \quad \phi_i(m^j) = 0 \text{ if } i \neq j
\]

Strictly speaking \(\partial \phi_i/\partial x, \partial \phi_i/\partial y\) contain Dirac distribution \(\rho \delta_{\partial T_k}\).

The numerical calculations will automatically ignore them. In [THOMASSET2012], there is a proof of the estimation

\[
\left(\sum_{k=1}^{n_v} \int_{T_k} |\nabla w - \nabla w_h|^2 dxdy\right)^{1/2} = O(h)
\]

The basis functions \(\phi_k\) have the following properties.

### 3.3. Finite element
1. For the bilinear form $a$ defined in Fig. 3.34(a) satisfy:

$$a(\phi_i, \phi_j) > 0, \quad a(\phi_i, \phi_j) \leq 0 \text{ if } i \neq j$$

$$\sum_{k=1}^{n_v} a(\phi_i, \phi_k) \geq 0$$

2. $f \geq 0 \Rightarrow u_h \geq 0$

3. If $i \neq j$, the basis function $\phi_i$ and $\phi_j$ are $L^2$-orthogonal:

$$\int_{\Omega} \phi_i \phi_j \, dx \, dy = 0 \quad \text{if } i \neq j$$

which is false for $P_1$-element.

See Fig. 3.34a for the projection of $f(x, y) = \sin(\pi x) \cos(\pi y)$ into $Vh(Th, P1nc)$.

### 3.3.9 Other FE-space

For each triangle $T_k \in T_h$, let $\lambda_{k_1}(x, y), \lambda_{k_2}(x, y), \lambda_{k_3}(x, y)$ be the area coordinate of the triangle (see Fig. 3.30), and put:

$$\beta_k(x, y) = 27 \lambda_{k_1}(x, y) \lambda_{k_2}(x, y) \lambda_{k_3}(x, y)$$

called bubble function on $T_k$. The bubble function has the feature: 1. $\beta_k(x, y) = 0 \quad \text{if } (x, y) \in \partial T_k$.

2. $\beta_k(q^{h_k}) = 1$ where $q^{h_k}$ is the barycenter $\frac{x^1 + x^2 + x^3}{3}$.

If we write:
\begin{verbatim}
Vh(Th, P1b); Vh fh = f(x,y);
\end{verbatim}

then:

\[ fh = f_h(x, y) = \sum_{i=1}^{n_1} f(q^i) \phi_i(x, y) + \sum_{k=1}^{n_2} f(q^k) \beta_k(x, y) \]

See Fig. 3.34b for the projection of \( f(x, y) = \sin(\pi x) \cos(\pi y) \) into \( \text{Vh}(\text{Th}, \text{P1b}) \).

### 3.3.10 Vector Valued FE-function

Functions from \( \mathbb{R}^2 \) to \( \mathbb{R}^N \) with \( N = 1 \) are called scalar functions and called \textit{vector valued} when \( N > 1 \). When \( N = 2 \)

\begin{verbatim}
fespace Vh(Th, [P0, P1]);
\end{verbatim}

makes the space

\[ V_h = \{(w_1, w_2) \mid w_1 \in \text{Vh}(\text{Th}, P_0), w_2 \in \text{Vh}(\text{Th}, P_1)\} \]

**Raviart-Thomas Element**

In the Raviart-Thomas finite element \( \text{RT}_0 \), the degrees of freedom are the fluxes across edges \( e \) of the mesh, where the flux of the function \( f : \mathbb{R}^2 \rightarrow \mathbb{R}^2 \) is \( \int_e f \cdot n_e \), \( n_e \) is the unit normal of edge \( e \).

This implies an orientation of all the edges of the mesh, for example we can use the global numbering of the edge vertices and we just go from small to large numbers.

To compute the flux, we use a quadrature with one Gauss point, the mid-point of the edge.

Consider a triangle \( T_k \) with three vertices \( (a, b, c) \).

Let's denote the vertices numbers by \( i_a, i_b, i_c \), and define the three edge vectors \( e^1, e^2, e^3 \) by \( \text{sgn}(i_b - i_c)(b - c) \), \( \text{sgn}(i_c - i_a)(c - a) \), \( \text{sgn}(i_a - i_b)(a - b) \).

We get three basis functions:

\[ \phi_1^k = \frac{\text{sgn}(i_b - i_c)}{2|T_k|}(x - a), \quad \phi_2^k = \frac{\text{sgn}(i_c - i_a)}{2|T_k|}(x - b), \quad \phi_3^k = \frac{\text{sgn}(i_a - i_b)}{2|T_k|}(x - c), \]

where \( |T_k| \) is the area of the triangle \( T_k \). If we write:

\begin{verbatim}
Vh(Th, RT0); Vh [f1h, f2h] = [f1(x, y), f2(x, y)];
\end{verbatim}

then:

\[ fh = f_h(x, y) = \sum_{k=1}^{6} \sum_{l=1}^{n_{ijl}} n_{ijl} |e| f_{ij}(m^l) \phi_{ijl} \]

where \( n_{ijl} \) is the \( j_l \)-th component of the normal vector \( n_{ij} \),

\[ \{m_1, m_2, m_3\} = \{\frac{b + c}{2}, \frac{a + c}{2}, \frac{b + a}{2}\} \]

and \( i_l = \{1, 1, 2, 2, 3, 3\}, j_l = \{1, 2, 1, 2, 1, 2\} \) with the order of \( l \).
Fig. 3.35: Normal vectors of each edge

```plaintext
// Mesh
mesh Th = square(2, 2);

// Fespace
fespace Xh(Th, P1);
Xh uh = x^2 + y^2, vh;

fespace Vh(Th, RT0);
Vh [Uxh, Uyh] = [sin(x), cos(y)]; //vectorial FE function

// Change the mesh
Th = square(5,5);
//Xh is unchanged
//Uxh = x; //error: impossible to set only 1 component
//of a vector FE function
vh = Uxh; //ok
//and now vh use the 5x5 mesh
//but the fespace of vh is always the 2x2 mesh

// Plot
plot(uh);

uh = uh; //do a interpolation of uh (old) of 5x5 mesh
//to get the new uh on 10x10 mesh
plot(uh);

vh([x-1/2, y]) = x^2 + y^2; //interpolate vh = ((x-1/2)^2 + y^2)
```

To get the value at a point $x = 1, y = 2$ of the FE function $uh$, or $[Uxh, Uyh]$, one writes:
To get the value of the array associated to the FE function \( uh \), one writes

```plaintext
real value = uh[2,4]; // get value = uh(2, 4)
value = Uxh(2, 4); // get value = Uxh(2, 4)
// OR
x = 1; y = 2;
value = uh; // get value = uh(1, 2)
value = Uxh; // get value = Uxh(1, 2)
value = Uyh; // get value = Uyh(1, 2)
```

3.3.11 A Fast Finite Element Interpolator

In practice, one may discretize the variational equations by the Finite Element method. Then there will be one mesh for \( \Omega_1 \) and another one for \( \Omega_2 \). The computation of integrals of products of functions defined on different meshes is difficult.

Quadrature formula and interpolations from one mesh to another at quadrature points are needed. We present below the interpolation operator which we have used and which is new, to the best of our knowledge.

Let \( T^0_h = \bigcup_k T^0_k, T^1_h = \bigcup_k T^1_k \) be two triangulations of a domain \( \Omega \). Let:

\[
V(T^i_h) = \{ C^0(\Omega^i_h) : f|_{T^i_h} \in P_0 \}, \quad i = 0, 1
\]

be the spaces of continuous piecewise affine functions on each triangulation.
Let $f \in V(T_h^0)$. The problem is to find $g \in V(T_h^1)$ such that:

$$g(q) = f(q) \quad \forall q \text{ vertex of } T_h^1$$

Although this is a seemingly simple problem, it is difficult to find an efficient algorithm in practice.

We propose an algorithm which is of complexity $N^1 \log N^0$, where $N^i$ is the number of vertices of $T_i^h$, and which is very fast for most practical 2D applications.

**Algorithm**

The method has 5 steps.

First a quadtree is built containing all the vertices of the mesh $T_h^0$ such that in each terminal cell there are at least one, and at most 4, vertices of $T_h^0$. For each $q^1$, vertex of $T_h^1$ do:

1. Find the terminal cell of the quadtree containing $q^1$.
2. Find the nearest vertex $q^0_j$ to $q^1$ in that cell.
3. Choose one triangle $T_k^0 \in T_h^0$ which has $q^0_j$ for vertex.
4. Compute the barycentric coordinates $\{\lambda_j\}_{j=1,2,3}$ of $q^1$ in $T_k^0$.
   - if all barycentric coordinates are positive, go to Step 5
   - otherwise, if one barycentric coordinate $\lambda_i$ is negative, replace $T_k^0$ by the adjacent triangle opposite $q^0_i$ and go to Step 4.
   - otherwise, if two barycentric coordinates are negative, take one of the two randomly and replace $T_k^0$ by the adjacent triangle as above.
5. Calculate $g(q^1)$ on $T_k^0$ by linear interpolation of $f$:

$$g(q^1) = \sum_{j=1,2,3} \lambda_j f(q^0_j)$$

**Fig. 3.37:** To interpolate a function at $q^0$, the knowledge of the triangle which contains $q^0$ is needed. The algorithm may start at $q^1 \in T_h^0$ and stall on the boundary (thick line) because the line $q^0q^1$ is not inside $\Omega$. But if the holes are triangulated too (doted line) then the problem does not arise.

Two problems need to be solved:
• What if: math:`q^1` is not in \( \Omega^0_h \)? Then Step 5 will stop with a boundary triangle.

So we add a step which tests the distance of \( q^1 \) with the two adjacent boundary edges and selects the nearest, and so on till the distance grows.

• What if \( \Omega^0_h \) is not convex and the marching process of Step 4 locks on a boundary? By construction Delaunay-Voronoï’s mesh generators always triangulate the convex hull of the vertices of the domain.

Therefore, we make sure that this information is not lost when \( T^0_h, T^1_h \) are constructed and we keep the triangles which are outside the domain on a special list.

That way, in step 5 we can use that list to step over holes if needed.

Note: Sometimes, in rare cases, the interpolation process misses some points, we can change the search algorithm through a global variable `searchMethod`

```c
searchMethod = 0; // default value for fast search algorithm
searchMethod = 1; // safe search algorithm, uses brute force in case of missing point
// (warning: can be very expensive in cases where a lot of points are outside of the domain)
searchMethod = 2; // always uses brute force. It is very computationally expensive.
```

Note: Step 3 requires an array of pointers such that each vertex points to one triangle of the triangulation.

Note: The operator = is the interpolation operator of FreeFEM, the continuous finite functions are extended by continuity to the outside of the domain.

Try the following example:

```c
// Mesh
mesh Ths = square(10, 10);
plot(Ths, wait=true);

// Fespace
fespace Ch(Ths, P2);
Ch us = (x-0.5)*(y-0.5);

fespace Dh(Ths, P2dc);
Dh vs = (x-0.5)*(y-0.5);

fespace Fh(Thg, P2dc);
Fh ug=us, vg=vs;

// Plot
plot(us, ug, wait=true);
plot(vs, vg, wait=true);
```

### 3.3.12 Keywords: Problem and Solve

For FreeFEM, a problem must be given in variational form, so we need a bilinear form \( a(u,v) \), a linear form \( \ell(f,v) \), and possibly a boundary condition form must be added.
FreeFEM Documentation, Release 4.6

Fig. 3.38: Extension of FE-function

```
problem P (u, v)
  = a(u,v) - l(f,v)
  + (boundary condition)
  ;
```

Note: When you want to formulate the problem and solve it in the same time, you can use the keyword `solve`.

**Weak Form and Boundary Condition**

To present the principles of Variational Formulations, also called weak form, for the Partial Differential Equations, let’s take a model problem: a Poisson equation with Dirichlet and Robin Boundary condition.

The problem: Find \( u \) a real function defined on a domain \( \Omega \) of \( \mathbb{R}^d \) (\( d = 2, 3 \)) such that:

\[
-\nabla \cdot (\kappa \nabla u) = f \quad \text{in } \Omega \\
au + \kappa \frac{\partial u}{\partial n} = b \quad \text{on } \Gamma_r \\
u = g \quad \text{on } \Gamma_d
\]

where:

- if \( d = 2 \) then \( \nabla \cdot (\kappa \nabla u) = \partial_x (\kappa \partial_x u) + \partial_y (\kappa \partial_y u) \) with \( \partial_x u = \frac{\partial u}{\partial x} \) and \( \partial_y u = \frac{\partial u}{\partial y} \)
- if \( d = 3 \) then \( \nabla \cdot (\kappa \nabla u) = \partial_x (\kappa \partial_x u) + \partial_y (\kappa \partial_y u) + \partial_z (\kappa \partial_z u) \) with \( \partial_x u = \frac{\partial u}{\partial x} \), \( \partial_y u = \frac{\partial u}{\partial y} \) and \( \partial_z u = \frac{\partial u}{\partial z} \)
- The border \( \Gamma = \partial \Omega \) is split in \( \Gamma_d \) and \( \Gamma_n \) such that \( \Gamma_d \cap \Gamma_n = \emptyset \) and \( \Gamma_d \cup \Gamma_n = \partial \Omega \),
- \( \kappa \) is a given positive function, such that \( \exists \kappa_0 \in \mathbb{R}, \ 0 < \kappa_0 \leq \kappa \).
- \( a \) a given non negative function,
- \( b \) a given function.

Note: This is the well known Neumann boundary condition if \( a = 0 \), and if \( \Gamma_d \) is empty.

In this case the function appears in the problem just by its derivatives, so it is defined only up to a constant (if \( u \) is a solution then \( u + c \) is also a solution).
Let \( v \), a regular test function, null on \( \Gamma_d \), by integration by parts we get:
\[
- \int_{\Omega} \nabla \cdot (\kappa \nabla u) v \, d\omega = \int_{\Omega} \kappa \nabla v \cdot \nabla u \, d\omega - \int_{\Gamma} v \kappa \frac{\partial u}{\partial n} \, d\gamma = \int_{\Omega} f v \, d\omega
\]
where if \( d = 2 \) the \( \nabla v, \nabla u = (\frac{\partial v}{\partial x} + \frac{\partial v}{\partial y}) \),
where if \( d = 3 \) the \( \nabla v, \nabla u = (\frac{\partial v}{\partial x} + \frac{\partial v}{\partial y} + \frac{\partial v}{\partial z}) \),
and where \( n \) is the unitary outer-pointing normal of the \( \Gamma \).

Now we note that \( \kappa \frac{\partial u}{\partial n} = -au + b \) on \( \Gamma_r \) and \( v = 0 \) on \( \Gamma_d \) and \( \Gamma = \Gamma_d \cup \Gamma_n \) thus:
\[
- \int_{\Gamma} v \kappa \frac{\partial u}{\partial n} = \int_{\Gamma_r} au v - \int_{\Gamma_r} bv
\]
The problem becomes:

Find \( u \in V_g = \{ w \in H^1(\Omega) / w = g \text{ on } \Gamma_d \} \) such that:
\[
\int_{\Omega} \kappa \nabla v \cdot \nabla u \, d\omega + \int_{\Gamma_r} au v \, d\gamma = \int_{\Omega} f v \, d\omega + \int_{\Gamma_r} bv \, d\gamma, \quad \forall v \in V_0
\]
where \( V_0 = \{ v \in H^1(\Omega) / v = 0 \text{ on } \Gamma_d \} \)

Except in the case of Neumann conditions everywhere, the problem (3.16) is well posed when \( \kappa \geq \kappa_0 > 0 \).

Note: If we have only the Neumann boundary condition, linear algebra tells us that the right hand side must be orthogonal to the kernel of the operator for the solution to exist.

One way of writing the compatibility condition is:
\[
\int_{\Omega} f \, d\omega + \int_{\Gamma_r} b \, d\gamma = 0
\]
and a way to fix the constant is to solve for \( u \in H^1(\Omega) \) such that:
\[
\int_{\Omega} (\varepsilon u v + \kappa \nabla v \cdot \nabla u) \, d\omega = \int_{\Omega} f v \, d\omega + \int_{\Gamma_r} bv \, d\gamma, \quad \forall v \in H^1(\Omega)
\]

where \( \varepsilon \) is a small parameter (\( \sim \kappa 10^{-10}|\Omega|^{\frac{1}{2}} \)).

Remark that if the solution is of order \( \frac{1}{2} \) then the compatibility condition is unsatisfied, otherwise we get the solution such that \( \int_{\Omega} u = 0 \), you can also add a Lagrange multiplier to solve the real mathematical problem like in the Lagrange multipliers example.

In FreeFEM, the bidimensional problem (3.16) becomes:

```plaintext
problem Pw (u, v)
1 = int2d(Th)( //int_(Omega) kappa nabla v . nabla u
2 kappa* * (dx(u)*dx(v) + dy(u)*dy(v))
3 )
4 + int1d(Th, gn)( //int_(Gamma_r) a u v
5 a * u*v
6 )
7 - int2d(Th)( //int_(Omega) f v
8 f*v
9 )
10 - int1d(Th, gn)( //int_(Gamma_r) b v
11 )
```

(continues on next page)
where \( \Theta \) is a mesh of the bi-dimensional domain \( \Omega \), and \( g_d \) and \( g_n \) are respectively the boundary labels of boundary \( \Gamma_d \) and \( \Gamma_n \).

And the three dimensional problem (3.16) becomes

```plaintext
macro Grad(u) [dx(u), dy(u), dz(u)] //

problem Pw (u, v)

= int3d(Th)( //int_{\Omega}) \kappa \nabla v . \nabla u
  \kappa \ast (\text{Grad}(u) \ast \text{Grad}(v))
+ int2d(Th, gn)( //int_{\Gamma_r} a u v
  a \ast u \ast v
)
- int3d(Th)( //int_{\Omega}) f v
  f \ast v
)
- int2d(Th, gn)( //int_{\Gamma_r} b v
  b \ast v
)
+ on(gd, u=g) //u = g on Gamma_d
;
```

where \( \Theta \) is a mesh of the three dimensional domain \( \Omega \), and \( g_d \) and \( g_n \) are respectively the boundary labels of boundary \( \Gamma_d \) and \( \Gamma_n \).

### 3.3.13 Parameters affecting solve and problem

The parameters are FE functions real or complex, the number \( n \) of parameters is even \((n = 2 \ast k)\), the \( k \) first function parameters are unknown, and the \( k \) last are test functions.

**Note:** If the functions are a part of vectorial FE then you must give all the functions of the vectorial FE in the same order (see Poisson problem with mixed finite element for example).

**Note:** Don’t mix complex and real parameters FE function.

**Warning:** Bug:

The mixing of multiple \textit{fespace} with different periodic boundary conditions are not implemented.

So all the finite element spaces used for tests or unknown functions in a problem, must have the same type of periodic boundary conditions or no periodic boundary conditions.

No clean message is given and the result is unpredictable.

The parameters are:

- \textbf{solver= }LU, CG, Crout, Cholesky, GMRES, sparsesolver, UMFPACK ...
The default solver is `sparsesolver` (it is equal to UMFPACK if no other sparse solver is defined) or is set to `LU` if no direct sparse solver is available.

The storage mode of the matrix of the underlying linear system depends on the type of solver chosen; for `LU` the matrix is sky-line non symmetric, for `Crout` the matrix is sky-line symmetric, for `Cholesky` the matrix is sky-line symmetric positive definite, for `CG` the matrix is sparse symmetric positive, and for `GMRES`, `sparsesolver` or UMFPACK the matrix is just sparse.

- **eps** = a real expression.

  $\varepsilon$ sets the stopping test for the iterative methods like `CG`.

  Note that if $\varepsilon$ is negative then the stopping test is:

  $$||Ax - b|| < |\varepsilon|$$

  if it is positive, then the stopping test is:

  $$||Ax - b|| < \frac{|\varepsilon|}{||Ax_0 - b||}$$

- **init** = boolean expression, if it is false or 0 the matrix is reconstructed.

  Note that if the mesh changes the matrix is reconstructed too.

- **precon** = name of a function (for example `P`) to set the preconditioner.

  The prototype for the function `P` must be:

  ```
  func real[int] P(real[int] xx);
  ```

- **tgv** = Huge value ($10^{30}$) used to implement Dirichlet boundary conditions.

- **tolpivot** = sets the tolerance of the pivot in UMFPACK ($10^{-1}$) and, `LU`, `Crout`, `Cholesky` factorisation ($10^{-20}$).

- **tolpivotsym** = sets the tolerance of the pivot sym in UMFPACK

- **strategy** = sets the integer UMFPACK strategy (0 by default).

### 3.3.14 Problem definition

Below $v$ is the unknown function and $w$ is the test function.

After the “=” sign, one may find sums of:

- **Identifier(s)**; this is the name given earlier to the variational form(s) (type `varf`) for possible reuse.

  Remark, that the name in the `varf` of the unknown test function is forgotten, we use the order in the argument list to recall names as in a C++ function,

- **The terms of the bilinear form itself**: if $K$ is a given function,

  - Bilinear part for 3D meshes $\text{Th}$
    
    $$\int_3 \text{Th}(K*v*w) = \sum_{T \in \text{Th}} \int_T K v w$$

    $$\int_3 \text{Th}, 1(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega} \int_T K v w$$

    $$\int_3 \text{Th}, \text{levelset} = \phi (K*v*w) = \sum_{T \in \text{Th}} \int_{T, \phi < 0} K v w$$
- \( \text{int3d}(\text{Th}, 1, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_l} \int_{T, \phi < 0} K v w \)

- \( \text{int2d}(\text{Th}, 2, 5)(K*v*w) = \sum_{T \in \text{Th}} \int_{(\partial T \cup \Gamma) \cap (\Gamma_2 \cup \Gamma_5)} K v w \)

- \( \text{int2d}(\text{Th}, 1)(K*v*w) = \sum_{T \in \text{th}, T \subset \Omega_1} \int_{T} K v w \)

- \( \text{int2d}(\text{Th}, 2, 5)(K*v*w) = \sum_{T \in \text{Th}} \int_{(\partial T \cup \Gamma) \cap (\Gamma_2 \cup \Gamma_5)} K v w \)

- \( \text{int2d}(\text{Th}, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}} \int_{T, \phi = 0} K v w \)

- \( \text{int2d}(\text{Th}, 1, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{T, \phi = 0} K v w \)

- \( \text{intallfaces}(\text{Th})(K*v*w) = \sum_{T \in \text{Th}} \int_{\partial T} K v w \)

- \( \text{intallfaces}(\text{Th}, 1)(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{\partial T} K v w \)

- They contribute to the sparse matrix of type \text{matrix} which, whether declared explicitly or not, is constructed by \text{FreeFEM}.

- Bilinear part for 2D meshes \( \text{Th} \)

  - \( \text{int2d}(\text{Th})(K*v*w) = \sum_{T \in \text{th}} \int_{T} K v w \)

  - \( \text{int2d}(\text{Th}, 1)(K*v*w) = \sum_{T \in \text{th}, T \subset \Omega_1} \int_{T} K v w \)

  - \( \text{int2d}(\text{Th}, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}} \int_{T, \phi < 0} K v w \)

  - \( \text{int2d}(\text{Th}, 1, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{T, \phi < 0} K v w \)

  - \( \text{int1d}(\text{Th}, 2, 5)(K*v*w) = \sum_{T \in \text{Th}} \int_{(\partial T \cup \Gamma) \cap (\Gamma_2 \cup \Gamma_5)} K v w \)

  - \( \text{int1d}(\text{Th}, 1)(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{T} K v w \)

  - \( \text{int1d}(\text{Th}, 2, 5)(K*v*w) = \sum_{T \in \text{Th}} \int_{(\partial T \cup \Gamma) \cap (\Gamma_2 \cup \Gamma_5)} K v w \)

  - \( \text{int1d}(\text{Th}, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}} \int_{T, \phi = 0} K v w \)

  - \( \text{int1d}(\text{Th}, 1, \text{levelset}=\text{phi})(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{T, \phi = 0} K v w \)
- \( \text{intalledges}(\text{Th})(K*v*w) = \sum_{T \in \text{Th}} \int_{\partial T} K \, v \, w \)

- \( \text{intalledges}(\text{Th}, 1)(K*v*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{\partial T} K \, v \, w \)

- They contribute to the sparse matrix of type matrix which, whether declared explicitly or not, is constructed by FreeFEM.

- The right hand-side of the Partial Differential Equation in 3D, the terms of the linear form: for given functions \( K, f \):

- \( \text{int3d}(\text{Th})(K*w) = \sum_{T \in \text{Th}} \int_{T} K \, w \)

- \( \text{int3d}(\text{Th}, 1)(K*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{T} K \, w \)

- \( \text{int3d}(\text{Th}, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi<0} K \, w \)

- \( \text{int3d}(\text{Th}, 1, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi<0} K \, w \)

- \( \text{int2d}(\text{Th}, 2, 5)(K*w) = \sum_{T \in \text{Th}} \int_{(\partial T \cup \Gamma) \cap (\Gamma_2 \cup \Gamma_5)} K \, w \)

- \( \text{int2d}(\text{Th}, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi=0} K \, w \)

- \( \text{int2d}(\text{Th}, 1, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi=0} K \, w \)

- \( \text{int1faces}(\text{Th})(f*w) = \sum_{T \in \text{Th}} \int_{\partial T} f \, w \)

- A vector of type real[int]

- The right hand-side of the Partial Differential Equation in 2D, the terms of the linear form: for given functions \( K, f \):

- \( \text{int2d}(\text{Th})(K*w) = \sum_{T \in \text{Th}} \int_{T} K \, w \)

- \( \text{int2d}(\text{Th}, 1)(K*w) = \sum_{T \in \text{Th}, T \subset \Omega_1} \int_{T} K \, w \)

- \( \text{int2d}(\text{Th}, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi<0} K \, w \)

- \( \text{int2d}(\text{Th}, 1, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi<0} K \, w \)

- \( \text{int1d}(\text{Th}, 2, 5)(K*w) = \sum_{T \in \text{Th}} \int_{(\partial T \cup \Gamma) \cap (\Gamma_2 \cup \Gamma_5)} K \, w \)

- \( \text{int1d}(\text{Th}, \text{levelset}=\phi)(K*w) = \sum_{T \in \text{Th}} \int_{T, \phi=0} K \, w \)
\[ \text{int1d}(\text{Th}, 1, \text{levelset} = \phi)(K \cdot w) = \sum_{T \in \text{Th}, T \subset \Omega} \int_{T, \phi = 0} K \cdot w \]

\[ \text{intalledges}(\text{Th})(f \cdot w) = \sum_{T \in \text{Th}} \int_{\partial T} f \cdot w \]

- a vector of type \text{real}[\text{int}]

- The boundary condition terms:
  - An “on” scalar form (for Dirichlet): \text{on}(1, u = g)
    Used for all degrees of freedom \( i \) of the boundary referred by “1”, the diagonal term of the matrix \( a_{ii} = tgv \) with the terrible giant value \( tgv = 10^{30} \) by default, and the right hand side \( b[i] = "(\Pi_h g)[i]" \times tgv \), where the “\( (\Pi_h g)[i] \)” is the boundary node value given by the interpolation of \( g \).
  - A linear form on \( \Gamma \) (for Neumann in 2d) \(-\text{int1d}(\text{Th})(f \cdot w)\) or \(-\text{int1d}(\text{Th}, 3)(f \cdot w)\)
  - A bilinear form on \( \Gamma \) or \( \Gamma_2 \) (for Robin in 2d) \text{int1d}(\text{Th})(K \cdot v \cdot w)\) or \text{int1d}(\text{Th}, 2)(K \cdot v \cdot w)\)
  - A linear form on \( \Gamma \) (for Neumann in 3d) \(-\text{int2d}(\text{Th})(f \cdot w)\) or \(-\text{int2d}(\text{Th}, 3)(f \cdot w)\)
  - A bilinear form on \( \Gamma \) or \( \Gamma_2 \) (for Robin in 3d) \text{int2d}(\text{Th})(K \cdot v \cdot w)\) or \text{int2d}(\text{Th}, 2)(K \cdot v \cdot w)\)

Note:
- An “on” vectorial form (for Dirichlet): \text{on}(1, u_1 = g_1, u_2 = g_2)
  If you have vectorial finite element like RT0, the 2 components are coupled, and so you have: \( b[i] = "(\Pi_h (g_1, g_2))[i]" \times tgv \), where \( \Pi_h \) is the vectorial finite element interpolant.
- An “on” vectorial form (for Dirichlet): \text{on}(u = g, tgv = \text{none positive value })
  if the value is equal to -2 (i.e \( tgv = \text{-2} \)) then we put to :math:`0` all term of the line and column \( i \) in the matrix, except diagonal term \( a_{ii} = 1 \), and \( b[i] = "(\Pi_h g)[i]" \) else if the value is equal to -20 (i.e \( tgv = \text{-20} \)) then we put to :math:`0` all term of the line and column \( i \) in the matrix, and \( b[i] = "(\Pi_h g)[i]" \) else if the value is equal to -10 (i.e \( tgv = \text{-10} \)) then we put to :math:`0` all term of the line the matrix, and \( b[i] = "(\Pi_h g)[i]" \) else (i.e \( tgv = \text{-1} \)) we put to :math:`0` all term of the line \( i \) in the matrix, except diagonal term \( a_{ii} = 1 \), and \( b[i] = "(\Pi_h g)[i]" \).
- If needed, the different kind of terms in the sum can appear more than once.
- The integral mesh and the mesh associated to test functions or unknown functions can be different in the case of \text{varf} form.
- \( N.x, N.y \) and \( N.z \) are the normal’s components.
- \( Ns.x, Ns.y \) and \( Ns.z \) are the normal’s components of the surface in case of \text{meshS} integral
- \( Tl.x, Tl.y \) and \( Tl.z \) are the tangent’s components of the line in case of \text{meshL} integral

\textbf{Warning:} It is not possible to write in the same integral the linear part and the bilinear part such as in \text{int1d}(\text{Th})(K \cdot v \cdot w \cdot f \cdot w)\).
3.3.15 Numerical Integration

Let $D$ be a $N$-dimensional bounded domain.

For an arbitrary polynomial $f$ of degree $r$, if we can find particular (quadrature) points $\xi_j$, $j = 1, \cdots, J$ in $D$ and (quadrature) constants $\omega_j$ such that

$$
\int_D f(x) = \sum_{\ell=1}^L c_{\ell} f(\xi_{\ell})
$$

then we have an error estimate (see [CROUSEIX1984]), and then there exists a constant $C > 0$ such that

$$
\left| \int_D f(x) - \sum_{\ell=1}^L \omega_{\ell} f(\xi_{\ell}) \right| \leq C |D|h^{r+1}
$$

for any function $r + 1$ times continuously differentiable $f$ in $D$, where $h$ is the diameter of $D$ and $|D|$ its measure (a point in the segment $[q'q'']$) is given as

$$
\{(x,y) | x = (1-t)q'_x + tq''_x, y = (1-t)q'_y + tq''_y, 0 \leq t \leq 1\}
$$

For a domain $\Omega_h = \bigcup_{k=1}^{n_1} T_k$, $\partial_h = \{ T_k \}$, we can calculate the integral over $\Gamma_h = \partial\Omega_h$ by:

$$
\int_{\Gamma_h} f(x)ds = \int(\text{Th}) (f) = \int(\text{Th, qfe}=\star) (f) = \int(\text{Th, qorder} = \star) (f)
$$

where * stands for the name of the quadrature formula or the precision (order) of the Gauss formula.

### Quadrature formula on an edge

<table>
<thead>
<tr>
<th>$L$</th>
<th>qfe</th>
<th>qforder</th>
<th>Point in $[q', q'']$</th>
<th>$\omega_\ell$</th>
<th>Exact on $P_k, k =$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>qf1pE</td>
<td>2</td>
<td>$1/2$</td>
<td>$|q'q''|$</td>
<td>1</td>
</tr>
<tr>
<td>2</td>
<td>qf2pE</td>
<td>3</td>
<td>$(1 \pm \sqrt{1/3})/2$</td>
<td>$|q'q''|/2$</td>
<td>3</td>
</tr>
<tr>
<td>3</td>
<td>qf3pE</td>
<td>6</td>
<td>$(1 \pm \sqrt{3/5})/2$</td>
<td>$|q'q''|/2$</td>
<td>5</td>
</tr>
<tr>
<td>4</td>
<td>qf4pE</td>
<td>8</td>
<td>$(1 \pm \sqrt{525+70\sqrt{30}}/45)/2$</td>
<td>$|q'q''|/2$</td>
<td>7</td>
</tr>
<tr>
<td>5</td>
<td>qf5pE</td>
<td>10</td>
<td>$(1 \pm \sqrt{245+14\sqrt{70}})/2$</td>
<td>$|q'q''|/2$</td>
<td>9</td>
</tr>
<tr>
<td></td>
<td>qf1pElump</td>
<td>0</td>
<td>$1$</td>
<td>$|q'q''|/2$</td>
<td>1</td>
</tr>
</tbody>
</table>

where $|q'q''|$ is the length of segment $q'q''$.

For a part $\Gamma_1$ of $\Gamma_h$ with the label “1”, we can calculate the integral over $\Gamma_1$ by:

$$
\int_{\Gamma_1} f(x, y)ds = \int(\text{Th, 1}) (f) = \int(\text{Th, 1, qfe}=\text{qf}2\text{pE}) (f)
$$

The integrals over $\Gamma_1, \Gamma_3$ are given by:

$$
\int_{\Gamma_1 \cup \Gamma_3} f(x, y)ds
$$

For each triangle $T_k = [q^k_1, q^k_2, q^k_3]$, the point $P(x, y)$ in $T_k$ is expressed by the area coordinate as $P(\xi, \eta)$:

$$
|T_k| = \frac{1}{2} \begin{vmatrix} 1 & q^k_1 & q^k_2 & q^k_3 \\ 1 & q^k_2 & q^k_3 & q^k_1 \\ 1 & q^k_3 & q^k_1 & q^k_2 \end{vmatrix}
\begin{array}{l}
\text{ } \\
\text{ } \\
\text{ } \\
\text{ } \\
\text{ } \\
\end{array}

D_1 = \begin{vmatrix} x & y \\
1 & q^k_1 \\
1 & q^k_2 \\
1 & q^k_3 \end{vmatrix}
\begin{array}{l} \\
\text{ } \\
\text{ } \\
\text{ } \\
\text{ } \\
\end{array}

D_2 = \begin{vmatrix} x & y \\
1 & q^k_1 \\
1 & q^k_2 \\
1 & q^k_3 \end{vmatrix}
\begin{array}{l} \\
\text{ } \\
\text{ } \\
\text{ } \\
\text{ } \\
\end{array}

D_3 = \begin{vmatrix} x & y \\
1 & q^k_1 \\
1 & q^k_2 \\
1 & q^k_3 \end{vmatrix}
\begin{array}{l} \\
\text{ } \\
\text{ } \\
\text{ } \\
\text{ } \\
\end{array}

\xi = \frac{1}{2}D_1/|T_k| 
\eta = \frac{1}{2}D_2/|T_k|
\text{ then } 1 - \xi - \eta = \frac{1}{2}D_3/|T_k|

3.3. Finite element
For a two dimensional domain or a border of three dimensional domain $\Omega_h = \sum_{k=1}^{n^i} T_k$, $\mathcal{T}_h = \{T_k\}$, we can calculate the integral over $\Omega_h$ by:

$$\int_{\Omega_h} f(x, y) \, \text{d}x \text{d}y = \text{int2d}(\mathcal{T}_h)(f) = \text{int2d}(\mathcal{T}_h, \text{qft} = \ast)(f) = \text{int2d}(\mathcal{T}_h, \text{qforder} = \ast)(f)$$

where * stands for the name of quadrature formula or the order of the Gauss formula.

### Quadrature formula on a triangle

<table>
<thead>
<tr>
<th>$L$</th>
<th>qft</th>
<th>qforder</th>
<th>$\omega_f$</th>
<th>Exact on $P_k$, $k =$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>qf1p1</td>
<td>2</td>
<td>$\left(\frac{1}{3}, \frac{1}{3}\right)$</td>
<td>$T_k$</td>
</tr>
<tr>
<td>3</td>
<td>qf2p1</td>
<td>3</td>
<td>$\left(\frac{1}{2}, \frac{1}{2}\right)$</td>
<td>$\left(\frac{1}{2}, 0\right)$</td>
</tr>
<tr>
<td>5</td>
<td>qf5p1</td>
<td>6</td>
<td>$\left(\frac{6 - \sqrt{15}}{21}, \frac{6 - \sqrt{15}}{21}\right)$</td>
<td>$\left(\frac{6 - \sqrt{15}}{21}, \frac{9 + 2\sqrt{15}}{21}\right)$</td>
</tr>
<tr>
<td>7</td>
<td>qf1pTlump</td>
<td>3</td>
<td>$\left(0, 0\right)$</td>
<td>$\left(1, 0\right)$</td>
</tr>
<tr>
<td>9</td>
<td>qf2P1</td>
<td>4</td>
<td>$\left(\frac{3}{4}, \frac{3}{4}\right)$</td>
<td>$\left(0, \frac{1}{4}\right)$</td>
</tr>
<tr>
<td>15</td>
<td>qf7p1</td>
<td>8</td>
<td>See [TAYLOR2005] for detail</td>
<td></td>
</tr>
<tr>
<td>21</td>
<td>qf9p1</td>
<td>10</td>
<td>See [TAYLOR2005] for detail</td>
<td></td>
</tr>
</tbody>
</table>

For a three dimensional domain $\Omega_h = \sum_{k=1}^{n^i} T_k$, $\mathcal{T}_h = \{T_k\}$, we can calculate the integral over $\Omega_h$ by:

$$\int_{\Omega_h} f(x, y) \, \text{d}x \text{d}y = \text{int3d}(\mathcal{T}_h)(f) = \text{int3d}(\mathcal{T}_h, \text{qf} = \ast)(f) = \text{int3d}(\mathcal{T}_h, \text{qforder} = \ast)(f)$$

where * stands for the name of quadrature formula or the order of the Gauss formula.
Quadrature formula on a tetrahedron

<table>
<thead>
<tr>
<th>L</th>
<th>qfV</th>
<th>qforder</th>
<th>Point in ( T_k \in \mathbb{R}^3 )</th>
<th>( \omega_\ell )</th>
<th>Exact on ( P_k ), ( k = )</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>qfV1</td>
<td>2</td>
<td>( (\frac{1}{3}, \frac{1}{3}, \frac{1}{3}) )</td>
<td>(</td>
<td>T_k</td>
</tr>
<tr>
<td>4</td>
<td>qfV2</td>
<td>3</td>
<td>( G^4(0.58\ldots, 0.13\ldots, 0.13\ldots) )</td>
<td>(</td>
<td>T_k</td>
</tr>
<tr>
<td>14</td>
<td>qfV5</td>
<td>6</td>
<td>( G^4(0.72\ldots, 0.092\ldots, 0.092\ldots) ) ( G^4(0.067\ldots, 0.31\ldots, 0.31\ldots) ) ( G^6(0.45\ldots, 0.045\ldots, 0.45\ldots) )</td>
<td>0.073... (</td>
<td>T_k</td>
</tr>
<tr>
<td>4</td>
<td>qfV1lump</td>
<td></td>
<td>( G^4(1, 0, 0) )</td>
<td>(</td>
<td>T_k</td>
</tr>
</tbody>
</table>

Where \( G^4(a, b, b) \) such that \( a + 3b = 1 \) is the set of the four point in barycentric coordinate:

\[
\{ (a, b, b), (b, a, b), (b, b, a), (b, b, b) \}
\]

and where \( G^6(a, b, b) \) such that \( 2a + 2b = 1 \) is the set of the six points in barycentric coordinate:

\[
\{ (a, a, b), (a, b, a), (b, b, a), (b, a, a), (b, a, b), (b, a, b) \}
\]

**Note:** These tetrahedral quadrature formulae come from [http://nines.cs.kuleuven.be/research/ecf/mtables.html](http://nines.cs.kuleuven.be/research/ecf/mtables.html)

**Note:** By default, we use the formula which is exact for polynomials of degree 5 on triangles or edges (in bold in three tables).

It is possible to add an own quadrature formulae with using plugin qf11to25 on segment, triangle or Tetrahedron.

The quadrature formulae in \( D \) dimension is a bidimensional array of size \( N_q \times (D + 1) \) such that the \( D + 1 \) value of on row \( i = 0, ..., N_p - 1 \) are \( w^i, \hat{x}_1^i, ..., \hat{x}_D^i \) where \( w^i \) is the weight of the quadrature point, and \( 1 - \sum_{k=1}^{D} \hat{x}_k^i, \hat{x}_1^i, ..., \hat{x}_D^i \) is the barycentric coordinate the quadrature point.

```plaintext
load "qf11to25"

// Quadrature on segment
real [int, int] qq1 = [
    [0.5, 0],
    [0.5, 1]
];
QF1 qf1(1, qq1); //def of quadrature formulae qf1 on segment
//remark:
//1 is the order of the quadrature exact for polynome of degree < 1

//Quadrature on triangle
real [int, int] qq2 = [
    [1./3., 0, 0],
    [1./3., 1, 0],
    [1./3., 0, 1]
];
QF2 qf2(1, qq2); //def of quadrature formulae qf2 on triangle
//remark:
//1 is the order of the quadrature exact for polynome of degree < 1
```

(continues on next page)
// so must have sum w^i = 1

// Quadrature on tetrahedron
real[int, int] qq3 = [1./4., 0, 0, 0], [1./4., 1, 0, 0], [1./4., 0, 1, 0], [1./4., 0, 0, 1];

QF3 qf3(1, qq3); // def of quadrature formulae qf3 on get
// remark:
// 1 is the order of the quadrature exact for polynomials of degree < 1)
// Verification in 1d and 2d
mesh Th = square(10, 10);

real I1 = int1d(Th, qfe=qf1)(x^2);
real I1l = int1d(Th, qfe=qf1pElump)(x^2);
real I2 = int2d(Th, qft=qf2)(x^2);
real I2l = int2d(Th, qft=qf1pTlump)(x^2);

cout << I1 << " == " << I1l << endl;
cout << I2 << " == " << I2l << endl;
assert(abs(I1-I1l) < 1e-10);
assert(abs(I2-I2l) < 1e-10);

The output is
1.67 == 1.67
0.335 == 0.335

3.3.16 Variational Form, Sparse Matrix, PDE Data Vector

In FreeFEM it is possible to define variational forms, and use them to build matrices and vectors, and store them to speed-up the script (4 times faster here).

For example let us solve the Thermal Conduction problem.

The variational formulation is in $L^2(0, T; H^1(\Omega))$; we shall seek $u^n$ satisfying:

$$
\forall w \in V_0; \quad \int_{\Omega} \frac{u^n - u^{n-1}}{\delta t} w + \kappa \nabla u^n \nabla w + \int_{\Gamma} \alpha(u^n - u_{\text{inc}}) w = 0
$$

where $V_0 = \{w \in H^1(\Omega)/w|_{\Gamma_{24}} = 0\}$.

So to code the method with the matrices $A = (A_{ij}), M = (M_{ij})$, and the vectors $u^n, b^n, b', b'', b_{i}$ (notation if $w$ is a vector then $w_i$ is a component of the vector).

$$
u^n = A^{-1}b^n, \quad b' = b_0 + Mu^{n-1}, \quad b'' = \frac{1}{\varepsilon} b_{\text{inc}}, \quad b^n_i = \begin{cases} b''_i & \text{if } i \in \Gamma_{24} \\ b'_i & \text{else if } i \notin \Gamma_{24} \end{cases}
$$
Where with $\frac{1}{\varepsilon} = tgv = 10^{30}$:

$$
A_{ij} = \begin{cases} 
\int_\Omega w_j w_i / dt + k(\nabla w_j, \nabla w_i) + \int_{\Gamma_{13}} \alpha w_j w_i & \text{if } i \in \Gamma_{24}, \text{and } j = i \\
\frac{1}{\varepsilon} & \text{if } i \in \Gamma_{24}, \text{and } j = i \\
\int_\Omega w_j w_i / dt & \text{else if } i \notin \Gamma_{24}, \text{or } j \neq i
\end{cases}
$$

$$
M_{ij} = \begin{cases} 
\int_\Omega w_j w_i / dt & \text{else if } i \notin \Gamma_{24}, \text{or } j \neq i
\end{cases}
$$

$$
b_{0,i} = \int_{\Gamma_{13}} \alpha w_i u_e w_i
$$

$$
b_{cl} = \int_\Omega w_i / \Omega^{30} \text{ the initial data}
$$

// Parameters
func fu0 = 10 + 90*x/6;
func k = 1.8*(y<0.5) + 0.2;
real ue = 25.;
real alpha = 0.25;
real T = 5;
real dt = 0.1;

// Mesh
mesh Th = square(30, 5, [6*x, y]);

// Fespace
fespace Vh(Th, P1);
Vh u0 = fu0, u = u0;

Create three variational formulation, and build the matrices $A,M$.

// Problem
varf vthermic (u, v) = int2d(Th)(
  u*v/dt
  + k*(dx(u)*dx(v) + dy(u)*dy(v))
  + int1d(Th, 1, 3)\{ 
    alpha*u*v \\
  \}
  + on(2, 4, u=1)
);

varf vthermic0 (u, v) = int1d(Th, 1, 3)\{ 
  alpha*ue*v
\}

varf vMass (u, v) = int2d(Th)(
  u*v/dt
  + on(2, 4, u=1)
);
real tgv = 1e30;
matrix A = vthermic(Vh, Vh, tgv=tgv, solver=CG);
matrix M = vMass(Vh, Vh);

Now, to build the right hand size we need 4 vectors.

3.3. Finite element
```
real[int] b0 = vthermic0(0, Vh); //constant part of the RHS
real[int] bcn = vthermic(0, Vh); //tgv on Dirichlet boundary node ( !=0 )
//we have for the node i : i in Gamma_24 -> bcn[i] != 0
real[int] bcl = tgv*u0[]; //the Dirichlet boundary condition part
```

**Note:** The boundary condition is implemented by penalization and vector `bcn` contains the contribution of the boundary condition \( u = 1 \), so to change the boundary condition, we have just to multiply the vector `bcn[]` by the current value \( f \) of the new boundary condition term by term with the operator `.*`.

*Uzawa model* gives a real example of using all this features.

And the new version of the algorithm is now:

```
// Time loop
ofstream ff("thermic.dat");
for(real t = 0; t < T; t += dt){
    // Update
    real[int] b = b0; //for the RHS
    b += M*u[]; //add the the time dependent part
    //lock boundary part:
    // Solve
    u[] = A^-1*b;
    // Save
    ff << t << " " << u(3, 0.5) << endl;
    // Plot
    plot(u);
}
// Display
for(int i = 0; i < 20; i++)
    cout << dy(u)(6.0*i/20.0, 0.9) << endl;
// Plot
plot(u, fill=true, wait=true);
```

**Note:** The functions appearing in the variational form are formal and local to the `varf` definition, the only important thing is the order in the parameter list, like in:

```
varf vb1([u1, u2], q) = int2d(Th)((dy(u1) + dy(u2))*q) + int2d(Th)(1*q);
varf vb2([v1, v2], p) = int2d(Th)((dy(v1) + dy(v2))*p) + int2d(Th)(1*p);
```

To build matrix \( A \) from the bilinear part the variational form \( a \) of type `varf` simply write:

```
A = a(Vh, Wh , [...]);
```

**Note:** \( \text{Vh} \) is "fespace" for the unknown fields with a correct number of component

**Note:** \( \text{Wh} \) is "fespace" for the test fields with a correct number of component

Possible named parameters in , , [...] are

- `solver= LU, CG, Crout, Cholesky, GMRES, sparsesolver, UMFPACK ...`
The default solver is GMRES.

The storage mode of the matrix of the underlying linear system depends on the type of solver chosen; for LU the matrix is sky-line non symmetric, for Crout the matrix is sky-line symmetric, for Cholesky the matrix is sky-line symmetric positive definite, for CG the matrix is sparse symmetric positive, and for GMRES, sparsesolver or UMFPACK the matrix is just sparse.

- **factorize** = If true then do the matrix factorization for LU, Cholesky or Crout, the default value is false.
- **eps**= A real expression. 
  
  \( \varepsilon \) sets the stopping test for the iterative methods like CG.

  Note that if \( \varepsilon \) is negative then the stopping test is:

  \[
  ||Ax - b|| < |\varepsilon|
  \]

  if it is positive then the stopping test is

  \[
  ||Ax - b|| < \frac{|\varepsilon|}{||Ax_0 - b||}
  \]

- **precon**= Name of a function (for example \( P \)) to set the preconditioner.

  The prototype for the function \( P \) must be:

  ```
  func real[int] P(real[int] & xx) ;
  ```

- **tgv**= Huge value (10^{30}) used to implement Dirichlet boundary conditions.
- **tolpivot**= Set the tolerance of the pivot in UMFPACK (10^{-1}) and, LU, Crout, Cholesky factorization (10^{-20}).
- **tolpivotsym**= Set the tolerance of the pivot sym in UMFPACK
- **strategy**= Set the integer UMFPACK strategy (0 by default).

**Note:** The line of the matrix corresponding to the space \( \mathcal{W}_h \) and the column of the matrix corresponding to the space \( \mathcal{V}_h \).

To build the dual vector \( b \) (of type real[int]) from the linear part of the variational form \( a \) do simply:

```
real b(Vh.ndof);
b = a(0, Vh);
```

A first example to compute the area of each triangle \( K \) of mesh \( T_h \), just do:

```
fespace Nh(Th, P0); //the space function constant / triangle
Nh areaK;
varf varea (unused, chiK) = int2d(Th)(chiK);
etaK[] = varea(0, Ph);
```

Effectively, the basic functions of space \( N_h \), are the characteristic function of the element of \( T_h \), and the numbering is the numeration of the element, so by construction:

\[
\eta_K[i] = \int_{K_i} 1 = \int_{K_i} 1;
\]

Now, we can use this to compute error indicators like in example *Adaptation using residual error indicator*.

First to compute a continuous approximation to the function \( h \) “density mesh size” of the mesh \( T_h \).
fespace Vh(Th, P1);
Vh h;
real[int] count(Th.nv);
varf vmeshsizen(u, v) = intalledges(Th, qfnbpE=1)(v);
varf vedgecount(u, v) = intalledges(Th, qfnbpE=1)(v/lenEdge);

// Computation of the mesh size
count = vedgecount(0, Vh); //number of edge / vertex
h[] = vmeshsizen(0, Vh); //sum length edge / vertex
h[] = h[]./count; //mean length edge / vertex

To compute error indicator for Poisson equation:

\[ \eta_K = \int_K h_K^2 |(f + \Delta u_h)|^2 + \int_{\partial K} h_e [\frac{\partial u_h}{\partial n}]^2 \]

where \( h_K \) is size of the longest edge (hTriangle), \( h_e \) is the size of the current edge (lenEdge), \( n \) the normal.

fespace Nh(Th, P0); // the space function constant / triangle
Nh etak;
varf vetaK(unused, chiK) = intalledges(Th) {
  chiK*lenEdge*square(jump(N.x*dx(u) + N.y*dy(u)));
} + int2d(Th) {
  chiK*square(hTriangle*(f + dxx(u) + dyy(u)));
};
etak[] = vetaK(0, Ph);

We add automatic expression optimization by default, if this optimization creates problems, it can be removed with the keyword optimize as in the following example:

varf a(u1, u2) = int2d(Th, optimize=0) {
  dx(u1)*dx(u2) + dy(u1)*dy(u2);
} + on(1, 2, 4, u1=0) + on(3, u1=1);

or you can also do optimization and remove the check by setting optimize=2.

Remark, it is all possible to build interpolation matrix, like in the following example:

mesh TH = square(3, 4);
mesh th = square(2, 3);
mesh Th = square(4, 4);
fespace VH(TH, P1);
fespace Vh(th, P1);
fespace Wh(Th, P1);
matrix B = interpolate(VH, Vh); //build interpolation matrix Vh->VH
matrix BB = interpolate(Wh, Vh); //build interpolation matrix Vh->Wh

and after some operations on sparse matrices are available for example:
3.3.17 Interpolation matrix

It is also possible to store the matrix of a linear interpolation operator from a finite element space $V_h$ to another $W_h$ to interpolate($W_h,V_h,...$) a function.

Note that the continuous finite functions are extended by continuity outside of the domain.

The named parameters of function interpolate are:

- **inside**= set true to create zero-extension.
- **t**= set true to get the transposed matrix
- **op**= set an integer written below
  - 0 the default value and interpolate of the function
  - 1 interpolate the $\partial_x$
  - 2 interpolate the $\partial_y$
  - 3 interpolate the $\partial_z$
- **U2Vc**= set the which is the component of $W_h$ come in $V_h$ in interpolate process in a int array so the size of the array is number of component of $W_h$, if the put $-1$ then component is set to 0, like in the following example:

  (by default the component number is unchanged).

```plaintext
fespace V4h(Th4, [P1, P1, P1, P1]);
fespace V3h(Th, [P1, P1, P1]);
int[int] u2vc = [1, 3, -1]; // -1 -> put zero on the component
matrix IV34 = interpolate(V3h, V4h, inside=0, U2Vc=u2vc); // V3h <- V4h
V4h [a1, a2, a3, a4] = [1, 2, 3, 4];
V3h [b1, b2, b3] = [10, 20, 30];
b1[] = IV34*a1[];
```

So here we have: freefem b1 == 2, b2 == 4, b3 == 0 ...
mesh Th4 = square(2, 2, [x*0.5, y*0.5]);
plot(Th, Th4, wait=true);

// Fespace
fespace Vh(Th, P1);
Vh v, vv;
fespace Vh4(Th4, P1);
Vh4 v4=x*y;

fespace Wh(Th, P0);
fespace Wh4(Th4, P0);

// Interpolation
matrix IV = interpolate(Vh, Vh4);  //here the function is extended by continuity
cout << "IV Vh<-Vh4 " << IV << endl;
v=v4;
vv[] = IV*v4[];  //here v == vv
real[int] diff= vv[] - v[];
cout << "|| v - vv || = " << diff.linfty << endl;
assert(diff.linfty <= 1e-6);

matrix IV0 = interpolate(Vh, Vh4, inside=1);  //here the function is extended by zero
cout << "IV Vh<-Vh4 (inside=1) " << IV0 << endl;

matrix IVt0 = interpolate(Vh, Vh4, inside=1, t=1);
cout << "IV Vh<-Vh4^t (inside=1) " << IVt0 << endl;

matrix IV4t0 = interpolate(Vh4, Vh);
cout << "IV Vh4<-Vh^t " << IV4t0 << endl;

matrix IW4 = interpolate(Wh4, Wh);
cout << "IV Wh4<-Wh " << IW4 << endl;

matrix IW4V = interpolate(Wh4, Vh);
cout << "IV Wh4<-Vh " << IW4V << endl;

Build interpolation matrix $A$ at a array of points $(xx[j], yy[j]), i = 0, 2$ here:

$$a_{i,j} = dop(w^c_i(xx[j], yy[j]))$$

where $w_i$ is the basic finite element function, $c$ the component number, $dop$ the type of diff operator like in op def.

real[int] xx = [.3, .4], yy = [.1, .4];
int c = 0, dop = 0;
matrix Ixx = interpolate(Vh, xx, yy, op=dop, composante=c);
cout << Ixx << endl;
Vh ww;
real[int] dd = [1, 2];
ww[] = Ixx*dd;

Tip: Schwarz

The following shows how to implement with an interpolation matrix a domain decomposition algorithm based on Schwarz method with Robin conditions.
Given a non-overlapping partition \( \bar{\Omega} = \bar{\Omega}_0 \cup \bar{\Omega}_1 \) with \( \Omega_0 \cap \Omega_1 = \emptyset, \Sigma := \bar{\Omega}_0 \cap \bar{\Omega}_1 \) the algorithm is:

\[
- \Delta u_i = f \text{ in } \Omega_i, \ i = 0, 1,
\]

\[
\frac{\partial (u_1 - u_0)}{\partial n} + \alpha (u_1 - u_0) = 0 \text{ on } \Sigma.
\]

The same in variational form is:

\[
- \int_{\Omega_i} \nabla u_i \cdot \nabla v + \int_{\Omega_i} \alpha u_i v = \int_{\Omega_i} f v
\]

\[
- \int_{\Omega_j} \nabla u_j \cdot \nabla v + \int_{\Sigma} \alpha u_j v, \ \forall v \in H^1_0(\Omega), i, j = [0, 1] \cup [1, 0]
\]

To discretize with the \( P_1 \) triangular Lagrangian finite element space \( V_h \) simply replace \( H^1_0(\Omega) \) by \( V_h(\Omega_0) \cup V_h(\Omega_1) \).

Then difficulty is to compute \( \int_{\Omega_j} \nabla u_j \cdot \nabla v \) when \( v \) is a basis function of \( V_h(\Omega_i), i \neq j \).

It is done as follows (with \( \Gamma = \partial \Omega \)):

```cpp
// Parameters
int n = 30;
int Gamma = 1;
int Sigma = 2;

func f = 1.;
real alpha = 1.;
int Niter = 50;

// Mesh
mesh[int] Th(2);
int[int] reg(2);

border a0(t=0, 1){x=t; y=0; label=Gamma;}
border a1(t=1, 2){x=t; y=0; label=Gamma;}
border b1(t=0, 1){x=2; y=t; label=Gamma;}
border c1(t=2, 1){x=t; y=1; label=Gamma;}
border c0(t=1, 0){x=t; y=1; label=Gamma;}
border b0(t=1, 0){x=0; y=t; label=Gamma;}
border d(t=0, 1){x=1; y=t; label=Sigma;}
plot(a0(n) + a1(n) + b1(n) + c1(n) + c0(n) + b0(n) + d(n));
mesh TH = buildmesh(a0(n) + a1(n) + b1(n) + c1(n) + c0(n) + b0(n) + d(n));
reg(0) = TH(0.5, 0.5).region;
reg(1) = TH(1.5, 0.5).region;
for(int i = 0; i < 2; i++) Th[i] = trunc(TH, region==reg(i));

// Fespace
fespace Vh0(Th[0], P1);
Vh0 u0 = 0;

fespace Vh1(Th[1], P1);
Vh1 ul = 0;

// Macro
macro grad(u) [dx(u), dy(u)] //

// Problem
int i;
varf a (u, v)

(continues on next page)
= int2d(Th[i]) (
    grad(u)'*grad(v)
) + int1d(Th[i], Sigma)(
    alpha*u*v
) + on(Gamma, u=0) ;

varf b (u, v)
= int2d(Th[i]) (f*v)
+ on(Gamma, u=0) ;

varf duldn (u, v)
= -int2d(Th[1]) (
    grad(u1)'*grad(v) - f*v
) + int1d(Th[1], Sigma)(
    alpha*u1*v
) + on(Gamma, u=0) ;

varf du0dn (u, v)
= -int2d(Th[0]) (
    grad(u0)'*grad(v) - f*v
) + int1d(Th[0], Sigma)(
    alpha*u0*v
) + on(Gamma, u=0) ;

matrix I01 = interpolate(Vh1, Vh0);
matrix I10 = interpolate(Vh0, Vh1);

matrix[int] A(2);
i = 0; A[i] = a(Vh0, Vh0);
i = 1; A[i] = a(Vh1, Vh1);

// Solving loop
for(int iter = 0; iter < Niter; iter++){
    // Solve on Th[0]
    i = 0;
    real[int] b0 = b(0, Vh0);
    real[int] Duldn = duldn(0, Vh1);
    real[int] Tduldn(Vh0.n dof); Tduldn = I01'*Duldn;
    b0 += Tduldn;
    u0[] = A[0]'-1*b0;
    // Solve on Th[1]
}
100  {  
101      i = 1;  
102      real[int] b1 = b(0, Vh1);  
103      real[int] Du0dn = du0dn(0, Vh0);  
104      real[int] Tdu0dn(Vh1.ndof); Tdu0dn = I10'*Du0dn;  
105      b1 += Tdu0dn;  
106      u1[] = A[1]^-1*b1;  
107  }
108  plot(u0, u1, cmm="iter="+iter);
109 }

3.3.18 Finite elements connectivity

Here, we show how to get informations on a finite element space $W_h(T_n, *)$, where "*" may be P1, P2, P1nc, etc.

- $W_h$.nt gives the number of element of $W_h$
- $W_h.ndof$ gives the number of degrees of freedom or unknown
- $W_h.ndofK$ gives the number of degrees of freedom on one element
- $W_h(k,i)$ gives the number of $i$th degrees of freedom of element $k$.

See the following example:

**Tip:** Finite element connectivity

```plaintext
// Mesh
mesh Th = square(5, 5);

// Fespace
fespace Wh(Th, P2);

cout << "Number of degree of freedom = " << Wh.ndof << endl;
cout << "Number of degree of freedom / ELEMENT = " << Wh.ndofK << endl;

int k = 2, kdf = Wh.ndofK; //element 2
cout << "Degree of freedom of element " << k << ":" << endl;
for (int i = 0; i < kdf; i++)
    cout << Wh(k,i) << ";
cout << endl;
```

The output is:

```
Number of degree of freedom = 121
Number of degree of freedom / ELEMENT = 6
Degree of freedom of element 2:
78 95 83 87 79 92
```

3.3. Finite element 215
3.4 Visualization

Results created by the finite element method can be a huge set of data, so it is very important to render them easy to grasp.

There are two ways of visualization in FreeFEM:

• One, the default view, which supports the drawing of meshes, isovalues of real FE-functions, and of vector fields, all by the command `plot` (see Plot section below). For publishing purpose, FreeFEM can store these plots as postscript files.

• Another method is to use external tools, for example, gnuplot (see Gnuplot section, medit section, Paraview section, Matlab/Octave section) using the command `system` to launch them and/or to save the data in text files.

3.4.1 Plot

With the command `plot`, meshes, isovalues of scalar functions, and vector fields can be displayed.

The parameters of the plot command can be meshes, real FE functions, arrays of 2 real FE functions, arrays of two double arrays, to plot respectively a mesh, a function, a vector field, or a curve defined by the two double arrays.

**Note:** The length of an arrow is always bound to be in [5%, 5%] of the screen size in order to see something.

The `plot` command parameters are listed in the Reference part.

The keyboard shortcuts are:

• `enter` tries to show plot

• `p` previous plot (10 plots saved)

• `?` shows this help

• `+,-` zooms in/out around the cursor 3/2 times

• `=` resets the view

• `r` refreshes plot

• `up, down, left, right` special keys to translate

• `3` switches 3d/2d plot keys:
  - `z,Z` focal zoom and zoom out
  - `H,h` increases or decreases the Z scale of the plot

• `mouse motion`:
  - `left button` rotates
  - `right button` zooms (ctrl+button on mac)
  - `right button +alt` translates (alt+ctrl+button on mac)

• `a,A` increases or decreases the arrow size

• `B` switches between showing the border meshes or not

• `i,I` updates or not: the min/max bound of the functions to the window

• `n,N` decreases or increases the number of iso value arrays
• b switches between black and white or color plotting
• g switches between grey or color plotting
• f switches between filling iso or iso line
• l switches between lighting or not
• v switches between show or not showing the numerical value of colors
• m switches between show or not showing the meshes
• w window dump in file ffglutXXXX.ppm
• * keep/drop viewpoint for next plot
• k complex data / change view type
• ESC closes the graphics process before version 3.22, after no way to close
• otherwise does nothing

For example:

```plaintext
1 real[int] xx(10), yy(10);
2 mesh Th = square(5,5);
3 fespace Vh(Th, P1);
4 //plot scalar and vectorial FE function
5 Vh uh=x*x+y*y, vh=-y^2+x^2;
6 plot(Th, uh, [uh, vh], value=true, ps="three.eps", wait=true);
7 //zoom on box defined by the two corner points [0.1,0.2] and [0.5,0.6]
8 plot(uh, [uh, vh], bb=[[0.1, 0.2], [0.5, 0.6]],
9     wait=true, grey=true, fill=true, value=true, ps="threeg.eps");
10 //compute a cut
11 for (int i = 0; i < 10; i++){
12     x = i/10.;
13     y = i/10.;
14     xx[i] = i;
15     yy[i] = uh; //value of uh at point (i/10., i/10.)
16 }
17 plot([xx, yy], ps="likegnu.eps", wait=true);
```

To change the color table and to choose the value of iso line you can do:

```plaintext
// from: \url{http://en.wikipedia.org/wiki/HSV_color_space}
// The HSV (Hue, Saturation, Value) model defines a color space
// in terms of three constituent components:
// HSV color space as a color wheel
// Hue, the color type (such as red, blue, or yellow):
// Ranges from 0-360 (but normalized to 0-100% in some applications, like here)
// Saturation, the "vibrancy" of the color: Ranges from 0-100%
// The lower the saturation of a color, the more "grayness" is present
// and the more faded the color will appear.
// Value, the brightness of the color: Ranges from 0-100%
mesh Th = square(10, 10, [2*x-1, 2*y-1]);
```

(continues on next page)
(a) Mesh, isovalue and vector  
(b) Enlargement in grey of isovalue and vector  
(c) Plots a cut of $u_h$. Note that a refinement of the same can be obtained in combination with gnuplot

**Fig. 3.39:** Plot
3.4.2 Link with gnuplot

Example *Membrane* shows how to generate a gnuplot from a *FreeFEM* file. Here is another technique which has the advantage of being online, i.e. one doesn’t need to quit *FreeFEM* to generate a gnuplot.

However, this works only if gnuplot is installed, and only on an Unix-like computer.

Add to the previous example:

```plaintext
//[ // file for gnuplot
ofstream gnu("plot.gp");
for (int i = 0; i < n; i++)
    gnu << xx[i] << " " << yy[i] << endl;

(continues on next page)
```

(continued from previous page)
// to call gnuplot command and wait 5 second (due to the Unix command)
// and make postscript plot
exec("echo 'plot "plot.gp" w l \n pause 5 \n set term postscript \n set output \n "gnuplot.eps" \n replot \n quit' | gnuplot");

Fig. 3.41: Plots a cut of uh with gnuplot

Note: See Plot example for the complete script.

3.4.3 Link with medit

As said above, medit is a freeware display package by Pascal Frey using OpenGL. Then you may run the following example.

Now medit software is included in FreeFEM under ffsedit name.

The medit command parameters are listed in the Reference part.

With version 3.2 or later

load "medit"

(continues on next page)
mesh Th = square(10, 10, [2*x-1, 2*y-1]);

fespace Vh(Th, P1);
Vh u=2-x*x-y*y;

medit("u", Th, u);

Before:

mesh Th = square(10, 10, [2*x-1, 2*y-1]);
fespace Vh(Th, P1);
Vh u=2-x*x-y*y;
savemesh(Th, "u", [x, y, u*.5]); //save u.points and u.faces file
// build a u.bb file for medit
{
  ofstream file("u.bb");
  file << "2 1 1 " << u[].n << " 2 \n";
  for (int j = 0; j < u[].n; j++)
    file << u[](j) << endl;
}
//call medit command
exec("ffmedit u");
//clean files on unix-like OS
exec("rm u.bb u.faces u.points");

Note: See Medit example for the complete script.

### 3.4.4 Link with Paraview

One can also export mesh or results in the .vtk format in order to post-process data using Paraview.
### 3.4.5 Link with Matlab© and Octave

In order to create a plot from a FreeFEM simulation in Octave and Matlab the mesh, the finite element space connectivity and the simulation data must be written to files:

```plaintext
#include "ffmatlib.idp"

mesh Th = square(10, 10, [2*x-1, 2*y-1]);
fespace Vh(Th, P1);
Vh u=2-x*x-y*y;
savemesh(Th,"export_mesh.msh");
ffSaveVh(Th,Vh,"export_vh.txt");
ffSaveData(u,"export_data.txt");
```

Within Matlab or Octave the files can be plot with the ffmatlib library:
```matlab
addpath('path to ffmatlib');
[p,b,t]=ffreadmesh('export_mesh.msh');
vh=ffreaddata('export_vh.txt');
u=ffreaddata('export_data.txt');
ffpdeplot(p,b,t,'VhSeq',vh,'XYData',u,'ZStyle','continuous','Mesh','on');
grid;
```

**Fig. 3.44:** Matlab / Octave plot

---

**Note:** For more Matlab / Octave plot examples have a look at the tutorial section *Matlab / Octave Examples* or visit the [ffmatlib library on github](https://github.com/).

### 3.5 Algorithms & Optimization

#### 3.5.1 Conjugate Gradient/GMRES

Suppose we want to solve the Euler problem (here $x$ has nothing to do with the reserved variable for the first coordinate in FreeFEM):

find $x \in \mathbb{R}^n$ such that

$$
\nabla J(x) = \left( \frac{\partial J}{\partial x_i}(x) \right) = 0
$$

(3.17)

where $J$ is a function (to minimize for example) from $\mathbb{R}^n$ to $\mathbb{R}$.

If the function is convex we can use the conjugate gradient algorithm to solve the problem, and we just need the function (named $dJ$ for example) which computes $\nabla J$, so the parameters are the name of that function with prototype

```matlab
func real[int] dJ(real[int] &xx);
```

which computes $\nabla J$, and a vector $x$ of type (of course the number 20 can be changed)

```matlab
real[int] x(20);
```

to initialize the process and get the result.

Given an initial value $x^{(0)}$, a maximum number $i_{\text{max}}$ of iterations, and an error tolerance $0 < \epsilon < 1$:

Put $x = x^{(0)}$ and write
NLCG(dJ, \mathbf{x}, \text{precon} = M, \text{nbiter} = \text{imax}, \text{eps} = \epsilon, \text{stop} = \text{stopfunc});

will give the solution of \mathbf{x} of \nabla J(\mathbf{x}) = 0. We can omit parameters precon, nbiter, eps, stop. Here \( M \) is the preconditioner whose default is the identity matrix.

The stopping test is

\[ \|\nabla J(\mathbf{x})\|_P \leq \epsilon \|\nabla J(\mathbf{x}(0))\|_P \]

Writing the minus value in \text{eps} = \epsilon, i.e.,

\[
\text{NLCG}(dJ, \mathbf{x}, \text{precon} = M, \text{nbiter} = \text{imax}, \text{eps} = -\epsilon);
\]

We can use the stopping test:

\[ \|\nabla J(\mathbf{x})\|_P^2 \leq \epsilon \]

The parameters of these three functions are:

- \text{nbiter} = \text{set the number of iteration (by default 100)}
- \text{precon} = \text{set the preconditioner function (P for example) by default it is the identity, note the prototype is func real[int] P(real[int] \& \mathbf{x})}
- \text{eps} = \text{set the value of the stop test} \( \epsilon = 10^{-6} \) by default if positive then relative test \( \|\nabla J(x)\|_P \leq \epsilon \|\nabla J(x(0))\|_P \), otherwise the absolute test is \( \|\nabla J(x)\|_P^2 \leq \epsilon \).
- \text{veps} = \text{set and return the value of the stop test, if positive, then relative test is} \( \|\nabla J(x)\|_P \leq \epsilon \|\nabla J(x(0))\|_P \), otherwise the absolute test is \( \|\nabla J(x)\|_P^2 \leq \epsilon \). The return value is minus the real stop test (remark: it is useful in loop).
- \text{stop} = \text{stopfunc} \text{ add your test function to stop before the eps criterion. The prototype for the function stopfunc is}

\[
\text{func bool stopfunc(int iter, real[int] u, real[int] g)}
\]

where \( u \) is the current solution, and \( g \), the current gradient, is not preconditioned.

**Tip:** *Algorithms.edp*

For a given function \( b \), let us find the minimizer \( u \) of the function

\[
J(u) = \frac{1}{2} \int_{\Omega} f(|\nabla u|^2) - \int_{\Omega} u b
\]

\[
f(x) = ax + x - \ln(1 + x), \quad f'(x) = a + \frac{x}{1+x}, \quad f''(x) = \frac{1}{(1+x)^2}
\]

under the boundary condition \( u = 0 \) on \( \partial \Omega \).

\[
fespace \text{Ph}(\text{Th}, \text{P0});
\]

\[
\text{Ph alpha; //store df(|nabla u|^2)}
\]

\[
// The functionn J
// J(u) = 1/2 \int_{\Omega} f(|\nabla u|^2) - \int_{\Omega} u b
\]

\[
fespace \text{Ph alpha; //store df(|nabla u|^2)}
\]

\[
// The functionn J
// J(u) = 1/2 \int_{\Omega} f(|\nabla u|^2) - \int_{\Omega} u b
\]

\[
\text{func real J (real[int] \& u)}
\]

\[
Vh w;
\]

\[
w[] = u;
\]

\[
\text{real r = int2d(Th) (0.5*integrate(dx(w)*dx(w) + dy(w)*dy(w)) - b*w)};
\]

\[
cout << "J(u) = " << r << " " << u.min << " " << u.max << endl;
\]

\[
\text{return r;}
\]

(continues on next page)
func real[int] dJ (real[int] & u) {
    Vh w;
    w[] = u;
    alpha = df(dx(w)*dx(w) + dy(w)*dy(w));
    varf au (uh, vh) = int2d(Th) {
        alpha*(dx(uh)*dx(vh) + dy(uh)*dy(vh)) - b*vh
    } + on(1, 2, 3, 4, uh=0) ;
    u = au(0, Vh);
    return u; //warning: no return of local array
}

We also want to construct a preconditioner $C$ with solving the problem:
find $u_h \in V_{0h}$ such that:
\[
\forall v_h \in V_{0h}, \quad \int_{\Omega} \alpha \nabla u_h \cdot \nabla v_h = \int_{\Omega} b v_h
\]
where $\alpha = f'(|\nabla u|^2)$.

alpha = df(dx(u)*dx(u) + dy(u)*dy(u));
varf alap (uh, vh) = int2d(Th) {
    alpha*(dx(uh)*dx(vh) + dy(uh)*dy(vh))
} + on(1, 2, 3, 4, uh=0) ;
varf amass(uh, vh) = int2d(Th) {
    uh*vh
} + on(1, 2, 3, 4, uh=0) ;
matrix Amass = amass(Vh, Vh, solver=CG);
matrix Alap= alap(Vh, Vh, solver=Cholesky, factorize=1);

// Preconditioner
func real[int] C(real[int] & u) {
    real[int] w = u;
    u = Alap^-1*w;
    return u; //warning: no return of local array variable
}

To solve the problem, we make 10 iterations of the conjugate gradient, recompute the preconditioner and restart the conjugate gradient:
```
int conv=0;
for(int i = 0; i < 20; i++) {
    conv = NLCG(dJ, u[], nbiter=10, precon=C, veps=eps, verbosity=5);
    if (conv) break;
    alpha = df(dx(u)*dx(u) + dy(u)*dy(u));
    Alap = alap(Vh, Vh, solver=Cholesky, factorize=1);
    cout << "Restart with new preconditioner " << conv << ", eps =" << eps << endl;
}
// Plot
plot (u, wait=true, cmm="solution with NLCG");
```

For a given symmetric positive matrix $A$, consider the quadratic form

$$J(x) = \frac{1}{2} x^T A x - b^T x$$

then $J(x)$ is minimized by the solution $x$ of $A x = b$. In this case, we can use the function `AffineCG`

```c
AffineCG(A, x, precon=M, nbiter=imax, eps=epsilon, stop=stp);
```

If $A$ is not symmetric, we can use GMRES (Generalized Minimum Residual) algorithm by

```c
AffineGMRES(A, x, precon=M, nbiter=imax, eps=epsilon);
```

Also, we can use the non-linear version of GMRES algorithm (the function $J$ is just convex)

```c
AffineGMRES(dJ, x, precon=M, nbiter=imax, eps=epsilon);
```

For the details of these algorithms, refer to [PIRONNEAU1998], Chapter IV, 1.3.

### 3.5.2 Algorithms for Unconstrained Optimization

Two algorithms of COOOL package are interfaced with the Newton Raphson method (called `Newton`) and the BFGS method. These two are directly available in FreeFEM (no dynamical link to load). Be careful with these algorithms, because their implementation uses full matrices. We also provide several optimization algorithms from the NLopt library as well as an interface for Hansen’s implementation of CMAES (a MPI version of this one is also available).

**Example of usage for BFGS or CMAES**

**Tip:** BFGS

```c
real[int] b(10), u(10);

func real J (real[int] & u) {
real s = 0;
for (int i = 0; i < u.n; i++)
    s += (i+1)*u[i]*u[i]*0.5 - b[i]*u[i];
if (debugJ)
    cout << "J = " << s << ", u = " << u[0] << " " << u[1] << endl;
return s;
```
It is almost the same as using the CMA evolution strategy except, that since it is a derivative free optimizer, the `dJ` argument is omitted and there are some other named parameters to control the behavior of the algorithm. With the same objective function as above, an example of utilization would be (see CMAES Variational inequality for a complete example):

```plaintext
load "ff-cmaes"
//define J, u, ...
real min = cmaes(J, u, stopTolFun=1e-6, stopMaxIter=3000);
cout << "minimum value is " << min << " for u = " << u << endl;
```

This algorithm works with a normal multivariate distribution in the parameters space and tries to adapt its covariance matrix using the information provided by the successive function evaluations (see NLopt documentation for more details). Therefore, some specific parameters can be passed to control the starting distribution, size of the sample generations, etc... Named parameters for this are the following:

- **seed**= Seed for random number generator (*val* is an integer). No specified value will lead to a clock based seed initialization.
- **initialStdDev**= Value for the standard deviations of the initial covariance matrix (*val* is a real). If the value $\sigma$ is passed, the initial covariance matrix will be set to $\sigma I$. The expected initial distance between initial $X$ and the $\text{argmin}$ should be roughly initialStdDev. Default is $0.3$.
- **initialStdDevs**= Same as above except that the argument is an array allowing to set a value of the initial standard deviation for each parameter. Entries differing by several orders of magnitude should be avoided (if it can’t be, try rescaling the problem).
- **stopTolFun**= Stops the algorithm if function value differences are smaller than the passed one, default is $10^{-12}$.  
- **stopTolFunHist**= Stops the algorithm if function value differences from the best values are smaller than the passed one, default is 0 (unused).  
- **stopTolX**= Stopping criteria is triggered if step sizes in the parameters space are smaller than this real value, default is 0.  
- **stopTolXFactor**= Stopping criteria is triggered when the standard deviation increases more than this value. The default value is $10^3$.  

---

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• **stopMaxFunEval** = Stops the algorithm when `stopMaxFunEval` function evaluations have been done. Set to $900(n + 3)^2$ by default, where $n$ is the parameters space dimension.

• **stopMaxIter** = Integer stopping the search when `stopMaxIter` generations have been sampled. Unused by default.

• **popsize** = Integer value used to change the sample size. The default value is $4 + \left\lfloor 3 \ln(n) \right\rfloor$. Increasing the population size usually improves the global search capabilities at the cost of, at most, a linear reduction of the convergence speed with respect to `popsize`.

• **paramFile** = This string type parameter allows the user to pass all the parameters using an extern file, as in Hansen’s original code. More parameters related to the CMA-ES algorithm can be changed with this file. Note that the parameters passed to the CMAES function in the FreeFEM script will be ignored if an input parameters file is given.

### 3.5.3 IPOPT

The `ff-Ipopt` package is an interface for the IPOPT [WÄCHTER2006] optimizer. IPOPT is a software library for large scale, non-linear, constrained optimization. It implements a primal-dual interior point method along with filter method based line searches.

IPOPT needs a direct sparse symmetric linear solver. If your version of FreeFEM has been compiled with the `--enable-downlad` tag, it will automatically be linked with a sequential version of MUMPS. An alternative to MUMPS would be to download the HSL subroutines (see Compiling and Installing the Java Interface JIPOPT) and place them in the `/ipopt/Ipopt-3.10.2/ThirdParty/HSL` directory of the FreeFEM downloads folder before compiling.

#### Short description of the algorithm

In this section, we give a very brief glimpse at the underlying mathematics of IPOPT. For a deeper introduction on interior methods for nonlinear smooth optimization, one may consult [FORSGREN2002], or [WÄCHTER2006] for more IPOPT specific elements. IPOPT is designed to perform optimization for both equality and inequality constrained problems. However, nonlinear inequalities are rearranged before the beginning of the optimization process in order to restrict the panel of nonlinear constraints to those of the equality kind. Each nonlinear inequality is transformed into a pair of simple bound inequalities and nonlinear equality constraints by the introduction of as many slack variables as is needed:

$$c_i(x) \leq 0 \text{ becomes } c_i(x) + s_i = 0 \text{ and } s_i \leq 0,$$

where $s_i$ is added to the initial variables of the problems $x_i$. Thus, for convenience, we will assume that the minimization problem does not contain any nonlinear inequality constraint. It means that, given a function $f : \mathbb{R}^n \mapsto \mathbb{R}$, we want to find:

$$x_0 = \underset{x \in V}{\text{argmin}} f(x)$$

with $V = \{ x \in \mathbb{R}^n \mid c(x) = 0 \text{ and } x_l \leq x \leq x_u \}$

Where $c : \mathbb{R}^n \to \mathbb{R}^m$ and $x_l, x_u \in \mathbb{R}^n$ and inequalities hold componentwise. The $f$ function as well as the constraints $c$ should be twice-continuously differentiable.

As a barrier method, interior points algorithms try to find a Karush-Kuhn-Tucker point for (3.18) by solving a sequence of problems, unconstrained with respect to the inequality constraints, of the form:

$$\text{for a given } \mu > 0, \text{ find } x_\mu = \underset{x \in \mathbb{R}^n \mid c(x) = 0}{\text{argmin}} B(x, \mu)$$

(3.19)

Where $\mu$ is a positive real number and

$$B(x, \mu) = f(x) - \mu \sum_{i=1}^{n} \ln(x_{u,i} - x_i) - \mu \sum_{i=1}^{m} \ln(x_i - x_{l,i})$$
The remaining equality constraints are handled with the usual Lagrange multipliers method. If the sequence of barrier parameters $\mu$ converge to 0, intuition suggests that the sequence of minimizers of (3.19) converge to a local constrained minimizer of (3.18). For a given $\mu$, (3.19) is solved by finding $(x_\mu, \lambda_\mu) \in \mathbb{R}^n \times \mathbb{R}^m$ such that:

$$\nabla B(x_\mu, \mu) + \sum_{i=1}^m \lambda_{\mu i} \nabla c_i(x_\mu) = \nabla B(x_\mu, \mu) + J_c(x_\mu)^T \lambda_\mu = 0$$

$$c(x_\mu) = 0$$

(3.20)

The derivations for $\nabla B$ only holds for the $x$ variables, so that:

$$\nabla B(x, \mu) = \nabla f(x) + \begin{pmatrix} \mu/(x_{u,1} - x_1) \\ \vdots \\ \mu/(x_{u,n} - x_n) \end{pmatrix} - \begin{pmatrix} \mu/(x_1 - x_{l,1}) \\ \vdots \\ \mu/(x_n - x_{l,n}) \end{pmatrix}$$

If we respectively call $z_u(x, \mu) = (\mu/(x_{u,1} - x_1), \ldots, \mu/(x_{u,n} - x_n))$ and $z_l(x, \mu)$ the other vector appearing in the above equation, then the optimum $(x_\mu, \lambda_\mu)$ satisfies:

$$\nabla f(x_\mu) + J_c(x_\mu)^T \lambda_\mu + z_u(x_\mu, \mu) - z_l(x_\mu, \mu) = 0 \quad \text{and} \quad c(x_\mu) = 0$$

(3.21)

In this equation, the $z_l$ and $z_u$ vectors seem to play the role of Lagrange multipliers for the simple bound inequalities, and indeed, when $\mu \to 0$, they converge toward some suitable Lagrange multipliers for the KKT conditions, provided some technical assumptions are fulfilled (see [FORSgren2002]).

Equation (3.21) is solved by performing a Newton method in order to find a solution of (3.20) for each of the decreasing values of $\mu$. Some order 2 conditions are also taken into account to avoid convergence to local maximizers, see [FORSgren2002] for details about them. In the most classic IP algorithms, the Newton method is directly applied to (3.20). This is in most case inefficient due to frequent computation of infeasible points. These difficulties are avoided in Primal-Dual interior point methods where (3.20) is transformed into an extended system where

$$\begin{pmatrix} \nabla f(x_\mu) + J_c(x_\mu)^T \lambda_\mu + z_u(x_\mu, \mu) - z_l(x_\mu, \mu) = 0 \\ c(x_\mu) = 0 \end{pmatrix}$$

(3.22)

Where if $a$ is a vector of $\mathbb{R}^n$, $A$ denotes the diagonal matrix $A = (a_i \delta_{ij})_{1 \leq i, j \leq n}$ and $e \in \mathbb{R}^n = (1, 1, \ldots, 1)$. Solving this nonlinear system by the Newton method is known as being the primal-dual interior point method. Here again, more details are available in [FORSgren2002]. Most actual implementations introduce features in order to globalize the convergence capability of the method, essentially by adding some line-search steps to the Newton algorithm, or by using trust regions. For the purpose of IPOPT, this is achieved by a filter line search methods, the details of which can be found in [Wächter2006].

More IPOPT specific features or implementation details can be found in [Wächter2006]. We will just retain that IPOPT is a smart Newton method for solving constrained optimization problems, with global convergence capabilities due to a robust line search method (in the sense that the algorithm will converge no matter the initializer). Due to the underlying Newton method, the optimization process requires expressions of all derivatives up to the order 2 of the fitness function as well as those of the constraints. For problems whose Hessian matrices are difficult to compute or lead to high dimensional dense matrices, it is possible to use a BFGS approximation of these objects at the cost of a much slower convergence rate.
IPOPT in FreeFEM

Calling the IPOPT optimizer in a FreeFEM script is done with the IPOPT function included in the ff-Ipopt dynamic library. IPOPT is designed to solve constrained minimization problems in the form:

\[
\begin{align*}
\text{find } x_0 &= \underset{x \in \mathbb{R}^n}{\text{argmin}} \, f(x) \\
\text{s.t. } \forall i \leq n, & \quad x_i^{lb} \leq x_i \leq x_i^{ub} \quad \text{(simple bounds)} \\
\forall i \leq m, & \quad c_i^{lb} \leq c_i(x) \leq c_i^{ub} \quad \text{(constraints functions)}
\end{align*}
\]

Where \( u^{lb} \text{ and } l^{lb} \) stand for “upper bound” and “lower bound”. If for some \( i, 1 \leq i \leq m \) we have \( c_i^{lb} = c_i^{ub} \), it means that \( c_i \) is an equality constraint, and an inequality one if \( c_i^{lb} < c_i^{ub} \).

There are different ways to pass the fitness function and constraints. The more general one is to define the functions using the keyword \textsc{func}. Any returned matrix must be a sparse one (type \texttt{matrix}, not a \texttt{real[int,int]}):

```cpp
1  func real J (real[int] &X) {...} //Fitness Function, returns a scalar
2  func real[int] gradJ (real[int] &X) {...} //Gradient is a vector
3  func real[int] C (real[int] &X) {...} //Constraints
4  func matrix jacC (real[int] &X) {...} //Constraints Jacobian
```

\textbf{Warning:} In the current version of FreeFEM, returning a \texttt{matrix} object that is local to a function block leads to undefined results. For each sparse matrix returning function you define, an extern matrix object has to be declared, whose associated function will overwrite and return on each call. Here is an example for \texttt{jacC}:

```cpp
1  matrix jacCBuffer; //just declare, no need to define yet
2  func matrix jacC (real[int] &X) {
3      ...//fill jacCBuffer
4      return jacCBuffer;
5  }
```

\textbf{Warning:} IPOPT requires the structure of each matrix at the initialization of the algorithm. Some errors may occur if the matrices are not constant and are built with the \texttt{matrix A = [I,J,C]} syntax, or with an intermediary full matrix (\texttt{real[int,int]}), because any null coefficient is discarded during the construction of the sparse matrix. It is also the case when making matrices linear combinations, for which any zero coefficient will result in the suppression of the matrix from the combination. Some controls are available to avoid such problems. Check the named parameter descriptions (\texttt{checkindex}, \texttt{structhess} and \texttt{structjac} can help). We strongly advice to use \texttt{varf} as much as possible for the matrix forging.

The Hessian returning function is somewhat different because it has to be the Hessian of the Lagrangian function:

\[
(x, \sigma, \lambda) \mapsto \sigma f''(x) + \sum_{i=1}^{m} \lambda_i c_i''(x) \quad \text{where } \lambda \in \mathbb{R}^m \text{ and } \sigma \in \mathbb{R}
\]

Your Hessian function should then have the following prototype:

```cpp
1  matrix hessianLBuffer; //Just to keep it in mind
2  func matrix hessianL (real[int] &X, real sigma, real[int] &lambda) {...}
```

If the constraints functions are all affine, or if there are only simple bound constraints, or no constraint at all, the Lagrangian Hessian is equal to the fitness function Hessian, one can then omit the \texttt{sigma} and \texttt{lambda} parameters:
matrix hessianJBuffer;
func matrix hessianJ (real[int] &X){...} //Hessian prototype when constraints are affine

When these functions are defined, IPOPT is called this way:

real[int] Xi = ...; //starting point
IPOPT(J, gradJ, hessianL, C, jacC, Xi, /*some named parameters*/);

If the Hessian is omitted, the interface will tell IPOPT to use the (L)BFGS approximation (it can also be enabled with a named parameter, see further). Simple bound or unconstrained problems do not require the constraints part, so the following expressions are valid:

IPOPT(J, gradJ, C, jacC, Xi, ... ); //IPOPT with BFGS
IPOPT(J, gradJ, hessianJ, Xi, ... ); //Newton IPOPT without constraints
IPOPT(J, gradJ, Xi, ... ); //BFGS, no constraints

Simple bounds are passed using the lb and ub named parameters, while constraint bounds are passed with the clb and cub ones. Unboundedness in some directions can be achieved by using the 1e19 and −1e19 values that IPOPT recognizes as +∞ and −∞:

real[int] xlb(n), xub(n), clb(m), cub(m);
//fill the arrays...
IPOPT(J, gradJ, hessianL, C, jacC, Xi, lb=xlb, ub=xub, clb=clb, cub=cub, /*some other named parameters*/);

P2 fitness function and affine constraints function: In the case where the fitness function or constraints function can be expressed respectively in the following forms:

\[ \forall x \in \mathbb{R}^n, f(x) = \frac{1}{2} \langle Ax, x \rangle + \langle b, x \rangle \quad (A, b) \in \mathcal{M}_{n,n}(\mathbb{R}) \times \mathbb{R}^n \]

or, \( C(x) = Ax + b \quad (A, b) \in \mathcal{M}_{m,n}(\mathbb{R}) \times \mathbb{R}^m \)

where \( A \) and \( b \) are constant, it is possible to directly pass the \( (A, b) \) pair instead of defining 3 (or 2) functions. It also indicates to IPOPT that some objects are constant and that they have to be evaluated only once, thus avoiding multiple copies of the same matrix. The syntax is:

// Affine constraints with "standard" fitness function
matrix A = ... ; //linear part of the constraints
real[int] b = ... ; //constant part of constraints
IPOPT([A, b], ..., Xi, /*bounds and named parameters*/);

//[b, A] would work as well.

Note that if you define the constraints in this way, they don’t contribute to the Hessian, so the Hessian should only take one real[int] as an argument.

// Affine constraints and P2 fitness func
matrix A = ... ; //bilinear form matrix
real[int] b = ... ; //linear contribution to f
matrix Ac = ... ; //linear part of the constraints
real[int] bc = ... ; //constant part of constraints
IPOPT([A, b], [Ac, bc], Xi, /*bounds and named parameters*/);

If both objective and constraint functions are given this way, it automatically activates the IPOPT mehrotra_algorithm option (better for linear and quadratic programming according to the documentation). Otherwise, this option can only be set through the option file (see the named parameters section).

A false case is the one of defining \( f \) in this manner while using standard functions for the constraints:

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Indeed, when passing \([A, b]\) in order to define \(f\), the Lagrangian Hessian is automatically built and has the constant \(x \mapsto A\) function, with no way to add possible constraint contributions, leading to incorrect second order derivatives. So, a problem should be defined like that in only two cases:

1. constraints are nonlinear but you want to use the BFGS mode (then add \(bfgs=1\) to the named parameter),
2. constraints are affine, but in this case, compatible to pass in the same way

Here are some other valid definitions of the problem (cases when \(f\) is a pure quadratic or linear form, or \(C\) a pure linear function, etc...):

```plaintext
// Pure quadratic f - A is a matrix
IPOPT(A, /*constraints arguments*/, Xi, /*bound and named parameters*/);
// Pure linear f - b is a real[int]
IPOPT(b, /*constraints arguments*/, Xi, /*bound and named parameters*/);
// Linear constraints - Ac is a matrix
IPOPT(/*fitness function arguments*/, Ac, Xi, /*bound and named parameters*/);
```

**Returned Value:** The IPOPT function returns an error code of type int. A zero value is obtained when the algorithm succeeds and positive values reflect the fact that IPOPT encounters minor troubles. Negative values reveal more problematic cases. The associated IPOPT return tags are listed in the table below. The IPOPT pdf documentation provides a more accurate description of these return statuses:

<table>
<thead>
<tr>
<th>Success</th>
<th>Failures</th>
</tr>
</thead>
<tbody>
<tr>
<td>0 Solve_Succeeded</td>
<td>-1 Maximum_Iterations_Exceeded</td>
</tr>
<tr>
<td>1 Solved_To_Acceptable_Level</td>
<td>-2 Restoration_Failed</td>
</tr>
<tr>
<td>2 Infeasible_Problem_Detected</td>
<td>-3 Error_In_Step_Computation</td>
</tr>
<tr>
<td>3 Search_Direction_Becomes_Too_Small</td>
<td>-4 Maximum_CpuTime_Exceeded</td>
</tr>
<tr>
<td>4 Diverging_Iterates</td>
<td></td>
</tr>
<tr>
<td>5 User_Requestened_Stop</td>
<td></td>
</tr>
<tr>
<td>6 Feasible_Point_Found</td>
<td></td>
</tr>
</tbody>
</table>

**Problem definition issues**

-10 Not_Enough_Degrees_Of_Freedom                                       
-11 Invalid_Problem_Definition                                         
-12 Invalid_Option                                                     
-13 Invalid_Number_Detected                                            

**Critical errors**

-100 Unrecoverable_Exception                                           
-101 NonIpopt_Exception_Thrown                                         
-102 Insufficient_Memory                                               
-199 Internal_Error                                                    

**Named Parameters:** The available named parameters in this interface are those we thought to be the most subject to variations from one optimization to another, plus a few that are interface specific. Though, as one could see at IPOPT Linear solver, there are many parameters that can be changed within IPOPT, affecting the algorithm behavior. These parameters can still be controlled by placing an option file in the execution directory. Note that IPOPT’s pdf documentation may provides more information than the previously mentioned online version for certain parameters. The in-script available parameters are:

- \(lb, ub: \text{real}[\text{int}]\) for lower and upper simple bounds upon the search variables must be of size \(n\) (search space dimension). If two components of the same index in these arrays are equal then the corresponding search variable is fixed. By default IPOPT will remove any fixed variable from the optimization process and always use the fixed value when calling functions. It can be changed using the \texttt{fixedVar} parameter.
- **clb, cub**: `real[int]` of size $m$ (number of constraints) for lower and upper constraints bounds. Equality between two components of the same index $i$ in `clb` and `cub` reflect an equality constraint.

- **struct jacc**: To pass the greatest possible structure (indexes of non null coefficients) of the constraint Jacobians under the form `[I, J]` where `I` and `J` are two integer arrays. If not defined, the structure of the constraint Jacobians, evaluated in $X_i$, is used (no issue if the Jacobian is constant or always defined with the same `varf`, hazardous if it is with a triplet array or if a full matrix is involved).

- **struct Hess**: Same as above but for the Hessian function (unused if $f$ is P2 or less and constraints are affine). Here again, keep in mind that it is the Hessian of the Lagrangian function (which is equal to the Hessian of $f$ only if constraints are affine). If no structure is given with this parameter, the Lagrangian Hessian is evaluated on the starting point, with $\sigma = 1$ and $\lambda = (1, 1, \ldots, 1)$ (it is safe if all the constraints and fitness function Hessians are constant or build with `varf`, and here again it is less reliable if built with a triplet array or a full matrix).

- **checkindex**: A boolean that triggers a dichotomic index search when matrices are copied from FreeFEM functions to IPOPT arrays. It is used to avoid wrong index matching when some null coefficients are removed from the matrices by FreeFEM. It will not solve the problems arising when a too small structure has been given at the initialization of the algorithm. Enabled by default (except in cases where all matrices are obviously constant).

- **warmstart**: If set to true, the constraints dual variables $\lambda$, and simple bound dual variables are initialized with the values of the arrays passed to `lm`, `lz` and `uz` named parameters (see below).

- **lm**: `real[int]` of size $m$, which is used to get the final values of the constraints dual variables $\lambda$ and/or initialize them in case of a warm start (the passed array is also updated to the last dual variables values at the end of the algorithm).

- **lz, uz**: `real[int]` of size $n$ to get the final values and/or initialize (in case of a warm start) the dual variables associated to simple bounds.

- **tol**: `real`, convergence tolerance for the algorithm, the default value is $10^{-8}$.

- **maxiter**: `int`, maximum number of iterations with 3000 as default value.

- **maxcputime**: `real` value, maximum runtime duration. Default is $10^6$ (almost 11 and a half days).

- **bfgs**: boolean enabling or not the (low-storage) BFGS approximation of the Lagrangian Hessian. It is set to false by default, unless there is no way to compute the Hessian with the functions that have been passed to IPOPT.

- **derivativetest**: Used to perform a comparison of the derivatives given to IPOPT with finite differences computation. The possible string values are: "none" (default), "first-order", "second-order" and "only-second-order". The associated derivative error tolerance can be changed via the option file. One should not care about any error given by it before having tried, and failed, to perform a first optimization.

- **dth**: Perturbation parameter for the derivative test computations with finite differences. Set by default to $10^{-8}$.

- **dttol**: Tolerance value for the derivative test error detection (default value unknown yet, maybe $10^{-5}$).

- **optfile**: `string` parameter to specify the IPOPT option file name. IPOPT will look for a `ipopt.opt` file by default. Options set in the file will overwrite those defined in the FreeFEM script.

- **printlevel**: An `int` to control IPOPT output print level, set to 5 by default, the possible values are from 0 to 12. A description of the output information is available in the PDF documentation of IPOPT.

- **fixedvar**: `string` for the definition of simple bound equality constraints treatment: use "make_parameter" (default value) to simply remove them from the optimization process (the functions will always be evaluated with the fixed value for those variables), "make_constraint" to treat them as any other constraint or "relax_bounds" to relax fixing bound constraints.

- **mustrategy**: `string` to choose the update strategy for the barrier parameter $\mu$. The two possible tags are "monotone", to use the monotone (Fiacco-McCormick) strategy, or "adaptive" (default setting).

---

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• **muinit**: real positive value for the barrier parameter initialization. It is only relevant when **mustrategy** has been set to **monotone**.

• **pivtol**: real value to set the pivot tolerance for the linear solver. A smaller number pivots for sparsity, a larger number pivots for stability. The value has to be in the \([0, 1]\) interval and is set to \(10^{-6}\) by default.

• **brf**: Bound relax factor: before starting the optimization, the bounds given by the user are relaxed. This option sets the factor for this relaxation. If it is set to zero, then the bound relaxation is disabled. This real has to be positive and its default value is \(10^{-8}\).

• **objvalue**: An identifier to a **real** type variable to get the last value of the objective function (best value in case of success).

• **mumin**: minimum value for the barrier parameter \(\mu\), a **real** with \(10^{-11}\) as default value.

• **linesearch**: A boolean which disables the line search when set to **false**. The line search is activated by default. When disabled, the method becomes a standard Newton algorithm instead of a primal-dual system. The global convergence is then no longer assured, meaning that many initializers could lead to diverging iterates. But on the other hand, it can be useful when trying to catch a precise local minimum without having some out of control process making the iterate caught by some other near optimum.

### 3.5.4 Some short examples using IPOPT

**Tip**: Ipopt variational inequality A very simple example consisting of, given two functions \(f\) and \(g\) (defined on \(\Omega \subset \mathbb{R}^2\)), minimizing \(J(u) = \frac{1}{2} \int_{\Omega} |\nabla u|^2 - \int_{\Omega} fu\), with \(u \leq g\) almost everywhere:

```plaintext
// Solve
//- Delta u = f
//u < g
//u = 0 on Gamma
load "ff-Ipopt";

// Parameters
int nn = 20;
func f = 1.; //rhs function
real r = 0.03, s = 0.1;
func g = r - r/2*exp(-0.5*(square(x-0.5) + square(y-0.5))/square(s));

// Mesh
mesh Th = square(nn, nn);

// Fespace
fespace Vh(Th, P2);
Vh u = 0;
Vh lb = -1.e19;
Vh ub = g;

// Macro
macro Grad(u) [dx(u), dy(u)]

// Problem
varf vP (u, v)
  = int2d(Th) (Grad(u)'*Grad(v)
              - int2d(Th) (continues on next page)
```

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Here we build the matrix and second member associated to the function to fully and finally minimize it. The $[A, b]$ syntax for the fitness function is then used to pass it to IPOPT.

```cpp
matrix A = vP(Vh, Vh, solver=CG);
real[int] b = vP(0, Vh);
```

We use simple bounds to impose the boundary condition $u = 0$ on $\partial \Omega$, as well as the $u \leq g$ condition.

```cpp
varf vGamma (u, v) = on(1, 2, 3, 4, u=1);
real[int] onGamma = vGamma(0, Vh);
ub[] = onGamma ? 0. : ub[];
lb[] = onGamma ? 0. : lb[];
IPOPT([A, b], u[], lb=lb[], ub=ub[]);
```

**Tip:** Ipopt variational inequality 2

Let $\Omega$ be a domain of $\mathbb{R}^2$. $f_1, f_2 \in L^2(\Omega)$ and $g_1, g_2 \in L^2(\partial \Omega)$ four given functions with $g_1 \leq g_2$ almost everywhere. We define the space:

$$V = \{(v_1, v_2) \in H^1(\Omega)^2; v_1|_{\partial \Omega} = g_1, v_2|_{\partial \Omega} = g_2, v_1 \leq v_2 \text{ a.e.}\}$$

as well as the function $J : H^1(\Omega)^2 \rightarrow \mathbb{R}$:

$$J(v_1, v_2) = \frac{1}{2} \int_{\Omega} |\nabla v_1|^2 - \int_{\Omega} f_1 v_1 + \frac{1}{2} \int_{\Omega} |\nabla v_2|^2 - \int_{\Omega} f_2 v_2$$

The problem entails finding (numerically) two functions $(u_1, u_2) = \arg\min_{(v_1, v_2) \in V} J(v_1, v_2)$.
Vh [lz, lz2] = [1, 1];
Vh [u1, u2] = [0, 0]; //starting point

fespace Wh(Th, [P1]);
Wh lm=1.;

// Macro
macro Grad(u) [dx(u), dy(u)] //

// Loop
int iter=0;
while (++iter){
  // Problem
  varf vP ([u1, u2], [v1, v2])
  = int2d(Th)(
      Grad(u1)'*Grad(v1)
      + Grad(u2)'*Grad(v2)
    )
    - int2d(Th)(
      f1*v1
      + f2*v2
    )
  ;

  matrix A = vP(Vh, Vh); //fitness function matrix
  real[int] b = vP(0, Vh); //and linear form

  int[int] II1 = [0], II2 = [1]; //Constraints matrix
  matrix C1 = interpolate (Wh, Vh, U2Vc =II1);
  matrix C2 = interpolate (Wh, Vh, U2Vc =II2);
  matrix CC = -1*C1 + C2; // u2 - u1 > 0
  Wh cl = 0; //constraints lower bounds (no upper bounds)

  //Boundary conditions
  varf vGamma ([u1, u2], [v1, v2]) = on(1, 2, 3, 4, u1=1, u2=1);
  real[int] onGamma = vGamma(0, Vh);
  Vh [ub1, ub2] = [g1, g2];
  Vh [lb1, lb2] = [g1, g2];
  ub1[] = onGamma ? ub1[] : 1e19; //Unbounded in interior
  lb1[] = onGamma ? lb1[] : -1e19;

  Vh [uzi, uzi2] = [uz, uz2], [lzi, lzi2] = [lz, lz2];
  Wh lmi = lm;
  Vh [ui1, ui2] = [u1, u2];

  // Solve
  IPOPT([b, A], CC, ui1[], lb=lb1[], clb=cl[]), ub=ub1[], warmstart=iter>1, uz=uzi[],
  lzi=lzi[]], lm=lmi[]);

  // Plot
  plot(ui1, ui2, wait=true, nbiso=60, dim=3);
  if(iter > 1) break;

  // Mesh adpatation
  Th = adaptmesh(Th, [ui1, ui2], err=0.004, nbvx=100000);
  [uz, uz2] = [uzi, uzi2];
(continues on next page)
3.5.5 3D constrained minimum surface with IPOPT

Area and volume expressions

This example is aimed at numerically solving some constrained minimum surface problems with the IPOPT algorithm. We restrain to $C^k$ ($k \geq 1$), closed, spherically parametrizable surfaces, i.e. surfaces $S$ such that:

$$\exists \rho \in C^k([0, 2\pi] \times [0, \pi])|S = \left\{ X = \begin{pmatrix} \rho(\theta, \phi) \\ 0 \\ 0 \end{pmatrix}, (\theta, \phi) \in [0, 2\pi] \times [0, \pi] \right\}$$

Where the components are expressed in the spherical coordinate system. Let’s call $\Omega$ the $[0, 2\pi] \times [0, \pi]$ angular parameters set. In order to exclude self crossing and opened shapes, the following assumptions upon $\rho$ are made:

$$\rho \geq 0 \quad \text{and} \quad \forall \phi, \rho(0, \phi) = \rho(2\pi, \phi)$$

For a given function $\rho$ the first fundamental form (the metric) of the defined surface has the following matrix representation:

$$G = \begin{pmatrix} \rho^2 \sin^2(\phi) + (\partial_\theta \rho)^2 & \partial_\theta \rho \partial_\phi \rho \\ \partial_\theta \rho \partial_\phi \rho & \rho^2 + (\partial_\phi \rho)^2 \end{pmatrix}$$

(3.23)

This metric is used to express the area of the surface. Let $g = \det(G)$, then we have:

$$A(\rho) = \int_{\Omega} \| \partial_\theta X \wedge \partial_\phi X \| = \int_{\Omega} \sqrt{g}$$

$$= \int_{\Omega} \sqrt{\rho^2(\partial_\theta \rho)^2 + \rho^4 \sin^2(\phi) + \rho^2(\partial_\phi \rho)^2 \sin^2(\phi)} d\theta d\phi$$

(3.24)

The volume of the space enclosed within the shape is easier to express:

$$V(\rho) = \int_{\Omega} \int_{0}^{\rho(\theta, \phi)} r^2 \sin(\phi) dr d\theta d\phi = \frac{1}{3} \int_{\Omega} \rho^4 \sin(\phi) d\theta d\phi$$

(3.25)
Derivatives

In order to use a Newton based interior point optimization algorithm, one must be able to evaluate the derivatives of $\mathcal{A}$ and $\mathcal{V}$ with respect to $\rho$. Concerning the area, we have the following result:

$$\forall v \in C^1(\Omega), \langle d\mathcal{A}(\rho), v \rangle = \int_{\Omega} \frac{1}{2} \frac{d\bar{g}(\rho)}{\sqrt{\bar{g}}} \, d\theta d\phi$$

Where $\bar{g}$ is the application mapping the $(\theta, \phi) \mapsto g(\theta, \phi)$ scalar field to $\rho$. This leads to the following expression, easy to transpose in a freefem script using:

$$\forall v \in C^1(\Omega), \langle d\mathcal{A}(\rho), v \rangle = \int_{\Omega} (2\rho^3 \sin^2(\phi) + \rho(\partial_{\theta}\rho)^2 + \rho(\partial_{\phi}\rho)^2 \sin^2(\phi)) \, v + \int_{\Omega} \rho^2 \partial_{\theta} \rho \partial_{\theta} v + \rho^2 \partial_{\phi} \rho \sin^2(\phi) \partial_{\phi} v$$

With a similar approach, one can derive an expression for second order derivatives. However, comporting no specific difficulties, the detailed calculus are tedious, the result is that these derivatives can be written using a $3 \times 3$ matrix $B$ whose coefficients are expressed in term of $\rho$ and its derivatives with respect to $\theta$ and $\phi$, such that:

$$\forall (w, v) \in C^1(\Omega), d^2\mathcal{A}(\rho)(w, v) = \int_{\Omega} \left( \begin{array}{c} v \\ \partial_{\theta} w \\ \partial_{\phi} w \end{array} \right) B \left( \begin{array}{c} v \\ \partial_{\theta} v \\ \partial_{\phi} v \end{array} \right) d\theta d\phi$$

Deriving the volume function derivatives is again an easier task. We immediately get the following expressions:

$$\forall v, \langle d\mathcal{V}(\rho), v \rangle = \int_{\Omega} \rho^2 \sin^2(\phi) \, v \, d\theta d\phi$$

$$\forall w, v, d^2\mathcal{V}(\rho)(w, v) = \int_{\Omega} 2\rho^2 \sin^2(\phi) w v \, d\theta d\phi$$

The problem and its script

The whole code is available in IPOPT minimal surface & volume example. We propose to solve the following problem:

**Tip:** Given a positive function $\rho_{\text{object}}$ piecewise continuous, and a scalar $\mathcal{V}_{\text{max}} > \mathcal{V}(\rho_{\text{object}})$, find $\rho_0$ such that:

$$\rho_0 = \arg\min_{\rho \in C^1(\Omega)} \mathcal{A}(\rho), \text{ s.t. } \rho_0 \geq \rho_{\text{object}} \text{ and } \mathcal{V}(\rho_0) \leq \mathcal{V}_{\text{max}}$$

If $\rho_{\text{object}}$ is the spherical parametrization of the surface of a 3-dimensional object (domain) $\mathcal{O}$, it can be interpreted as finding the surface with minimum area enclosing the object with a given maximum volume. If $\mathcal{V}_{\text{max}}$ is close to $\mathcal{V}(\rho_{\text{object}})$, so should be $\rho_0$ and $\rho_{\text{object}}$. With higher values of $\mathcal{V}_{\text{max}}$, $\rho$ should be closer to the unconstrained minimum surface surrounding $\mathcal{O}$ which is obtained as soon as $\mathcal{V}_{\text{max}} \geq \frac{1}{3} \pi \|\rho_{\text{object}}\|_{\infty}^3$ (sufficient but not necessary).

It also could be interesting to solve the same problem with the constraint $\mathcal{V}(\rho_0) \geq \mathcal{V}_{\text{min}}$ which leads to a sphere when $\mathcal{V}_{\text{min}} \geq \frac{1}{6} \pi \text{diam}(\mathcal{O})^3$ and moves toward the solution of the unconstrained problem as $\mathcal{V}_{\text{min}}$ decreases.

We start by meshing the domain $[0, 2\pi] \times [0, \pi]$, then a periodic P1 finite elements space is defined.

```plaintext
// Parameters
int nadapt = 3;
real alpha = 0.9;
int np = 30;
real regtest;
```

(continues on next page)
int shapeswitch = 1;
real sigma = 2*pi/40.;
real treshold = 0.1;
real e = 0.1;
real r0 = 0.25;
real rr = 2-r0;
real E = 1./(e*e);
real RR = 1./(rr*rr);

// Mesh
mesh Th = square(2*np, np, [2*pi*x, pi*y]);

// Fespace
fespace Vh(Th, P1, periodic=[[2, y], [4, y]]);

//Initial shape definition
//outside of the mesh adaptation loop to initialize with the previous optimal shape
//found on further iterations
Vh startshape = 5;

We create some finite element functions whose underlying arrays will be used to store the values of dual variables associated to all the constraints in order to reinitialize the algorithm with it in the case where we use mesh adaptation. Doing so, the algorithm will almost restart at the accuracy level it reached before mesh adaptation, thus saving many iterations.

Vh uz = 1., lz = 1.;
real[int] lm = [1];

Then, follows the mesh adaptation loop, and a rendering function, Plot3D, using 3D mesh to display the shape it is passed with medit (the movemesh23 procedure often crashes when called with ragged shapes).

for(int kkk = 0; kkk < nadapt; ++kkk){
    int iter=0;
    func sin2 = square(sin(y));

    // A function which transform Th in 3d mesh (r=rho)
    //a point (theta,phi) of Th becomes ( r(theta,phi)*cos(theta)*sin(phi) , r(theta, phi)*sin(theta)*sin(phi) , r(theta,phi)*cos(phi) )
    //then displays the resulting mesh with medit
    func int Plot3D(real[int] &rho, string cmm, bool ffplot){
        Vh rhoo;
        rhoo[,] = rho;
        //mesh sTh = square(np, np/2, [2*pi*x, pi*y]);
        //fespace sVh(sTh, P1);
        //Vh rhoplot = rhoo;
        try{
            mesh3 Sphere = movemesh23(Th, transfo=[rhoo(x,y)*cos(x)*sin(y), rhoo(x, y)*sin(x)*sin(y), rhoo(x,y)*cos(y)]);
            if(ffplot)
                plot(Sphere);
            else
                medit(cmm, Sphere);
        }
        catch(...){
            cout << "PLOT ERROR" << endl;
        }
    return 1;
    }
}

(continues on next page)
Here are the functions related to the area computation and its shape derivative, according to equations (3.24) and (3.26):

```plaintext
// Surface computation
//Maybe is it possible to use movemesh23 to have the surface function less complicated
//However, it would not simplify the gradient and the hessian
func real Area (real[int] &X)
{
    Vh rho;
    rho[] = X;
    Vh rho2 = square(rho);
    Vh rho4 = square(rho2);
    real res = int2d(Th)(sqrt(rho4*sin2 + rho2*square(dx(rho)) +
        rho2*sin2*square(dy(rho))));
    ++iter;
    if(1)
        plot(rho, value=true, fill=true, cmm="rho(theta,phi) on [0,2pi]x[0,pi] - S=
            res, dim=3);
    else
        Plot3D(rho[], "shape_evolution", 1);
    return res;
}

func real[int] GradArea (real[int] &X)
{
    Vh rho, rho2;
    rho[] = X;
    rho2[] = square(X);
    Vh sqrtPsi, alpha;
    {
        Vh dxrho2 = dx(rho)*dx(rho), dyrho2 = dy(rho)*dy(rho);
        sqrtPsi = sqrt(rho2*rho2*sin2 + rho2*dxrho2 + rho2*dyrho2*sin2);
        alpha = 2.*rho2*rho*sin2 + rho*dxrho2 + rho*dyrho2*sin2;
    }
    varf dArea (u, v)
        = int2d(Th)(
            1./sqrtPsi * (alpha*v + rho2*dx(rho)*dx(v) + rho2*dy(rho)*sin2*dy(v))
        );
    real[int] grad = dArea(0, Vh);
    return grad;
}
```

The function returning the hessian of the area for a given shape is a bit blurry, thus we won’t show here all of equation (3.27) coefficients definition, they can be found in the edp file.

```plaintext
matrix hessianA;
func matrix HessianArea (real[int] &X)
{
    Vh rho, rho2;
    rho[] = X;
    rho2 = square(rho);
    Vh sqrtPsi, sqrtPsi3, C00, C01, C02, C11, C12, C22, A;
    {
        Vh C0, C1, C2;
    }
}
```
Vh dxrho2 = dx(rho)*dx(rho), dyrho2 = dy(rho)*dy(rho);
sqrtPsi = sqrt( rho2*rho2*sin2 + rho2*dxrho2 + rho2*dyrho2*sin2);
sqrtPsi3 = (rho2*rho2*sin2 + rho2*dxrho2 + rho2*dyrho2*sin2)*sqrtPsi;
C0 = 2*rho2*rho*sin2 + rho*dxrho2 + rho*dyrho2*sin2;
C1 = rho2*dx(rho);
C2 = rho2*dyrho2*sin2;
C00 = square(C0);
C01 = C0*dx(rho);
C02 = C0*dy(rho);
C11 = square(C1);
C12 = C1*dx(rho);
C22 = square(C2);
A = 6.*rho2*rho*sin2 + dxrho2 + dyrho2*sin2;
}

varf d2Area (w, v)
  = int2d(Th) (1./sqrtPsi * {
    A*w*v
   + 2*rho*dx(rho)*dx(w)*v
   + 2*rho*dx(rho)*w*dx(v)
   + 2*rho*dy(rho)*sin2*dy(w)*v
   + 2*rho*dy(rho)*sin2*w*dy(v)
   + rho2*dx(w)*dx(v)
   + rho2*dy(w)*dy(v)
})
  + 1./sqrtPsi3 * {
    C00*w*v
   + C01*dx(w)*v
   + C01*w*dx(v)
   + C02*dy(w)*v
   + C02*w*dy(v)
   + C11*dx(w)*dx(v)
   + C12*dx(w)*dy(v)
   + C12*dy(w)*dx(v)
   + C22*dy(w)*dy(v)
}
  ;

hessianA = d2Area(Vh, Vh);

return hessianA;
}

And the volume related functions:

// Volume computation
func real Volume (real[int] &X)
  { 
    Vh rho;
    rho[] = X;
    Vh rho3 = rho*rho*rho;
    real res = 1./3.*int2d(Th) (rho3*sin(y));
    return res;
  }

func real[int] GradVolume (real[int] &X)
  {
    Vh rho;
    rho[] = X;
    varf dVolume(u, v) = int2d(Th) (rho*rho*sin(y)*v);
}
real[int] grad = dVolume(0, Vh);
return grad;
}

matrix hessianV;
func matrix HessianVolume(real[int] &X) {
    Vh rho;
    rho[] = X;
    varf d2Volume(w, v) = int2d(Th)(2*rho*sin(y)*v*w);
    hessianV = d2Volume(Vh, Vh);
    return hessianV;
}

If we want to use the volume as a constraint function we must wrap it and its derivatives in some FreeFEM functions returning the appropriate types. It is not done in the above functions in cases where one wants to use it as a fitness function. The lagrangian hessian also has to be wrapped since the Volume is not linear with respect to \( \rho \), it has some non-null second order derivatives.

func real[int] ipVolume (real[int] &X) {
    real[int] vol = [Volume(X)];
    return vol;
}

matrix mdV;
func matrix ipGradVolume (real[int] &X) {
    real[int],int dvol(1,Vh.ndof);
    dvol(0,:) = GradVolume(X);
    mdV = dvol;
    return mdV;
}

matrix HLagrangian;
func matrix ipHessianLag (real[int] &X, real objfact, real[int] &lambda) {
    HLagrangian = objfact*HessianArea(X) + lambda[0]*HessianVolume(X);
    return HLagrangian;
}

The ipGradVolume function could pose some troubles during the optimization process because the gradient vector is transformed in a sparse matrix, so any null coefficient will be discarded. Here we create the IPOPT structure manually and use the checkindex named-parameter to avoid bad indexing during copies. This gradient is actually dense, there is no reason for some components to be constantly zero:

int[int] gvi(Vh.ndof), gvj=0:Vh.ndof-1;
gvi = 0;

These two arrays will be passed to IPOPT with structjacc=[gvi,gvj]. The last remaining things are the bound definitions. The simple lower bound must be equal to the components of the P1 projection of \( \rho_{\text{object}} \). And we choose \( \alpha \in [0,1] \) to set \( V_{\text{max}}(1-\alpha)V(\rho_{\text{object}}) + \alpha \frac{4}{3}\pi \|\rho_{\text{object}}\|_\infty^3 \):

func disc1 = sqrt(1./(RR+(E-RR)*cos(y)*cos(y)))*(1+0.1*cos(7*x));
func disc2 = sqrt(1./(RR+(E-RR)*cos(x)*cos(x)*sin2));

if(1){
    lb = r0;
    for (int q = 0; q < 5; ++q){
        func f = rr*Gaussian(x, y, 2*q*pi/5., pi/3.);
        func g = rr*Gaussian(x, y, 2*q*pi/5.+pi/5., 2.*pi/3.);
        lb = max(max(lb, f), g);
    }
    lb = max(lb, rr*Gaussian(x, y, 2*pi, pi/3));
}

lb = max(lb, max(disc1, disc2));
real Vobj = Volume(lb[]);
real Vnvc = 4./3.*pi*pow(lb[].linfty, 3);
\begin{verbatim}
if (1)
    Plot3D(lb[], "object_inside", 1);
real[int] clb = 0., cub = [(1-alpha)*Vobj + alpha*Vnvc];

Calling IPOPT:

   int res = IPOPT(Area, GradArea, ipHessianLag, ipVolume, ipGradVolume,
   rc[], ub=ub[], lb=lb[], clb=clb, cub=cub, checkindex=1, maxiter=kkk<nadapt-1 ?
   40:150, 
   warmstart=kkk, lm=lm, uz=uz[], lz=lz[], toI=0.00001, structjacc=[gvi,gvj]);
   cout << "IPOPT: res =" << res << endl ;

   // Plot
   Plot3D(rc[], "Shape_at_"+kkk, 1);
   Plot3D(GradArea(rc[]), "ShapeGradient", 1);

Finally, before closing the mesh adaptation loop, we have to perform the said adaptation. The mesh is adaptated with
respect to the $X = (\rho, 0, 0)$ (in spherical coordinates) vector field, not directly with respect to $\rho$, otherwise the true
curvature of the 3D-shape would not be well taken into account.

   if (kkk < nadapt-1){
      Th = adaptmesh(Th, rc*cos(x)*sin(y), rc*sin(x)*sin(y), rc*cos(y),
      nbvx=50000, periodic=[[2, y], [4, y]]);
      plot(Th, wait=true);
      startshape = rc;
      uz = uz;
      lz = lz;
   }

Here are some pictures of the resulting surfaces obtained for decreasing values of $\alpha$ (and a slightly more complicated
object than two orthogonal discs). We return to the enclosed object when $\alpha = 0$:

\end{verbatim}

\section{3.5.6 The \texttt{nlOpt} optimizers}

The \texttt{ff-NLopt} package provides a \texttt{FreeFEM} interface to the free/open-source library for nonlinear optimization,
easing the use of several different free optimization (constrained or not) routines available online along with the PDE
solver. All the algorithms are well documented in \texttt{NLopt} documentation, therefore no exhaustive information
concerning their mathematical specificities will be found here and we will focus on the way they are used in a \texttt{FreeFEM}
script. If needing detailed information about these algorithms, visit the website where a description of each of them is
given, as well as many bibliographical links.

Most of the gradient based algorithms of \texttt{NLopt} uses a full matrix approximation of the Hessian, so if you’re planning
to solve a large scale problem, use the IPOPT optimizer which definitely surpass them.

All the \texttt{NLopt} features are identified that way:

\begin{verbatim}
load "ff-NLopt"
   //define J, u, and maybe grad(J), some constraints etc...
   real min = nloptXXXXXX(J, u, //Unavoidable part
   grad=<name of grad(J)>, //if needed
   lb= //Lower bounds array
   ub= //Upper bounds array
   ... //Some optional arguments:
   //Constraints functions names,
\end{verbatim}
$0.4 \leq \alpha \leq 1$

$\alpha = 0.3$

$\alpha = 0.2$

$\alpha = 0.1$

$\alpha = 0.05$

$\alpha = 0$
//Stopping criteria,
//Algorithm specific parameters,
//Etc...

XXXXXX refers to the algorithm tag (not necessarily 6 characters long). \( u \) is the starting position (a \( \text{real}[\text{int}] \) type array) which will be overwritten by the algorithm, the value at the end being the found \( \arg\min \). And as usual, \( J \) is a function taking a \( \text{real}[\text{int}] \) type array as argument and returning a \( \text{real} \). \( \text{grad}, \text{lb} \) and \( \text{ub} \) are “half-optional” arguments, in the sense that they are obligatory for some routines but not all.

The possible optionally named parameters are the following, note that they are not used by all algorithms (some do not support constraints, or a type of constraints, some are gradient-based and others are derivative free, etc...). One can refer to the table after the parameters description to check which are the named parameters supported by a specific algorithm. Using an unsupported parameter will not stop the compiler work, seldom breaks runtime, and will just be ignored. When it is obvious you are missing a routine, you will get a warning message at runtime (for example if you pass a gradient to a derivative free algorithm, or set the population of a non-genetic one, etc...). In the following description, \( n \) stands for the dimension of the search space.

**Half-optional parameters :**

- \( \text{grad} \): The name of the function which computes the gradient of the cost function (prototype should be \( \text{real}[\text{int}] \to \text{real}[\text{int}] \), both argument and result should have the size \( n \)). This is needed as soon as a gradient-based method is involved, which is ignored if defined in a derivative free context.

- \( \text{lb}/\text{ub} \): Lower and upper bounds arrays (\( \text{real}[\text{int}] \) type) of size \( n \). Used to define the bounds within which the search variable is allowed to move. Needed for some algorithms, optional, or unsupported for others.

- \( \text{subOpt} \): Only enabled for the Augmented Lagrangian and MLSL methods who need a sub-optimizer in order to work. Just pass the tag of the desired local algorithm with a string.

**Constraints related parameters (optional - unused if not specified):**

- \( \text{IConst}/\text{EConst} \): Allows to pass the name of a function implementing some inequality (resp. equality) constraints on the search space. The function type must be \( \text{real}[\text{int}] \to \text{real}[\text{int}] \) where the size of the returned array is equal to the number of constraints (of the same type - it means that all of the constraints are computed in one vectorial function). In order to mix inequality and equality constraints in a same minimization attempt, two vectorial functions have to be defined and passed. See example (3.29) for more details about how these constraints have to be implemented.

- \( \text{gradIConst}/\text{gradEConst} \): Use to provide the inequality (resp. equality) constraints gradient. These are \( \text{real}[\text{int}] \to \text{real}[\text{int},\text{int}] \) type functions. Assuming we have defined a constraint function (either inequality or equality) with \( p \) constraints, the size of the matrix returned by its associated gradient must be \( p \times n \) (the \( i \)-th line of the matrix is the gradient of the \( i \)-th constraint). It is needed in a gradient-based context as soon as an inequality or equality constraint function is passed to the optimizer and ignored in all other cases.

- \( \text{tolIConst}/\text{tolEConst} \): Tolerance values for each constraint. This is an array of size equal to the number of inequality (resp. equality) constraints. Default value is set to \( 10^{-12} \) for each constraint of any type.

**Stopping criteria :**

- \( \text{stopFuncValue} \): Makes the algorithm end when the objective function reaches this \( \text{real} \) value.

- \( \text{stopRelXTol} \): Stops the algorithm when the relative moves in each direction of the search space is smaller than this \( \text{real} \) value.

- \( \text{stopAbsXTol} \): Stops the algorithm when the moves in each direction of the search space is smaller than the corresponding value in this \( \text{real}[\text{int}] \) array.

- \( \text{stopRelFTol} \): Stops the algorithm when the relative variation of the objective function is smaller than this \( \text{real} \) value.
• stopAbsFTol: Stops the algorithm when the variation of the objective function is smaller than this real value.
• stopMaxFEval: Stops the algorithm when the number of fitness evaluations reaches this integer value.
• stopTime: Stops the algorithm when the optimization time in seconds exceeds this real value. This is not a strict maximum: the time may exceed it slightly, depending upon the algorithm and on how slow your function evaluation is.

Note that when an AUGLAG or MLSL method is used, the meta-algorithm and the sub-algorithm may have different termination criteria. Thus, for algorithms of this kind, the following named parameters has been defined (just adding the SO prefix - for Sub-Optimizer) to set the ending condition of the sub-algorithm (the meta one uses the ones above): SOStopFuncValue, SOStopRelXTol, and so on... If these are not used, the sub-optimizer will use those of the master routine.

Other named parameters:
• popSize: integer used to change the size of the sample for stochastic search methods. Default value is a peculiar heuristic to the chosen algorithm.
• SOPopSize: Same as above, but when the stochastic search is passed to a meta-algorithm.
• nGradStored: The number (integer type) of gradients to “remember” from previous optimization steps: increasing this increases the memory requirements but may speed convergence. It is set to a heuristic value by default. If used with AUGLAG or MLSL, it will only affect the given subsidiary algorithm.

The following table sums up the main characteristics of each algorithm, providing the more important information about which features are supported by which algorithm and what are the unavoidable arguments they need. More details can be found in NLopt documentation.

Tip: Variational inequality

Let $\Omega$ be a domain of $\mathbb{R}^2$, $f_1, f_2 \in L^2(\Omega)$ and $g_1, g_2 \in L^2(\partial \Omega)$ four given functions with $g_1 \leq g_2$ almost everywhere.

We define the space:

$$V = \{(v_1, v_2) \in H^1(\Omega)^2; v_1|_{\partial \Omega} = g_1, v_2|_{\partial \Omega} = g_2, v_1 \leq v_2 \text{ a.e.}\}$$

as well as the function $J : H^1(\Omega)^2 \rightarrow \mathbb{R}$:

$$J(v_1, v_2) = \frac{1}{2} \int_{\Omega} |\nabla v_1|^2 - \int_{\Omega} f_1 v_1 + \frac{1}{2} \int_{\Omega} |\nabla v_2|^2 - \int_{\Omega} f_2 v_2$$

(3.29)

The problem consists in finding (numerically) two functions $(u_1, u_2) = \arg\min_{(v_1, v_2) \in V} J(v_1, v_2)$.

This can be interpreted as finding $u_1, u_2$ as close as possible (in a certain sense) to the solutions of the Laplace equation with respectively $f_1, f_2$ second members and $g_1, g_2$ Dirichlet boundary conditions with the $u_1 \leq u_2$ almost everywhere constraint.

Here is the corresponding script to treat this variational inequality problem with one of the NLOpt algorithms.

```plaintext
1 // A brief script to demonstrate how to use the freefemm interfaced nlopt routines
2 // The problem consist in solving a simple variational inequality using one of the
3 // optimization algorithm of nlopt. We restart the algorithm a few times after
4 // performing some mesh adaptation to get a more precise output
5
6 load "ff-NLopt"
7
8 // Parameters
9 int kas = 3; // choose of the algorithm
```
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<th>Stochastic</th>
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Legend: ✓ Supported and optional
✓ Should be supported and optional, may lead to weird behaviour though.
✓/✓ Intrinsic characteristic of the algorithm which then need one or more unavoidable parameter to work (for stochastic algorithm, the population size always have a default value, they will then work if it is omitted)
✓ For routines with subsidiary algorithms only, indicates that the corresponding feature will depend on the chosen sub-optimizer.
```
int NN = 10;
func f1 = 1.;
func f2 = -1.;
func g1 = 0.1;
func g2 = 0.1;
int iter = 0;
int nadapt = 2;
real starttol = 1e-6;
real bctol = 6.e-12;

// Mesh
mesh Th = square(NN, NN);

// Fespace
fespace Vh(Th, P1);
Vh oldu1, oldu2;

// Adaptation loop
for (int al = 0; al < nadapt; ++al) {
    varf BVF (v, w) = int2d(Th)(0.5*DX(v)*DX(w) + 0.5*DY(v)*DY(w));
    varf LVF1 (v, w) = int2d(Th)(f1*w);
    varf LVF2 (v, w) = int2d(Th)(f2*w);
    matrix A = BVF(Vh, Vh);
    real[int] b1 = LVF1(0, Vh), b2 = LVF2(0, Vh);
    varf Vbord (v, w) = on(1, 2, 3, 4, v=1);
    Vh In, Bord;
    Bord[] = Vbord(0, Vh, tgv=1);
    In[] = Bord[] ? 0:1;
    Vh gh1 = Bord*g1, gh2 = Bord*g2;

    func real J (real[int] &X) {
        Vh u1, u2;
        u1[] = X(0:Vh.ndof-1);
        u2[] = X(Vh.ndof:2*Vh.ndof-1);
        iter++;
        real[int] Au1 = A*u1[], Au2 = A*u2[];
        Au1 -= b1;
        Au2 -= b2;
        real val = u1[]'*Au1 + u2[]'*Au2;
        if (iter%10 == 9)
            plot(u1, u2, nbiso=30, fill=1, dim=3, cmm="adapt level "+al+" - iteration "+"+iter+" - J = "+val, value=1);
        return val;
    }

    varf dBFV (v, w) = int2d(Th)(dx(v)*dx(w)+dy(v)*dy(w));
    matrix dA = dBFV(Vh, Vh);
    func real[int] dJ (real[int] &X) {
        Vh u1, u2;
        u1[] = X(0:Vh.ndof-1);
        u2[] = X(Vh.ndof:2*Vh.ndof-1);
        real[int] grad1 = dA*u1[], grad2 = dA*u2[];
        grad1 -= b1;
        grad2 -= b2;
    }
```

(continues on next page)
real[int] Grad(X.n);
Grad(0:Vh.ndof-1) = grad1;
Grad(Vh.ndof:2*Vh.ndof-1) = grad2;
return Grad;
}

func real[int] IneqC (real[int] &X){
real[int] constraints(Vh.ndof);
for (int i = 0; i < Vh.ndof; ++i) constraints[i] = X[i] - X[i+Vh.ndof];
return constraints;
}

func real[int, int] dIneqC (real[int] &X){
real[int, int] dconst(Vh.ndof, 2*Vh.ndof);
dconst = 0;
for (int i = 0; i < Vh.ndof; ++i){
dconst(i, i) = 1.;
dconst(i, i+Vh.ndof) = -1.;
}
return dconst;
}

real[int] BordIndex(Th.nbe); //Indexes of border d.f.
{
    int k = 0;
    for (int i = 0; i < Bord.n; ++i) if (Bord[i][i]) { BordIndex[k] = i; ++k; }
}

func real[int] BC (real[int] &X){
real[int] bc(2*Th.nbe);
for (int i = 0; i < Th.nbe; ++i){
    int I = BordIndex[i];
    bc[i] = X[I] - gh1[I];
    bc[i+Th.nbe] = X[I+Th.nv] - gh2[I];
}
return bc;
}

func real[int, int] dBC(real[int] &X){
real[int, int] dbc(2*Th.nbe,2*Th.nv);
dbc = 0.;
for (int i = 0; i < Th.nbe; ++i){
    int I = BordIndex[i];
    dbc(i, I) = 1.;
    dbc(i+Th.nbe, I+Th.nv) = 1.;
}
return dbc;
}

real[int] start(2*Vh.ndof), up(2*Vh.ndof), lo(2*Vh.ndof);
if (al == 0){
    start(0:Vh.ndof-1) = 0.;
    start(Vh.ndof:2*Vh.ndof-1) = 0.01;
} else{
    start(0:Vh.ndof-1) = oldu1[];
}
3.5.7 Optimization with MPI

The only quick way to use the previously presented algorithms on a parallel architecture lies in parallelizing the used cost function (which is in most real life cases, the expensive part of the algorithm). Somehow, we provide a parallel version of the CMA-ES algorithm. The parallelization principle is the trivial one of evolving/genetic algorithms: at each iteration the cost function has to be evaluated \( N \) times without any dependence at all; these \( N \) calculus are then equally distributed to each process. Calling the MPI version of CMA-ES is nearly the same as calling its sequential version (a complete example of use can be found in the CMAES MPI variational inequality example):

```
load "mpi-cmaes"
... // Define J, u and all here
real min = cmaesMPI(J, u, stopTolFun=1e-6, stopMaxIter=3000);
cout << "minimum value is " << min << " for u = " << u << " end1;
```

If the population size is not changed using the popsize parameter, it will use the heuristic value slightly changed
to be equal to the closest greatest multiple of the size of the communicator used by the optimizer. The FreeFEM mpicommworld is used by default. The user can specify his own MPI communicator with the named parameter comm=, see the MPI section of this manual for more information about communicators in FreeFEM.

3.6 Parallelization

A first attempt of parallelization of FreeFEM is made here with MPI. An extended interface with MPI has been added to FreeFEM version 3.5, (see the MPI documentation for the functionality of the language).

3.6.1 MPI

MPI Keywords

The following keywords and concepts are used:

- mpiComm to defined a communication world
- mpiGroup to defined a group of processors in the communication world
- mpiRequest to defined a request to wait for the end of the communication

MPI Constants

- mpisize The total number of processes,
- mpirank the id-number of my current process in \{0, \ldots, mpisize-1\},
- mpiUndefined The MPI_Undefined constant,
- mpiAnySource The MPI_ANY_SOURCE constant,
- mpiCommWorld The MPI_COMM_WORLD constant,
- [... ] and all the keywords of MPI_Op for the reduce operator: mpiMAX, mpiMIN, mpiSUM, mpiPROD, mpiLAND, mpiLOR, mpiLXOR, mpiBAND, mpiBXOR.

MPI Constructor

```cpp
// Parameters
int[int] proc1 = [1, 2], proc2 = [0, 3];
int color = 1;
int key = 1;

// MPI ranks
cout << "MPI rank = " << mpirank << endl;

// MPI
mpiComm comm(mpiCommWorld, 0, 0); //set a MPI_Comm to MPI_COMM_WORLD
mpiGroup grp(proc1); //set MPI_Group to proc 1,2 in MPI_COMM_WORLD
mpiGroup grp1(comm, proc1); //set MPI_Group to proc 1,2 in comm
mpiComm ncomm1(mpiCommWorld, grp); //set the MPI_Comm form grp
```

(continues on next page)
mpiComm ncomm2(comm, color, key); //MPI_Comm_split(MPI_Comm comm, int color, int key,(MPI_Comm *ncomm)

mpiRequest rq; //defined an MPI_Request
mpiRequest[int] arq(10); //defined an array of 10 MPI_Request

MPI Functions

mpiComm Comm(mpiCommWorld, 0, 0);

int MPICommSize = mpiSize(Comm);
int MPIRank = mpiRank(Comm);
if (MPIRank == 0) cout << "MPI Comm size = " << MPICommSize << endl;
cout << "MPI rank in Comm = " << mpiRank(Comm) << endl;

mpiRequest Req;
mpiRequest[int] ReqArray(10);
for (int i = 0; i < MPICommSize; i++){
    //return processor i with no Resquest in MPI_COMM_WORLD
    processor(i);
    //return processor any source with no Resquest in MPI_COMM_WORLD
    processor(mpiAnySource);
    //return processor i with no Resquest in Comm
    processor(i, Comm);
    //return processor i with Resquest rq in Comm
    processor(i, Req, Comm);
    //return processor i with Resquest rq in MPI_COMM_WORLD
    processor(i, Req);
    //return processor i in MPI_COMM_WORLD in block mode for synchronously
    processorblock(i);
    //return processor any source in MPI_COMM_WORLD in block mode for synchronously
    processorblock(mpiAnySource);
    //return processor i in in Comm in block mode
    processorblock(i, Comm);
}

mpiBarrier(Comm); //do a MPI_BARRIER on communicator Comm
mpiWaitAny(ReqArray); //wait add of Request array,
mpiWait(Req); //wait on a Request
real t = mpiWtime(); //return MPI_Wtime in second
real tick = mpiWtick(); //return MPI_Wtick in second

where a processor is just a integer rank, pointer to a MPI_comm and pointer to a MPI_Request, and processorblock with a special MPI_Request.
MPI Communicator operator

```cpp
int status; //to get the MPI status of send / recv
real a, b;

mpiComm comm(mpiCommWorld, 0, 0);
mpiRequest req;

//send a,b asynchronously to the process 1
processor(1) << a << b;
//receive a,b synchronously from the process 10
processor(10) >> a >> b;

//broadcast from processor of comm to other comm processor
// broadcast(processor(10, comm), a);
//send synchronously to the process 10 the data a
status = Send(processor(10, comm), a);
//receive synchronously from the process 10 the data a
status = Recv(processor(10, comm), a);

//send asynchronously to the process 10 the data a without request
status = Isend(processor(10, comm), a);
//send asynchronously to the process 10 the data a with request
status = Isend(processor(10, comm, req), a);
//receive asynchronously from the process 10 the data a
status = Irecv(processor(10, req), a);
//Error asynchronously without request.
// status = Irecv(processor(10), a);

where the data type of a can be of type of int, real, complex, int[int], real[int], complex[int],
int[int,int], double[int,int], complex[int,int], mesh, mesh3, mesh[int], mesh3[int],
matrix, matrix<complex>

//send asynchronously to the process 10 the data a with request
processor(10, req) << a;
//receive asynchronously from the process 10 the data a with request
processor(10, req) >> a;
```

If a, b are arrays or full matrices of int, real, or complex, we can use the following MPI functions:

```cpp
mpiAlltoall(a, b, [comm]);
mpiAllgather(a, b, [comm]);
mpiGather(a, b, processor(..));
mpiScatter(a, b, processor(..));
mpiReduce(a, b, processor(..), mpiMAX);
mpiAllReduce(a, b, comm, mpiMAX);
```

Thank you to Guy-Antoine Atenekeng Kahou for his help to code this interface.

Schwarz example in parallel

This example is a rewriting of example Schwarz overlapping.

```bash
ff-mpirun -np 2 SchwarzParallel.edp
# OR
mpirun -np 2 FreeFem++-mpi SchwarzParallel.edp
```

3.6. Parallelization
if (mpisize != 2) {
  cout << " sorry, number of processors !=2 " << endl;
  exit(1);
}

// Parameters
verbosity = 0;
int interior = 2;
int exterior = 1;
n = 4;

// Mesh
border a(t=1, 2) { x=t; y=0; label=exterior; }
border b(t=0, 1) { x=2; y=t; label=exterior; }
border c(t=2, 0) { x=t; y=1; label=exterior; }
border d(t=1, 0) { x=1-t; y=t; label=interior; }
border e(t=0, pi/2) { x=cos(t); y=sin(t); label=interior; }
border e1(t=pi/2, 2*pi) { x=cos(t); y=sin(t); label=exterior; }

mesh[int] Th[mpisize];
if (mpirank == 0)
  Th[0] = buildmesh(a(5*n) + b(5*n) + c(10*n) + d(5*n));
else
  Th[1] = buildmesh(e(5*n) + e1(25*n));

broadcast(processor(0), Th[0]);
broadcast(processor(1), Th[1]);

// Fespace
fespace Vh(Th[mpirank], P1);
Vh u = 0, v;

fespace Vhother(Th[1-mpirank], P1);
Vhother U = 0;

//Problem
int i = 0;
problem pb { u, v, init=i, solver=Cholesky }
  = int2d(Th[mpirank]){
    dx(u) * dx(v)
    + dy(u) * dy(v)
  } - int2d(Th[mpirank]){
    v
  }
  + on(interior, u=U)
  + on(exterior, u= 0 };

// Loop
for (i = 0; i < 20; i++){
  cout << " - Loop " << i << endl;
 解决问题;
  //send u to the other proc, receive in U
  processor(1-mpirank) << u[];
  processor(1-mpirank) >> U[];
}
Todo: script freeze in the loop

True parallel Schwarz example

Thank you to F. Nataf

This is a explanation of the two examples MPI-GMRES 2D and MPI-GMRES 3D, a Schwarz parallel with a complexity almost independent of the number of process (with a coarse grid preconditioner).

To solve the following Poisson problem on domain $\Omega$ with boundary $\Gamma$ in $L^2(\Omega)$:

\[-\Delta u = f \text{ in } \Omega\]
\[u = g \text{ on } \Gamma\]

where $f$ and $g$ are two given functions of $L^2(\Omega)$ and of $H^\frac{1}{2}(\Gamma)$.

Let's introduce $(\pi_i)_{i=1,...,N_p}$ a regular partition of the unity of $\Omega$, q-e-d:

\[\pi_i \in C^0(\Omega): \quad \pi_i \geq 0 \text{ and } \sum_{i=1}^{N_p} \pi_i = 1.\]

Denote $\Omega_i$ the sub domain which is the support of $\pi_i$ function and also denote $\Gamma_i$ the boundary of $\Omega_i$.

The parallel Schwarz method is:

Let $\ell = 0$ the iterator and a initial guest $u^0$ respecting the boundary condition (i.e. $u^0|_{\Gamma_i} = g$).

\[\forall i = 1,..,N_p: \]
\[-\Delta u^\ell_i = f \text{ in } \Omega_i\]
\[u^\ell_i = u^\ell \text{ on } \Gamma_i \setminus \Gamma\]
\[u^\ell_i = g \text{ on } \Gamma_i \cap \Gamma\]

\[u^{\ell+1} = \sum_{i=1}^{N_p} \pi_i u^\ell_i\] \hspace{1cm} (3.30)

After discretization with the Lagrange finite element method, with a compatible mesh $T_{h_i}$ of $\Omega_i$, i.e., the exist a global mesh $T_h$ such that $T_{h_i}$ is include in $T_h$.

Let us denote:

- $V_{h_i}$ the finite element space corresponding to domain $\Omega_i$, ...

3.6. Parallelization
\[ \mathcal{N}_{h_i} \] is the set of the degree of freedom \( \sigma_i^k \),
\[ \mathcal{N}^{T_i}_{h_i} \] is the set of the degree of freedom of \( V_{h_i} \) on the boundary \( \Gamma_i \) of \( \Omega_i \),
\( \sigma_i^k(v_h) \) is the value the degree of freedom \( k \),
\[ V_{h_i} = \{ v_h \in V_{h_i} : \forall k \in \mathcal{N}^{T_i}_{h_i}, \quad \sigma_i^k(v_h) = 0 \}. \]

The conditional expression \( a ? b : c \) is defined like in \texttt{C} of \texttt{C++} language by
\[
\begin{align*}
    a ? b : c & \equiv \\
    & \begin{cases} 
        \text{if } a \text{ is true then return } b \\
        \text{else return } c
    \end{cases}
\end{align*}
\]

Note: We never use finite element space associated to the full domain \( \Omega \) because it is too expensive.

We have to defined to operator to build the previous algorithm:

We denote \( u_{h|\Gamma}^\ell \) the restriction of \( u_h^\ell \) on \( V_{h|\Gamma} \), so the discrete problem on \( \Omega_i \) of problem (3.30) is find \( u_{h|\Gamma}^\ell \in V_{h|\Gamma} \) such that:
\[
\forall v_{h|\Gamma} \in V_{h|\Gamma}, \quad \int_{\Omega_i} \nabla v_{h|\Gamma} \cdot \nabla u_{h|\Gamma}^\ell = \int_{\Omega_i} f v_{h|\Gamma}, \quad \forall k \in \mathcal{N}_{h_i}^{T_i} : \quad \sigma_i^k(u_{h|\Gamma}^\ell) = \begin{cases} 
    g_i^k & (k \in \Gamma) \\
    0 & (k \notin \Gamma)
\end{cases}
\]

where \( g_i^k \) is the value of \( g \) associated to the degree of freedom \( k \in \mathcal{N}_{h_i}^{T_i} \).

In FreeFEM, it can be written has with \( U \) is the vector corresponding to \( u_{h|\Gamma}^\ell \), and the vector \( U1 \) is the vector corresponding to \( u_{h|\Gamma}^\ell \) is the solution of:

```plaintext
1  real[int] U1(Ui.n);
2  real[int] b = onG .* U;
3  b = onG ? b : Bi ;
4  U1 = Ai^-1*b;
```

where \( onG[i] = (i \in \Gamma_i \setminus \Gamma)\)?1 : 0, and \( Bi \) the right of side of the problem, are defined by

```plaintext
1  // Fespace
2  fespace Whi(Thi, P2);
3  // Problem
4  varf vPb (U, V) = int3d(Thi) (grad(U)'*grad(V))
5    + int3d(Thi) (F*V + on(1, U=g) + on(10, U=G));
6  varf vPbon (U, V) = on(10, U=1) + on(1, U=0);
7  matrix Ai = vPb (Whi, Whi, solver=sparsesolver);
8  real[int] onG = vPbon(0, Whi);
9  real[int] Bi=vPb(0, Whi);
```

where the FreeFEM label of \( \Gamma \) is 1 and the label of \( \Gamma_i \setminus \Gamma \) is 10.

---

**Note:** We never use finite element space associated to the full domain \( \Omega \) because it is too expensive.

We have to defined to operator to build the previous algorithm:

We denote \( u_{h|\Gamma}^\ell \) the restriction of \( u_h^\ell \) on \( V_{h|\Gamma} \), so the discrete problem on \( \Omega_i \) of problem (3.30) is find \( u_{h|\Gamma}^\ell \in V_{h|\Gamma} \) such that:

\[
\forall v_{h|\Gamma} \in V_{h|\Gamma}, \quad \int_{\Omega_i} \nabla v_{h|\Gamma} \cdot \nabla u_{h|\Gamma}^\ell = \int_{\Omega_i} f v_{h|\Gamma}, \quad \forall k \in \mathcal{N}_{h_i}^{T_i} : \quad \sigma_i^k(u_{h|\Gamma}^\ell) = \begin{cases} 
    g_i^k & (k \in \Gamma) \\
    0 & (k \notin \Gamma)
\end{cases}
\]

where \( g_i^k \) is the value of \( g \) associated to the degree of freedom \( k \in \mathcal{N}_{h_i}^{T_i} \).

In FreeFEM, it can be written has with \( U \) is the vector corresponding to \( u_{h|\Gamma}^\ell \), and the vector \( U1 \) is the vector corresponding to \( u_{h|\Gamma}^\ell \) is the solution of:

```plaintext
1  real[int] U1(Ui.n);
2  real[int] b = onG .* U;
3  b = onG ? b : Bi ;
4  U1 = Ai^-1*b;
```

where \( onG[i] = (i \in \Gamma_i \setminus \Gamma)\)?1 : 0, and \( Bi \) the right of side of the problem, are defined by

```plaintext
1  // Fespace
2  fespace Whi(Thi, P2);
3  // Problem
4  varf vPb (U, V) = int3d(Thi) (grad(U)'*grad(V))
5    + int3d(Thi) (F*V + on(1, U=g) + on(10, U=G));
6  varf vPbon (U, V) = on(10, U=1) + on(1, U=0);
7  matrix Ai = vPb (Whi, Whi, solver=sparsesolver);
8  real[int] onG = vPbon(0, Whi);
9  real[int] Bi=vPb(0, Whi);
```

where the FreeFEM label of \( \Gamma \) is 1 and the label of \( \Gamma_i \setminus \Gamma \) is 10.
To build the transfer/update part corresponding to (3.30) equation on process \( i \), let us call \( n_{\text{part}} \) the number the neighborhood of domain of \( \Omega_i \) (i.e. \( \pi_j \) is none 0 of \( \Omega_i \)), we store in an array \( j_{\text{part}} \) of size \( n_{\text{part}} \) all this neighborhood.

Let us introduce two array of matrix, \( S_{mj[j]} \) to defined the vector to send from \( i \) to \( j \) a neighborhood process, and the matrix \( r_{Mj[j]} \) to after to reduce owith neighborhood \( j \) domain.

So the tranfert and update part compute \( u_i = \pi_i u_i + \sum_{j \in J_i} \pi_j u_j \) and can be write the FreeFEM function Update:

```cpp
func bool Update (real[int] &ui, real[int] &vi){
    int n = jpart.n;
    for (int j = 0; j < njpart; ++j) Usend[j][j] = Smj[j]*ui;
    mpiRequest[int] rq(n*2);
    for (int j = 0; j < n; ++j) Irecv(processor(jpart[j], comm, rq[j]), Ri[j][j]);
    for (int j = 0; j < n*2; ++j) Srecv[j][j] = Smj[j]*ui;
    // apply the unity local partition
    vi = Pii*ui; //set to pi_i u_i
    for (int j = 0; j < njpart; ++j) vi += rMj[j]*Vrecv[j][j]; //add pi_j u_j
    return true;
}
```

where the buffer are defined by:

```cpp
InitU(njpart, Whij, Thij, aThij, Usend) //defined the send buffer
InitU(njpart, Whij, Thij, aThij, Vrecv) //defined the revc buffer
```

with the following macro definition:

```cpp
macro InitU(n, Vh, Th, aTh, U) Vh[int] U(n); for (int j = 0; j < n; ++j){Th = aTh[j]; U[j] = 0;}
```

**First GMRES algorithm:** you can easily accelerate the fixed point algorithm by using a parallel GMRES algorithm after the introduction the following affine \( A_i \) operator sub domain \( \Omega_i \).

```cpp
func real[int] DJ0 (real[int] & U){
    real[int] V(U.n), b = onG .* U;
    b = onG ? b : Bi ;
    V = A^-1* b;
    Update(V, U);
    V -= U;
    return V;
}
```

Where the parallel MPIGMRES or MPICG algorithm is just a simple way to solve in parallel the following \( A_i x_i = b_i, i = 1, \ldots, N_p \) by just changing the dot product by reduce the local dot product of all process with the following MPI code:

```cpp
template<class R> R ReduceSum1(R s, MPI_Comm *comm){
    R r = 0;
    MPI_Allreduce(&s, &r, 1, MPI_TYPE<R>::TYPE(), MPI_SUM, *comm );
    return r;
}
```

This is done in MPIGC dynamics library tool.

**Second GMRES algorithm:** Use scharwz algorithm as a preconditioner of basic GMRES method to solving the parallel problem.

### 3.6. Parallelization
Third GMRES algorithm: Add a coarse solver to the previous algorithm

First build a coarse grid on processor 0, and the

```cpp
matrix AC, Rci, Pci;
if (mpiRank(comm) == 0)
    AC = vPbC(VhC, VhC, solver=sparsesolver); //the coarse problem
Pci = interpolate(Whi, VhC); //the projection on coarse grid
Rci = Pci'*Pii; //the restriction on Process i grid with the partition pi_i
func bool CoarseSolve (real[int]& V, real[int]& U, mpiComm& comm)
    // solving the coarse problem
    real[int] Uc(Rci.n), Bc(Uc.n);
    Uc = Rci*U;
    mpiReduce(Uc, Bc, processor(0, comm), mpiSUM);
    if (mpiRank(comm) == 0)
        Uc = AC^-1*Bc;
    broadcast(processor(0, comm), Uc);
    V = Pci*Uc;
}
```

The New preconditioner

```cpp
func real[int] PDJC (real[int]& U){
    // Idea: F. Nataf.
    // 0 ~ (I C1A)(I-C2A) => I ~ - C1AC2A +C1A +C2A
    // New Prec P= C1+C2 - C1AC2 = C1(I- A C2) +C2
    // ( C1(I- A C2) +C2 ) Uo
    // V = - C2*Uo
    // ....
    real[int] V(U.n);
    CoarseSolve(V, U, comm);
    V = -V; // -C2*Uo
    U += Ai*V; //U = (I-A C2) Uo
    real[int] b = onG ? 0. : U;
    Update(V, U);
    return U;
}
```
The code of the 4 algorithms:

```plaintext
real epss = 1e-6;
int rgmres = 0;
if (gmres == 1){
    rgmres = MPIAffineGMRES(DJ0, u[], veps=epss, nbiter=300,
                           comm=comm, dimKrylov=100, verbosity=ipart?0: 50);
    real[int] b = onG .* u[];
    b = onG ? b : Bi ;
    v[] = Ai^-1*b;
    Update(v[], u[]);
}
else if (gmres == 2)
    rgmres = MPILinearGMRES(DJ, precon=PDJ, u[], Bi,
                           veps=epss, nbiter=300, comm=comm, dimKrylov=100,
                           verbosity=ipart?0: 50);
else if (gmres == 3)
    rgmres = MPILinearGMRES(DJ, precon=PDJC, u[], Bi, veps=epss,
                           nbiter=300, comm=comm, dimKrylov=100, verbosity=ipart?0: 50);
else //algo Shwarz for demo
    for (int iter = 0; iter < 10; ++iter)
    ...
}
```

We have all ingredient to solve in parallel if we have et the partitions of the unity. To build this partition we do:

The initial step on process 1 to build a coarse mesh, $T_h^*$ of the full domain, and build the partition $\pi$ function constant equal to $i$ on each sub domain $\mathcal{O}_i$, $i = 1, \ldots, N_p$, of the grid with the metis graph partitioner [KARYPIS1995] and on each process $i$ in $1, \ldots, N_p$ do:

1. Broadcast from process 1, the mesh $T_h^*$ (call Thii in FreeFEM script), and $\pi$ function,
2. remark that the characteristic function $\mathbf{1}_{\mathcal{O}_i}$ of domain $\mathcal{O}_i$, is defined by $(\pi = i)?1 : 0$,
3. Let us call $\Pi^2_{V_h}$ (resp. $\Pi^2_{V_h}$) the $L^2$ on $P_h^*$ the space of the constant finite element function per element on $T_h^*$ (resp. $V_h^*$ the space of the affine continuous finite element per element on $T_h^*$) and build in parallel the $\pi_i$ and $\Omega_i$ such that $\mathcal{O}_i \subset \Omega_i$ where $\mathcal{O}_i = \text{supp}(\Pi^2_{V_h}^* \mathbf{1}_{\mathcal{O}_i})$, and $m$ is the overlaps size on the coarse mesh (generally one), (this is done in function AddLayers(Thi, suppii[], nlayer, phi[]); We choose a function $\pi^*_i = (\Pi^2_{V_h}^* \mathbf{1}_{\mathcal{O}_i})$ so the partition of the unity is simply defined by

$$
\pi_i = \frac{\pi^*_i}{\sum_{j=1}^{N_p} \pi^*_j}
$$

The set $J_i$ of neighborhood of the domain $\mathcal{O}_i$, and the local version on $V_{hi}$ can be defined the array jpart and njpart with:

```plaintext
Vhi[pii] = piistar;
Vhi[int] pij(npij); //local partition of 1 = pii + sum_j pij[j]
int[int] jpart(npart);
int njpart = 0;
Vhi sumphi = piistar;
for (int i = 0; i < npart; ++i)
    if (i != ipart){
        if (int3d(Thi)(pijstar,j) > 0){
            pij[njpart] = pijstar;
            sumphi[] += pij[njpart][];
            jpart[njpart++] = i;
        }
    }
}
```

(continues on next page)
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(continued from previous page)

14  pii[] = pii[] ./ sumphi[];
15  for (int j = 0; j < njpart; ++j)
16  pij[j][] = pij[j][] ./ sumphi[];
17  jpart.resize(njpart);

4. We call \( T_{hi}^* \) the sub mesh part of \( T_{hi} \) where \( \pi_j \) are none zero. And thanks to the function `trunc` to build this array,

1  for(int jp = 0; jp < njpart; ++jp)
2  aThij[jp] = trunc(Thi, pij[jp] > 1e-10, \text{label}=10);

5. At this step we have all on the coarse mesh, so we can build the fine final mesh by splitting all meshes: \( \text{Thi}, \text{Thij[j]}, \text{Thij[j]} \) with FreeFEM `trunc` mesh function which do restriction and slipping.

6. The construction of the send/recv matrices \( sMj \) and `freefem`:`rMj`: can done with this code:

\begin{verbatim}
1  mesh3 Thij = Thi;
2  fespace Whij(Thij, Pk);
3  matrix Pii; Whi wpii = pii; Pii = wpii[]; //Diagonal matrix corresponding \( \pi_i \)
4  matrix[int] sMj(njpart), rMj(njpart); //M send/recive case
5  for (int jp = 0; jp < njpart; ++jp){
6    int j = jpart[jp];
7    Thij = aThij[jp]; //change mesh to change Whij, Whij
8    matrix I = interpolate(Whij, Whi); //Whij <- Whi
9    sMj[jp] = I*Pii; //Whi -> s Whij
10   rMj[jp] = interpolate(Whij, Whi, \text{t}=1); //Whij -> Whi
11 }
\end{verbatim}

To build a not too bad application, all variables come from parameters value with the following code

\begin{verbatim}
1  include "getARGV.idp"
2  verbosity = getARGV("-vv", 0);
3  int vdebug = getARGV("-d", 1);
4  int ksplit = getARGV("-k", 10);
5  int nloc = getARGV("-n", 25);
6  string sff = getARGV("-p", ");
7  int gmres = getARGV("-gmres", 3);
8  bool dplot = getARGV("-dp", 0);
9  int nC = getARGV("-N", \text{max}(nloc/10, 4));
\end{verbatim}

And small include to make graphic in parallel of distribute solution of vector \( u \) on mesh \( T_h \) with the following interface:

\begin{verbatim}
1  include "MPIplot.idp"
2  func bool plotMPIall(mesh &Th, real[int] &u, string cm) {
3    PLOTMPIALL(mesh, Pk, Th, u, \{cmm=cm, nbiso=20, fill=1, dim=3, value=1\});
4    return 1;
5  }
\end{verbatim}

Note: The \text{cmm}=cm, \ldots in the macro argument is a way to quote macro argument so the argument is \text{cmm}=cm, \ldots.
3.6.2 Parallel sparse solvers

Parallel sparse solvers use several processors to solve linear systems of equation. Like sequential, parallel linear solvers can be direct or iterative. In FreeFEM both are available.

Using parallel sparse solvers in FreeFEM

We recall that the solver parameters are defined in the following commands: solve, problem, set (setting parameter of a matrix) and in the construction of the matrix corresponding to a bilinear form. In these commands, the parameter solver must be set to sparsesolver for parallel sparse solver. We have added specify parameters to these command lines for parallel sparse solvers. These are:

- lparams: vector of integer parameters (l is for the C++ type long)
- dparams: vector of real parameters
- sparms: string parameters
- datafilename: name of the file which contains solver parameters

The following four parameters are only for direct solvers and are vectors. These parameters allow the user to preprocess the matrix (see the section on sparse direct solver for more information).

- permr: row permutation (integer vector)
- permc: column permutation or inverse row permutation (integer vector)
- scaler: row scaling (real vector)
- scalec: column scaling (real vector)

There are two possibilities to control solver parameters. The first method defines parameters with lparams, dparams and sparms in .edp file.

The second one reads the solver parameters from a data file. The name of this file is specified by datafilename. If lparams, dparams, sparms or datafilename is not provided by the user, the solver’s default values are used.

To use parallel solver in FreeFEM, we need to load the dynamic library corresponding to this solver. For example to use MUMPS solver as parallel solver in FreeFEM, write in the .edp file load "MUMPS_FreeFem".

If the libraries are not loaded, the default sparse solver will be loaded (default sparse solver is UMFPACK). The Table 3.2 gives this new value for the different libraries.

<table>
<thead>
<tr>
<th>Libraries</th>
<th>Default sparse solver</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>real</td>
</tr>
<tr>
<td>MUMPS_FreeFem</td>
<td>mumps</td>
</tr>
<tr>
<td>real_SuperLU_DIST_FreeFem</td>
<td>SuperLU_DIST</td>
</tr>
<tr>
<td>complex_SuperLU_DIST_FreeFem</td>
<td>previous solver</td>
</tr>
<tr>
<td>real_pastix_FreeFem</td>
<td>PaStiX</td>
</tr>
<tr>
<td>complex_pastix_FreeFem</td>
<td>previous solver</td>
</tr>
<tr>
<td>hips_FreeFem</td>
<td>hips</td>
</tr>
<tr>
<td>hypre_FreeFem</td>
<td>hypre</td>
</tr>
<tr>
<td>parms_FreeFem</td>
<td>parms</td>
</tr>
</tbody>
</table>

We also add functions (see Table 3.3) with no parameter to change the default sparse solver in the .edp file. To use these functions, we need to load the library corresponding to the solver. An example of using different parallel sparse solvers for the same problem is given in Direct solvers example.
Table 3.3: Functions that allow to change the default sparse solver for real and complex arithmetics and the result of these functions

<table>
<thead>
<tr>
<th>Function</th>
<th>default sparse solver</th>
</tr>
</thead>
<tbody>
<tr>
<td>defaulttoMUMPS()</td>
<td>defaulttoMUMPS()</td>
</tr>
<tr>
<td>realdefaulttoSuperLUDist()</td>
<td>SuperLU_DIST</td>
</tr>
<tr>
<td>complexdefaulttoSuperLUDist()</td>
<td>SuperLU_DIST</td>
</tr>
<tr>
<td>realdefaulttopastix()</td>
<td>pastix</td>
</tr>
<tr>
<td>complexdefaulttopastix()</td>
<td>pastix</td>
</tr>
<tr>
<td>defaulttohips()</td>
<td>hips</td>
</tr>
<tr>
<td>defaulttohypre()</td>
<td>hypre</td>
</tr>
<tr>
<td>defaulttoparms()</td>
<td>arms</td>
</tr>
</tbody>
</table>

Tip: Test direct solvers

```c
load "MUMPS_FreeFem"
//default solver: real-> MUMPS, complex -> MUMPS
load "real_SuperLU_DIST_FreeFem"
//default solver: real-> SuperLU_DIST, complex -> MUMPS
load "real_pastix_FreeFem"
//default solver: real-> pastix, complex -> MUMPS

// Solving with pastix
{
  matrix A =
  [[1, 2, 2, 2, 1, 2],
   [1, 2, 0, 2, 0, 0],
   [1, 2, 0, 0, 0, 0]];
  real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
  b = A*xx;
  cout << "b = " << b << endl; cout << "xx = " << xx << endl;
  set(A, solver=sparse, datafilename="fpastix_iparm_dparm.txt");
  cout << "solve" << endl;
  x = A^(-1)*b;
  cout << "b = " << b << endl;
  cout << "x = " << endl;
  cout << x << endl;
  di = xx - x;
  if (mpirank == 0){
    cout << "x-xx = " << endl;
    cout << "Linfty = " << di.linfty << ", L2 = " << di.l2 << endl;
  }
}

// Solving with SuperLU_DIST
realdefaulttoSuperLUDist();
//default solver: real-> SuperLU_DIST, complex -> MUMPS
{
  matrix A =
  [[1, 2, 2, 1, 1],
   [1, 2, 0, 1, 0, 2],
   [1, 2, 0, 0, 2, 0],
   [1, 2, 0, 0, 0, 0]];
```
[ 2, 12, 0, 10, 10],
[ 2, 0, 1, 0, 2],
[ 1, 10, 0, 22, 0.],
[ 1, 10, 2, 0., 22]);

real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
b = A*xx;
cout << "b = " << b << endl;
cout << "xx = " << xx << endl;

set(A, solver=sparsesolver, datafilename="ffsuperlu_dist_fileparam.txt");
cout << "solve" << endl;
x = A^-1*b;
cout << "b = " << b << endl;
cout << "x = " << endl;
di = xx - x;
if (mpirank == 0){
    cout << "x-xx = " << endl;
    cout << "Linf = " << di.linfty << " , L2 = " << di.l2 << endl;
}
}

// Solving with MUMPS
defaulttoMUMPS();
//default solver: real -> MUMPS, complex -> MUMPS
{
    matrix A =
    [[1, 2, 2, 1, 1],
     [2, 12, 0, 10, 10],
     [2, 0, 1, 0, 2],
     [1, 10, 0, 22, 0.],
     [1, 10, 2, 0., 22]];

real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
b = A*xx;
cout << "b = " << b << endl;
cout << "xx = " << xx << endl;

set(A, solver=sparsesolver, datafilename="ffmumps_fileparam.txt");
cout << "solving solution" << endl;
x = A^-1*b;
cout << "b = " << b << endl;
cout << "x = " << endl;
di = xx - x;
if (mpirank == 0){
    cout << "x-xx = " << endl;
    cout << "Linf = " << di.linfty << " , L2 = " << di.l2 << endl;
}
}

Sparse direct solver

In this section, we present the sparse direct solvers interfaced with FreeFEM.
MUMPS solver

MUltifrontal Massively Parallel Solver (MUMPS) is an open-source library.

This package solves linear system of the form \( A x = b \) where \( A \) is a square sparse matrix with a direct method. The square matrix considered in MUMPS can be either unsymmetric, symmetric positive definite or general symmetric.

The method implemented in MUMPS is a direct method based on a multifrontal approach. It constructs a direct factorization \( A = LU \) or \( A = L^t D L \) depending of the symmetry of the matrix \( A \).

MUMPS uses the following libraries:

- BLAS,
- BLACS,
- ScaLAPACK.

**Warning:** MUMPS does not solve linear system with a rectangular matrix.

MUMPS parameters:

There are four input parameters in MUMPS. Two integers sym and par, a vector of integer of size 40 inctl and a vector of real of size 15 cntl.

The first parameter gives the type of the matrix: 0 for unsymmetric matrix, 1 for symmetric positive matrix and 2 for general symmetric.

The second parameter defined if the host processor work during the factorization and solves steps: par=1 host processor working and par=0 host processor not working.

The parameter inctl and cntl is the control parameter of MUMPS. The vectors icntl and cntl in MUMPS becomes with index 1 like vector in Fortran. For more details see the MUMPS user's guide.

We describe now some elements of the main parameters of icntl for MUMPS.

- **Input matrix parameter** The input matrix is controlled by parameters ICNTL(5) and ICNTL(18).
  
  The matrix format (resp. matrix pattern and matrix entries) are controlled by ICNTL(5) (resp. ICNTL(18)).

  The different values of ICNTL(5) are 0 for assembled format and 1 for element format. In the current release of FreeFEM, we consider that FE matrix or matrix is storage in assembled format. Therefore, ICNTL(5) is treated as 0 value.

  The main option for ICNTL(18): INCLTL(18)=0 centrally on the host processor, ICNTL(18)=3 distributed the input matrix pattern and the entries (recommended option for distributed matrix by developer of MUMPS). For other values of ICNTL(18) see the MUMPS user's guide. These values can be used also in FreeFEM.

  The default option implemented in FreeFEM are ICNTL(5)=0 and ICNTL(18)=0.

- **Preprocessing parameter** The preprocessed matrix \( A_p \) that will be effectively factored is defined by

  \[
  A_p = P D_r A Q_c D_c P^t
  \]

  where \( P \) is the permutation matrix, \( Q_c \) is the column permutation, \( D_r \) and \( D_c \) are diagonal matrix for respectively row and column scaling.

  The ordering strategy to obtain \( P \) is controlled by parameter ICNTL(7). The permutation of zero free diagonal \( Q_c \) is controlled by parameter ICNTL(6). The row and column scaling is controlled...
by parameter ICNTL(18). These option are connected and also strongly related with ICNTL(12)
(see the MUMPS user’s guide for more details).

The parameters permr, scaler, and scalec in FreeFEM allow to give permutation matrix($P$),
row scaling ($D_r$) and column scaling ($D_c$) of the user respectively.

Calling MUMPS in FreeFEM

To call MUMPS in FreeFEM, we need to load the dynamic library MUMPS_freefem.dylib (MacOSX),
MUMPS_freefem.so (Unix) or MUMPS_freefem.dll (Windows).

This is done in typing load "MUMPS_FreeFem" in the .edp file. We give now the two methods to give the option
of MUMPS solver in FreeFEM.

- Solver parameters is defined in .edp file: In this method, we need to give the parameters lparams and dparams.
  These parameters are defined for MUMPS by :
  
  - lparams[0] = SYM, lparams[1] = PAR,
  - ∀i = 1,…,40, lparams[i+1] = ICNTL(i)
  - ∀i = 1,…,15, dparams[i-1] = CNTL(i)

- Reading solver parameters on a file:

  The structure of data file for MUMPS in FreeFEM is : first line parameter SYM and second line
  parameter PAR and in the following line the different value of vectors ICNTL and CNTL. An example
  of this parameter file is given in ffmumpsfileparam.txt.

```
0 /* SYM :: 0 for non symmetric matrix, 1 for symmetric definite positive,
   matrix and 2 general symmetric matrix*/
1 /* PAR :: 0 host not working during factorization and solves steps, 1
   host working during factorization and solves steps*/
-1 /* ICNTL(1) :: output stream for error message */
-1 /* ICNTL(2) :: output for diagnostic printing, statics and warning
   message */
-1 /* ICNTL(3) :: for global information */
0 /* ICNTL(4) :: Level of printing for error, warning and diagnostic
   message */
0 /* ICNTL(5) :: matrix format : 0 assembled format, 1 elemental format.
   */
7 /* ICNTL(6) :: control option for permuting and/or scaling the matrix
   in analysis phase */
3 /* ICNTL(7) :: pivot order strategy : AMD, AMF, metis, pord scotch*/
77 /* ICNTL(8) :: Row and Column scaling strategy */
1 /* ICNTL(9) :: 0 solve Ax = b, 1 solve the transposed system A^t x = b
   : parameter is not considered in the current release of FreeFEM*/
0 /* ICNTL(10) :: number of steps of iterative refinement*/
0 /* ICNTL(11) :: statics related to linear system depending on ICNTL(9)
   */
1 /* ICNTL(12) :: constrained ordering strategy for general symmetric
   matrix */
0 /* ICNTL(13) :: method to control splitting of the root frontal matrix
   */
20 /* ICNTL(14) :: method for given : matrix pattern and matrix entries :
   */
0 /* ICNTL(15) :: not used in this release of MUMPS */
0 /* ICNTL(16) :: not used in this release of MUMPS */
0 /* ICNTL(17) :: not used in this release of MUMPS */
3 /* ICNTL(18) :: method for given : matrix pattern and matrix entries :
   */
```
/* ICNTL(19) :: method to return the Schur complement matrix */
/* ICNTL(20) :: right hand side form ( 0 dense form, 1 sparse form) : parameter will be set to 0 for FreeFEM */
/* ICNTL(21) :: 0, 1 kept distributed solution : parameter is not considered in the current release of FreeFEM */
/* ICNTL(22) :: controls the in-core/out-of-core (OOC) facility */
/* ICNTL(23) :: maximum size of the working memory in Megabyte than MUMPS can allocate per working processor */
/* ICNTL(24) :: control the detection of null pivot */
/* ICNTL(25) :: control the computation of a null space basis */
/* ICNTL(26) :: This parameter is only significant with Schur option (ICNTL(19) not zero). : parameter is not considered in the current release of FreeFEM */
/* ICNTL(27) (Experimental parameter subject to change in next release of MUMPS) :: control the blocking factor for multiple righthand side during the solution phase : parameter is not considered in the current release of FreeFEM */
/* ICNTL(28) :: not used in this release of MUMPS*/
/* ICNTL(29) :: not used in this release of MUMPS*/
/* ICNTL(30) :: not used in this release of MUMPS*/
/* ICNTL(31) :: not used in this release of MUMPS*/
/* ICNTL(32) :: not used in this release of MUMPS*/
/* ICNTL(33) :: not used in this release of MUMPS*/
/* ICNTL(34) :: not used in this release of MUMPS*/
/* ICNTL(35) :: not used in this release of MUMPS*/
/* ICNTL(36) :: not used in this release of MUMPS*/
/* ICNTL(37) :: not used in this release of MUMPS*/
/* ICNTL(38) :: not used in this release of MUMPS*/
/* ICNTL(39) :: not used in this release of MUMPS*/
/* ICNTL(40) :: not used in this release of MUMPS*/
0.01 /* CNTL(1) :: relative threshold for numerical pivoting */
1e-8 /* CNTL(2) :: stopping criteria for iterative refinement */
-1 /* CNTL(3) :: threshold for null pivot detection */
-1 /* CNTL(4) :: determine the threshold for partial pivoting */
0.0 /* CNTL(5) :: fixation for null pivots */
/* CNTL(6) :: not used in this release of MUMPS */
/* CNTL(7) :: not used in this release of MUMPS */
/* CNTL(8) :: not used in this release of MUMPS */
/* CNTL(9) :: not used in this release of MUMPS */
/* CNTL(10) :: not used in this release of MUMPS */
/* CNTL(11) :: not used in this release of MUMPS */
/* CNTL(12) :: not used in this release of MUMPS */
/* CNTL(13) :: not used in this release of MUMPS */
/* CNTL(14) :: not used in this release of MUMPS */
/* CNTL(15) :: not used in this release of MUMPS */

If no solver parameter is given, we used default option of MUMPS solver.

**Tip:** MUMPS example

A simple example of calling MUMPS in FreeFEM with this two methods is given in the Test solver MUMPS example.
SuperLU distributed solver

The package SuperLU_DIST solves linear systems using LU factorization. It is a free scientific library.

This library provides functions to handle square or rectangular matrix in real and complex arithmetics. The method implemented in SuperLU_DIST is a supernodal method. New release of this package includes a parallel symbolic factorization. This scientific library is written in C and MPI for communications.

SuperLU_DIST parameters:

We describe now some parameters of SuperLU_DIST. The SuperLU_DIST library use a 2D-logical process group. This process grid is specified by nprow (process row) and npcol (process column) such that \( N_p = nprow \times npcol \) where \( N_p \) is the number of all process allocated for SuperLU_DIST.

The input matrix parameters is controlled by “matrix=” in sparams for internal parameter or in the third line of parameters file. The different value are

- \( \text{matrix=assembled} \) global matrix are available on all process
- \( \text{matrix=distributedglobal} \) The global matrix is distributed among all the process
- \( \text{matrix=distributed} \) The input matrix is distributed (not yet implemented)

The option arguments of SuperLU_DIST are described in the section Users-callable routine of the SuperLU users’ guide.

The parameter Fact and TRANS are specified in FreeFEM interfaces to SuperLU_DIST during the different steps. For this reason, the value given by the user for this option is not considered.

The factorization LU is calculated in SuperLU_DIST on the matrix \( A_p \).

\[
A_p = P_c \ P_r \ D_r \ A \ D_c \ P_r^t
\]

where \( P_c \) and \( P_r \) is the row and column permutation matrix respectively, \( D_r \) and \( D_c \) are diagonal matrix for respectively row and column scaling.

The option argument RowPerm (resp. ColPerm) control the row (resp. column) permutation matrix. \( D_r \) and \( D_c \) is controlled by the parameter DiagScale.

The parameter permr, permc, scaler, and scalec in FreeFEM is provided to give row permutation, column permutation, row scaling and column scaling of the user respectively.

The other parameters for LU factorization are ParSymFact and ReplaceTinyPivot. The parallel symbolic factorization works only on a power of two processes and need the ParMetis ordering. The default option argument of SuperLU_DIST are given in the file ffsuperlu_dist_fileparam.txt.

Calling SuperLU_DIST in FreeFEM

To call SuperLU_DIST in FreeFEM, we need to load the library dynamic correspond to interface. This done by the following line load "real_superlu_DIST_FreeFem" (resp. load "complex_superlu_DIST_FreeFem") for real (resp. complex) arithmetics in the file .edp.

Solver parameters is defined in .edp file:

To call SuperLU_DIST with internal parameter, we used the parameters sparams. The value of parameters of SuperLU_DIST in sparams are defined by :

- nprow=1,
- npcol=1,
- matrix= distributedglobal,
- Fact= DOFACT,
FreeFEM Documentation, Release 4.6

- Equil=NO,
- ParSymbFact=NO,
- ColPerm= MMD_AT_PLUS_A,
- RowPerm= LargeDiag,
- DiagPivotThresh=1.0,
- IterRefine=DOUBLE,
- Trans=NOTRANS,
- ReplaceTinyPivot=NO,
- SolveInitialized=NO,
- PrintStat=NO,
- DiagScale=NOEQUIL

This value correspond to the parameter in the file ffsuperlu_dist_fileparam.txt. If one parameter is not specified by the user, we take the default value of SuperLU_DIST.

Reading solver parameters on a file: The structure of data file for SuperLU_DIST in FreeFEM is given in the file ffsuperlu_dist_fileparam.txt (default value of the FreeFEM interface).

```
1 /* nprow : integer value */
2 /* npcol : integer value */
3 distributedglobal /* matrix input : assembled, distributedglobal, distributed */
4 DOFACT /* Fact : DOFACT, SamePattern, SamePattern_SameRowPerm, FACTORED */
5 NO /* Equil : NO, YES */
6 NO /* ParSymbFact : NO, YES */
7 MMD_AT_PLUS_A /* ColPerm : NATURAL, MMD_AT_PLUS_A, MMD_ATA, METIS_AT_PLUS_A, PARMETIS, 
→ MY_PERMC */
8 LargeDiag /* RowPerm : NOROWPERM, LargeDiag, MY_PERMR */
9 1.0 /* DiagPivotThresh : real value */
10 NOTRANS /* Trans : NOTRANS, TRANS, CONJ */
11 NO /* ReplaceTinyPivot : NO, YES */
12 NO /* SolveInitialized : NO, YES */
13 NO /* RefineInitialized : NO, YES */
14 NO /* PrintStat : NO, YES */
15 NOEQUIL /* DiagScale : NOEQUIL, ROW, COL, BOTH */
```

If no solver parameter is given, we used default option of SuperLU_DIST solver.

Tip: A simple example of calling SuperLU_DIST in FreeFEM with this two methods is given in the Solver superLU_DIST example.

PaStiX solver

PaStiX (Parallel Sparse matrix package) is a free scientific library under CECILL-C license. This package solves sparse linear system with a direct and block ILU(k) iterative methods. His solver can be applied to a real or complex matrix with a symmetric pattern.

PaStiX parameters:
The input matrix parameter of FreeFEM depend on PaStiX interface. matrix = assembled for non distributed matrix. It is the same parameter for SuperLU_DIST.

There are four parameters in PaStiX: iparm, dparm, perm and invp. These parameters are respectively the integer parameters (vector of size 64), real parameters (vector of size 64), permutation matrix and inverse permutation matrix respectively. iparm and dparm vectors are described in PaStiX RefCard.

The parameters permr and permc in FreeFEM are provided to give permutation matrix and inverse permutation matrix of the user respectively.

**Solver parameters defined in .edp file:**

To call PaStiX in FreeFEM in this case, we need to specify the parameters lparams and dparams. These parameters are defined by:

\[ \forall i = 0, \ldots, 63, \text{lparams}[i] = \text{iparm}[i]. \]
\[ \forall i = 0, \ldots, 63, \text{dparams}[i] = \text{dparm}[i]. \]

**Reading solver parameters on a file:**

The structure of data file for PaStiX parameters in FreeFEM is: first line structure parameters of the matrix and in the following line the value of vectors iparm and dparm in this order.

```plaintext
assembled /* matrix input :: assembled, distributed global and distributed */
iparm[0]
iparm[1]
...
iparm[63]
dparm[0]
dparm[1]
...
dparm[63]
```

An example of this file parameter is given in ffpastix_iparm_dparm.txt with a description of these parameters. This file is obtained with the example file iparm.txt and dparm.txt including in the PaStiX package.

If no solver parameter is given, we use the default option of PaStiX solver.

**Tip:** A simple example of calling PaStiX in FreeFEM with this two methods is given in the Solver PaStiX example.

In Table 3.4, we recall the different matrix considering in the different direct solvers.

<table>
<thead>
<tr>
<th>direct solver</th>
<th>square matrix</th>
<th>rectangular matrix</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>sym</td>
<td>sym pattern</td>
</tr>
<tr>
<td>SuperLU_DIST</td>
<td>yes</td>
<td>yes</td>
</tr>
<tr>
<td>MUMPS</td>
<td>yes</td>
<td>yes</td>
</tr>
<tr>
<td>Pastix</td>
<td>yes</td>
<td>yes</td>
</tr>
</tbody>
</table>

**Parallel sparse iterative solver**

Concerning iterative solvers, we have chosen pARMS, HIPS and Hypre.

Each software implements a different type of parallel preconditioner.
So, pARMS implements algebraic domain decomposition preconditioner type such as additive Schwartz [CAI1989] and interface method; while HIPS implement hierarchical incomplete factorization and finally HYPRE implements multilevel preconditioner are AMG(Algebraic MultiGrid) and parallel approximated inverse.

To use one of these programs in FreeFEM, you have to install it independently of FreeFEM. It is also necessary to install the MPI communication library which is essential for communication between the processors and, in some cases, software partitioning graphs like METIS or Scotch.

All this preconditioners are used with Krylov subspace methods accelerators.

Krylov subspace methods are iterative methods which consist in finding a solution \( x \) of linear system \( Ax = b \) inside the affine space \( x_0 + K_m \) by imposing that \( b - Ax \perp L_m \), where \( K_m \) is Krylov subspace of dimension \( m \) defined by \( K_m = \{ r_0, Ar_0, A^2r_0, ..., A^{m-1}r_0 \} \) and \( L_m \) is another subspace of dimension \( m \) which depends on type of Krylov subspace. For example in GMRES, \( L_m = AK_m \).

We realized an interface which is easy to use, so that the call of these different softwares in FreeFEM is done in the same way. You just have to load the solver and then specify the parameters to apply to the specific solvers. In the rest of this chapter, when we talk about Krylov subspace methods we mean one among GMRES, CG and BICGSTAB.

**pARMS solver**

pARMS (parallel Algebraic Multilevel Solver) is a software developed by Youssef Saad and al at University of Minnesota.

This software is specialized in the resolution of large sparse non symmetric linear systems of equation. Solvers developed in pARMS are of type "Krylov’s subspace".

It consists of variants of GMRES like FGMRES (Flexible GMRES), DGMRES (Deflated GMRES) [SAAD2003] and BICGSTAB. pARMS also implements parallel preconditioner like RAS (Restricted Additive Schwarz) [CAI1989] and Schur Complement type preconditioner.

All these parallel preconditioners are based on the principle of domain decomposition. Thus, the matrix \( A \) is partitioned into sub matrices \( A_i(i = 1, ..., p) \) where \( p \) represents the number of partitions one needs. The union of \( A_i \) forms the original matrix. The solution of the overall system is obtained by solving the local systems on \( A_i \) (see [SMITH1996]). Therefore, a distinction is made between iterations on \( A \) and the local iterations on \( A_i \).

To solve the local problem on \( A_i \) there are several preconditioners as ilut (Incomplete LU with threshold), iluk (Incomplete LU with level of fill in) and ARMS (Algebraic Recursive Multilevel Solver).

**Tip:** Default parameters

```plaintext
load "parms_FreeFem" //Tell FreeFem that you will use pARMS

// Mesh
border C(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
mesh Th = buildmesh (C(50));

// Fespace
fespace Vh(Th, P2); Vh u, v;

// Function
func f= x*y;

// Problem
problem Poisson (u, v, solver=sparsesolver)
    = int2d(Th) (dx(u) dx(v)
```

(continues on next page)
In line 1, the pARMS dynamic library is loaded with interface FreeFEM. After this, in line 15 we specify that the bilinear form will be solved by the last sparse linear solver load in memory which, in this case, is pARMS.

The parameters used in pARMS in this case are the default one since the user does not have to provide any parameter.

**Note:** In order to see the plot of a parallel script, run the command FreeFem++-mpi -glut ffglut script.edp

Here are some default parameters:

- solver=FGMRES,
- Krylov dimension=30,
- Maximum of Krylov=1000,
- Tolerance for convergence=1e-08 (see book [SAAD2003] to understand all this parameters),
- preconditionner=Restricted Additive Schwarz [CAI1989],
- Inner Krylov dimension=5,
- Maximum of inner Krylov dimension=5,
- Inner preconditionner=ILUK.

To specify the parameters to apply to the solver, the user can either give an integer vector for integer parameters and real vectors for real parameters or provide a file which contains those parameters.

**Tip:** User specifies parameters inside two vectors

Let us consider Navier-Stokes example. In this example we solve linear systems coming from discretization of Navier-Stokes equations with pARMS. Parameters of solver is specified by user.
We need two vectors to specify the parameters of the linear solver. In line 5-6 of the example, we have declared these vectors(int[int] iparm(16); real[int] dparm(6));. In line 7-10 we have initialized these vectors by negative values. We do this because all parameters values in pARMS are positive and if you do not change the negative values of one entry of this vector, the default value will be set.

In Table 3.7 and Table 3.8, we have the meaning of different entries of these vectors. We run this example on a cluster paradent of Grid5000 and report results in Table 3.5.

<table>
<thead>
<tr>
<th>np</th>
<th>add(iluk)</th>
<th>time</th>
<th>nit</th>
<th>time</th>
</tr>
</thead>
<tbody>
<tr>
<td>4</td>
<td>230</td>
<td>637.57</td>
<td>21</td>
<td>557.8</td>
</tr>
<tr>
<td>8</td>
<td>240</td>
<td>364.12</td>
<td>22</td>
<td>302.25</td>
</tr>
<tr>
<td>16</td>
<td>247</td>
<td>212.07</td>
<td>24</td>
<td>167.5</td>
</tr>
<tr>
<td>32</td>
<td>261</td>
<td>111.16</td>
<td>25</td>
<td>81.5</td>
</tr>
</tbody>
</table>
Table 3.6: Legend of Table 3.5

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>n</td>
<td>matrix size</td>
</tr>
<tr>
<td>nnz</td>
<td>number of non null entries inside matrix</td>
</tr>
<tr>
<td>nit</td>
<td>number of iteration for convergence</td>
</tr>
<tr>
<td>time</td>
<td>Time for convergence</td>
</tr>
<tr>
<td>Te</td>
<td>Time for constructing finite element matrix</td>
</tr>
<tr>
<td>np</td>
<td>number of processor</td>
</tr>
</tbody>
</table>

In this example, we fix the matrix size (in term of finite element, we fix the mesh) and increase the number of processors used to solve the linear system. We saw that, when the number of processors increases, the time for solving the linear equation decreases, even if the number of iteration increases. This proves that, using pARMS as solver of linear systems coming from discretization of partial differential equation in FreeFEM can decrease drastically the total time of simulation.

Table 3.7: Meaning of iparams corresponding variables

<table>
<thead>
<tr>
<th>Entries of iparm</th>
<th>Significations of each entries</th>
</tr>
</thead>
<tbody>
<tr>
<td>iparm[0]</td>
<td>Krylov subspace methods</td>
</tr>
<tr>
<td></td>
<td>Different values for this parameters are specify on Table 3.9</td>
</tr>
<tr>
<td>iparm[1]</td>
<td>Preconditionner</td>
</tr>
<tr>
<td></td>
<td>Different preconditioners for this parameters are specify on Table 3.10</td>
</tr>
<tr>
<td>iparm[4]</td>
<td>Number of level in arms when used</td>
</tr>
<tr>
<td>iparm[7]</td>
<td>Symmetric(=1 for symmetric) or unsymmetric matrix: default value 0(unsymmetric matrix)</td>
</tr>
<tr>
<td>iparm[8]</td>
<td>Overlap size between different subdomain: default value 0(no overlap)</td>
</tr>
<tr>
<td>iparm[9]</td>
<td>Scale the input matrix or not: Default value 1 (Matrix should be scaled)</td>
</tr>
<tr>
<td>iparm[10]</td>
<td>Block size in arms when used: default value 20</td>
</tr>
<tr>
<td>iparm[12]</td>
<td>ilfil for Schur complement const : default value 20</td>
</tr>
<tr>
<td>iparm[14]</td>
<td>Multicoloring or not in ILU when used : default value 1</td>
</tr>
<tr>
<td>iparm[15]</td>
<td>Inner iteration : default value 0</td>
</tr>
<tr>
<td>iparm[16]</td>
<td>Print message when solving: default 0 (no message print)</td>
</tr>
<tr>
<td></td>
<td>• 0: no message is print,</td>
</tr>
<tr>
<td></td>
<td>• 1: Convergence informations like number of iteration and residual,</td>
</tr>
<tr>
<td></td>
<td>• 2: Timing for a different step like preconditioner,</td>
</tr>
<tr>
<td></td>
<td>• 3 : Print all informations</td>
</tr>
</tbody>
</table>

3.6. Parallelization
### Interfacing with HIPS

**HIPS (Hierarchical Iterative Parallel Solver)** is a scientific library that provides an efficient parallel iterative solver for very large sparse linear systems. HIPS is available as free software under the CeCILL-C licence.

HIPS implements two solver classes which are the iteratives class (GMRES, PCG) and the Direct class. Concerning preconditionners, HIPS implements a type of multilevel ILU. For further informations on those preconditionners see the HIPS documentation.

**Tip:** Laplacian 3D solved with HIPS

Let us consider the 3D Laplacian example inside **FreeFEM** package where after discretization we want to solve the linear equation with HIPS.

The following example is a Laplacian 3D using Hips as linear solver. We first load Hips solver at line 2. From line 7 to 18 we specify the parameters for the Hips solver and in line 82 we set these parameters in the linear solver.
In Table 3.11 results of running on Cluster Paradent of Grid5000 are reported. We can see in this running example the efficiency of parallelism.

```plaintext
load "msh3"
load "hips_FreeFem" //load Hips library

// Parameters
int nn = 10;
real zmin = 0, zmax = 1;
int[int] iparm(14);
real[int] dparm(6);
for (int iii = 0; iii < 14; iii++)
    iparm[iii] = -1;
for (int iii = 0; iii < 6; iii++)
    dparm[iii] = -1;

//use iterative solver
iparm[0] = 0; //use iterative solver
iparm[1] = 1; //PCG as Krylov method
iparm[4] = 0; //Matrix are symmetric
iparm[5] = 1; //Pattern are also symmetric
iparm[9] = 1; //Scale matrix
dparm[0] = 1e-13; //Tolerance to convergence
dparm[1] = 5e-4; //Threshold in ILUT
dparm[2] = 5e-4; //Threshold for Schur preconditionner

// Functions
func ue = 2*x*x + 3*y*y + 4*z*z + 5*x*y + 6*x*z + 1;
func uex = 4*x + 5*y + 6*z;
func uey = 6*y + 5*x;
func uez = 8*z + 6*x;
func f = -18.;

// Mesh
mesh Th2 = square(nn, nn);
int[int] rup = [0,2], rdown=[0, 1];
int[int] rmid=[1, 1, 2, 1, 3, 1, 4, 1];

mesh3 Th = buildlayers(Th2, nn, zbound=[zmin, zmax], reffacemid=rmid,
                          reffaceup = rup, reffacelow = rdown);

// Fespace
fespace Vh2(Th2, P2);
Vh2 ux, uz, p2;

fespace Vh(Th, P2);
Vh uhe = ue;
cout << "uhe min =" << uhe[].min << ", max =" << uhe[].max << endl;
Vh u, v;
Vh F;

// Macro
macro Grad3(u) [dx(u), dy(u), dz(u)] //

// Problem
varf va (u, v) = int3d(Th)(
    Grad3(v)' * Grad3(u)
);
```

(continues on next page)
\begin{verbatim}
  + int2d(Th, 2) (
    u*v
  )
  - int3d(Th) (
    f*v
  )
  - int2d(Th, 2) (
    ue*v + (uex*N.x + uey*N.y + uez*N.z)*v
  )
  + on(1, u=ue);

  \textbf{varf l} (\textit{unused}, v) = \textbf{int3d}(Th)(f*v);

  \textbf{real cpu=\textbf{clock}();}

  \textbf{matrix Aa = \textbf{va}(Vh, Vh);}

  \textbf{F[]} = \textbf{va}(0, Vh);

  \textbf{if} (mpirank == 0) {
    \textbf{cout} << \texttt{"Size of A ="} \textbf{Aa.n} \textbf{\texttt{\textless} \texttt{endl;}}
    \textbf{cout} << \texttt{"Non zero coefficients ="} \textbf{Aa.nbcoef} \textbf{\texttt{\textless} \texttt{endl;}}
    \textbf{cout} << \texttt{"CPU TIME FOR FORMING MATRIX ="} \textbf{\texttt{\textless} clock()-\textbf{cpu} \texttt{\textless} \texttt{endl;}}
  }

  \textbf{set(Aa, solver=sparsesolver, dparams=\textbf{dparm}, lparams=\textbf{iparm}); //Set hips as linear}

  // Solve
  u[] = Aa^{-1}*F[];

  // Plot
  plot(u);
\end{verbatim}

\textbf{Table 3.11:} Legend of this table are give in \textbf{Table 3.6}

<table>
<thead>
<tr>
<th>np</th>
<th>nit</th>
<th>time</th>
</tr>
</thead>
<tbody>
<tr>
<td>8</td>
<td>190</td>
<td>120.34</td>
</tr>
<tr>
<td>16</td>
<td>189</td>
<td>61.08</td>
</tr>
<tr>
<td>32</td>
<td>186</td>
<td>31.70</td>
</tr>
<tr>
<td>64</td>
<td>183</td>
<td>23.44</td>
</tr>
</tbody>
</table>

\textbf{Tip:}

\textbf{Table 3.12:} Significations of \texttt{lparams corresponding to HIPS interface}
Entries of iparm

<table>
<thead>
<tr>
<th>iparm[0]</th>
<th>Strategy use for solving (Iterative=0 or Hybrid=1 or Direct=2). Defaults values are: Iterative</th>
</tr>
</thead>
<tbody>
<tr>
<td>iparm[1]</td>
<td>Krylov methods. If iparm[0]=0, give type of Krylov methods: 0 for GMRES, 1 for PCG</td>
</tr>
<tr>
<td>iparm[4]</td>
<td>Symmetric(=0 for symmetric) and 1 for unsymmetric matrix: default value 1 (unsymmetric matrix)</td>
</tr>
<tr>
<td>iparm[5]</td>
<td>Pattern of matrix are symmetric or not: default value 0</td>
</tr>
<tr>
<td>iparm[6]</td>
<td>Partition type of input matrix: default value 0</td>
</tr>
<tr>
<td>iparm[7]</td>
<td>Number of level that use the HIPS locally consistent fill-in: Default value 2</td>
</tr>
<tr>
<td>iparm[8]</td>
<td>Numbering in indices array will start at 0 or 1: Default value 0</td>
</tr>
<tr>
<td>iparm[9]</td>
<td>Scale matrix. Default value 1</td>
</tr>
<tr>
<td>iparm[10]</td>
<td>Reordering use inside subdomains for reducing fill-in: Only use for iterative. Default value 1</td>
</tr>
<tr>
<td>iparm[11]</td>
<td>Number of unknowns per node in the matrix non-zero pattern graph: Default value 1</td>
</tr>
<tr>
<td>iparm[12]</td>
<td>This value is used to set the number of time the normalization is applied to the matrix: Default 2.</td>
</tr>
<tr>
<td>iparm[14]</td>
<td>HIPS_DOMSIZE Subdomain size</td>
</tr>
</tbody>
</table>

Table 3.13: Significations of dparams corresponding to HIPS interface

<table>
<thead>
<tr>
<th>dparm[0]</th>
<th>HIPS_PREC: Relative residual norm: Default=1e-9</th>
</tr>
</thead>
<tbody>
<tr>
<td>dparm[1]</td>
<td>HIPS_DROPTOL0: Numerical threshold in ILUT for interior domain (important : set 0.0 in HYBRID: Default=0.005)</td>
</tr>
<tr>
<td>dparm[2]</td>
<td>HIPS_DROPTOL1: Numerical threshold in ILUT for Schur preconditioner: Default=0.005</td>
</tr>
<tr>
<td>dparm[3]</td>
<td>HIPS_DROPTOLE: Numerical threshold for coupling between the interior level and Schur: Default 0.005</td>
</tr>
<tr>
<td>dparm[4]</td>
<td>HIPS_AMALG: Numerical threshold for coupling between the interior level and Schur: Default=0.005</td>
</tr>
<tr>
<td>dparm[5]</td>
<td>HIPS_DROPSCHUR: Numerical threshold for coupling between the interior level and Schur: Default=0.005</td>
</tr>
</tbody>
</table>

Interfacing with HYPRE

Hypre (High Level Preconditioner) is a suite of parallel preconditioner developed at Lawrence Livermore National Lab.

There are two main classes of preconditioners developed in HYPRE: AMG (Algebraic MultiGrid) and Parasails (Parallel Sparse Approximate Inverse).

Now, suppose we want to solve $Ax = b$.

At the heart of AMG there is a series of progressively coarser (smaller) representations of the matrix $A$. Given an approximation $\tilde{x}$ to the solution $x$, consider solving the residual equation $Ae = r$ to find the error $e$, where $r = b - A\tilde{x}$. A fundamental principle of AMG is that it is an algebraically smooth error. To reduce the algebraically smooth errors
further, they need to be represented by a smaller defect equation (coarse grid residual equation) \( A_c e_c = r_c \), which is cheaper to solve. After solving this coarse equation, the solution is then interpolated in fine grid represented here by matrix \( A \). The quality of AMG depends on the choice of coarsening and interpolating operators.

The *sparse approximate inverse* approximates the inverse of a matrix \( A \) by a sparse matrix \( M \). A technical idea to construct matrix \( M \) is to minimize the Frobenius norm of the residual matrix \( I - MA \). For more details on this preconditioner technics see [CHOW1997].

HYPRE implement three Krylov subspace solvers: GMRES, PCG and BiCGStab.

**Tip:** Laplacian 3D solved with HYPRE

Let us consider again the 3D Laplacian example inside FreeFEM package where after discretization we want to solve the linear equation with HYPRE. The following example is a Laplacian 3D using HYPRE as linear solver. This is the same example as Hips one, so we just show here the lines where we set some HYPRE parameters.

We first load the HYPRE solver at line 2. From line 6 to 18 we specifies the parameters to set to HYPRE solver and in line 22 we set parameters to HYPRE solver.

It should be noted that the meaning of the entries of these vectors is different from those of Hips. In the case of HYPRE, the meaning of different entries of vectors \( iparm \) and \( dparm \) are given in Table 3.1.4 to Table 3.18.

In Table 3.19 the results of running on Cluster Paradent of Grid5000 are reported. We can see in this running example the efficiency of parallelism, in particular when AMG are used as preconditioner.

```plaintext
load "msh3"
load "hipre_FreeFem" //Load Hipre librairy

// Parameters
int nn = 10;
int[int] iparm(20);
real[int] dparm(6);
for (int iii = 0; iii < 20; iii++)
  iparm[iii] = -1;
for (int iii = 0; iii < 6; iii++)
  dparm[iii] = -1;
iparm[0] = 2; //PCG as krylov method
iparm[1] = 0; //AMG as preconditioner 2: if ParaSails
iparm[7] = 7; //Interpolation
iparm[9] = 6; //AMG Coarsen type
iparm[10] = 1; //Measure type
iparm[16] = 2; //Additive schwarz as smoother
dparm[0] = 1e-13; //Tolerance to convergence
dparm[1] = 5e-4; //Threshold
dparm[2] = 5e-4; //Truncation factor
...
set(Aa, solver=sparseSolver, dparams=dparm, lparams=iparm);
```

**Table 3.14:** Definitions of common entries of \( iparms \) and \( dparms \) vectors for every preconditioner in HYPRE
### iparms[0]
Solver identification:
- 0: BiCGStab, 1: GMRES, 2: PCG. Default=1

### iparms[1]
Preconditioner identification:
- 0: BOOMER AMG, 1: PILUT, 2: Parasails, 3: Schwartz Default=0

### iparms[2]
Maximum of iteration: Default=1000

### iparms[3]
Krylov subspace dim: Default=40

### iparms[4]
Solver print info level: Default=2

### iparms[5]
Solver log: Default=1

### iparms[6]
Solver stopping criteria only for BiCGStab: Default=1

### dparms[0]
Tolerance for convergence: Default=:math:`1.0e-11`

**Table 3.15:** Definitions of other entries of `iparms` and `dparms` if preconditioner is BOOMER AMG

<table>
<thead>
<tr>
<th>iparms[7]</th>
<th>AMG interpolation type: Default=6</th>
</tr>
</thead>
<tbody>
<tr>
<td>iparms[8]</td>
<td>Specifies the use of GSMG - geometrically smooth coarsening and interpolation: Default=1</td>
</tr>
<tr>
<td>iparms[9]</td>
<td>AMG coarsen type: Default=6</td>
</tr>
<tr>
<td>iparms[10]</td>
<td>Defines whether local or global measures are used: Default=1</td>
</tr>
<tr>
<td>iparms[11]</td>
<td>AMG cycle type: Default=1</td>
</tr>
<tr>
<td>iparms[12]</td>
<td>AMG Smoother type: Default=1</td>
</tr>
<tr>
<td>iparms[13]</td>
<td>AMG number of levels for smoothers: Default=3</td>
</tr>
<tr>
<td>iparms[14]</td>
<td>AMG number of sweeps for smoothers: Default=2</td>
</tr>
<tr>
<td>iparms[15]</td>
<td>Maximum number of multigrid levels: Default=25</td>
</tr>
<tr>
<td>iparms[16]</td>
<td>AMG Smoother type: Default=1</td>
</tr>
<tr>
<td>iparms[17]</td>
<td>Size of the system of PDEs: Default=1</td>
</tr>
<tr>
<td>iparms[18]</td>
<td>Overlap for the Schwarz method: Default=1</td>
</tr>
<tr>
<td>iparms[19]</td>
<td>Type of domain used for the Schwarz method</td>
</tr>
<tr>
<td></td>
<td>0: each point is a domain</td>
</tr>
<tr>
<td></td>
<td>1: each node is a domain (only of interest in “systems” AMG)</td>
</tr>
<tr>
<td></td>
<td>2: each domain is generated by agglomeration (default)</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>dparms[1]</th>
<th>AMG strength threshold: Default=0.25</th>
</tr>
</thead>
<tbody>
<tr>
<td>dparms[3]</td>
<td>Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix: Default=0.9</td>
</tr>
<tr>
<td>dparms[4]</td>
<td>Defines a smoothing parameter for the additive Schwartz method. Default=1</td>
</tr>
</tbody>
</table>

**Table 3.16:** Definitions of other entries of `iparms` and `dparms` if preconditioner is PILUT

<table>
<thead>
<tr>
<th>iparms[7]</th>
<th>Row size in Parallel ILUT: Default=1000</th>
</tr>
</thead>
<tbody>
<tr>
<td>iparms[8]</td>
<td>Set maximum number of iterations: Default=30</td>
</tr>
<tr>
<td>dparms[1]</td>
<td>Drop tolerance in Parallel ILUT: Default=1e-5</td>
</tr>
</tbody>
</table>

**Table 3.17:** Definitions of other entries of `iparms` and `dparms` if preconditioner is ParaSails

3.6. Parallelization
**Table 3.18:** Definitions of other entries of `iparms` and `dparms` if preconditioner is Schwartz

<table>
<thead>
<tr>
<th><code>iparms[7]</code></th>
<th>Defines which variant of the Schwartz method is used:</th>
</tr>
</thead>
<tbody>
<tr>
<td>0: hybrid multiplicative Schwartz method (no overlap across processor boundaries)</td>
<td></td>
</tr>
<tr>
<td>1: hybrid additive Schwartz method (no overlap across processor boundaries)</td>
<td></td>
</tr>
<tr>
<td>2: additive Schwartz method</td>
<td></td>
</tr>
<tr>
<td>3: hybrid multiplicative Schwartz method (with overlap across processor boundaries)</td>
<td></td>
</tr>
<tr>
<td>Default=1</td>
<td></td>
</tr>
</tbody>
</table>

| `iparms[8]` | Overlap for the Schwartz method: Default=1 |

<table>
<thead>
<tr>
<th><code>iparms[9]</code></th>
<th>Type of domain used for the Schwartz method</th>
</tr>
</thead>
<tbody>
<tr>
<td>0: each point is a domain</td>
<td></td>
</tr>
<tr>
<td>1: each node is a domain (only of interest in “systems” AMG)</td>
<td></td>
</tr>
<tr>
<td>2: each domain is generated by agglomeration (default)</td>
<td></td>
</tr>
</tbody>
</table>

**Table 3.19:** Convergence and time for solving linear system

<table>
<thead>
<tr>
<th>$n = 4 \times 10^6$</th>
<th>$nnz = 13 \times 10^6$</th>
<th>$Te = 571.29$</th>
</tr>
</thead>
<tbody>
<tr>
<td>np</td>
<td>AMG</td>
<td>nit</td>
</tr>
<tr>
<td>-----</td>
<td>-----</td>
<td>-----</td>
</tr>
<tr>
<td>8</td>
<td>6</td>
<td>1491.83</td>
</tr>
<tr>
<td>16</td>
<td>5</td>
<td>708.49</td>
</tr>
<tr>
<td>32</td>
<td>4</td>
<td>296.22</td>
</tr>
<tr>
<td>64</td>
<td>4</td>
<td>145.64</td>
</tr>
</tbody>
</table>

**Conclusion**

With the different runs presented here, we wanted to illustrate the gain in time when we increase the number of processors used for the simulations. We saw that in every case the time for the construction of the finite element matrix is constant. This is normal because until now this phase is sequential in **FreeFEM**. In contrast, phases for solving the linear system are parallel. We saw on several examples presented here that when we increase the number of processors, in general we decrease the time used for solving the linear systems. But this is not true in every case. In several cases, when we increase the number of processors the time to convergence also increases. There are two main reasons for this. First, the increase of processors can lead to the increase of volume of exchanged data across processors consequently increasing the time for solving the linear systems.

Furthermore, in decomposition domain type preconditioners, the number of processors generally corresponds to the number of sub domains. In subdomain methods, generally when we increase the number of subdomains we decrease convergence quality of the preconditioner. This can increase the time used for solving linear equations.

To end this, we should note that good use of the preconditioners interfaced in **FreeFEM** is empiric, because it is difficult to know what is a good preconditioner for some type of problems. Although, the efficiency of preconditioners sometimes depends on how its parameters are set. For this reason we advise the user to pay attention to the meaning of the parameters in the user guide of the iterative solvers interfaced in **FreeFEM**.
Domain decomposition

In the previous section, we saw that the phases to construct a matrix are sequential. One strategy to construct the matrix in parallel is to divide geometrically the domain into subdomains. In every subdomain we construct a local submatrix and after that we assemble every submatrix to form the global matrix.

We can use this technique to solve PDE directly in domain $\Omega$. In this case, in every subdomains you have to define artificial boundary conditions to form consistent equations in every subdomains. After this, you solve equation in every subdomains and define a strategy to obtain the global solution.

In terms of parallel programming for FreeFEM, with MPI, this means that the user must be able to divide processors available for computation into subgroups of processors and also must be able to realize different type of communications in FreeFEM script. Here is a wrapper of some MPI functions.

Communicators and groups

Groups

`mpiGroup grpe(mpiGroup gp, KN_<long>)`: Create MPI_Group from existing group `gp` by given vector.

Communicators

Communicators is an abstract MPI object which allows MPI user to communicate across group of processors. Communicators can be Intra-communicators (involves a single group) or Inter-communicators (involves two groups). When we not specify type of communicator it will be Intra-communicators

`mpiComm cc(mpiComm comm, mpiGroup gp)`: Creates a new communicator.

`mpiComm cc(mpiGroup gp)`: Same as previous constructor but default `comm` here is MPI_COMM_WORLD.

`mpiComm cc(mpiComm comm, int color, int key)`: Creates new communicators based on colors and key. This constructor is based on MPI_Comm_split routine of MPI.

`mpiComm cc(MPIrank p, int key)`: Same constructor than the last one.

Tip: Split communicator

```
mpiComm comm(mpiCommWorld, 0, 0);
int color = mpiRank(comm)%2;
mpiComm ccc(processor(color, comm), 0);
mpiComm qpp(comm, 0, 0);
mpiComm cp(ccc, color, 0);
```

`mpiComm cc(mpiComm comm, int high)`: Creates an intracomunicator from an intercommunicator. `comm` intercommunicator, `high`.

Used to order the groups within `comm` (logical) when creating the new communicator. This constructor is based on MPI_Intercomm_merge routine of MPI.

`mpiComm cc(MPIrank p1, MPIrank p2, int tag)`: This constructor creates an intercommunicator from two intra-communicators. `p1` defined local (intra)communicator and rank in local_comm of leader (often 0) while `p2` defined remote communicator and rank in peer_comm of remote leader (often 0). `tag` Message tag to use in constructing intercommunicator. This constructor is based on MPI_Intercomm_create.
Tip: Merge

```plaintext
mpiComm comm, cc;
int color = mpiRank(comm) % 2;
int rk = mpiRank(comm);
int size = mpiSize(comm);
cout << "Color values: " << color << endl;
mpiComm ccc(processor((rk < size / 2), comm), rk);
mpiComm cp(cc, color, 0);
int rleader;
if (rk == 0) { rleader = size / 2; }
else if (rk == size / 2) { rleader = 0; }
else { rleader = 3; }
mpiComm qqp(processor(0, ccc), processor(rleader, comm), 12345);
int aaa = mpiSize(ccc);
cout << "Number of processor: " << aaa << endl;
```

Process

In FreeFEM we wrap MPI process by function call processor which create internal FreeFEM object call MPIrank. This mean that do not use MPIrank in FreeFEM script.

`processor(int rk): Keep process rank inside object MPIrank. Rank is inside MPI_COMM_WORLD.`
`processor(int rk, mpiComm cc) and processor(mpiComm cc, int rk) process rank inside communicator cc.`

`processorblock(int rk): This function is exactly the same than processor(int rk) but is use in case of blocking communication.`
`processorblock(int rk, mpiComm cc): This function is exactly the same as processor(int rk, mpiComm cc) but uses a synchronization point.`

Points to Points communicators

In FreeFEM you can call MPI points to points communications functions.

`Send(processor(int rk, mpiComm cc), Data D) : Blocking send of Data D to processor of rank rk inside communicator cc. Note that Data D can be: int, real, complex, int[int], real[int], complex[int], Mesh, Mesh3, Matrix.`
`Recv(processor(int rk, mpiComm cc), Data D): Receive Data D from process of rank rk in communicator cc. Note that Data D can be: int, real, complex, int[int], real[int], complex[int], Mesh, Mesh3, Matrix and should be the same type than corresponding send.`
`Isend(processor(int rk, mpiComm cc), Data D) : Non blocking send of Data D to processor of rank rk inside communicator cc. Note that Data D can be: int, real, complex, int[int], real[int], complex[int], mesh, mesh3, matrix.}`
Global operations

In **FreeFEM** you can call MPI global communication functions.

```plaintext
broadcast(processor(int rk, mpiComm cc), Data D): Process rk Broadcast Data D to all process inside communicator cc. Note that Data D can be: int, real, complex, int[int], real[int], complex[int], Mesh, Mesh3, Matrix.

broadcast(processor(int rk), Data D): Process rk Broadcast Data D to all processes inside communicator MPI_COMM_WORLD. Note that Data D can be: int, real, complex, int[int], real[int], complex[int], Mesh, Mesh3, Matrix.

mpiAlltoall(Data a, Data b): Sends data a from all to all processes. Receive buffer is Data b. This is done inside communicator MPI_COMM_WORLD.

mpiAlltoall(Data a, Data b, mpiComm cc): Sends data a from all to all processes. Receive buffer is Data b. This is done inside communicator cc.

mpiGather(Data a, Data b, processor(mpiComm, int rk)): Gathers together values Data a from a group of processes. Process of rank rk get data on communicator rk. This function is like MPI_Gather.

mpiAllgather(Data a, Data b): Gathers Data a from all processes and distribute it to all in Data b. This is done inside communicator MPI_COMM_WORLD. This function is like MPI_Allgather.

mpiAllgather(Data a, Data b, mpiComm cc): Gathers Data a from all processes and distribute it to all in Data b. This is done inside communicator cc. This function is like MPI_Allgather.

mpiScatter(Data a, Data b, processor(int rk, mpiComm cc)): Sends Data a from one process with rank rk to all other processes in group represented by communicator mpiComm cc.

mpiReduce(Data a, Data b, processor(int rk, mpiComm cc), MPI_Op op): Reduces values Data a on all processes to a single value Data b on process of rank rk and communicator cc.

Operation use in reduce is: MPI_Op op which can be: mpiMAX, mpiMIN, mpiSUM, mpiPROD, mpiLAND, mpiLOR, mpiLXOR, mpiBAND, mpiBXOR, mpiMAXLOC, mpiMINLOC.

Note that, for all global operations, only int[int] and real[int] are data type take in account in **FreeFEM**.

### 3.7 PETSc and SLEPc

**FreeFEM** is interfaced with **PETSc** and **SLEPc** which offer a wide range of sequential or parallel linear or nonlinear solvers, time steppers, optimizers, and eigensolvers. In particular, it gives access transparently (without much changes to user code) to: distributed and multithreaded direct solvers (**PARDISO, MUMPS, SuperLU**), multigrid solvers (**hypre, GAMG**), domain decomposition methods (**block Jacobi, ASM, HPDDM**). For a detailed introduction to these tools, interested readers are referred to the tutorial Introduction to FreeFEM with an emphasis on parallel computing.

In most of the scripts listed below, the following standard procedure is used.

- Load an initial sequential mesh (in 2D or 3D).
- Partition the mesh and generate connectivity information according to the number of processes.
- Provide these information to PETSc so that subsequent computations may be done in a distributed fashion.

Combining the power and flexibility of PETSc with the ease-of-use of FreeFEM may help design multiphysics solvers, e.g., for **Navier–Stokes equations**, advanced matrix-free discretizations, and such.
### 3.7.1 Examples

#### Linear problems

<table>
<thead>
<tr>
<th>Filename</th>
<th>Comments (preconditioners, numerical schemes)</th>
</tr>
</thead>
<tbody>
<tr>
<td>diffusion-2d-PETSc.edp</td>
<td>Distributed LU/Cholesky, domain decomposition and multigrid methods</td>
</tr>
<tr>
<td>diffusion-2d-PETSc-complex.edp</td>
<td></td>
</tr>
<tr>
<td>heat-2d-PETSc.edp</td>
<td>Transient diffusion equation, same as above</td>
</tr>
<tr>
<td>diffusion-periodic-2d-PETSc.edp</td>
<td>Periodic boundary conditions, multigrid methods</td>
</tr>
<tr>
<td>diffusion-periodic-balanced-2d-PETSc.edp</td>
<td>Better load balancing than above example</td>
</tr>
<tr>
<td>diffusion-substructuring-2d-PETSc.edp</td>
<td>Balancing Domain Decomposition with Constraints</td>
</tr>
<tr>
<td>diffusion-3d-PETSc.edp</td>
<td>Three-dimensional problem, domain decomposition and multigrid methods</td>
</tr>
<tr>
<td>diffusion-mg-2d-PETSc.edp</td>
<td>Geometric non-nested multigrid methods</td>
</tr>
<tr>
<td>diffusion-mg-3d-PETSc.edp</td>
<td>Geometric nested multigrid methods</td>
</tr>
<tr>
<td>helmholtz-2d-PETSc-complex.edp</td>
<td>Domain decomposition methods with optimized boundary conditions</td>
</tr>
<tr>
<td>helmholtz-mg-2d-PETSc-complex.edp</td>
<td>Geometric multigrid methods</td>
</tr>
<tr>
<td>laplace-RT-2d-PETSc.edp</td>
<td>Vectorial two-dimensional problem with a block preconditioner (fieldsplit)</td>
</tr>
<tr>
<td>laplace-adapt-3d-PETSc.edp</td>
<td>Three-dimensional problem with $h$ adaptivity, multigrid methods using Mmg</td>
</tr>
<tr>
<td>laplace-adapt-dist-3d-PETSc.edp</td>
<td>Three-dimensional problem with fully-distributed $h$ adaptivity using ParMmg</td>
</tr>
<tr>
<td>laplace-lagrange-PETSc.edp</td>
<td>Laplace equation with constraints and a block preconditioner (fieldsplit)</td>
</tr>
<tr>
<td>elasticity-2d-PETSc.edp</td>
<td>Vectorial problem, domain decomposition (GenEO) and multigrid methods</td>
</tr>
<tr>
<td>elasticity-3d-PETSc.edp</td>
<td></td>
</tr>
<tr>
<td>stokes-2d-PETSc.edp</td>
<td>Distributed LU/Cholesky</td>
</tr>
<tr>
<td>stokes-3d-PETSc.edp</td>
<td></td>
</tr>
<tr>
<td>stokes-block-2d-PETSc.edp</td>
<td>Stokes equation defined as a block system with four matrices (fieldsplit)</td>
</tr>
<tr>
<td>stokes-fieldsplit-2d-PETSc.edp</td>
<td>Block preconditioner (fieldsplit)</td>
</tr>
<tr>
<td>stokes-fieldsplit-3d-PETSc.edp</td>
<td></td>
</tr>
<tr>
<td>oseen-2d-PETSc.edp</td>
<td>Oseen problem preconditioned by Pressure Convection–Diffusion (PCD)</td>
</tr>
<tr>
<td>maxwell-2d-PETSc.edp</td>
<td>Direct LU/Cholesky</td>
</tr>
<tr>
<td>maxwell-3d-PETSc.edp</td>
<td>Multigrid method</td>
</tr>
<tr>
<td>maxwell-mg-3d-PETSc-complex.edp</td>
<td>Two-grid preconditioner</td>
</tr>
<tr>
<td>helmholtz-3d-surf-PETSc-complex.edp</td>
<td>BEM with hierarchical matrices from Htool on surfaces</td>
</tr>
<tr>
<td>helmholtz-3d-line-PETSc-complex.edp</td>
<td>BEM with hierarchical matrices from Htool on lines</td>
</tr>
<tr>
<td>helmholtz-coupled-2d-PETSc-complex.edp</td>
<td>FEM coupled with BEM in a MatNest using a block preconditioner (fieldsplit)</td>
</tr>
<tr>
<td>PtAP-2d-PETSc.edp</td>
<td>Parallel interpolation on non-matching grids + $P'$ A P operation</td>
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3.8 Plugins

### 3.8.1 gsl

The interface with gsl spline is available in **FreeFEM**, the seven kind of spline are

- 0. gslinterpcspline: default type of spline
- 1. gslinterpakima
- 2. gslinterpsteffen
- 3. gslinterlinear
- 4. gslinterppolynomial
- 5. gslinterpcsplineperiodic
- 6. gslinterpakimaperiodic

A brief wing example given all the syntax:

```plaintext
load "gsl"

// Parameters
int n = 10;
real[int, int] dspline(2, n+1); //data points to define the spline
for(int i = 0; i <= n; ++i){ //set data points
    real xx = square(real(i)/n);
    real yy = sin(xx*pi*2);
    dspline(0, i) = xx;
    dspline(1, i) = yy;
}

// GSL splines
glspline spline1(gslinterpcspline, dspline); //define the spline1
spline11(dspline);

//define the spline1

//evaluate the function spline1 at t
real t = 1.;
real s1 = spline1(t);  cout << "spline1(t) = " << s1 << endl;

//evaluate the derivative of function spline1 at t
real ds1 = spline1.d(t); cout << "spline1.d(t) = " << ds1 << endl;

//evaluate the second derivative of function spline1 at t
real dds1 = spline1.dd(t); cout << "spline1.dd(t) = " << dds1 << endl;
```

This can be usefull to build function from data value.
The list of all gsl functions and the FreeFEM equivalent is available in the Language references (same names without _).

### 3.8.2 ffrandom

Plugin to linux random functions.

The range of the random generator is from 0 to \((2^{31}) - 1\).

```cpp
1
load "ffrandom"
2
3
srandomdev(); // set a true random seed
4
// warning: under window this command
5
// change the seed by randinit(random())) so all
6
// FreeFEM random function are changed
7
8
int maxrang = 2^31 - 1;
9
cout " max range " maxrang endl;
10
11
cout random() endl;
12
cout random() endl;
13
cout random() endl;
14
15
srandom(10);
16
cout random() endl;
17
cout random() endl;
18
cout random() endl;
```

### 3.8.3 mmap / semaphore

The idea is just to use Interprocess communication using POSIX Shared Memory in Linux.

We build a small library libff-mmap-semaphore.c and libff-mmap-semaphore.h to easily interface.

- mmap - allocate memory, or map files or devices into memory
- semaphore - allow processes and threads to synchronize their actions

A semaphore is an integer whose value is never allowed to fall below zero. Two operations can be performed on semaphores: increment the semaphore value by one \(\text{sem_post}\); and decrement the semaphore value by one \(\text{sem_wait}\).

If the value of a semaphore is currently zero, then a \text{sem_wait} operation will block until the value becomes greater than zero.

The functions of library

First the semaphore interface to make synchronization:

- typedef struct FF_P_sem *ff_Psem; the pointer to data structure
- ff_Psem ffsem_malloc(); malloc an empty data structure
- void ffsem_del(ff_Psem sem); clean and free the pointer
- void ffsem_destroy(ff_Psem sem); clean, close the data structure
- void ffsem_init0(ff_Psem sem); make a correct empty of the data structure
• void ffsem_init(ff_Psem sem, const char *nmm, int crea); create or use a new semaphore

• long ffsem_post(ff_Psem sem); nlocked, the value of the semaphore is incremented, and all threads which are waiting on the semaphore are awakened

• long ffsem_wait(ff_Psem sem); the semaphore referenced by sem is locked. When calling sem.wait(), if the semaphore’s value is zero, the calling thread will block until the lock is acquired or until the call is interrupted by a signal.

Alternatively, the sem.trywait() function will fail if the semaphore is already locked, rather than blocking on the semaphore

• long ffsem_trywait(ff_Psem p);

Secondly, the mmap functions:

• typedef struct FF_P_mmap *ff_Pmmap; the pointer to data structure

• ff_Psem ffmmap_malloc(); malloc an empty data structure

• void ffmmap_del(ff_Pmmap p); clean and free the pointer

• void ffmmap_destroy(ff_Pmmap p); clean, close the data structure

• void ffmmap_init0(ff_Pmmap p); make a correct empty of the data structure

• long ffmmap_msync(ff_Pmmap p, long off, long ln); call writes modified whole pages back to the filesystem and updates the file modification time. Only those pages containing addr and len-1 succeeding locations will be examined.

• void ffmmap_init(ff_Pmmap p, const char *nmm, long len); allocate memory, or map files or devices into memory.

• long ffmmap_read(ff_Pmmap p, void *t, size_t n, size_t off); read n bytes from the mmap at memory offset in pointer t.

• long ffmmap_write(ff_Pmmap p, void *t, size_t n, size_t off); write n bytes to the mmap at memory offset in pointer t.

The FreeFEM corresponding functions:

• Pmmap sharedata(filename, 1024); new type to store the mmap informations of name store in string filename with 1024 is the size the sharedata zone and file.

• Psemaphore smff("ff-slave", creat); new type to store the semaphore of name ff-slave where creat is a boolean to create or use a existing semaphore.

• Wait(sem) the semaphore referenced by sem is locked. When calling Wait(sem), if the semaphore’s value is zero, the calling thread will block until the lock is acquired or until the call is interrupted by a signal. Alternatively, the trywait(sem) function will fail if the semaphore is already locked, rather than blocking on the semaphore.

• Post(sem) the semaphore referenced by sem is unlocked, the value of the semaphore is incremented, and all threads which are waiting on the semaphore are awakened.

• Read(sharedata, offset, data); read the variable data from the place offset in sharedata mmap.

• Write(sharedata, offset, data); write the variable data at the place offset in sharedata mmap.

The full example:

The FFMaster.c file:
```c
#include "libff-mmap_semaphore.h"
#include <unistd.h>
#include <stdlib.h>
#include <stdio.h>

ff_Psem sem_ff, sem_c; //the semaphore for mutex

int main(int argc, const char ** argv)
{
    int debug = 0;
    ff_Pmmap shd;
    double cff, rff;
    long status;
    int i;
    if (argc > 1) debug = atoi(argv[1]);
    ff_mmap_sem_verb = debug;

    sem_ff = ffsem_malloc();
    sem_c = ffsem_malloc();
    shd = ffmmap_malloc();

    ffsem_init(sem_ff, "ff-slave1", 1);
    ffsem_init(sem_c, "ff-master1", 1);
    ffmmap_init(shd, "shared-data", 1024);

    status = 1;
    ffmmap_write(shd, &status, sizeof(status), 8);
    ffmmap_msync(shd, 0, 32);

    char ff[1024];
    sprintf(ff, "FreeFem++ FFSlave.edp -nw -ns -v %d&", debug);
    system(ff); //lauch FF++ in batch no graphics
    if(debug) printf("cc: before wait
");

    if(debug) printf("cc: before wait 0 ff\n");
    ffsem_wait(sem_ff);

    for (i = 0; i < 10; ++i){
        printf(" iter : %d 
", i);
        cff = 10+i;
        ffmmap_write(shd, &cff, sizeof(cff), 0);
        ffsem_post(sem_c);

        if(debug) printf(" cc: before wait 2\n");
        ffsem_wait(sem_ff);
        ffmmap_read(shd, &rff, sizeof(rff), 16);
        printf(" iter = %d rff= %f\n", i, rff);
    }

    status = 0; //end
    ffmmap_write(shd, &status, sizeof(status), 8);
    ffsem_post(sem_c);
    printf("End Master \n");
    ffsem_wait(sem_ff);
    ffsem_del(sem_ff);
    ffsem_del(sem_c);
    ffmmap_del(shd);
    return 0;
}
```
The FFSlave.edp file:

```cpp
load "ff-mmap-semaphore"

Psemaphore smff("ff-slave1", 0);
Psemaphore smc("ff-master1", 0);
Pmmap sharedata("shared-data", 1024);
if (verbosity < 4) verbosity = 0;

// Mesh
mesh Th = square(10, 10);
int[int] Lab = [1, 2, 3, 4];

// Fespace
fespace Vh(Th, P1);
Vh u, v;

// Macro
macro grad(u) [dx(u), dy(u)] //
int status = 1;
cout << " FF status = " << status << endl;
real cff, rff;

// Problem
problem Pb (u, v)
  = int2d(Th)(
    grad(u)'*grad(v)
  )
  - int2d(Th)(
    cff*v
  )
  + on(Lab, u=0);

if (verbosity > 9) cout << " FF: before FF post

Post(smff); //unlock master end init

while (1){
  if (verbosity > 9) cout << " FF: before FF wait \n";
  Wait(smc); //wait from cint write ok
  Read(sharedata, 0, cff);
  Read(sharedata, 8, status);
  cout << " After wait .. FF " << cff << " " << status << endl;
  if(status <= 0) break;

  // Solve
  Pb;
  rff = int2d(Th)(u*u);
  cout << " ** FF " << cff << " " << rff << endl;

  // Write
  Write(sharedata, 16, rff);
  Post(smff); //unlock cc
}
Post(smff); //wait from cint
```

(continues on next page)
cout << " End FreeFEM " << endl;

To test this example of coupling C program and FreeFEM script:

```plaintext
cc -c libff-mmap-semaphore.c
cc FFMaster.c -o FFMaster libff-mmap-semaphore.o -g -pthread
ff-c++ -auto ff-mmap-semaphore.cpp
./FFMaster
```

The output:

```plaintext
len 1024 size 0
len 1024 size 1024
FF status = 1
iter : 0
After wait .. FF 10 1
** FF 10 0.161797
iter = 0 rff= 0.161797
iter : 1
After wait .. FF 11 1
** FF 11 0.195774
iter = 1 rff= 0.195774
iter : 2
After wait .. FF 12 1
** FF 12 0.232987
iter = 2 rff= 0.232987
iter : 3
After wait .. FF 13 1
** FF 13 0.273436
iter = 3 rff= 0.273436
iter : 4
After wait .. FF 14 1
** FF 14 0.317121
iter = 4 rff= 0.317121
iter : 5
After wait .. FF 15 1
** FF 15 0.364042
iter = 5 rff= 0.364042
iter : 6
After wait .. FF 16 1
** FF 16 0.414199
iter = 6 rff= 0.414199
iter : 7
After wait .. FF 17 1
** FF 17 0.467592
iter = 7 rff= 0.467592
iter : 8
After wait .. FF 18 1
** FF 18 0.524221
iter = 8 rff= 0.524221
iter : 9
After wait .. FF 19 1
** FF 19 0.584086
iter = 9 rff= 0.584086
End Master
After wait .. FF 19 0
```
3.9 Developers

3.9.1 File formats

Mesh file data structure

The mesh data structure, output of a mesh generation algorithm, refers to the geometric data structure and in some case to another mesh data structure.

In this case, the fields are

```plaintext
MeshVersionFormatted 0
Dimension [DIM] (int)
Vertices
[Number of vertices](int)
X_1(double) Y_1(double) (Z_1(double)) Ref_1(int)
...
X_nv(double) Y_nv(double) (Z_nv(double)) Ref_nv(int)
Edges
[Number of edges](int)
Vertex1_1(int) Vertex2_1(int) Ref_1(int)
...
Vertex1_ne(int) Vertex2_ne(int) Ref_ne(int)
Triangles
[Number of triangles](int)
Vertex1_1(int) Vertex2_1(int) Vertex3_1(int) Ref_1(int)
...
Vertex1_nt(int) Vertex2_nt(int) Vertex3_nt(int) Ref_nt(int)
Quadrilaterals
[Number of Quadrilaterals](int)
Vertex1_1(int) Vertex2_1(int) Vertex3_1(int) Vertex4_1(int) Ref_1(int)
...
Vertex1_nq(int) Vertex2_nq(int) Vertex3_nq(int) Vertex4_nq(int) Ref_nq(int)
Geometry
[File name of geometric support](char*)
VertexOnGeometricVertex
[Number of vertex on geometric vertex](int)
Vertex_1(int) VertexGeometry_1(int)
...
Vertex_nvg(int) VertexGeometry_nvg(int)
EdgeOnGeometricEdge
[Number of geometric edge](int)
Edge_1(int) EdgeGeometry_1(int)
...
Edge_neg(int) EdgeGeometry_neg(int)
CrackedEdges
[Number of cracked edges](int)
Edge1_1(int) Edge2_1(int)
```

(continues on next page)
When the current mesh refers to a previous mesh, we have in addition

```c
MeshSupportOfVertices
[File name of mesh support](char*)

VertexOnSupportVertex
[Number of vertex on support vertex](int)
Vertex_1(int) VertexSupport_1(int)
...
Vertex_nvsv(int) VertexSupport_nvsv(int)

VertexOnSupportEdge
[Number of vertex on support edge](int)
Vertex_1(int) EdgeSupport_1(int) USupport_1(double)
...
Vertex_nvse(int) EdgeSupport_nvse(int) USupport_nvse(double)

VertexOnSupportTriangle
[Number of vertex on support triangle](int)
Vertex_1(int) TriangleSupport_1(int) USupport_1(double) VSupport_1(double)
...
Vertex_nvst(int) TriangleSupport_nvst(int) USupport_nvst(double) VSupport_nvst(double)

VertexOnSupportQuadrilaterals
[Number of vertex on support quadrilaterals]
Vertex_1(int) TriangleSupport_1(int) USupport_1(double) VSupport_1(double)
...
Vertex_nvsq(int) TriangleSupport_nvsq(int) USupport_nvsq(double) VSupport_nvsq(double)
```

- `nv` means the number of vertices
- `ne` means the number of edges
- `nt` means the number of triangles
- `nq` means the number of quadrilaterals
- `nvg` means the number of vertex on geometric vertex
- `neg` means the number of edges on geometric edge
- `nce` means the number of cracked edges

**bb file type to Store Solutions**

The file is formatted such that:

```c
[Number of solutions](int) [Number of vertices](int) 2
U_1_1(double) ... U_ns_1(double)
...
U_1_nv(double) ... U_ns_nv(double)
```
• \( ns \) means the number of solutions
• \( nv \) means the number of vertices
• \( U_{i,j} \) is the solution component \( i \) at the vertex \( j \) on the associated mesh.

**BB file type to store solutions**

The file is formatted such that:

```
1  [Number of solutions](int)  [Type 1](int) ... [Type ns](int)  [Number of
→vertices](int) 2
2  U_{1,1,1}(double) ... U_{(type_k),1,1}(double)
3  ...  
4  U_{1,1,1}(double) ... U_{(type_k),nbv,1}(double)
5  ...  
6  U_{1,1,ns}(double) ... U_{(type_k),1,ns}(double)
7  ...  
8  U_{1,nbv,ns}(double) ... U_{(type_k),nbv,ns}(double)
```

• \( ns \) means the number of solutions
• \( type_k \) mean the type of solution \( k \):
  - 1: the solution is scalar (1 value per vertex)
  - 2: the solution is vectorial (2 values per vertex)
  - 3: the solution is a \( 2 \times 2 \) symmetric matrix (3 values per vertex)
  - 4: the solution is a \( 2 \times 2 \) matrix (4 values per vertex)
• \( nbv \) means the number of vertices
• \( U_{i,j,k} \) is the value of the component \( i \) of the solution \( k \) at vertex \( j \) on the associated mesh

**Metric file**

A metric file can be of two types, isotropic or anisotropic.

The isotropic file is such that

```
1  [Number of vertices](int) 1
2  h_0(double)
3  ...
4  h_nv(double)
```

• \( nv \) is the number of vertices
• \( h_i \) is the wanted mesh size near the vertex \( i \) on associated mesh.

The metric is \( M_i = h_i^{-2} I \) where \( I \) is the identity matrix.

The anisotropic file is such that
• $\text{nv}$ is the number of vertices

• $a_{11,i}, a_{21,i}$ and $a_{22,i}$ represent metric $M_i = \begin{pmatrix} a_{11,i} & a_{12,i} \\ a_{12,i} & a_{22,i} \end{pmatrix}$ which define the wanted size in a vicinity of the vertex $i$ such that $h$ in direction $u \in \mathbb{R}^2$ is equal to $\frac{|u|}{\sqrt{u \cdot M_i u}}$, where $\cdot$ is the dot product in $\mathbb{R}^2$, and $|\cdot|$ is the classical norm.

List of AM_FMT, AMDBA Meshes

The mesh is only composed of triangles and can be defined with the help of the following two integers and four arrays:

• $\text{nbt}$ the number of triangles

• $\text{nbv}$ the number of vertices

• $\text{nu}(1:3, 1:\text{nbt})$ an integer array giving the three vertex numbers counterclockwise for each triangle

• $\text{c}(1:2, 1:\text{nbv})$ a real array giving the two coordinates of each vertex

• $\text{refs}(1:\text{nbv})$ an integer array giving the reference numbers of the vertices

• $\text{reft}(1:\text{nbt})$ an integer array giving the reference numbers of the triangles

AM_FMT Files

In Fortran the am_fmt files are read as follows:

```fortran
1 open (1, file='xxx.am_fmt', form='formatted', status='old')
2 read (1, *) nbv, nbt
3 read (1, *) ((nu(i, j), i=1, 3), j=1, nbt)
4 read (1, *) ((c(i, j), i=1, 2), j=1, nbv)
5 read (1, *) ( reft(i), i=1, nbt)
6 read (1, *) ( refs(i), i=1, nbv)
7 close(1)
```

AM Files

In Fortran the am files are read as follows:

```fortran
1 open (1, file='xxx.am', form='unformatted', status='old')
2 read (1, *) nbv, nbt
3 read (1) ((nu(i, j), i=1, 3), j=1, nbt),
4 & ((c(i, j), i=1, 2), j=1, nbv),
5 & (reft(i), i=1, nbt),
6 & (refs(i), i=1, nbv)
7 close(1)
```

AMDBA Files

In Fortran the amdba files are read as follows:

```fortran
1 open (1, file='xxx.amdba', form='formatted', status='old')
2 read (1, *) nbv, nbt
3 read (1, *) (k, (c(i, k), i=1, 2), refs(k), j=1, nbv)
```

(continues on next page)
msh Files

First, we add the notions of boundary edges

- \( nbbe \) the number of boundary edge
- \( nube(1:2, 1:nbbe) \) an integer array giving the two vertex numbers of boundary edges
- \( refbe(1:nbbe) \) an integer array giving the reference numbers of boundary edges

In Fortran the msh files are read as follows:

```fortran
open (1, file='xxx.msh', form='formatted', status='old')
read (1, *) nbv, nbt, nbbe
read (1, *) ((c(i, k), i=1, 2), refs(k), j=1, nbv)
read (1, *) ((nu(i, k), i=1, 3), reft(k), j=1, nbt)
read (1, *) ((ne(i, k), i=1, 2), refbe(k), j=1, nbbe)
close(1)
```

ftq Files

In Fortran the ftq files are read as follows:

```fortran
open(1, file='xxx.ftq', form='formatted', status='old')
read (1, *) nbv, nbbe, nbv, nbe
read (1, *) (k(j), (nu(i, j), i=1, k(j)), reft(j), j=1, nbe)
read(1, *) ((c(i, k), i=1, 2), refs(k), j=1, nbv)
close(1)
```

where if \( k(j) = 3 \) when the element \( j \) is a triangle and \( k(j) = 4 \) when the the element \( j \) is a quadrilateral.

sol and solb files

With the keyword `savesol`, we can store a scalar functions, a scalar finite element functions, a vector fields, a vector finite element fields, a symmetric tensor and a symmetric finite element tensor.

Such format is used in medit.

Extension file .sol

The first two lines of the file are:

- MeshVersionFormatted 0
- Dimension [DIM](int)

The following fields begin with one of the following keyword: SolAtVertices, SolAtEdges, SolAtTriangles, SolAtQuadrilaterals, SolAtTetrahedra, SolAtPentahedra, SolAtHexahedra.

In each field, we give then in the next line the number of elements in the solutions (SolAtVertices: number of vertices, SolAtTriangles: number of triangles, ...). In other lines, we give the number of solutions, the type of solution (1: scalar, 2: vector, 3: symmetric tensor). And finally, we give the values of the solutions on the elements.

The file must be ended with the keyword End.
The real element of symmetric tensor:

\[
ST^{3d} = \begin{pmatrix}
ST^{3d}_{xx} & ST^{3d}_{xy} & ST^{3d}_{xz} \\
ST^{3d}_{yx} & ST^{3d}_{yy} & ST^{3d}_{yz} \\
ST^{3d}_{zx} & ST^{3d}_{zy} & ST^{3d}_{zz}
\end{pmatrix}
\]

\[
ST^{2d} = \begin{pmatrix}
ST^{2d}_{xx} & ST^{2d}_{xy} \\
ST^{2d}_{yx} & ST^{2d}_{yy}
\end{pmatrix}
\]

(3.31)

stored in the extension .sol are respectively \(ST^{3d}_{xx}, ST^{3d}_{xy}, ST^{3d}_{xz}, ST^{3d}_{yx}, ST^{3d}_{yy}, ST^{3d}_{yz}, ST^{3d}_{zx}, ST^{3d}_{zy}, ST^{3d}_{zz}\) and \(ST^{2d}_{xx}, ST^{2d}_{yx}, ST^{2d}_{yy}\).

An example of field with the keyword SolAtTetrahedra:

```plaintext
SolAtTetrahedra
[Number of tetrahedra](int) [Number of solutions](int) [Type of solution 1](int) ... [Type of solution nt](int)
U_1_l_1(double) ... U_nrs_l_1(double)
... U_l ns_1(double) ... U_(nrs_k)_ns_1(double)
... U_l ns_l(double) ... U_nrs_l ns_1(double)
U_1_l nt(double) ... U_nrs_l nt(double)
... U_l ns nt(double) ... U_(nrs_k)_ns nt(double)
```

- \(ns\) is the number of solutions
- \(typesol_k\), type of the solution number \(k\)
  - \(typesol_k = 1\) the solution \(k\) is scalar
  - \(typesol_k = 2\) the solution \(k\) is vectorial
  - \(typesol_k = 3\) the solution \(k\) is a symmetric tensor or symmetric matrix
- \(nrs_k\) is the number of real to describe solution \(k\)
  - \(nrs_k = 1\) if the solution \(k\) is scalar
  - \(nrs_k = \text{dim}\) if the solution \(k\) is vectorial (\(\text{dim}\) is the dimension of the solution)
  - \(nrs_k = \text{dim}*(\text{dim}+1)/2\) if the solution \(k\) is a symmetric tensor or symmetric matrix
- \(U_i_j_^k\) is a real equal to the value of the component \(i\) of the solution \(k\) at tetrahedron \(j\) on the associated mesh

The format .solb is the same as format .sol but in binary (read/write is faster, storage is less).

A real scalar functions \(f_1\), a vector fields \(\Phi = [\Phi_1, \Phi_2, \Phi_3]\) and a symmetric tensor \(ST^{3d}\) (3.31) at the vertices of the three dimensional mesh Th3 is stored in the file f1PhiTh3.sol using:

```plaintext
savesol("f1PhiST3dTh3.sol", Th3, f1, [Phi(1), Phi(2), Phi(3)], VV3, order=1);
```

where \(VV3 = [ST^{3d}_{xx}, ST^{3d}_{yx}, ST^{3d}_{xy}, ST^{3d}_{yy}, ST^{3d}_{yz}, ST^{3d}_{zy}, ST^{3d}_{zx}, ST^{3d}_{zy}, ST^{3d}_{zz}]\).

For a two dimensional mesh Th, A real scalar functions \(f_2\), a vector fields \(\Psi = [\Psi_1, \Psi_2]\) and a symmetric tensor \(ST^{2d}\) (3.31) at triangles is stored in the file f2PsiST2dTh3.solb using:

```plaintext
savesol("f2PsiST2dTh3.solb", Th, f2, [Psi(1), Psi(2)], VV2, order=0);
```

where \(VV2 = [ST^{2d}_{xx}, ST^{2d}_{yx}, ST^{2d}_{yy}]\).

The arguments of savesol functions are the name of a file, a mesh and solutions. These arguments must be given in this order.
The parameters of this keyword are:

- **order** = 0 is the solution is given at the center of gravity of elements. 1 is the solution is given at the vertices of elements.

In the file, solutions are stored in this order: scalar solutions, vector solutions and finally symmetric tensor solutions.

### 3.9.2 Adding a new finite element

#### Some notations

For a function $f$ taking value in $\mathbb{R}^N$, $N = 1, 2, \ldots$, we define the finite element approximation $\Pi_h f$ of $f$.

Let us denote the number of the degrees of freedom of the finite element by $N_{\text{DoF}}$. Then the $i$-th base $\omega^K_i$ ($i = 0, \cdots, N_{\text{DoF}} - 1$) of the finite element space has the $j$-th component $\omega^K_{ij}$ for $j = 0, \cdots, N - 1$.

The operator $\Pi_h$ is called the interpolator of the finite element.

We have the identity $\omega^K_i = \Pi_h \omega^K_i$.

Formally, the interpolator $\Pi_h$ is constructed by the following formula:

$$\Pi_h f = \sum_{k=0}^{kP1-1} \alpha_k f_{jk}(P_k) \omega^K_{ik}$$  \hspace{1cm} (3.32)

where $P_k$ is a set of $nP Pi$ points.

In the formula (3.32), the list $p_k$, $j_k$, $i_k$ depend just on the type of finite element (not on the element), but the coefficient $\alpha_k$ can be depending on the element.

**Tip:** Classical scalar Lagrange finite element

With the classical scalar Lagrange finite element, we have $kPi = npPi = NbOfNode$ and

- $P_p$ is the point of the nodal points.
- the $\alpha_k = 1$, because we take the value of the function at the point $P_k$.
- $p_k = k$, $j_k = k$ because we have one node per function.
- $j_k = 0$ because $N = 1$.

**Tip:** The Raviart-Thomas finite element

$$RT0_h = \{ v \in H(div) / \forall K \in T_h \quad v|_K(x,y) = [\frac{\alpha^K_{\beta}}{\beta^K} + \gamma^K | y] \}$$  \hspace{1cm} (3.33)

The degrees of freedom are the flux through an edge $e$ of the mesh, where the flux of the function $f : \mathbb{R}^2 \to \mathbb{R}^2$ is $\int_e f \cdot n_e$, $n_e$ is the unit normal of edge $e$ (this implies a orientation of all the edges of the mesh, for example we can use the global numbering of the edge vertices and we just go to small to large number).

To compute this flux, we use a quadrature formula with one point, the middle point of the edge. Consider a triangle $T$ with three vertices $(a, b, c)$.

Let denote the vertices numbers by $i_a, i_b, i_c$, and define the three edge vectors $e^0, e^1, e^2$ by $sgn(i_b - i_c)(b - c)$, $sgn(i_c - i_a)(c - a)$, $sgn(i_a - i_b)(a - b)$.

The three basis functions are:

$$\omega^K_0 = \frac{sgn(i_b - i_a)}{2|T|} (x - a), \quad \omega^K_1 = \frac{sgn(i_c - i_a)}{2|T|} (x - b), \quad \omega^K_2 = \frac{sgn(i_a - i_b)}{2|T|} (x - c),$$
where \(|T|\) is the area of the triangle \(T\).

So we have \(N = 2\), \(k\Pi = 6\); \(np\Pi = 3\); and:

- \(P_p = \left\{ \frac{b+c}{2}, \frac{a+c}{2}, \frac{b+a}{2} \right\}\)
- \(\alpha_0 = -e_2, \alpha_1 = e_1, \alpha_2 = -e_1, \alpha_3 = e_1, \alpha_4 = -e_2, \alpha_5 = e_2\) (effectively, the vector \((-e_2, e_1)\) is orthogonal to the edge \(e^m = (e_1^m, e_2^m)\) with a length equal to the side of the edge or equal to \(\int e^m 1\)).
- \(i_k = \{0, 0, 1, 1, 2, 2\}\),
- \(p_k = \{0, 0, 1, 1, 2, 2\}\), \(j_k = \{0, 1, 0, 1, 0, 0\}\).

### Which class to add?

Add file `FE_ADD.cpp` in directory `FreeFem-sources/src/femlib` for example first to initialize:

```cpp
#include "error.hpp"
#include "rgraph.hpp"
using namespace std;
#include "RNM.hpp"
#include "fem.hpp"
#include "FESpace.hpp"
#include "AddNewFE.h"

namespace Fem2D { ... }
```

Then add a class which derive for public `TypeOfFE` like:

```cpp
class TypeOfFE_RTortho : public TypeOfFE { public:
    static int Data[]; //some numbers
    TypeOfFE_RTortho():
        TypeOfFE(0+3+0, //nb degree of freedom on element
            2, //dimension N of vectorial FE (1 if scalar FE)
            Data, //the array data
            1, //nb of subdivision for plotting
            1, //nb of sub finite element (generaly 1)
            6, //number kPi of coef to build the interpolator
            3, //number npPi of integration point to build interpolator
            0 //an array to store the coef \alpha_k to build interpolator
            //here this array is no constant so we have
            //to rebuilt for each element
            )
    {
        const R2 Pt[] = {R2(0.5, 0.5), R2(0.0, 0.5), R2(0.5, 0.0) }; // the set of Point in hat{K}
        for (int p = 0; p < 3; p++){
            P_Pi_h[p] = Pt[p];
            for (int j = 0; j < 2; j++)
                pij_alpha[kk++] = IPJ(p, p, j);
        }
    } //definition of i_k, p_k, j_k in interpolator

    void FB(const bool *watdd, const Mesh &Th, const Triangle &K,
            const R2 &PHat, RNMK_ &val) const;
```

(continues on next page)
where the array data is formed with the concatenation of five array of size \( N_{\text{DoF}} \) and one array of size \( N \).

This array is:

```c
int TypeOfFE_RTortho::Data[] = {
  // for each df 0, 1, 3:
  3, 4, 5, // the support of the node of the df
  0, 0, 0, // the number of the df on the node
  0, 1, 2, // the node of the df
  0, 0, 0, // the df come from which FE (generally 0)
  0, 1, 2, // which are the df on sub FE
  0, 0
}; // for each component j=0, N-1 it give the sub FE associated
```

where the support is a number 0, 1, 2 for vertex support, 3, 4, 5 for edge support, and finally 6 for element support.

The function to defined the function \( \omega^K_i \), this function return the value of all the basics function or this derivatives in array \( \text{val} \), computed at point \( \hat{P} \) on the reference triangle corresponding to point \( R^2 \hat{P} = K(\hat{P}) \); on the current triangle \( K \).

The index \( i, j, k \) of the array \( \text{val}(i, j, k) \) correspond to:

- \( i \) is the basic function number on finite element \( i \in [0, NoF[ \)
- \( j \) is the value of component \( j \in [0, N[ \)
- \( k \) is the type of computed value \( f(P), dx(f)(P), dy(f)(P), ... i \in [0, \text{last_operatortype}] \).

Note: For optimization, this value is computed only if \( \text{whatd}[k] \) is true, and the numbering is defined with

```c
enum operatortype {
  op_id = 0,
  op_dx = 1, op_dy = 2,
  op_dxx = 3, op_dyy = 4,
  op_dyx = 5, op_dxy = 5,
  op_dz = 6,
  op_dzz = 7,
  op_dzx = 8, op_dxz = 8,
  op_dzy = 9, op_dyz = 9
};
const int last_operatortype = 10;
```

The shape function:

```c
void TypeOfFE_RTortho::FB(const bool *whatd, const Mesh &Th, const Triangle & K, const R2 &P, RNMK_ &val) const
{
    R2 P(K(PHat));
    R2 A(K[0]), B(K[1]), C(K[2]);
    R 10 = 1 - P.x-P.y;
    R 11 = P.x, 12 = P.y;
    assert(val.N() >= 3);
    assert(val.M() == 2);
    val = 0;
}
```

(continues on next page)
The function to defined the coefficient \( \alpha_k \):

```cpp
void TypeOfFE_RT::Pi_h_alpha(const baseFElement &K, KN_<double> &v) const
{
    const Triangle &T(K.T);
    for (int i = 0, k = 0; i < 3; i++){
        R2 E(T.Edge(i));
        R signe = T.EdgeOrientation(i) ;
        v[k++] = signe*E.y;
        v[k++] = -signe*E.x;
    }
}
```

Now, we just need to add a new key work in FreeFEM.

Two way, with static or dynamic link so at the end of the file, we add:

**With dynamic link** it is very simple (see section *Dynamical link*), just add before the end of `FEM2d` namespace:
static TypeOfFE_RTortho The_TypeOfFE_RTortho;
static AddNewFE("RT0Ortho", The_TypeOfFE_RTortho);

Try with ./load.link command in examples++-load/ and see BernardiRaugel.cpp or Morley.cpp new finite element examples.

Otherwise with static link (for expert only), add

To enforce in loading of this new finite element, we have to add the two new lines close to the end of files src/femlib/FESpace.cpp like:

and now you have to change the makefile.

First, create a file FE_ADD.cpp containing all this code, like in file src/femlib/Element_P2h.cpp, after modify the Makefile.am by adding the name of your file to the variable EXTRA_DIST like:

and do in the FreeFEM root directory
autoreconf
./reconfigure
make

For codewarrior compilation add the file in the project an remove the flag in panel PPC linker FreeFm++ Setting Dead-strip Static Initialization Code Flag.

3.9.3 Dynamical link

Now, it’s possible to add built-in functionalities in FreeFEM under the three environments Linux, Windows and MacOS X 10.3 or newer.

It is a good idea to first try the example load.edp in directory example++-load.

You will need to install a compiler (generally gcc/g++ compiler) to compile your function.

- Windows Install the cygwin environment or the mingw one
- MacOs Install the developer tools Xcode on the apple DVD
- Linux/Unix Install the correct compiler (gcc for instance)

Now, assume that you are in a shell window (a cygwin window under Windows) in the directory example++-load.

Note: In the sub directory include, they are all the FreeFEM include file to make the link with FreeFEM.

Note: If you try to load dynamically a file with command load "xxx" - Under Unix (Linux or MacOs), the file xxx.so will be loaded so it must be either in the search directory of routine dlopen (see the environment variable $LD_LIBRARY_PATH) or in the current directory, and the suffix " .so" or the prefix " ./" is automatically added.

- Under Windows, the file xxx.dll will be loaded so it must be in the loadLibrary search directory which includes the directory of the application.

Compilation of your module:

The script ff-c++ compiles and makes the link with FreeFEM, but be careful, the script has no way to know if you try to compile for a pure Windows environment or for a cygwin environment so to build the load module under cygwin you must add the -cygwin parameter.

A first example myfunction.cpp

The following defines a new function call myfunction with no parameter, but using the x, y current value.
using namespace Fem2D;

double myfunction(Stack stack) {
    // to get FreeFEM data
    MeshPoint &mp = *MeshPointStack(stack); // the struct to get x, y, normal, value
    double x = mp.P.x; // get the current x value
    double y = mp.P.y; // get the current y value
    // cout << "x = " << x << " y=" << y << endl;
    return sin(x) * cos(y);
}

Now the Problem is to build the link with FreeFEM, to do that we need two classes, one to call the function myfunction.

All FreeFEM evaluable expression must be a C++ struct/class which derivate from E_F0. By default this expression does not depend of the mesh position, but if they derivate from E_F0mps the expression depends of the mesh position, and for more details see [HECHT2002].

// A class build the link with FreeFEM
// generaly this class are already in AFunction.hpp
// but unfortunately I have no simple function with no parameter
// in FreeFEM depending of the mesh
template<class R>
class OneOperator0s : public OneOperator {
    // the class to define and evaluate a new function
    // It must devive from E_F0 if it is mesh independent
    // or from E_F0mps if it is mesh dependent
    class E_F0_F : public E_F0mps {
        public:
            typedef R (*func)(Stack stack);
            func f; // the poineuter to the function myfunction
            E_F0_F(func ff) : f(ff) {}  
            // the operator evaluation in FreeFEM
            AnyType operator() (Stack stack) const { return SetAny<R>(f(stack)); }
    };

typedef R (*func)(Stack);
    func f;

    public:
        // the function which build the FreeFEM byte code
        E_F0 *code(const basicAC_F0 &) const { return new E_F0_F(f); }  
        // the constructor to say ff is a function without parameter
        // and returning a R
        OneOperator0s(func ff) : OneOperator(map_type[typeid(R).name()],f(ff){
    };

To finish we must add this new function in FreeFEM table, to do that include:

```cpp
void init() {
    Global.Add("myfunction", ", new OneOperator0s<double>(myfunction));
}
```

It will be called automatically at load module time.

To compile and link, use the ff-c++ script:

```
ff-c++ myfunction.cpp
g++ -c -g -Iinclude myfunction.cpp
```
To try the simple example under Linux or MacOS, do `FreeFem++-nw load.edp`

The output must be:

```plaintext
-- FreeFem++ v *.****** (date *** ** *** ****, **:**:** (UTC+0*00))
Load: lg_fem lg_mesh lg_mesh3 eigenvalue
1 : // Example of dynamic function load
2 : // --------------------------------
3 : // $Id$
4 :
5 : load "myfunction"
6 : // dumtable(cout);
7 : mesh Th=square(5,5);
8 : fespace Vh(Th,P1);
9 : Vh uh= myfunction(); // warning do not forget ()
10 : cout << uh[][.min << " " << uh[].max << endl;
11 : cout << " test io ( " << endl;
12 : testio();
13 : cout << ") end test io .. " << endl; sizestack + 1024 =1416 ( 392 )
-- Square mesh : nb vertices =36 , nb triangles = 50 , nb boundary edges 20
0 0.841471
test io ( test cout 3.14159
test cout 512
test cerr 3.14159
test cerr 512 ) end test io ..
times: compile 0.012854s, execution 0.000313s, mpirank:0
CodeAlloc : nb ptr 2715, size :371104 mpirank: 0
Ok: Normal End
```

Under Windows, launch `FreeFEM` with the mouse (or ctrl O) on the example.

**Example: Discrete Fast Fourier Transform**

This will add FFT to `FreeFEM`, taken from `FFTW`. To download and install under `download/include` just go in `download/fftw` and try make.

The 1D dfft (fast discrete fourier transform) for a simple array \( f \) of size \( n \) is defined by the following formula:

\[
dfft(f, \varepsilon)_k = \sum_{j=0}^{n-1} f_j e^{2\pi i k j / n}
\]

The 2D DFFT for an array of size \( N = n \times m \) is:

\[
dfft(f, m, \varepsilon)_{k+nl} = \sum_{j'=0}^{m-1} \sum_{j=0}^{n-1} f_{i+nj} e^{2\pi i (kj/n + lj'/m)}
\]

**Note:** The value \( n \) is given by \( size(f)/m \), and the numbering is row-major order.

So the classical discrete DFFT is \( \hat{f} = dfft(f, -1)/\sqrt{n} \) and the reverse DFFT \( f = dfft(\hat{f}, 1)/\sqrt{n} \).
Note: The 2D Laplace operator is

\[ f(x, y) = \frac{1}{\sqrt{N}} \sum_{j'=0}^{m-1} \sum_{j=0}^{n-1} \hat{f}_{i+j_j} e^{2\pi i (x_j + y_{j'})} \]

and we have

\[ f_{k+nl} = f(k/n, l/m) \]

So

\[ \hat{\Delta} f_{kl} = -((2\pi)^2((\hat{k})^2 + (\hat{l})^2)) \hat{f}_{kl} \]

where \( \hat{k} = k \) if \( k \leq n/2 \) else \( \hat{k} = k - n \) and \( \hat{l} = l \) if \( l \leq m/2 \) else \( \hat{l} = l - m \).

And to have a real function we need all modes to be symmetric around zero, so \( n \) and \( m \) must be odd.

---

### Compile to build a new library

```bash
ff-c++ dfft.cpp ../download/install/lib/libfftw3.a -I../download/install/include
export MACOSX_DEPLOYMENT_TARGET=10.3
g++ -c -Iinclude -I../download/install/include dfft.cpp
g++ -bundle -undefined dynamic_lookup dfft.o -o ./dfft.dylib ../download/install/lib/
```

To test, try *FFT example*.

---

### Load Module for Dervieux P0-P1 Finite Volume Method

The associated edp file is `examples++-load/convect_dervieux.edp`.

See `mat_dervieux.cpp`.

---

### More on Adding a new finite element

First read the *Adding a new finite element section*, we add two new finite elements examples in the directory `examples++-load`.

---

### The Bernardi-Raugel Element

The Bernardi-Raugel finite element is meant to solve the Navier Stokes equations in \( u, p \) formulation; the velocity space \( P_K^{br} \) is minimal to prove the inf-sup condition with piecewise constant pressure by triangle.

The finite element space \( V_h \) is

\[ V_h = \{ u \in H^1(\Omega)^2; \quad \forall K \in T_h, u_{|K} \in P_K^{br} \} \]

where

\[ P_K^{br} = \text{span}\{\lambda^K_i e_k\}_{i=1,2,3,k=1,2} \cup \{\lambda^K_{i+1} n^K_{i+1,2}\}_{i=1,2,3} \]
with notation $4 = 1, 5 = 2$ and where $\lambda^K_i$ are the barycentric coordinates of the triangle $K$, $(e_k)_{k=1,2}$ the canonical basis of $\mathbb{R}^2$ and $n^K_k$ the outer normal of triangle $K$ opposite to vertex $k$.

See BernardiRaugel.cpp.

A way to check the finite element

```cpp
load "BernardiRaugel"
// Macro
// a macro to compute the numerical derivative
macro DD(f, hx, hy) ((f(x1+hx, y1+hy) - f(x1-hx, y1-hy))/(2*(hx+hy)) )
// Mesh
mesh Th = square(1, 1, [10*(x+y/3), 10*(y-x/3)]);
// Parameters
real x1 = 0.7, y1 = 0.9, h = 1e-7;
int it1 = Th(x1, y1).nuTriangle;
// Fespace
fespace Vh2(Th, P2BR);
Vh2 [a1, a2], [b1, b2], [c1, c2];
for (int i = 0; i < Vh2.ndofK; ++i)
    cout << i << " " << Vh2(0,i) << endl;
for (int i = 0; i < Vh2.ndofK; ++i)
{
    a1[] = 0;
    int j = Vh(it1, i);
    a1[][j] = 1;
    plot([a1, a2], wait=1);
    [b1, b2] = [a1, a2]; // do the interpolation
    c1[] = a1[] - b1[];
    cout << " ---------" << i << " " << c1[].max << " " << c1[].min << endl;
    cout << " a = " << a1[] << endl;
    cout << " b = " << b1[] << endl;
    assert(c1[].max < 1e-9 && c1[].min > -1e-9); // check if the interpolation is correct
    // check the derivative and numerical derivative
    cout << " dx(a1)(x1, y1) = " << dx(a1)(x1, y1) << " == " << DD(a1, h, 0) << endl;
    assert( abs(dx(a1)(x1, y1) - DD(a1, h, 0) ) < 1e-5);
    assert( abs(dx(a2)(x1, y1) - DD(a2, h, 0) ) < 1e-5);
    assert( abs(dy(a1)(x1, y1) - DD(a1, 0, h) ) < 1e-5);
    assert( abs(dy(a2)(x1, y1) - DD(a2, 0, h) ) < 1e-5);
}
```

A real example using this finite element, just a small modification of the Navier-Stokes P2-P1 example, just the beginning is change to

```cpp
load "BernardiRaugel"
real s0 = clock();
mesh Th = square(10, 10);
fespace Vh2(Th, P2BR);
```
6 \texttt{fespace Vh(Th, P0);}
7 Vh2 [u1, u2], [up1, up2];
8 Vh2 [v1, v2];

And the plot instruction is also changed because the pressure is constant, and we cannot plot isovalue of piecewise constant functions.

\textbf{The Morley Element}

See the example \texttt{bilapMorley.edp}.

\section{3.10 ffddm}

In the acronym \texttt{ffddm}, \texttt{ff} stands for FreeFEM and \texttt{ddm} for domain decomposition methods. The idea behind \texttt{ffddm} is to simplify the use of parallel solvers in FreeFEM: distributed direct methods and domain decomposition methods.

Parallelism is an important issue because, since about 2004, the clock speed of cores stagnates at 2-3 GHz. The increase in performance is almost entirely due to the increase in the number of cores per processor. All major processor vendors are producing multicore chips and now every machine is a parallel machine. Waiting for the next generation machine does not guarantee anymore a better performance of a software. To keep doubling performance parallelism must double. It implies a huge effort in algorithmic development.

Thanks to \texttt{ffddm}, FreeFEM users have access to high-level functionalities for specifying and solving their finite element problems in parallel. The first task handled by \texttt{ffddm} is the data distribution among the processors. This is done via an overlapping domain decomposition and a related distributed linear algebra. Then, solving a linear system is possible either via an interface to the parallel MUMPS solver or by using domain decomposition methods as preconditioners to the GMRES Krylov method. The \texttt{ffddm} framework makes it easy to use scalable Schwarz methods enhanced by a coarse space correction built either from a coarse mesh or a GenEO (Generalized Eigenvalue in the Overlap) coarse space, see also the book \textit{An Introduction to Domain Decomposition Methods: algorithms, theory, and parallel implementation}. State-of-the-art three level methods are also implemented in \texttt{ffddm}.

The \texttt{ffddm} framework is entirely written in the FreeFEM language and the ‘\texttt{idp}’ scripts can be found here (‘\texttt{ffddm*.idp}’ files). It makes it also a very good tool for learning and prototyping domain decomposition methods without compromising efficiency.

\texttt{ffddm} can also act as a wrapper for the \texttt{HPDDM} library. \texttt{HPDDM} is an efficient implementation of various domain decomposition methods and a variety of Krylov subspace algorithms, with advanced block and recycling methods for solving sequences of linear systems with multiple right-hand sides: GMRES and Block GMRES, CG, Block CG, and Breakdown-Free Block CG, GCRO-DR and Block GCRO-DR. For more details on how to use \texttt{HPDDM} within \texttt{ffddm}, see the \textit{ffddm documentation}.

\textbf{Getting Started}

\begin{verbatim}
macro dimension 2 // EOM    // 2D or 3D
include "ffddm.idp"

mesh Th = square(50,50); // global mesh

// Step 1: Decompose the mesh
ffddmbuildDmesh( M , Th , mpiCommWorld )

// Step 2: Define your finite element
macro def(u) u // EOM
macro init(u) u // EOM

// Step 3: Define your problem
ffddmbuildDfespace( FE , M , real , def , init , P2 )
\end{verbatim}

(continues on next page)
This example solves a Laplace problem in 2D in parallel with a two-level GenEO domain decomposition method. To try this example, just copy and paste the script above in a file ‘test.edp’ and run it on 2 cores with

```
ff-mpirun -np 2 test.edp -wg
```

### 3.10.1 Domain Decomposition (DD)

When the size of a three dimensional problem is large (whatever it means), it is necessary to distribute data among several processors especially for solving linear systems. A natural way is to do it via domain decomposition.

#### Mesh Decomposition

The starting point is a collection of \( N \) sub-meshes \( (T_{h_i})_{i=1}^{N} \) that together form a global mesh

\[
T_h := \bigcup_{i=1}^{N} T_{h_i}.
\]

These meshes may be overlapping or not. This decomposition induces a natural decomposition of the global finite element space \( V_h \) on \( T_h \) into \( N \) local finite element spaces \( (V_{h_i})_{i=1}^{N} \) each of them defined on \( T_{h_i} \).

**Note** By global, we mean that the corresponding structure can be referred to in the code (most often only) by its local values. In computer science term, it corresponds to a distributed data where each piece of data is stored by a MPI process.

#### Distributed Linear Algebra

For a given finite element space \( V_h \), the domain decomposition induces a natural decomposition of the set of the global degrees of freedom (d.o.f.) of \( V_h \) into the \( N \) subsets of d.o.f.’s \( (\mathcal{N}_i)_{i=1}^{N} \) each associated with the local finite element space \( V_{h_i} \). We have thus

\[
\mathcal{N} = \bigcup_{i=1}^{N} \mathcal{N}_i,
\]

but with duplications of some of the d.o.f.’s.

Associated with this decomposition of the set of d.o.f.’s \( \mathcal{N} \), a *distributed vector* is a collection of local vectors \((\mathbf{V}_i)_{i \leq i \leq N}\) so that the values on the duplicated d.o.f.’s are the same.
Note: In mathematical terms, it can be described as follows for a real valued problem. Let $R_i$ be the restriction operator from $\mathbb{R}^{\# \mathcal{N}}$ to $\mathbb{R}^{\# \mathcal{N}_i}$, where $\# \mathcal{N}_i$ denotes the number of elements of $\mathcal{N}_i$. A collection of local vectors $(V_i)_{1 \leq i \leq N} \in \prod_{i=1}^{N} \mathbb{R}^{\# \mathcal{N}_i}$ is a distributed vector iff there exists a global vector $V \in \mathbb{R}^{\# \mathcal{N}}$ such that for all subset $1 \leq i \leq N$, we have:

$$V_i = R_i V.$$ 

We will also say that the collection of local vectors $(V_i)_{1 \leq i \leq N}$ is consistent. For a complex valued problem, simply replace $\mathbb{R}$ with $\mathbb{C}$.

### Partition of Unity Matrices (POUM)

Let $(D_i)_{1 \leq i \leq N}$ be square diagonal matrices of size $\# \mathcal{N}_i$ which form a partition of unity in the sense that:

$$Id = \sum_{i=1}^{N} R_i^T D_i R_i \text{ in } \mathbb{R}^{\# \mathcal{N} \times \# \mathcal{N}}.$$ 

For instance if a degree of freedom is shared by $k$ subdomains defining the corresponding entry of the diagonal matrix $D$ to be $1/k$ yields partition of unity matrices. The matrices $R_i$ and $D_i$ are the heart of distributed linear algebra.

### Distributed scalar product

For two global vectors $U$ and $V$ of size $\# \mathcal{N}$, the formula for the scalar product $V^T U = (U, V)$ in terms of their distributed vector counterparts makes use of the partition of unity matrices $(D_i)_{1 \leq i \leq N}$ introduced above:

$$(U, V) = \left(U, \sum_{i=1}^{N} R_i^T D_i R_i V\right) = \sum_{i=1}^{N} (R_i U, D_i R_i V) = \sum_{i=1}^{N} (U_i, D_i V_i).$$

Local scalar products are performed concurrently. Thus, the implementation is parallel except for the sum which corresponds to a MPI_Reduce call across the $N$ MPI processes. Note also that the implementation relies on the knowledge of a partition of unity so that the FreeFEM syntax is $dscalprod(D_i,u,v)$ or equivalently $myFEprefix#scalprod(u,v)$ where $myFEprefix$ is a user defined prefix for the finite element space decomposition, see the ffddm documentation.

### Update

From a collection of local vectors $(U_i)_{1 \leq i \leq N}$, it is possible ensure consistency of the duplicated data by modifying the distributed vector $(U_i)_{1 \leq i \leq N}$ by calling the function $myFEprefix#update(Ui, TRUE)$ where $myFEprefix$ is the user defined prefix that refers to the finite element space decomposition. This function performs the following operation for all $1 \leq i \leq N$:

$$U_i \leftarrow R_i \sum_{j=1}^{N} R_j^T D_j U_j$$

Note: The implementation corresponds to

$$U_i \leftarrow R_i \sum_{j=1}^{N} R_j^T D_j U_j = D_i U_i + \sum_{j \in \mathcal{O}(i)} R_i R_j^T D_j U_j$$

where $\mathcal{O}(i)$ is the set of neighbors of subdomain $i$. Therefore, the matrix vector product is computed in three steps:
• concurrent computing of $D_j U_j$ for all $1 \leq j \leq N$;
• neighbor to neighbor MPI-communications from subdomain $j$ to subdomain $i \left( R_i R_i^T \right)$;
• concurrent sum of neighbor contributions.

### Distributed Matrix and Vector resulting from a variational formulation

The discretization of a variational formulation on the global mesh $\mathcal{T}_h$ yields a global matrix $A$ and a global right hand side $\text{RHS}$. Thanks to the sparsity of finite element matrices for partial differential equations and thanks to the overlap between subdomains, the knowledge of the local matrix $R_i A R_i^T$ on each subdomain $1 \leq i \leq N$ is sufficient to perform the matrix-vector product $A \times U$ for any global vector $U$. Once the problem has been set up by a call to `ffddmssetupOperator(myprefix, myFEprefix, myVarf)`, the matrix-vector product is performed by calling the function `myprefix#A(Ui)` where `myprefix` is a user defined prefix that refers to the problem at hand which itself implicitly refers to the triplet (domain decomposition, finite element, variational formulation). See more on problem definition in this documentation and more on distributed linear algebra in chapter 8 of “An Introduction to Domain Decomposition Methods: algorithms, theory and parallel implementation” SIAM 2015.

### Distributed Linear Solvers

In many cases, we are interested in the solution of the problem in terms of the vector of d.o.f.’s $X$ that satisfies:

$$A X = \text{RHS}.$$  

`ffddm` offers two parallel solvers: direct factorization and iterative preconditioned solvers via Schwarz domain decomposition methods.

### Distributed Direct Solvers

In order to benefit from the sparsity of the matrix arising from a finite element discretization of a partial differential equation, a variant of Gauss elimination, the frontal method, that automatically avoids a large number of operations involving zero terms was developed. A frontal solver builds a $LU$ or Cholesky decomposition of a sparse matrix given as the assembly of element matrices by eliminating equations only on a subset of elements at a time. This subset is called the **front** and it is essentially the transition region between the part of the system already finished and the part not touched yet. These methods are basically sequential since the unknowns are processed the one after another or one front after another. In order to benefit from multicore processors, a multifrontal solver is an improvement of the frontal solver that uses several independent fronts at the same time. The fronts can be worked on by different processors, which enables parallel computing. `ffddm` provides an interface to the parallel sparse direct solver MUMPS. These methods have the advantage to be very robust and to have a predictable cost. The main drawback is the memory requirement which can be prohibitive especially for three-dimensional problems.

### Schwarz methods

These methods are part of the large family of preconditioned iterative solvers. When considering the solve of the equation $A X = \text{RHS}$, a preconditioner is a linear operator that approximates the inverse of $A$ and whose cost of the associated matrix-vector product is much cheaper than solving the original linear system. It enables to accelerate the solution of the latter with Krylov type methods such as the conjugate gradient (in the symmetric positive definite case), GMRES or BiCGSTAB in the general case. Two options are possible.

Left preconditioning: the preconditioner is applied to the left of the equation

$$M^{-1} A X = M^{-1} \text{RHS}.$$  

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and the Krylov method is applied to the left preconditioned system with a residual that is preconditioner dependent.

Right preconditioning: the preconditioner is inserted on the right of the operator:

\[ A M^{-1} Y = \text{RHS} \quad \text{where} \quad X = M^{-1} Y. \]

and the Krylov method is applied to the right preconditioned system with a residual that is preconditioner independent.

In both cases, if the preconditioner is efficient the number of iterations to get a converged solution is much smaller than the number of iterations of the Krylov method applied to the original equation \( A X = \text{RHS} \). Although right preconditioning seems more intricate, it is much safer to use since the convergence is checked on a residual that does not depend on the preconditioner.

In the sequel, we consider the solution of the equation \( A X = \text{RHS} \) preconditioned by domain decomposition methods and with a flexible GMRES Krylov method which is thus necessarily right preconditioned.

**Restricted Additive Schwarz (RAS)**

The RAS preconditioner reads:

\[ M_{\text{RAS}}^{-1} := \sum_{j=1}^{N} R_j^T D_j (R_j A R_j^T)^{-1} R_j, \]

where for each subdomain \( j \) the restriction matrix \( R_j \) and the partition of unity matrix \( D_j \) to a global right hand side \( \text{RHS} \) is detailed below. Recall that this global vector is distributed among processes via the local vectors \( (\text{RHS}_i)_{i=1}^{N} \). Let \( A_j \) denote the local matrix \( (R_j A R_j^T) \). The local vector in subdomain \( i \) resulting from the matrix vector product \( M_{\text{RAS}}^{-1} \text{RHS} \) consists in computing:

\[ R_i \sum_{j=1}^{N} R_j^T D_j A_j^{-1} \text{RHS}_j = D_i A_i^{-1} \text{RHS}_i + \sum_{j \in \mathcal{O}(i)} (R_i R_j^T) D_j A_j^{-1} \text{RHS}_j. \]

This task is performed by first solving concurrently on all subdomains a linear system for \( Y_j \) for all \( 1 \leq j \leq N \):

\[ A_j Y_j = \text{RHS}_j. \]

Each local vector \( Y_j \) is weighted by the partition of unity matrix \( D_j \). Then data transfers between neighboring subdomains implement the \( R_i R_j^T D_j Y_j \) formula. The contribution from neighboring subdomains are summed locally. This pattern is very similar to that of the **update** procedure.

**Optimized Restricted Additive Schwarz (ORAS)**

The ORAS preconditioner may be seen as a variant of the RAS preconditioner. It reads:

\[ M_{\text{ORAS}}^{-1} := \sum_{j=1}^{N} R_j^T D_j B_j^{-1} R_j, \]

where \( B_j \) are local matrices of size \( \#N_j \times \#N_j \) for \( 1 \leq j \leq N \). This variant is very useful when dealing with wave propagation phenomena such as Helmholtz problems in acoustics or Maxwell system in the frequency domain for electromagnetism. Defining \( B_j \) as the discretization of the physical equation with impedance conditions on the boundary of the subdomain has been proved to be a good choice.
Two level methods

The RAS and ORAS methods are called a one-level method in the sense that sub-domains only interact with their direct neighbors. For some problems such as Darcy problems or static elasticity problems and when the number of subdomains is large, such one-level methods may suffer from a slow convergence. The fix is to add to the preconditioner an auxiliary coarse problem that couples all subdomains at each iteration and is inexpensive to calculate.

In mathematical terms, we first choose a full rank rectangular matrix $Z \in \mathbb{R}^{\#N \times NC}$ where $NC \ll \#N$ denotes the dimension of the coarse space spanned by the columns of $Z$. We also pick a coarse matrix $A_C \in \mathbb{R}^{NC \times NC}$. A generic one-level method preconditioner $M_1^{-1}$ is enriched by a solve on the coarse space. The simplest correction formula is additive:

$$M_2^{-1} := Z A_C^{-1} Z^T + M_1^{-1}$$

Other correction formulas are given in documentation.

We consider two ways to build $Z$ and thus the coarse space and the coarse problem $A_C$, see below Coarse Mesh and GenEO.

Coarse Mesh

A first possibility is to discretize the problem on a coarse mesh, following the same principle as multi-grid methods. For 3-D problems, a coarsening of the mesh size by a factor 2, reduces by a factor $2^3 = 8$ the size of the coarse problem which is then easier to solve by a direct method. Then, $Z$ is the interpolation matrix from the coarse finite element space to the fine one.

GenEO

For highly heterogeneous or anisotropic problems, two level methods based on coarse meshes might fail and a more sophisticated construction must be used. A provable robust coarse space called GenEO is built by first solving the following local generalized eigenvalue problem in parallel for each subdomain $1 \leq i \leq N$, where $A_{i}^{\text{Neu}}$ denotes the local matrix resulting from the variational formulation:

$$D_i A_i D_i V_{i,k} = \lambda_{i,k} A_{i}^{\text{Neu}} V_{i,k}$$

The eigenvectors selected to enter the coarse space correspond to eigenvalues $\lambda_{i,k} \geq \tau$, where the threshold parameter $\tau$ is user-defined. The precise formulas are given in this documentation. From a mathematical point of view, it has been proved that for a symmetric positive definite matrix $A$, the spectrum of the preconditioned by the two-level method with a GenEO coarse space lies in the interval $[\frac{1}{1 + k_1 \tau}, k_0]$.

Note A heuristic that justifies this construction is as follows. We first introduce the Additive Schwarz method (ASM) which can be seen as a symmetrized variant of the RAS preconditioner:

$$M_{ASM}^{-1} := \sum_{j=1}^{N} R_j^T A_j^{-1} R_j .$$

It can be proved that the lower bound for the eigenvalue of $M_{ASM}^{-1}$ $A$ is close to zero (which is bad for convergence) whereas the upper bound depends only on the number of neighbors of a subdomain (which is good for convergence).

Second, we also introduce the following preconditioner $M_{NN}^{-1}$:

$$M_{NN}^{-1} := \sum_{1 \leq j \leq N} D_i (A_j^{\text{Neu}})^{-1} D_j .$$
We have a very good lower bound for the preconditioned operator $M^{-1}_{NN} A$ that does not depend on the number of subdomains but only on the maximum multiplicity of intersections $k_1$ (which is good for convergence). But the upper bound for this preconditioner is very large (which is bad for convergence).

Now, if we compare formulas for $M^{-1}_{NN}$ and $M^{-1}_{ASM}$, we may suspect that vectors $V_{ik}$ for which $D_i (A_i^{Neu})^{-1} D_i V_{ik}$ and $A_i^{-1} V_{ik}$ have very different values are responsible for the slow convergence and should contribute to the coarse space. This is a way to interpret the above generalized eigenvalue problem which controls the lower bound of the two-level preconditioned system.

### 3.10.2 ffddm documentation

#### Minimal example

```plaintext
// EOM // 2D or 3D
macro dimension 3

include "ffddm.idp"

load "msh3"

int[int] LL = [2,2, 1,2, 2,2];
mesh3 ThGlobal = cube(10, 10, 10, [x, y, z], label = LL); // global mesh

macro grad(u) [dx(u), dy(u), dz(u)] // three-dimensional gradient

macro Varf(varfName, meshName, VhName)
    varf varfName(u,v) = int3d(meshName)(grad(u)'* grad(v)) + int3d(meshName)(v) + on(1, u = 1.0);

// Domain decomposition
ffddmbuildDmesh( LapMesh, ThGlobal, mpiCommWorld )

macro def(i)i // scalar field definition
macro init(i)i // scalar field initialization
ffddmbuildDfespace( LapFE, LapMesh, real, def, init, P1 )

ffddmsetupOperator( Lap, LapFE, Varf )

real[int] rhsi(0);
ffddmbuildrhs( Lap, Varf, rhsi )
LapFEVhi def(ui);

//Direct solve
ui[] = Lapdirectsolve(rhsi);
Lapwritesummary
ffddmplot(LapFE,ui,"u")
```

#### Overlapping mesh decomposition

```plaintext
ffddmbuildDmesh(prmesh,Th,comm)
```
decomposes the mesh $\text{Th}$ into overlapping submeshes. The mesh will be distributed over the mpi ranks of communicator $\text{comm}$. This will create and expose variables whose names will be prefixed by $\text{prmesh}$, see below ($\#$ is the concatenation operator). The way the initial mesh $\text{Th}$ is partitioned depends on the value of $\text{ffddmpartitioner}$.

The size of the overlap between subdomains (its width in terms of number of mesh elements) is given by $\text{ffddmoverlap}$.

The level of refinement of the resulting submeshes with respect to the input mesh $\text{Th}$ is given by $\text{ffddmsplit}$.

If $\text{ffddmexclude} \neq 0$, the first $\text{ffddmpCS}$ mpi ranks of $\text{comm}$ will be excluded from the spatial domain decomposition, in order to dedicate them later to the coarse problem (for two-level preconditioners).

The label of the new border of the submeshes (the interface between the subdomains) is given by $\text{ffddminterfacelabel}$.

defines:

- $\text{int prmesh\#npart}$ number of subdomains for this decomposition; should be equal to $\text{mpiSize} (\text{comm}) - \text{ffddmexclude} * \text{ffddmpCS}$
- $\text{int prmesh\#pCS}$ equal to $\text{ffddmpCS}$
- $\text{int prmesh\#exclude}$ equal to $\text{ffddmexclude}$
- $\text{int prmesh\#excluded}$ true if $\text{ffddmexclude}$ is true ($\neq 0$) and $\text{mpiRank}(\text{comm}) < \text{prmesh\#pCS}$. In this case, this mpi rank will be excluded from the spatial domain decomposition and will only work on the coarse problem.
- $\text{mpiComm prmesh\#commddm}$ mpi communicator for ranks participating in the spatial domain decomposition (ranks 0 to $\text{prmesh\#npart-1}$ in $\text{comm}$ if $\text{prmesh\#exclude}$ is false, ranks $\text{prmesh\#pCS}$ to $\text{prmesh\#pCS}$+$\text{prmesh\#npart-1}$ otherwise)
- $\text{mpiComm prmesh\#commCS}$ mpi communicator for ranks participating in the assembly and resolution of the coarse problem for two-level preconditioners (ranks 0 to $\text{prmesh\#pCS}$-1 in $\text{comm}$)
- $\text{mpiComm prmesh\#commself}$ self mpi communicator (this mpi rank only), used for factorizing local matrices
- $\text{meshN[int] prmesh\#aTh}$ array (size $\text{prmesh\#npart}$) of local meshes of the subdomains. In the standard parallel case, only the local mesh for this mpi rank $\text{prmesh\#aTh}[\text{mpiRank}(\text{prmesh\#commddm})]$ is defined (unless this mpi rank is excluded from the spatial domain decomposition, i.e. $\text{prmesh\#excluded} = 1$, see below). In the sequential case, all local meshes are defined.
- $\text{meshN prmesh\#Thi}$ the local mesh of the subdomain for this mpi rank, i.e. $\text{prmesh\#aTh}[\text{mpiRank}(\text{prmesh\#commddm})]$ in the parallel case
- $\text{int prmesh\#numberIntersection}$ the number of neighbors for this mpi rank
- $\text{int[int] prmesh\#arrayIntersection}$ the list of neighbor ranks in $\text{prmesh\#commddm}$ for this mpi rank

Remark for sequential use (see `-seqddm`):

- $\text{meshN[int] prmesh\#aTh}$ array (size $\text{prmesh\#npart}$) of local meshes of the subdomains

Local finite element spaces

```cpp
ffddmbuildDfespace(prfe,prmesh,scalar,def,\textit{init},Pk)
```

builds the local finite element spaces and associated distributed operators on top of the mesh decomposition $\text{prmesh}$. This will create and expose variables whose names will be prefixed by $\text{prfe}$, see below. It is assumed that $\text{ffddmbuildDmesh}$ has already been called with prefix $\text{prmesh}$ in order to build the mesh decomposition.

The local finite element spaces of type $\text{Pk}$ (where $\text{Pk}$ is the type of finite element: $\text{P1}$, $\text{[P2,P2,P1]}$, ...) are defined on the local meshes of the subdomains based on the mesh decomposition previously created with prefix $\text{prmesh}$.  

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**scalar** determines the type of data for this finite element: *real* or *complex*.

Two macros, **def** and **init**, are needed: **def** specifies how to define a finite element function in the finite element space \( P_k \), and **init** specifies how to interpolate a scalar function onto the (possibly multiple) components of \( P_k \). Two examples are given below:

For scalar \( P_2 \) finite elements and complex-valued problems:

```plaintext
1 macro def(u) u // EOM
2 macro init(u) u // EOM
3 ffddmbuildDfespace(myFEprefix,mymeshprefix,complex,def,init,P2)
```

For vectorial \([P_2,P_2,P_1]\) finite elements and real-valued problems:

```plaintext
1 macro def(u) [u, u#, u#B, u#C] // EOM
2 macro init(u) [u, u, u] // EOM
3 ffddmbuildDfespace(myFEprefix,mymeshprefix,real,def,init,[P2,P2,P1])
```

In practice, this builds the necessary distributed operators associated to the finite element space: the local partition of unity functions \( (D_i)_{i=1,\ldots,N} \) (see prfe#Dk and prfe#Dih below) as well as the function prfe#update (see below) which synchronizes local vectors \( (u_i)_{i=1,\ldots,N} \) between neighboring subdomains, performing the equivalent of

\[
u_i = R_i (\sum_{j=1}^{N} R_T^{ij} u_j) \quad \text{or} \quad u_i = R_i (\sum_{j=1}^{N} R_T^{ij} D_j u_j)
\]

in a distributed parallel environment.

prfe#scalprod (see below) performs the parallel scalar product for vectors defined on this finite element.

**defines:**

- **prfe#prmesh** macro, saves the parent prefix **prmesh** of the mesh decomposition
- **prfe#K** macro, saves the type of data **scalar** for this finite element space (real or complex)
- **func prfe#fPk** saves the type of finite element **Pk**, e.g. **P1, [P2,P2,P1], ...**
- **fespace prfe#Vhi** the local finite element space for this mpi rank, defined on the local mesh **prmesh#Thi**
- **int prfe#Ndofglobal** the total number of degrees of freedom \( n \) for this finite element discretization
- **prfe#mdef** macro, saves the macro **def** giving the definition of a finite element function in the finite element space **Pk**
- **prfe#minit** macro, saves the macro **init** specifying how to interpolate a scalar function onto the (possibly multiple) components of a finite element function of **Pk**. This is used to create the local partition of unity function in **prfe#Vhi**, by interpolating the local P1 partition of unity function onto the components of **prfe#Vhi**. For non Lagrange finite element spaces (e.g. **RT0, Edge03d, ...**), see **ffddmbuildDfespaceEdge**.
- **prfe#K[1][1] prfe#Dk** array (size **prmesh#npart**) of local partition of unity vectors in the subdomains, equivalent to \( (D_i)_{i=1,\ldots,N} \). In the standard parallel case, only the local partition of unity vector for this mpi rank \( prfe#Dk[mpiRank(prmesh#commddm)] \) is defined (unless this mpi rank is excluded from the spatial domain decomposition, i.e. \( prmesh#excluded = 1 \)). In the sequential case, all local partition of unity vectors are defined.
- **matrix<prfe#K>[int] prfe#Dih** array (size **prmesh#npart**) similar to **prfe#Dk** but in matrix form, allowing for easier matrix-matrix multiplications. **prfe#Dih[i]** is a diagonal matrix, with the diagonal equal to **prfe#Dk[i]**.
- **fespace prfe#Vhglob** the global finite element space defined on the global mesh **prmesh#Thglob**. Defined only if **-noGlob** is not used.
- **matrix<prfe#K>[int] prfe#Rih** array (size **prmesh#npart**) of restriction matrices from the global finite element space to the local finite element spaces on the local submeshes of the subdomains. In the standard parallel case, only the restriction matrix for this mpi rank **prfe#Rih[mpiRank(prmesh#commdm)]** is
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defined (unless this mpi rank is excluded from the spatial domain decomposition, i.e., prmesh#excluded = 1). In the sequential case, all restriction matrices are defined. The restriction matrices prfe#Rih are defined only if -noGlob is not used.

- **func int prfe#update(scalar[int] ui, bool scale)** The function prfe#update synchronizes the local vector ui between subdomains by exchanging the values of ui shared with neighboring subdomains (in the overlap region) using point-to-point MPI communications. If scale is true, ui is multiplied by the local partition of unity beforehand. This is equivalent to $u_i = R_i \left( \sum_{j=1}^{N} R_j^T u_j \right)$ when scale is false and $u_i = R_i \left( \sum_{j=1}^{N} R_j^T D_j u_j \right)$ when scale is true.

- **func scalar prfe#scalprod(scalar[int] ai, scalar[int] bi)** The function prfe#scalprod computes the global scalar product of two vectors whose local restriction to the subdomain of this mpi rank are ai and bi. The result is computed as $\sum_{j=1}^{N} (D_j a_j, b_j)$.

**Define the problem to solve**

```plaintext
ffddmsetupOperator(pr,prfe,Varf)
```

builds the distributed operator associated to the variational problem given by Varf, on top of the distributed finite element space prfe. This will create and expose variables whose names will be prefixed by pr, see below. It is assumed that ffddmbuildDfespace has already been called with prefix prfe in order to define the distributed finite element space.

In practice, this builds the so-called local ‘Dirichlet’ matrices $A_i = R_i A R_i^T$, the restrictions of the global operator $A$ to the subdomains (see pr#aRd below). The matrices correspond to the discretization of the bilinear form given by the macro Varf, which represents the abstract variational form of the problem. These matrices are then used to implement the action of the global operator $A$ on a local vector (the parallel matrix-vector product with $A$), see pr#A below.

At this point, we already have the necessary data to be able to solve the problem with a parallel direct solver (MUMPS), which is the purpose of the function pr#directsolve (see below). See ffddmbuildrhs for building the right-hand side.

The macro Varf is required to have three parameters: the name of the variational form, the mesh, and the finite element space. The variational form given in this ‘abstract’ format will then be used by ffddm to assemble the discrete operators by setting the appropriate mesh and finite element space as parameters. An example is given below:

```plaintext
macro myVarf(varfName, meshName, VhName)
  varf varfName(u,v) = int3d(meshName)(grad(u)'* grad(v)) + on(1, u = 1.0);
  // EOM
ffddmsetupOperator(myprefix,myFEprefix,myVarf)
```

**Remark** In this simple example, the third parameter VhName is not used. However, for more complex cases such as non-linear or time dependent problems where the problem depends on a solution computed at a previous step, it is useful to know for which discrete finite element space the variational form is being used. See for example TODO defines:

- **pr#prfe macro**, saves the parent prefix prfe of the finite element space
- **int pr#verbosity** the level of verbosity for this problem, initialized with the value of ffddmverbosity
- **pr#writesummary macro**, prints a summary of timings for this problem, such as the time spent to assemble local matrices or solve the linear system.
- **matrix<prfe#K> pr#Aglobal** the global matrix $A$ corresponding to the discretization of the variational form given by the macro Varf on the global finite element space prfe#Vhglob. Defined only in the sequential case.
• matrix<prfe#K>[int] pr#aRd array (size prfe#prmesh#npart) of so-called local ‘Dirichlet’ matrices in the subdomains; these are the restrictions of the global operator to the subdomains, equivalent to $A_i = R_i A R_i^T$ with $A$ the global matrix corresponding to the discretization of the variational form given by the macro Varf on the global finite element space. In the standard parallel case, only the local matrix for this mpi rank pr#aRd[mpiRank(prmesh#commddm)] is defined (unless this mpi rank is excluded from the spatial domain decomposition, i.e. prmesh#excluded = 1). In the sequential case, all local matrices are defined.

• func prfe#K[int] pr#A(prfe#K[int] &ui) The function pr#A computes the parallel matrix-vector product, i.e. the action of the global operator $A$ on the local vector $u_i$. The computation is equivalent to $R_i (\sum_{j=1}^N R_j^T D_j A_j u_j)$ and is performed in parallel using local matrices pr#aRd and the function prfe#update. In the sequential case, the global matrix pr#Aglobal is used instead.

• func prfe#K[int] pr#AT(prfe#K[int] &ui) Similarly to pr#A, The function pr#AT computes the action of $A^T$, the transpose of the global operator $A$, on $u_i$.

• func prfe#K[int] pr#directsolve(prfe#K[int]& rhsi) The function pr#directsolve allows to solve the linear system $Ax = b$ in parallel using the parallel direct solver MUMPS. The matrix is given to MUMPS in distributed form through the local matrices pr#aRd. The input rhsi is given as a distributed vector (rhsi is the restriction of the global right-hand side $b$ to the subdomain of this mpi rank, see ffddmbuildrhs) and the returned vector is local as well.

**Remark: rectangular operators**

It is possible to define a non-square distributed operator where the variational form takes two different finite element spaces of unknown and test functions. This is done through macro ffddmsetupOperatorRect which takes two FE prefixes (which must be defined on the same mesh prefix), see below:

```cpp
macro myVarf(varName, meshName, VhName)
  varf varName([u, uB, uC], [q]) = int3d(meshName)(div(u) * q);
// EOM
ffddmsetupOperatorRect(myprefix,myFEprefixV,myFEprefixP,myVarf)
```

```cpp
ffddmbuildrhs(pr,Varfrhs,rhs)
```

builds the right-hand side associated to the variational form given by Varfrhs for the problem corresponding to prefix pr. The resulting right-hand side vector rhs corresponds to the discretization of the abstract linear form given by the macro Varfrhs (see ffddmsetupOperator for more details on how to define the abstract variational form as a macro).

The input vector rhs is resized and contains the resulting local right-hand side $R_i b$, the restriction of the global right-hand side $b$ to the subdomain of this mpi rank. In the sequential case, the global right-hand side vector $b$ is assembled instead.

An example is given below:

```cpp
macro myVarfrhs(varName, meshName, VhName)
  varf varName(u,v) = intN(meshName)(v) + on(1, u = 1.0);
// EOM
real[int] rhsi(0);
ffddmbuildrhs(myprefix,myVarfrhs,rhsi)
```

**One level preconditioners**

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ffddmsetupPrecond(pr, VarfPrec)

builds the one level preconditioner for problem pr. This will create and expose variables whose names will be prefixed by pr, see below. It is assumed that ffddmsetupOperator has already been called with prefix pr in order to define the problem to solve.

In practice, this builds and performs the factorization of the local matrices used in the one level preconditioner. The local matrices depend on the choice of ffddmprecond and VarfPrec, see pr#ARbelow.

defines:

- string pr#prec equal to ffddmprecond. Sets the type of one level preconditioner $M^{-1}$ to be used: “asm” (Additive Schwarz), “ras” (Restricted Additive Schwarz), “oras” (Optimized Restricted Additive Schwarz), “soras” (Symmetric Optimized Restricted Additive Schwarz) or “none” (no preconditioner).
- matrix<pr#prfe#K>[int] pr#AR array (size pr#ARmesh#npart) of local matrices used for the one level preconditioner. Each mpi rank of the spatial domain decomposition performs the $LU$ (or $LDL^T$) factorization of the local matrix corresponding to its subdomain using the direct solver MUMPS.

- If VarfPrec is not a previously defined macro (just put null for example), the matrices pr#AR are set to be equal to the so-called local ‘Dirichlet’ matrices pr#ARd (see ffddmsetupOperator). This is for the classical ASM preconditioner as $M^{-1} = M^{-1}_ASM = \sum_{i=1}^{N} R_i^T A_i^{-1} R_i$, or classical RAS preconditioner $M^{-1} = M^{-1}_RAS = \sum_{i=1}^{N} R_i^T D_i A_i^{-1} R_i$ (it is assumed that ffddmprecond is equal to “asm” or “ras”).

- If VarfPrec is a macro, it is assumed that VarfPrec defines an abstract bilinear form (see ffddmsetupOperator for more details on how to define the abstract variational form as a macro).

  - If ffddmprecond is equal to “asm” or “ras”, the matrices pr#AR will be assembled as local ‘Dirichlet’ matrices in the same manner as pr#ARd, but using the bilinear form defined by VarfPrec instead. This defines the ASM preconditioner as $M^{-1} = M^{-1}_ASM = \sum_{i=1}^{N} R_i^T (A_i^{Prec})^{-1} R_i$ and the RAS preconditioner as $M^{-1} = M^{-1}_RAS = \sum_{i=1}^{N} R_i^T D_i (A_i^{Prec})^{-1} R_i$, where $A_i^{Prec} = R_i A_i^{Prec} R_i^T$.

  - If ffddmprecond is equal to “oras” or “soras”, the matrices pr#AR will correspond to the discretization of the variational form VarfPrec in the subdomains $\Omega_i$. In particular, various boundary conditions can be imposed at the interface between subdomains (corresponding to mesh boundary of label ffddminterfacelabel set by the parent call to ffddmbuildDmesh), such as Optimized Robin boundary conditions. We note the ORAS preconditioner as $M^{-1} = M^{-1}_ORAS = \sum_{i=1}^{N} R_i^T D_i (B_i^{Prec})^{-1} R_i$ and the SORAS preconditioner as $M^{-1} = M^{-1}_SORAS = \sum_{i=1}^{N} R_i^T D_i (B_i^{Prec})^{-1} D_i R_i$.

- func pr#AR[K>[int] pr#PREC1(pr#AR[K>[int] &ui) The function pr#PREC1 computes the parallel application of the one level preconditioner $M^{-1}$, i.e. the action of $M^{-1}$ on the local vector $u_i$. In the sequential case, it computes the action of $M^{-1}$ on a global vector. The action of the inverse of local matrices pr#ARd is computed by forward-backward substitution using their $LU$ (or $LDL^T$) decomposition.

- func pr#AR[K>[int] pr#PREC(pr#AR[K>[int] &ui) The function pr#PREC corresponds to the action of the preconditioner $M^{-1}$ for problem pr. It coincides with the one level preconditioner pr#PREC1 after the call to ffddmsetupPrecond. If a second level is subsequently added (see the next section about Two level preconditioners), it will then coincide with the two level preconditioner $M^{-2}$ (see pr#PREC2level).

- func pr#AR[K>[int] pr#FGMRES(pr#AR[K]> x0i, pr#AR[K]>b, real eps, int nbiter, string sprec) The function pr#FGMRES allows to solve the linear system $Ax = b$ in parallel using the flexible GMRES method preconditioned by $M^{-1}$. The action of the global operator $A$ is given by pr#, the action of the preconditioner $M^{-1}$ is given by pr#PREC and the scalar products are computed by pr#scalprod. More details are given in the section Solving the linear system.

Two level preconditioners

The main ingredient of a two level preconditioner is the so-called ‘coarse’ space matrix $Z$. 
$Z$ is a rectangular matrix of size $n \times n_c$, where usually $n_c \ll n$.

$Z$ is used to build the 'coarse space operator' $E = Z^T A Z$, a square matrix of size $n_c \times n_c$. We can then define the 'coarse space correction operator' $Q = Z E^{-1} Z^T = Z (Z^T A Z)^{-1} Z^T$, which can then be used to enrich the one level preconditioner through a correction formula. The simplest one is the additive coarse correction: $M_2^{-1} = M_1^{-1} + Q$. See pr\#corr below for all other available correction formulas.

There are multiple ways to define a relevant coarse space $Z$ for different classes of problems. \texttt{ffddmgeneosetup} defines a coarse space correction operator by building the GenEO coarse space, while \texttt{ffddmcoarsemeshsetup} builds the coarse space using a coarse mesh.

After a call to either \texttt{ffddmgeneosetup} or \texttt{ffddmcoarsemeshsetup}, the following variables and functions are set up:

1. \texttt{int pr\#ncoarsespace} the size of the coarse space $n_c$.
2. \texttt{string pr\#corr} initialized with the value of \texttt{ffddmcorrection}. Specifies the type of coarse correction formula to use for the two level preconditioner. The possible values are:
   
   - "AD": Additive, $M^{-1} = M_2^{-1} = M_1^{-1} + Q$
   - "BNN": Balancing Neumann-Neumann, $M^{-1} = M_2^{-1} = (I - QA) M_1^{-1} (I - AQ) + Q$
   - "ADEF1": Adapted Deflation Variant 1, $M^{-1} = M_2^{-1} = M_1^{-1} (I - AQ) + Q$
   - "ADEF2": Adapted Deflation Variant 2, $M^{-1} = M_2^{-1} = (I - QA) M_1^{-1} + Q$
   - "RBNN1": Reduced Balancing Variant 1, $M^{-1} = M_2^{-1} = (I - QA) M_1^{-1}$
   - "RBNN2": Reduced Balancing Variant 2, $M^{-1} = M_2^{-1} = (I - AQ) M_1^{-1}$
   - "none": no coarse correction, $M^{-1} = M_2^{-1} = M_1^{-1}$

   Note that AD, ADEF1 and RBNN2 only require one application of $Q$, while BNN, ADEF2 and RBNN1 require two. The default coarse correction is \texttt{ADEF1}, which is cheaper and generally as robust as BNN or ADEF2.

3. \texttt{func pr\#prfe\#K[int] pr\#Q(pr\#prfe\#K[int] &ui)} The function \texttt{pr\#Q} computes the parallel application of the coarse correction operator $Q$, i.e. the action of $Q = Z E^{-1} Z^T$ on the local vector $u_i$. In the sequential case, it computes the action of $Q$ on a global vector. The implementation differs depending on the method used to build the coarse space (with GenEO or using a coarse mesh), but the idea is the same: the action of the transpose of the distributed operator $Z$ on the distributed vector $u_i$ is computed in parallel, with the contribution of all subdomains being gathered in a vector of size $n_c$ in the mpi process of rank 0. The action of the inverse of the coarse space operator $E$ is then computed by forward-backward substitution using its $LU$ (or $LDL^T$) decomposition previously computed by the first \texttt{func pr\#prfe\#prmesh\#pCS} ranks of the mpi communicator. The result is then sent back to all subdomains to perform the last application of $Q$ and obtain the resulting local vector in each subdomain.

4. \texttt{func pr\#prfe\#K[int] pr\#PREC2level(pr\#prfe\#K[int] &ui)} The function \texttt{pr\#PREC2level} computes the parallel application of the two level preconditioner $M_2^{-1}$, i.e. the action of $M_2^{-1}$ on the local vector $u_i$. In the sequential case, it computes the action of $M_2^{-1}$ on a global vector. The two level preconditioner depends on the choice of the coarse correction formula which is determined by \texttt{pr\#corr}, see above.

### Building the GenEO coarse space

```cpp
ffddmgeneosetup(pr, Varf)
```

This builds the GenEO coarse space for problem \texttt{pr}. This will create and expose variables whose names will be prefixed by \texttt{pr}, see below. It is assumed that \texttt{ffddmsetupPrecond} has already been called for prefix \texttt{pr} in order to define the one level preconditioner for problem \texttt{pr}. The GenEO coarse space is $Z = \{ R_i^T D_i V_{i,k} \}_{i=1,...,N, \lambda_{i,k} \geq \tau}$, where $V_{i,k}$ are eigenvectors corresponding to eigenvalues $\lambda_{i,k}$ of the following local generalized eigenvalue problem in subdomain $i$:

$$D_i A_i^\text{Neu} V_{i,k} = \lambda_{i,k} A_i^\text{Neu} V_{i,k},$$
where $A_{i}^{\text{Neu}}$ is the local Neumann matrix of subdomain $i$ (with Neumann boundary conditions at the subdomain interface).

In practice, this builds and factorizes the local Neumann matrices $A_{i}^{\text{Neu}}$ corresponding to the abstract bilinear form given by the macro $\text{Varf}$ (see $\text{ffddmsetupOperator}$ for more details on how to define the abstract variational form as a macro). In the GenEO method, the abstract bilinear form $\text{Varf}$ is assumed to be the same as the one used to define the problem $\text{pr}$ through the previous call to $\text{ffddmsetupOperator}$. The local generalized eigenvalue problem is then solved in each subdomain to find the eigenvectors $V_{i,k}$ corresponding to the largest eigenvalues $\lambda_{i,k}$ (see $\text{pr#Z}$ below). The number of computed eigenvectors $\nu$ is given by $\text{ffddmnu}$. The eigenvectors selected to enter $Z$ correspond to eigenvalues $\lambda_{i,k}$ larger than $\tau$, where the threshold parameter $\tau$ is given by $\text{ffddmtau}$. If $\text{ffddmtau} = 0$, all $\text{ffddmnu}$ eigenvectors are selected. Finally, the coarse space operator $E = Z^{T}AZ$ is assembled and factorized (see $\text{pr#E}$ below).

defines:

- $\text{pr#prfe#K[\text{int}][\text{int}]} \text{pr#Z}$ array of local eigenvectors $Z_{i,k} = D_{i}V_{i,k}$ obtained by solving the local generalized eigenvalue problem above in the subdomain of this mpi rank using Arpack. The number of computed eigenvectors $\nu$ is given by $\text{ffddmnu}$. The eigenvectors selected to enter $Z$ correspond to eigenvalues $\lambda_{i,k}$ larger than $\tau$, where the threshold parameter $\tau$ is given by $\text{ffddmtau}$. If $\text{ffddmtau} = 0$, all $\text{ffddmnu}$ eigenvectors are selected.

- $\text{matrix}<\text{pr#prfe#K}> \text{pr#E}$ the coarse space operator $E = Z^{T}AZ$. The matrix $\text{pr#E}$ is assembled in parallel and is factorized by the parallel direct solver MUMPS using the first $\text{pr#prfe#prmesh#pCS}$ ranks of the mpi communicator, with mpi rank 0 as the master process. The number of mpi processes dedicated to the coarse problem is set by the underlying mesh decomposition of problem $\text{pr}$, which also specifies if these mpi ranks are excluded from the spatial decomposition or not. These parameters are set by $\text{ffddmpCS}$ and $\text{ffddmexclude}$ when calling $\text{ffddmbuildDmesh}$ (see $\text{ffddmbuildDmesh}$ for more details).

Building the coarse space from a coarse mesh

```
ffddmcoarsemeshsetup(pr,Thc,VarfEprec,VarfAprec)
```

builds the coarse space for problem $\text{pr}$ from a coarse problem which corresponds to the discretization of a variational form on a coarser mesh $\text{Thc}$ of $\Omega$. This will create and expose variables whose names will be prefixed by $\text{pr}$, see below. It is assumed that $\text{ffddmsetupPrecond}$ has already been called for prefix $\text{pr}$ in order to define the one level preconditioner for problem $\text{pr}$. The abstract variational form for the coarse problem can differ from the original problem $\text{pr}$ and is given by macro $\text{VarfEprec}$ (see $\text{ffddmsetupOperator}$ for more details on how to define the abstract variational form as a macro). For example, absorption can be added in the preconditioner for wave propagation problems, see examples for Helmholtz and Maxwell equations in the Examples section.

The coarse space $Z$ corresponds to the interpolation operator from the coarse finite element space to the original finite element space of the problem. Thus, the coarse space operator $E = Z^{T}A^{\text{Eprec}}Z$ corresponds to the matrix of the problem given by $\text{VarfEprec}$ discretized on the coarse mesh $\text{Thc}$ and is assembled as such.

Similarly, $\text{VarfAprec}$ specifies the global operator involved in multiplicative coarse correction formulas. For example, $M_{2}^{-1}A_{DEFI} = M_{1}^{-1}(I - A^{\text{Aprec}}Q) + Q$ (where $Q = ZE^{-1}Z^{T}$). $A^{\text{Aprec}}$ defaults to $A$ if $\text{VarfAprec}$ is not a valid macro (you can put $\text{null}$ for example).

defines:

- $\text{meshN} \text{pr#ThCoarse}$ the coarse mesh $\text{Thc}$
- $\text{fespace} \text{pr#VhCoarse}$ the coarse finite element space of type $\text{pr#prfe#fPk}$ defined on the coarse mesh $\text{pr#ThCoarse}$
- $\text{matrix}<\text{pr#prfe#K}> \text{pr#AglobEprec}$ the global matrix $A^{\text{Aprec}}$ corresponding to the discretization of the variational form given by the macro $\text{VarfAprec}$ on the global finite element space $\text{pr#prfe#Vhglob}$. 

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Defined only in the sequential case. `pr#AglobEprec` is equal to `pr#Aglobal` if `VarfAprec` is not a valid macro.

- `matrix<pr#prfe#K> pr#aRdEprec` the local ‘Dirichlet’ matrix corresponding to `VarfAprec`; it is the local restriction of the global operator $A^{Aprec}$ to the subdomain, equivalent to $A^{Aprec}_i = R_i A^{Aprec} R_i^T$ with $A^{Aprec}$ the global matrix corresponding to the discretization of the variational form given by the macro `VarfAprec` on the global finite element space. Defined only if this mpi rank is not excluded from the spatial domain decomposition, i.e. `prmesh#excluded = 0`. `pr#aRdEprec` is equal to `pr#aRd[mpiRank(prmesh#commddm)]` if `VarfAprec` is not a valid macro.

- `func pr#prfe#K[int] pr#AEprec(pr#prfe#K[int] &ui)` The function `pr#AEprec` computes the parallel matrix-vector product, i.e. the action of the global operator $A^{Aprec}$ on the local vector $u_i$. The computation is equivalent to $R_i (\sum_{j=1}^{N} R_j^T D_j A^{Aprec}_j u_j)$ and is performed in parallel using local matrices `pr#aRdEprec` and the function `pr#prfe#update`. In the sequential case, the global matrix `pr#AglobEprec` is used instead.

- `matrix<pr#prfe#K> pr#ZCM` the interpolation operator $Z$ from the coarse finite element space `pr#VhCoarse` to the global finite element space `pr#Vhglob`. Defined only in the sequential case.

- `matrix<pr#prfe#K> pr#ZCMi` the local interpolation operator $Z_i$ from the coarse finite element space `pr#VhCoarse` to the local finite element space `pr#Vhi`. Defined only if this mpi rank is not excluded from the spatial domain decomposition, i.e. `prmesh#excluded = 0`. `pr#ZCMi` is used for the parallel application of $Z$ and $Z^T$.

- `matrix<pr#prfe#K> pr#ECM` the coarse space operator $E = Z^T A^{Aprec} Z$. The matrix `pr#ECM` is assembled by discretizing the variational form given by `VarfEprec` on the coarse mesh and factorized by the parallel direct solver `MUMPS` using the first `pr#prfe#prmesh#pcs` ranks of the mpi communicator, with mpi rank 0 as the master process. The number of mpi processes dedicated to the coarse problem is set by the underlying mesh decomposition of problem `pr`, which also specifies if these mpi ranks are excluded from the spatial decomposition or not. These parameters are set by `ffddmpcs` and `ffddmexclude` when calling `ffddmbuildDmesh` (see `ffddmbuildDmesh` for more details).

### Solving the linear system

```cpp
func pr#prfe#K[int] pr#fGMRES(pr#prfe#K[int] & x0i, pr#prfe#K[int] & bi, real eps, int sp, int maxit, string sp)
```
solves the linear system for problem `pr` using the flexible GMRES algorithm with preconditioner $M^{-1}$ (corresponding to `pr#PREC`). Returns the local vector corresponding to the restriction of the solution to `pr#Vhi`. `x0i` and `bi` are local distributed vectors corresponding respectively to the initial guess and the right-hand side (see `ffddmbuildrhs`). `eps` is the stopping criterion in terms of the relative decrease in residual norm. If `eps < 0`, the residual norm itself is used instead. `maxit` sets the maximum number of iterations. `sp` selects between the "left" or "right" preconditioning variants: left preconditioned GMRES solves $M^{-1}Ax = M^{-1}b$, while right preconditioned GMRES solves $AM^{-1}y = b$ for $y$, with $x = M^{-1}y$.

### Using HPDDDM within ffddm

`ffddm` allows you to use HPDDDM to solve your problem, effectively replacing the `ffddm` implementation of all parallel linear algebra computations. `ffddm` can then be viewed as a finite element interface for HPDDDM.

You can use HPDDDM features unavailable in `ffddm` such as advanced Krylov subspace methods implementing block and recycling techniques.

To switch to HPDDDM, simply define the macro `pr#withhpddm` before using `ffddmsetupOperator`. You can then pass HPDDDM options with command-line arguments or directly to the underlying HPDDDM operator `pr#hpddmOP`. Options need to be prefixed by the operator prefix.

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You can also choose to replace only the Krylov solver, by defining the macro pr\#withhpddmKrylov before using \texttt{ffddmsetupOperator}. Doing so, a call to pr\#fGMRES will call the HPDDM Krylov solver, with \texttt{ffddm} providing the operator and preconditioner through pr\#A and pr\#PREC. You can then pass HPDDM options to the Krylov solver through command-line arguments:

```plaintext
macro PBwithhpddmkrylov() // EOM
ffddmsetupOperator( PB , FE , Varf )
```

For example, using restarted GCRO-DR(40) and recycling 10 Ritz vectors at each restart:

```plaintext
ff-mpirun -np 4 test.edp -wg -hpddm_krylov_method gcrodr -hpddm_recycle 10 -ffddm_gmres_restart 40
```

An example can be found in \texttt{Helmholtz-2d-HPDDM-BGMRES.edp}, see the \textit{Examples} section.

### Advanced use

#### Local finite element spaces for non Lagrange finite elements

For Lagrange finite elements, the partition of unity \( (D_i)_{i=1,...,N} \) (see \texttt{prfe\#Dk} and \texttt{prfe\#Dih}) is built by interpolating the local P1 partition of unity function onto the components of the Pk finite element space \texttt{prfe\#Vhi}. For non Lagrange finite element spaces, such as Raviart–Thomas or Nédélec edge elements, the definition of the degrees of freedom can be more involved, and interpolating the P1 partition of unity functions directly is inappropriate. The idea is then to use a “pseudo” finite element \texttt{Pkpart} derived from Pk which is suitable for interpolating the P1 partition of unity, in the sense that it will produce a partition of unity for Pk.

For example, for first-order Nédélec edge elements (\texttt{Edge03d}), whose degrees of freedom are the circulations along the edges, we define the “pseudo” finite element \texttt{Edge03ds0} which can be seen as a scalar Lagrange counterpart: the numbering of the degrees of freedom is the same, but they correspond to the value at the edge midpoints.

For Lagrange finite elements, the distributed finite element spaces are built using \texttt{ffddmbuildDfespace}. Here you must use \texttt{ffddmbuildDfespaceEdge}, which builds the distributed finite element space using a “pseudo” finite element to build the partition of unity:

```plaintext
ffddmbuildDfespaceEdge(prfe,prmesh,scalar,def,\texttt{init},Pk,defpart,initpart,Pkpart)
```

where macros \texttt{defpart} and \texttt{initpart} specify how to define and interpolate a function in the ‘pseudo’ finite element space \texttt{Pkpart}, similar to \texttt{def} and \texttt{init} for \texttt{Pk}.

An example with first-order Nédélec edge elements (\texttt{Edge03d} + \texttt{Edge03ds0}) for Maxwell equations can be found in \texttt{Maxwell-3d-simple.edp}, see the \textit{Examples} section.

#### Inexact coarse solves for two level methods

We have seen in the \textit{Two level preconditioners section} that two level methods produce a ‘coarse space operator’ \( E \) that needs to be inverted at each iteration. By default the coarse space operator matrix is factorized by the direct solver \texttt{MUMPS}. This can become a bottleneck and hinder scalability for large problems, where \( E \) can become too large to be factorized efficiently. To remedy this, we can instead opt to use an iterative method to solve the coarse problem at each iteration. Moreover, in order to retain robustness, a DD preconditioner can be used to solve the inner coarse problem more efficiently.
Coarse mesh and inexact coarse solve

When the coarse problem comes from a coarse mesh discretization, a natural way to do inexact coarse solve is to use a one level domain decomposition method on the coarse problem, with the same subdomain partitioning for the coarse and fine meshes. This means that each processor is associated to one spatial subdomain and hosts the two local (nested) coarse and fine submeshes corresponding to this subdomain, as well as the corresponding local matrices for the two discretizations. This natural choice offers interesting benefits:

- We naturally recover a load-balanced parallel implementation, provided that the initial partitioning is balanced.
- The communication pattern between neighboring subdomains is the same for the coarse and fine discretizations.
- The assembly and the application of the interpolation operator $Z$ (and $Z^T$) between the fine and the coarse spaces can be computed locally in each subdomain and require no communication.

In `ffddm`, the first step is to build the two nested mesh decompositions using `ffddmbuildDmeshNested`:

```plaintext
ffddmbuildDmeshNested(prmesh, Thc, s, comm)
```

decomposes the coarse mesh $Thc$ into overlapping submeshes and creates the fine decomposition by locally refining submeshes by a factor of $s$, i.e. splitting each mesh element into $s^d$ elements, $s \geq 1$. This will create and expose variables corresponding to both decompositions, prefixed by `prmesh` for the fine mesh and by `prmesh#Coarse` for the coarse mesh (see `ffddmbuildDmesh`). It also sets the integer variable `prmesh#binexactCS` to 1, which specifies that any two level method defined on mesh prefix `prmesh` will use inexact coarse solves.

The distributed finite element spaces, operators and preconditioners can then be defined for both decompositions. Here is an example where the coarse problem is solved using a one level method:

```plaintext
ffddmbuildDmeshNested(M, Thc, 3, mpiCommWorld)
ffddmbuildDfespace(FE, M, real, def, init, Pk)
ffddmbuildDfespace(FECoarse, MCoarse, real, def, init, Pk)

// coarse operator (Varf of E):
ffddmsetupOperator(PBCoarse, FECoarse, VarfEprec)
// one level preconditioner for the coarse problem:
ffddmsetupPrecond(PBCoarse, VarfPrecC)

// operator for the fine problem:
ffddmsetupOperator(PB, FE, Varf)
// one level preconditioner for the fine problem:
ffddmsetupPrecond(PB, VarfPrec)

// add the second level:
ffddmcoarsemeshsetup(PB, Thc, VarfEprec, null)

[...]
u[] = PBfGMRES(x0, rhs, 1.e-6, 200, "right");
```

Remarks:

- Note that the different prefixes need to match: prefixes for the coarse decomposition have to be those of the fine decomposition, appended with `Coarse`.
- The operator and preconditioner for the coarse problem have to be defined before those of the fine problem, because the `pr#Q` function is actually defined by `ffddmsetupPrecond` and involves a call to `pr#CoarsefGMRES` (which is defined by `ffddmsetupPrecond` for the coarse problem) for the iterative solution of the coarse problem if `pr#prfe#prmesh#binexactCS \neq 0`. 
In this case, `ffddmcoarsemeshsetup` does not use `Thc` or `VarfEprec` and only builds the local interpolation matrices between fine and coarse local finite element spaces `pr#prfe#Vhi` and `pr#prfe#CoarseVhi` to be able to apply $Z$ and $Z^T$.

The GMRES tolerance for the inner solution of the coarse problem is set by `ffddminexactCStol` and is equal to 0.1 by default.

In practice, these methods can give good results for wave propagation problems, where the addition of artificial absorption in the preconditioner helps with the convergence of the one level method for the inner solution of the coarse problem. You can find an example for Maxwell equations in `Maxwell_CobraCavity.edp`, see the Examples section. More details can be found here and in


### 3.10.3 Parameters

**Command-line arguments**

- `-ffddm_verbosity N`, the level of verbosity of `ffddm`, see `ffddmverbosity` (default 3).
- `-seqddm N` use `ffddm` in sequential mode, with N the number of subdomains.
- `-noGlob` if present, do not define any global quantity (such as saving the global mesh for plotting or building the global restriction matrices). Cannot be used in sequential mode or with plotting.
- `-ffddm_partitioner N` specifies how to partition the initial domain, see `ffddmpartitioner` (default 1, `metis`).
- `-ffddm_overlap N` specifies the width of the overlap region between subdomains, see `ffddmoverlap` (default 1).
- `-ffddm_master_p N` number of master processes for the coarse problem (for two level preconditioners), see `ffddmpCS` (default 1).
- `-ffddm_master_exclude 0|1` exclude master processes from the domain decomposition, see `ffddmexclude` (default 0).
- `-ffddm_split N`, level of refinement of the local submeshes with respect to the initial global mesh, see `ffddmsplit` (default 1).
- `-ffddm_schwarz_method S`, specifies the type of one level preconditioner $M_I^{-1}$: “asm” (Additive Schwarz), “ras” (Restricted Additive Schwarz), “oras” (Optimized Restricted Additive Schwarz), “soras” (Symmetric Optimized Restricted Additive Schwarz) or “none” (no preconditioner), see `ffddmprecond` (default “ras”).
- `-ffddm_geneo_nu N`, number of local eigenvectors to compute in each subdomain when solving the local generalized eigenvalue problem for the GenEO method, see `ffddmnu` (default 20).
- `-ffddm_geneo_threshold R`, threshold parameter for selecting local eigenvectors when solving the local generalized eigenvalue problems for the GenEO method, see `ffddmtau` (default 0.5). If the command-line parameter `-ffddm_geneo_nu N` is used, then `ffddmtau` is initialized to 0.
- `-ffddm_schwarz_coarse_correction S`, specifies the coarse correction formula to use for the two level preconditioner: “AD” (Additive), “BNN” (Balancing Neumann-Neumann), “ADEF1” (Adapted Deflation Variant 1), “ADEF2” (Adapted Deflation Variant 2), “RBNN1” (Reduced Balancing Variant 1), “RBNN2” (Reduced Balancing Variant 2) or “none” (no coarse correction), see `ffddmcorrection` (default “ADEF1”).
- `-ffddm_inexactCS_tol R`, specifies the GMRES tolerance for the inner solution of the coarse problem when using a two level method with approximate coarse solves, see `ffddminexactCStol` (default 0.1).
Global parameters

- **ffddmverbosity** initialized by command-line argument `-ffddmverbosity N`, specifies the level of verbosity of `ffddm` (default 3).

- **ffddmpartitioner** initialized by command-line argument `-ffddmpartitioner N`, specifies how to partition the initial domain:
  - N=0: user-defined partition through the definition of a macro, see `ffddmbuildDmesh`
  - N=1: use the automatic graph partitioner `metis` (default)
  - N=2: use the automatic graph partitioner `scotch`

- **ffddmoverlap** initialized by command-line argument `-ffddmoverlap N`, specifies the number of layers of mesh elements in the overlap region between subdomains N >= 1 (default 1). **Remark** The actual width of the overlap region between subdomains is 2N, since each subdomain is extended by N layers of elements in a symmetric way.

- **ffddminterfacelabel** the label of the new border of the subdomain meshes (the interface between the subdomains) (default 10). Used for imposing problem-dependent boundary conditions at the interface between subdomains for the preconditioner, for example optimized Robin boundary conditions (see ORAS).

- **ffddmpCSC** initialized by command-line argument `-ffddmpCSC N`, number of mpi processes used for the assembly and resolution of the coarse problem for two level preconditioners (default 1).

- **ffddmexclude** initialized by command-line argument `-ffddmexclude N`, 0 or 1 (default 0). If true, mpi ranks participating in the assembly and resolution of the coarse problem for two level preconditioners will be excluded from the spatial domain decomposition and will only work on the coarse problem.

- **ffddmsplit** initialized by command-line argument `-ffddmsplit N`, level of refinement of the local submeshes with respect to the initial global mesh (default 1). This is useful for large problems, where we want to avoid working with a very large global mesh. The idea is to start from a coarser global mesh, and generate finer local meshes in parallel during the mesh decomposition step in order to reach the desired level of refinement for the subdomains. For example, calling `ffddmbuildDmesh` with `ffddmsplit = 3` will generate local submeshes where each mesh element of the initial mesh is split into $3^d$ elements.

- **ffddmprecond** initialized by command-line argument `-ffddm_schwarz_method S`, specifies the type of one level preconditioner $M_l^{-1}$ to build when calling `ffddmsetupPrecond`: “asm” (Additive Schwarz), “ras” (Restricted Additive Schwarz), “oras” (Optimized Restricted Additive Schwarz), “soras” (Symmetric Optimized Restricted Additive Schwarz) or “none” (no preconditioner). Default is “ras”. See `ffddmsetupPrecond` for more details.

- **ffddmnu** initialized by command-line argument `-ffddm_geneo_nu N`, number of local eigenvectors to compute in each subdomain when solving the local generalized eigenvalue problem for the GenEO method (default 20). See `ffddmsetupGenEOset` for more details.

- **ffddmtau** initialized by command-line argument `-ffddm_geneo_threshold R`, threshold parameter for selecting local eigenvectors when solving the local generalized eigenvalue problems for the GenEO method (default 0.5). If the command-line parameter `-ffddm_geneo_nu N` is used, then `ffddmtau` is initialized to 0. See `ffddmsetupGenEOset` for more details.

- **ffddmcurrence** initialized by command-line argument `-ffddm_schwarz_coarse_correction S`, specifies the coarse correction formula to use for the two level preconditioner: “AD” (Additive), “BNN” (Balancing Neumann-Neumann), “ADEF1” (Adapted Deflation Variant 1), “ADEF2” (Adapted Deflation Variant 2), “RBNN1” (Reduced Balancing Variant 1), “RBNN2” (Reduced Balancing Variant 2) or “none” (no coarse correction). Default is “ADEF1”. See the section about Two level preconditioners for more details.

- **ffddminexactCStol** initialized by command-line argument `-ffddminexactCS_tol R`, GMRES tolerance for the inner solution of the coarse problem when using a two level method with approximate coarse solves (default 0.1). See the section about Approximate coarse solves for two level methods for more details.
3.10.4 Tutorial

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What is ffddm?

- ffddm implements a class of parallel solvers in FreeFEM: overlapping Schwarz domain decomposition methods.
- The entire ffddm framework is written in the FreeFEM language ffddm aims at simplifying the use of parallel solvers in FreeFEM. You can find the ffddm scripts here (‘ffddm*.idp’ files) and examples here.
- ffddm provides a set of high-level macros and functions to:
  - handle data distribution: distributed meshes and linear algebra
  - build DD preconditioners for your variational problems
  - solve your problem using preconditioned Krylov methods
- ffddm implements scalable two level Schwarz methods, with a coarse space correction built either from a coarse mesh or a GenEO coarse space. Ongoing research: approximate coarse solves and three level methods.
- ffddm can also act as a wrapper for the HPDDM library. HPDDM is an efficient C++11 implementation of various domain decomposition methods and Krylov subspace algorithms with advanced block and recycling techniques. More details on how to use HPDDM within ffddm here.

Why Domain Decomposition Methods?

How can we solve a large sparse linear system \( Au = b \in \mathbb{R}^n \)?

- Memory consumption
- Robustness
- Parallelizable

Direct Methods

Iterative Methods

DDM “Hybrid” methods

Step 1: Decompose the mesh

See documentation

Build a collection of \( N \) overlapping sub-meshes \( (Th_i)_{i=1}^N \) from the global mesh \( Th \)
freeFEM documentation, release 4.6

1. \texttt{ffddmbuildDmesh( prmesh, ThGlobal, comm )}

- mesh distributed over the MPI processes of communicator \texttt{comm}
- initial mesh \texttt{ThGlobal} partitioned with \textit{metis} by default
- size of the overlap given by \texttt{ffddmoverlap} (default 1)

\texttt{prmesh#Thi} is the local mesh of the subdomain for each mpi process

```
macro dimension 2// EOM    // 2D or 3D
include "ffddm.idp"

mesh ThGlobal = square(100,100);    // global mesh

// Step 1: Decompose the mesh
ffddmbuildDmesh( M, ThGlobal, mpiCommWorld )
medit("Th"+mpirank, MThi);
```

Copy and paste this to a file `test.edp` and run it:
```
ff-mpirun -np 2 test.edp -wg
```

### Step 2: Define your finite element

See documentation
```
ffddmbuildDfespace( prfe, prmesh, scalar, def, init, Pk )
```

builds the local finite element spaces and associated distributed operators on top of the mesh decomposition \texttt{prmesh}

- \textbf{scalar}: type of data for this finite element: \textit{real} or \textit{complex}
- \textbf{Pk}: your type of finite element: P1, [P2,P2,P1], ...
- \textbf{def, init}: macros specifying how to define and initialize a \textbf{Pk} FE function

\texttt{prfe#Vhi} is the local FE space defined on \texttt{prmesh#Thi} for each mpi process

Example for P2 complex:

```
3.10. ffddm
```

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Distributed vectors and restriction operators

Natural decomposition of the set of d.o.f.’s \( \mathcal{N} \) of \( Vh \) into the \( N \) subsets of d.o.f.’s \( (\mathcal{N}_i)_{i=1}^N \) each associated with the local FE space \( Vh_i \)

\[ \mathcal{N} = \bigcup_{i=1}^N \mathcal{N}_i, \]

but with duplications of the d.o.f.’s in the overlap

**Definition.** A distributed vector is a collection of local vectors \( (\mathbf{V}_i)_{1 \leq i \leq N} \) so that the values on the duplicated d.o.f.’s are the same:

\[ \mathbf{V}_i = R_i \mathbf{V}, \quad i = 1, \ldots, N \]

where \( \mathbf{V} \) is the corresponding global vector and \( R_i \) is the restriction operator from \( \mathcal{N} \) into \( \mathcal{N}_i \)

**Remark** \( R_i^T \) is the extension operator: extension by 0 from \( \mathcal{N}_i \) into \( \mathcal{N} \)
Partition of unity

\[
I = \sum_{i=1}^{N} R_i^T D_i R_i
\]

\((D_i)_{1\leq i \leq N}\) are square diagonal matrices of size \(\#N_i\)

\[
V = \sum_{i=1}^{N} R_i^T D_i R_i V = \sum_{i=1}^{N} R_i^T D_i V_i
\]

Data exchange between neighbors

```cpp
func prfe#update(K[int] vi, bool scale)
```

synchronizes local vectors \(V_i\) between subdomains ⇒ exchange the values of mathbf\(V_i\) shared with neighbors in the overlap region

\[
V_i \leftarrow R_i \left( \sum_{j=1}^{N} R_j^T D_j V_j \right) = D_i V_i + \sum_{j \in \mathcal{O}(i)} R_i R_j^T D_j V_j
\]

where \(\mathcal{O}(i)\) is the set of neighbors of subdomain \(i\). Exchange operators \(R_i R_j^T\) correspond to neighbor-to-neighbor MPI communications

```cpp
FUpdate(vi, false);
```

```cpp
V_i \leftarrow R_i \left( \sum_{j=1}^{N} R_j^T V_j \right)
```

```cpp
FUpdate(vi, true);
```

\[
V_i \leftarrow R_i \left( \sum_{j=1}^{N} R_j^T D_j V_j \right)
\]
### Step 3: Define your problem

See documentation

```plaintext
ffddmsetupOperator( pr, prfe, Varf )
```

builds the distributed operator associated to your variational form on top of the distributed FE `prfe`

**Varf** is a macro defining your abstract variational form

```plaintext
macro Varf(varfName, meshName, VhName)
    varf varfName(u,v) = int2d(meshName)(grad(u)'* grad(v)) + int2d(meshName)(f*v) + on(1, u = 0); // EOM
```

⇒ assemble local ‘Dirichlet’ matrices $A_i = R_i A R_i^T$

$$A = \sum_{i=1}^{N} R_i^T D_i A_i R_i$$

**Warning:** only true because $D_i R_i A = D_i A_i R_i$ due to the fact that $D_i$ vanishes at the interface !!

`pr#A` applies $A$ to a distributed vector: $U_i \leftarrow R_i \sum_{j=1}^{N} R_j^T D_j A_j V_j$

⇒ multiply by $A_i + prfe#update$
```plaintext
macro dimension 2 // EOM // 2D or 3D

include "ffddm.idp"

mesh ThGlobal = square(100,100); // global mesh

// Step 1: Decompose the mesh
ffddmbuildDmesh( M , ThGlobal , mpiCommWorld )

// Step 2: Define your finite element
macro def(u) u // EOM
macro init(u) u // EOM
ffddmbuildDfespace( FE , M , real , def , init , P2 )

// Step 3: Define your problem
macro grad(u) [dx(u), dy(u)] // EOM
macro Varf(varfName, meshName, VhName)
  varf varfName(u,v) = int2d(meshName)(grad(u)'* grad(v))
                 + int2d(meshName)(1*v) + on(1, u = 0); // EOM
ffddmsetupOperator( PB , FE , Varf )

FEVhi ui, bi;
ffddmbuildrhs( PB , Varf , bi[] )
ui[] = PBA(bi[]);
ffddmplot(FE, ui, "A*b")
```

Summary so far: translating your sequential **FreeFEM** script

**Step 1: Decompose the mesh**

See documentation

```plaintext
mesh Th = square(100,100);
```

**Step 2: Define your finite element**

See documentation

```plaintext
fespace Vh(Th, P1);
```

**Step 3: Define your problem**

See documentation

```plaintext
varf Pb(u, v) = ...;
matrix A = Pb(Vh, Vh);
```
FreeFEM Documentation, Release 4.6

```c
macro Varf(varfName, meshName, VhName)
  varf varfName(u,v) = ... // EOM
  ffddmsetupOperator(PB, FE, Varf)
```

Solve the linear system

See documentation

```c
u[] = A^-1 * b[];
ui[] = PBdirectsolve(bi[]);
```

Solve the linear system with the parallel direct solver **MUMPS**

See documentation

```c
func K[int] pr#directsolve(K[int]& bi)

We have $A$ and $b$ in distributed form, we can solve the linear system $Au = b$ using the parallel direct solver **MUMPS**

```c
// Solve the problem using the direct parallel solver MUMPS
ui[] = PBdirectsolve(bi[]);
ffddmplot(FE, ui, "u")
```

**Step 4: Define the one level DD preconditioner**

See documentation

```c
ffddmsetupPrecond( pr, VarfPrec )
```

builds the one level preconditioner for problem `pr`.

By default it is the *Restricted Additive Schwarz (RAS)* preconditioner:

$$M_i^{-1} = M_{RAS}^{-1} = \sum_{i=1}^{N} R_i^T D_i A_i^{-1} R_i \quad \text{with} \quad A_i = R_i A R_i^T$$

**Setup step**: compute the $LU$ (or $LDL^T$) factorization of local matrices $A_i$

`pr#PRECI` applies $M_i^{-1}$ to a distributed vector: $U_i \leftarrow R_i \sum_{j=1}^{N} R_j^T D_j A_j^{-1} V_i$

⇒ apply $A_i^{-1}$ (forward/backward substitutions) + `prfe#update`

**Step 5: Solve the linear system with preconditioned GMRES**

See documentation

```c
func K[int] pr#fGMRES(K[int]& x0i, K[int]& bi, real eps, int itmax, string sp)
```

solves the linear system with flexible GMRES with DD preconditioner $M^{-1}$
• \(x_0\): initial guess  
• \(b_i\): right-hand side  
• \(\varepsilon\): relative tolerance  
• \(\text{itmax}\): maximum number of iterations  
• \(\text{sp}\): “left” or “right” preconditioning

**left preconditioning**

solve \(M^{-1}Ax = M^{-1}b\)

**right preconditioning**

solve \(AM^{-1}y = b\)

\[\Rightarrow x = M^{-1}y\]

```plaintext
macro dimension 2 // EOM // 2D or 3D
include "ffddm.idp"

mesh ThGlobal = square(100,100); // global mesh
ffddmbuildDmesh( M , ThGlobal , mpiCommWorld ) // Step 2: Define your finite element
macro def(u) u // EOM
macro init(u) u // EOM
ffddmbuildDfespace( FE , M , real , def , init , P2 ) // Step 2: Define your finite element
macro grad(u) [dx(u), dy(u)] // EOM
macro Varf(varfName, meshName, VhName)
    varf varfName(u,v) = int2d(meshName)(grad(u)'* grad(v))
    + int2d(meshName)(1*v) + on(1, u = 0); // EOM
ffddmsetupOperator( PB , FE , Varf )

FEVhi ui, bi;
ffddmbuildrhs( PB , Varf , bi[] )
// Step 4: Define the one level DD preconditioner
ffddmsetupPrecond( PB , Varf )
// Step 5: Solve the linear system with GMRES
FEVhi x0i = 0;
ui[] = PBfGMRES(x0i[], bi[], 1.e-6, 200, "right");
ffddmplot(FE, ui, "u")
PBwritesummary
```

**Define a two level DD preconditioner**

See documentation

**Goal** improve scalability of the one level method  
\[\Rightarrow\] enrich the one level preconditioner with a *coarse problem* coupling all subdomains

Main ingredient is a rectangular matrix \(Z\) of size \(n \times n_c\), where \(n_c \ll n\)  
\(Z\) is the *coarse space* matrix

The *coarse space operator* \(E = Z^T A Z\) is a square matrix of size \(n_c \times n_c\)
The simplest way to enrich the one level preconditioner is through the *additive coarse correction* formula:

\[ M_2^{-1} = M_1^{-1} + Z E^{-1} Z^T \]

How to choose \( Z \)?

**Build the GenEO coarse space**

See documentation

```
ffddmgeneosetup( pr, Varf )
```

The *GenEO* method builds a robust coarse space for highly heterogeneous or anisotropic *SPD* problems

⇒ solve a local generalized eigenvalue problem in each subdomain

\[ D_i A_i D_i V_{i,k} = \lambda_{i,k} A_{i,Neu} V_{i,k} \]

with \( A_{i,Neu} \) the local Neumann matrices built from \( \text{Varf} \) (same \( \text{Varf} \) as *Step 3*)

The GenEO coarse space is

\[ Z = (R_i^T D_i V_{i,k})_{i=1,...,N} \]

The eigenvectors \( V_{i,k} \) selected to enter the coarse space correspond to eigenvalues \( \lambda_{i,k} \geq \tau \), where \( \tau \) is a threshold parameter

**Theorem** the spectrum of the preconditioned operator lies in the interval \([\frac{1}{1+k_1\tau}, k_0]\) where \( k_0 - 1 \) is the # of neighbors and \( k_1 \) is the multiplicity of intersections ⇒ \( k_0 \) and \( k_1 \) do not depend on \( N \) nor on the PDE

```
macro dimension 2 // EOM             // 2D or 3D
include "ffddm.idp"

mesh ThGlobal = square(100,100); // global mesh
// Step 1: Decompose the mesh
ffddmbuildDmesh( M, ThGlobal, mpiCommWorld )
// Step 2: Define your finite element
macro def(u) u // EOM
macro init(u) u // EOM
ffddmbuildDfespace( FE, M, real, def, init, P2 )
// Step 3: Define your problem
macro grad(u) [dx(u), dy(u)] // EOM
macro Varf(varfName, meshName, VhName)
    varf varfName(u,v) = int2d(meshName)(grad(u)'* grad(v))
    + int2d(meshName)(1*v) + on(1, u = 0); // EOM
ffddmsetupOperator( PB, FE, Varf )

FEVhi ui, bi;
ffddmbuildrhs( PB, Varf, bi[] )
// Step 4: Define the one level DD preconditioner
ffddmsetupPrecond( PB, Varf )
// Build the GenEO coarse space
ffddmgeneosetup( PB, Varf )
// Step 5: Solve the linear system with GMRES
FEVhi x0i = 0;
ui[] = PBfGMRES(x0i[], bi[], 1.e-6, 200, "right");
```
Build the coarse space from a coarse mesh

See documentation

```
ffddmcoarsemeshsetup( pr , Thc , VarfEprec , VarfAprec )
```

For non SPD problems, an alternative is to build the coarse space by discretizing the PDE on a coarser mesh \( \text{Th}_c \)

\( Z \) will be the interpolation matrix from the coarse FE space \( V_{h_c} \) to the original FE space \( V_h \)

\[ E = Z^T A Z \]

is the matrix of the problem discretized on the coarse mesh

The variational problem to be discretized on \( \text{Th}_c \) is given by macro \text{VarfEprec}

\text{VarfEprec} can differ from the original \text{Varf} of the problem

**Example:** added absorption for wave propagation problems

Similarly, \text{VarfAprec} specifies the global operator involved in multiplicative coarse correction formulas. It defaults to \( A \) if \text{VarfAprec} is not defined

```
macro dimension 2 // EOM
// 2D or 3D
include "ffddm.idp"

mesh ThGlobal = square(100,100); // global mesh
// Step 1: Decompose the mesh
ffddmbuildDmesh( M , ThGlobal , mpiCommWorld )
// Step 2: Define your finite element
macro def(u) u // EOM
macro init(u) u // EOM
ffddmbuildDfespace( FE , M , real , def , init , P2 )
// Step 3: Define your problem
macro grad(u) [dx(u), dy(u)] // EOM
macro Varf(varfName, meshName, VhName)
    varf varfName(u,v) = int2d(meshName)(grad(u)'* grad(v))
    int2d(meshName)(1*v) + on(1, u = 0); // EOM
ffddmsetupOperator( PB , FE , Varf )

FEVhi ui, bi;
ffddmbuildrhs( PB , Varf , bi[] )
// Step 4: Define the one level DD preconditioner
ffddmsetupPrecond( PB , Varf )
// Build the coarse space from a coarse mesh
mesh Thc = square(10,10);
ffddmcoarsemeshsetup( PB , Thc , Varf , null )
// Step 5: Solve the linear system with GMRES
FEVhi x0i = 0;
ui[] = PBfGMRES(x0i[], bi[], 1.e-6, 200, "right");
```

Use HPDDM within ffddm

See documentation

\text{ffddm} allows you to use \text{HPDDM} to solve your problem, effectively replacing the \text{ffddm} implementation of all parallel linear algebra computations

\[ \Rightarrow \text{define your problem with \text{ffddm}}, \text{solve it with \text{HPDDM}} \]
ffddm acts as a finite element interface for HPDDM

You can use HPDDM features unavailable in ffddm such as advanced Krylov subspace methods implementing block and recycling techniques

To switch to HPDDM, simply define the macro \texttt{pr\#withhpddm} before using \texttt{ffddmsetupOperator (Step 3)}. You can then pass HPDDM options with command-line arguments or directly to the underlying HPDDM operator. Options need to be prefixed by the operator prefix:

\begin{verbatim}
macro PBwithhpddm()1 // EOM
ffddmsetupOperator( PB , FE , Varf )
set(PBhpddmOP,sparams="-hpddm_PB_krylov_method gcrodr -hpddm_PB_recycle 10");
\end{verbatim}

Or, define \texttt{pr\#withhpddmkrylov} to use HPDDM only for the Krylov method

Example here: Helmholtz problem with multiple rhs solved with Block GMRES

Some results: Heterogeneous 3D elasticity with GenEO

Heterogeneous 3D linear elasticity equation discretized with P2 FE solved on 4096 MPI processes $n \approx 262$ million

Some results: 2-level DD for Maxwell equations, scattering from the COBRA cavity

$f = 10$ GHz
Some results: 2-level DD for Maxwell equations, scattering from the COBRA cavity

- order 2 Nedelec edge FE
- fine mesh: 10 points per wavelength
- coarse mesh: 3.33 points per wavelength
- two level ORAS preconditioner with added absorption
- $f = 10$ GHz: $n \approx 107$ million, $n_c \approx 4$ million
- $f = 16$ GHz: $n \approx 198$ million, $n_c \approx 7.4$ million
  
  \[ \rightarrow \text{coarse problem too large for a direct solver} \Rightarrow \text{inexact coarse solve: GMRES + one level ORAS preconditioner} \]

<table>
<thead>
<tr>
<th>$f$</th>
<th>$N$</th>
<th># it</th>
<th>Total # inner it</th>
<th>Total times (seconds)</th>
</tr>
</thead>
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<td>10GHz</td>
<td>1536</td>
<td>32</td>
<td>1527</td>
<td>Total: 515.8</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td></td>
<td>Setup: 383.2</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td></td>
<td>GMRES: 132.6</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td></td>
<td>inner: 61.8</td>
</tr>
<tr>
<td>10GHz</td>
<td>3072</td>
<td>33</td>
<td>2083</td>
<td>Total: 285.0</td>
</tr>
<tr>
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<td></td>
<td></td>
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<td>Setup: 201.6</td>
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<tr>
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<td></td>
<td></td>
<td>GMRES: 83.4</td>
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<td></td>
<td></td>
<td></td>
<td>inner: 40.6</td>
</tr>
<tr>
<td>16GHz</td>
<td>3072</td>
<td>43</td>
<td>3610</td>
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</tr>
<tr>
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<td>Setup: 336.8</td>
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<tr>
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<td></td>
<td></td>
<td></td>
<td>GMRES: 212.4</td>
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<tr>
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<td></td>
<td></td>
<td></td>
<td>inner: 118.6</td>
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<tr>
<td>16GHz</td>
<td>6144</td>
<td>46</td>
<td>4744</td>
<td>Total: 363.0</td>
</tr>
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<td></td>
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<td>Setup: 210.5</td>
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<tr>
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<td></td>
<td></td>
<td></td>
<td>GMRES: 152.5</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td></td>
<td>inner: 96.8</td>
</tr>
</tbody>
</table>
speedup of 1.81 from 1536 to 3072 cores at 10GHz
1.51 from 3072 to 6144 cores at 16GHz
You can find the script here

## 3.10.5 Examples

<table>
<thead>
<tr>
<th>File name</th>
<th>$M_1^{-1}$</th>
<th>$M_2^{-1}$</th>
<th>inexact CS</th>
<th>comments</th>
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<td>RAS</td>
<td>GenEO</td>
<td></td>
<td>direct solver MUMPS</td>
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<td></td>
<td>comparison with direct solver</td>
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<td>RAS</td>
<td>GenEO</td>
<td>RAS + GenEO</td>
<td>comparison with direct solver</td>
</tr>
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<td></td>
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<tr>
<td>elasticity-3d-thirdlevelgeneo.edp</td>
<td>RAS</td>
<td>GenEO</td>
<td>RAS + GenEO</td>
<td></td>
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<tr>
<td>Helmholtz-2d-simple.edp</td>
<td>ORAS</td>
<td>Coarse Mesh/DtN</td>
<td></td>
<td>for the DtN coarse space see this paper</td>
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<td>ORAS</td>
<td>Coarse Mesh</td>
<td></td>
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<tr>
<td>Helmholtz-3d-simple.edp</td>
<td>ORAS</td>
<td>Coarse Mesh</td>
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<tr>
<td>Helmholtz-3d-overthrust.edp</td>
<td>ORAS</td>
<td>Coarse Mesh</td>
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<td>ORAS</td>
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<td></td>
<td>multi-rhs Block GMRES with HPDDM</td>
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<td>Coarse Mesh</td>
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<tr>
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<td>ORAS</td>
<td>Coarse Mesh</td>
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<td></td>
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<tr>
<td>Maxwell_Cobracavity.edp</td>
<td>ORAS</td>
<td>Coarse Mesh</td>
<td>ORAS</td>
<td></td>
</tr>
<tr>
<td>natural_convection.edp</td>
<td>ORAS</td>
<td>Coarse Mesh</td>
<td></td>
<td>nonlinear</td>
</tr>
<tr>
<td>natural_convection_3D_obstacle.edp</td>
<td>ORAS</td>
<td>Coarse Mesh</td>
<td></td>
<td>nonlinear</td>
</tr>
<tr>
<td>Richards-2d.edp</td>
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<td></td>
<td>nonlinear time dependent mesh adaptation</td>
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<tr>
<td>heat-torus-3d-surf.edp</td>
<td>RAS</td>
<td>GenEO</td>
<td></td>
<td>3d surface time dependent</td>
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</tbody>
</table>
In essence **FreeFEM** is a compiler: its language is typed, polymorphic, with exception and reentrant. Every variable must be declared of a certain type, in a declarative statement; each statement are separated from the next by a semicolon ;.

The language allows the manipulation of basic types integers (int), reals (real), strings (string), arrays (example: real[int]), bi-dimensional (2D) finite element meshes (mesh), 2D finite element spaces (fespace), analytical functions (func), arrays of finite element functions (func[basic_type]), linear and bilinear operators, sparse matrices, vectors, etc. For example:

```cpp
int i, n = 20; //i, n are integer
real[int] xx(n), yy(n); //two array of size n
for (i = 0; i < 20; i++) { //which can be used in statements such as
    xx[i] = cos(i*pi/10);
    yy[i] = sin(i*pi/10);
}
```

The life of a variable is the current block {...}, except the fespace variable, and the variables local to a block are destroyed at the end of the block as follows.

**Tip:** Example

```cpp
real r = 0.01;
mesh Th = square(10, 10); //unit square mesh
fespace Vh(Th, P1); //P1 Lagrange finite element space
Vh u = x + exp(y);
func f = z*x + r*log(y);
plot(u, wait=true);
{
    real r = 2; //not the same r
    fespace Vh(Th, P1); //error because Vh is a global name
} // end of block
//here r back to 0.01
```

The type declarations are mandatory in **FreeFEM**; in the end this feature is an asset because it is easy to make bugs in a language with many implicit types.

The variable name is just an alphanumeric string, the underscore character _ is not allowed, because it will be used as an operator in the future.
4.1 Types

4.1.1 Standard types

int

Integer value (equivalent to long in C++).

```cpp
int i = 0;
```

bool

Boolean value.

```cpp
bool b = true;
```

**Tip:** The result of a comparison is a boolean

```cpp
bool b = (1 < 2);
```

real

Real value (equivalent to double in C++).

```cpp
real r = 0.;
```

complex

Complex value (equivalent to two double or complex<double> in C++).

```cpp
complex c = 0. + 1i;
```

The imaginary number $i$ is defined as $1i$

**Tip:** Example

```cpp
complex a = 1i, b = 2 + 3i;
cout << "a + b = " << a + b << endl;
cout << "a - b = " << a - b << endl;
cout << "a*b = " << a*b << endl;
cout << "a/b = " << a/b << endl;
```

The output of this script is:

```
a + b = (2,4)
a - b = (-2,-2)
a*b = (-3,2)
a/b = (0.230769,0.153846)
```
Note: See Complex example for a detailed example.

string

String value.

```c
string s = "this is a string";
```

Note: string value is enclosed within double quotes.

Other types can be concatenate to a string, like:

```c
int i = 1;
real r = 1.;
string s = "the int i = " + i +", the real r = " + r +", the complex z = " + (1. +
˓→1i);
```

To append a string in a string at position 4:

```c
s(4:3) = "+++";
```

To copy a substring in an other string:

```c
string s2 = s1(5:10);
```

See String Example for a complete example.

4.1.2 Mesh design

border

Border type.

```c
border b(t=0., 1.){x=cos(2.*pi*t); y=sin(2.*pi*t);}
```

Define the 2D geometrical border in parametric coordinates.

Note: Label

A label can be defined with the border:

```c
border b(t=0., 1.){x=cos(2.*pi*t); y=sin(2.*pi*t); label=1;}
```

Note: Inner variable

An inner variable can be defined inside a border:
\begin{verbatim}
border b(t=0., 1.){real tt=2.*p1*t; x=cos(tt); y=sin(tt);}
\end{verbatim}

**Note:** From vector
A border can be defined from two vectors using \( P.x \) and \( P.y \):

\begin{verbatim}
border b(t=0, vectorX.n-1){P.x=vectorX[t]; P.y=vectorY[t];}
\end{verbatim}

**mesh**

2D Mesh type (see *Mesh Generation*).

\begin{verbatim}
mesh Th;
\end{verbatim}

**mesh3**

3D mesh type (see *Mesh Generation*).

\begin{verbatim}
mesh3 Th;
\end{verbatim}

### 4.1.3 Finite element space design

**fespace**

Finite element space type (see *Finite Element*).

\begin{verbatim}
fespace Uh(Th, P1);
fespace UPh(Th, [P2, P2, P1]);
\end{verbatim}

A finite element space is based on a mesh (Th) with an element definition, scalar (P1) or vector ([P2, P2, P1]).

**Available finite element space:**

Generic:

- P0 / P03d
- P0Edge
- P1 / P13d
- P1dc
- P1b / P1b3d
- P1bl / P1bl3d
- P1nc
- P2 / P23d
- P2b
- P2dc
• P2h
• RT0 / RT03d
• RT0Ortho
• Edge03d

Using *Element_P3*:
• P3

Using *Element_P3dc*:
• P3dc

Using *Element_P4*:
• P4

Using *Element_P4dc*:
• P4dc

Using *Element_PkEdge*:
• P1Edge
• P2Edge
• P3Edge
• P4Edge
• P5Edge

Using *Morlay*:
• P2Morley

Using *HCT*:
• HCT

Using *BernardiRaugel*:
• P2BR

Using *Element_Mixte*:
• RT1
• RT1Ortho
• RT2
• RT2Ortho
• BDM1
• BDM1Ortho

Using *Element_Mixte3d*:
• Edge13d
• Edge23d

Using *Element_QF*:
• FEQF
A finite element function is defined as follow:

```plaintext
fespace Uh(Th, P1);
Uh u;

fespace UPh(Th, [P2, P2, P1]);
UPh [Ux, Uy, p];
```

### 4.1.4 Macro design

**macro**

Macro type.

```plaintext
macro vU() [Ux, Uy] //
macro grad(u) [dx(u), dy(u)] //
```

Macro ends with `//`.

**Note:** Macro concatenation

You can use the C concatenation operator `##` inside a macro using `#`.

If $U_x$ and $U_y$ are defined as finite element function, you can define:

```plaintext
macro Grad(U) [grad(U##x), grad(U##y)] // End of macro
```

See Macro example

**NewMacro / EndMacro**

**Warning:** In development - Not tested

Set and end a macro

```plaintext
NewMacro grad(u) [dx(u), dy(u)] EndMacro
```

**IFMACRO**

Check if a macro exists and check its value.

```plaintext
IFMACRO(AA) //check if macro AA exists
...
ENDIFMACRO

IFMACRO(AA, tt) //check if macro exists and is equall to tt
...
ENDIFMACRO
```
4.1.5 Functions design

**func**

Function type.

Function without parameters ($x$, $y$ and $z$ are implicitly considered):

```plaintext
func f = x^2 + y^2;
```

**Note:** Function’s type is defined by the expression’s type.

Function with parameters:

```plaintext
func real f (real var){
    return x^2 + y^2 + var^2;
}
```

**Elementary functions**

Class of basic functions (polynomials, exponential, logarithmic, trigonometric, circular) and the functions obtained from those by the four arithmetic operations

$$f(x) + g(x), f(x) - g(x), f(x)g(x), \frac{f(x)}{g(x)}$$

and by composition $f(g(x))$, each applied a finite number of times.

In FreeFEM, all elementary functions can thus be created. The derivative of an elementary function is also an elementary function; however, the indefinite integral of an elementary function cannot always be expressed in terms of elementary functions.

See [Elementary function example](#) for a complete example.

**Random functions**

FreeFEM includes the Mersenne Twister random number generator. It is a very fast and accurate random number generator of period $2^{2^19937} - 1$.

See `randint32()`, `randint64()`, `randreal1()`, `randreal2()`, `randreal3()`, `randres53()`, `randinit(seed)`.

In addition, the ffrandom plugin interface `random`, `srandom` and `srandomdev` functions of the Unix libc library. The range is $0 - 2^{31} - 1$.

**Note:** If `srandomdev` is not defined, a seed based on the current time is used.

gsl plugin equally allows usage of all random functions of the gsllib, see [gsl external library](#).
FE-functions

Finite element functions are also constructed like elementary functions by an arithmetic formula involving elementary functions.

The difference is that they are evaluated at declaration time and FreeFEM stores the array of its values at the places associated with he degree of freedom of the finite element type. By opposition, elementary functions are evaluated only when needed. Hence FE-functions are not defined only by their formula but also by the mesh and the finite element which enter in their definitions.

If the value of a FE-function is requested at a point which is not a degree of freedom, an interpolation is used, leading to an interpolation error, while by contrast, an elementary function can be evaluated at any point exactly.

1
2
3
4
5
6

```plaintext
func f = x^2*(1+y)^3 + y^2;
mesh Th = square(20, 20, [-2+4*x, -2+4*y]); // ]-2, 2[^2
fespace Vh(Th, P1);
Vh fh=f; //fh is the projection of f to Vh (real value)
func zf = (x^2*(1+y)^3 + y^2)*exp(x + 1i*y);
Vh<complex> zh = zf; //zh is the projection of zf to complex value Vh space
```

The construction of \( fh = f \) is explained in *Finite Element*.

**Warning:** The `plot` command only works for real or complex FE-functions, not for elementary functions.

4.1.6 Problem design

**problem**

Problem type.

```plaintext
problem Laplacian (u, uh) = ... 
```

FreeFEM needs the variational form in the problem definition.

In order to solve the problem, just call:

```plaintext
Laplacian; 
```

**Note:** Solver

A solver can be specified in the problem definition:

```plaintext
problem Laplacian(u, uh, solver=CG) = ... 
```

The default solver is `sparsesolver` or `LU` if any direct sparse solver is available.

Solvers are listed in the *Global variables* section.

**Note:** Stop test

A criterion \( \varepsilon \) can be defined for iterative methods, like CG for example:

```plaintext
problem Laplacian(u, uh, solver=CG, eps=1.e-6) = ... 
```
If $\varepsilon > 0$, the stop test is:

$$||Ax - b|| < \varepsilon$$

Else, the stop test is:

$$||Ax - b|| < \frac{\varepsilon}{||Ax_0 - b||}$$

---

**Note:** Reconstruction

The keyword `init` controls the reconstruction of the internal problem matrix.

If `init` is set to `false` or `0`, the matrix is reconstructed at each problem call (or after a mesh modification), else the previously constructed matrix is used.

```plaintext
problem Laplacian(u, uh, init=1) = ...
```

---

**Note:** Preconditioning

A preconditioner can be specified in the problem definition:

```plaintext
problem Laplacian(u, uh, precon=P) = ...
```

The preconditioning function must have a prototype like:

```plaintext
func real[int] P(real[int] &xx);
```

---

**Note:** “Très grande valeur”

The “Très grand valeur” `tgv` (or **Terrible giant value**) used to implement the Dirichlet conditions can be modified in the problem definition:

```plaintext
problem Laplacian(u, uh, tgv=1e30) = ...
```

Refer to *Problem definition* for a description of the Dirichlet condition implementation.

---

**Note:** Pivot tolerance

The tolerance of the pivot in UMFPACK, LU, Crout, Cholesky factorization can be modified in the problem definition:

```plaintext
problem Laplacian(u, uh, solver=LU, tolpivot=1e-20) = ...
```

---

**Note:** UMFPACK

Two specific parameters for the UMFPACK can be modified:

- Tolerance of the pivot sym
- strategy
problem Laplacian(u, uh, solver=LU, tolpivotsym=le-1, strategy=0) = ...

Refer to the UMFPACK website for more informations.

Note: dimKrylov
Dimension of the Krylov space

Usage of problem is detailed in the tutorials.

solve
Solve type.
Identical to problem but automatically solved.
Usage of solve is detailed in the tutorials.

varf
Variational form type.

varf vLaplacian (u, uh) = ...

Directly define a variational form.
This is the other way to define a problem in order to directly manage matrix and right hand side.
Usage of varf is detailed in the tutorial.

4.1.7 Array

An array stores multiple objects, and there are 2 kinds of arrays:
- the first is similar to vector, i.e. array with integer indices
- the second is array with string indices

In the first case, the size of the array must be known at execution time, and implementation is done with the KN<> class and all the vector operator of KN<> are implemented.

Arrays can be set like in Matlab or Scilab with the operator ::, the array generator of a:c is equivalent to a:1:c, and the array set by a:b:c is set to size ⌊(b − a)/c⌋ + 1 and the value i is set by a + i(b − a)/c.

There are int, real, complex array with, in the third case, two operators (.im, .re) to generate the real and imaginary real array from the complex array (without copy).

Note: Quantiles are points taken at regular intervals from the cumulative distribution function of a random variable. Here the array values are random.

This statistical function a.quantile(q) computes v from an array a of size n for a given number q ∈[0, 1[ such that:

$$\#\{i/a[i] < v\} \sim q * n$$
it is equivalent to \( v = a[q \cdot n] \) when the array \( a \) is sorted.

For example, to declare, fill and display an array of real of size \( n \):

```cpp
int n = 5;
real[int] Ai(n);
for (int i = 0; i < n; i++)
    Ai[i] = i;
cout << Ai << endl;
```

The output of this script is:

```
5
0 1 2 3 4
```

See the Array example for a complete example.

**Array index**

Array index can be int or string:

```cpp
real[int] Ai = [1, 1, 0, 0];
real[string] As = [1, 1, 0, 0];
```

**Array size**

The size of an array is obtained using the keyword \( n \):

```cpp
int ArraySize = Ai.n;
```

**Array sort**

To sort an array:

```cpp
Ai.sort;
```

**Double array**

A double array (matrix) can be defined using two indexes:

```cpp
real[int, int] Aii = {{1, 1}, {0, 0}};
```

The two sizes are obtained using the keywords \( n \) and \( m \):

```cpp
int ArraySize1 = Aii.n;
int ArraySize2 = Aii.m;
```

The minimum and maximum values of an array (simple or double) can be obtained using:

```cpp
real ArrayMin = Aii.min;
real ArrayMax = Aii.max;
```
Th minimum and maximum position of an array can be obtained using:

```c
int mini = Aii.imin;
int minj = Aii.jmin;
int maxi = Aii.imax;
int maxj = Aii.jmax;
```

**Tip:** An array can be obtained from a finite element function using:

```c
real[int] aU = U[];
```

where U is a finite element function.

### Array of FE functions

It is also possible to make an array of FE functions, with the same syntax, and we can treat them as vector valued function if we need them.

The syntax for space or vector finite function is

```c
int n = 100; // size of the array.
Vh[int] wh(n); // real scalar case
Vh[int] [uh,vh](n); // real vectorial case
Vh<complex>[int] cwh(n); // complex scalar case
Vh<complex>[int] [cuh, cvh](n); // complex vectorial case
[cuh[2], cvh[2]] = [x, y]; // set interpolation of index 2
```

**Tip:** Example

In the following example, Poisson’s equation is solved for 3 different given functions $f = 1, \sin(\pi x)\cos(\pi y), |x - 1||y - 1|$, whose solutions are stored in an array of FE function.

```c
// Mesh
mesh Th = square(20, 20, [2*x, 2*y]);

// Fespace
fespace Vh(Th, P1);
Vh u, v, f;

// Problem
problem Poisson (u, v) = int2d(Th) (dx(u)*dx(v) + dy(u)*dy(v)) + int2d(Th) (-f*v)
```

(continues on next page)
See FE array example.

Map arrays

```cpp
real[string] map; // a dynamic array
map["1"] = 2.0;
map[2] = 3.0; // is automatically cast to the string "2"
cout << "map["1"] = " << map["1"] << endl;
```

It is just a map of the standard template library so no operations on vector are allowed, except the selection of an item.

4.1.8 matrix

Defines a sparse matrix.

Matrices can be defined like vectors:

```cpp
matrix A = [[1, 2, 3],
            [4, 5, 6],
            [7, 8, 9]];
```

or using a variational form type (see Finite Element):

```cpp
matrix Laplacian = vLaplacian(Uh, Uh);
```

or from block of matrices:

```cpp
matrix A1, ..., An;
matrix A = [[A1, ...], ... , [... , An]];
```
or using sparse matrix set:

```plaintext
A = [I, J, C];
```

**Note:** I and J are `int[int]` and C is `real[int]`. The matrix is defined as:

\[
A = \sum_k C[k] M_{I[k], J[k]}
\]

where \( M_{a,b} = (\delta_{ia} \delta_{jb})_{ij} \)

I, J and C can be retrieved using \([I, J, C] = A\) (arrays are automatically resized).

The size of the matrix is \( n = I.\text{max}; m = J.\text{max}; \).

Matrices are designed using templates, so they can be real or complex:

```plaintext
matrix<real> A = ...  
matrix<complex> A = ...  
```

**Note:** Solver

See *problem*.

The default solver is *GMRES*.

```plaintext
matrix A = vLaplacian(Uh, Uh, solver=sparseresolver);
```

or

```plaintext
set(A, solver=sparseresolver);
```

**Note:** Factorize

If `true`, the factorization is done for LU, Cholesky or Crout.

```plaintext
matrix A = vLaplacian(Uh, Uh, solver=LU, factorize=1);
```

or

```plaintext
set(A, solver=LU, factorize=1);
```

**Note:** Stop test

See *problem*.

**Note:** Très grande valeur

See *problem*.

**Note:** Preconditioning
See `problem`.

**Note:** Pivot tolerance
See `problem`.

**Note:** UMFPACK
See `problem`.

**Note:** dimKrylov
See `problem`.

**Note:** datafilename
Name of the file containing solver parameters, see *Parallel sparse solvers*

**Note:** lparams
Vector of integer parameters for the solver, see *Parallel sparse solvers*

**Note:** dparams
Vector of real parameters for the solver, see *Parallel sparse solvers*

**Note:** sparams
String parameters for the solver, see *Parallel sparse solvers*

**Tip:** To modify the solver, the stop test,… after the matrix construction, use the `set` keyword.

### Matrix size

The size of a matrix is obtain using:

```c
int NRows = A.n;
int NColumns = A.m;
```

### Matrix resize

To resize a matrix, use:
\begin{verbatim}
A.resize(n, m);
\end{verbatim}

\textbf{Warning:} When resizing, all new terms are set to zero.

\section*{Matrix diagonal}

The diagonal of the matrix is obtained using:
\begin{verbatim}
real[int] Aii = A.diag;
\end{verbatim}

\section*{Matrix renumbering}

\begin{verbatim}
int[int] I(15), J(15);
matrix B = A;
B = A(I, J);
B = A(I^{-1}, J^{-1});
\end{verbatim}

\section*{Complex matrix}

Use \texttt{.im} and \texttt{.re} to get the imaginary and real part of a complex matrix, respectively:
\begin{verbatim}
matrix<complex> C = ...;
matrix R = C.re;
matrix I = C.im;
\end{verbatim}

\section*{Dot product / Outer product}

The dot product of two matrices is realized using:
\begin{verbatim}
real d = A' * B;
\end{verbatim}

The outer product of two matrices is realized using:
\begin{verbatim}
matrix C = A * B';
\end{verbatim}

See \textit{Matrix operations example} for a complete example.

\section*{Matrix inversion}

See \textit{Matrix inversion example}.

\section*{4.2 Global variables}

\subsection*{4.2.1 area}

Area of the current triangle.
4.2.2 ARGV

Array that contains all the command line arguments.

```
for (int i = 0; i < ARGV.n; i++)
    cout << ARGV[i] << endl;
```

See *Command line arguments example* for a complete example.

4.2.3 BoundaryEdge

Return 1 if the current edge is on a boundary, 0 otherwise.

```
real B = int2d(Th)(BoundaryEdge);
```

4.2.4 CG

Conjugate gradient solver.

Usable in *problem* and *solve* definition

```
problem Laplacian (U, V, solver=CG) = ...  
```

Or in *matrix* construction

```
matrix A = vLaplacian(Uh, Uh, solver=CG);
```

Or in *set function*

```
set(A, solver=CG);
```

4.2.5 Cholesky

Cholesky solver.

4.2.6 Crout

Crout solver.

4.2.7 edgeOrientation

Sign of $i - j$ if the current edge is $[q_i, q_j]$.

```
real S = int1d(Th, 1)(edgeOrientation);
```
4.2.8 false

False boolean value.

```cpp
bool b = false;
```

4.2.9 GMRES

GMRES solver (Generalized minimal residual method).

4.2.10 hTriangle

Size of the current triangle.

```cpp
fespace Vh(Th, P0);
Vh h = hTriangle;
```

4.2.11 include

Include an external library.

```cpp
include "iovtk"
```

4.2.12 InternalEdge

Return 0 if the current edge is on a boundary, 1 otherwise.

```cpp
real I = int2d(Th) (InternalEdge);
```

4.2.13 label

Label number of a boundary if the current point is on a boundary, 0 otherwise.

```cpp
int L = Th(xB, yB).label;
```

4.2.14 lenEdge

Length of the current edge.

For an edge $[q_i, q_j]$, return $|q_i - q_j|$.

```cpp
real L = int1d(Th, 1) (lenEdge);
```
4.2.15 load
Load a script.

```plaintext
load "Element_P3"
```

4.2.16 LU
LU solver.

4.2.17 N
Outward unit normal at the current point if it is on a curve defined by a border. \( N.x, \ N.y, \ N.z \) are respectively the \( x, \ y \) and \( z \) components of the normal.

```plaintext
func Nx = N.x;
func Ny = N.y;
func Nz = N.z;
```

4.2.18 nTonEdge
Number of adjacent triangles of the current edge.

```plaintext
real nTE = int2d(Th)(nTonEdge);
```

4.2.19 nuEdge
Index of the current edge in the triangle.

```plaintext
real nE = int2d(Th)(nuEdge);
```

4.2.20 nuTriangle
Index of the current triangle.

```plaintext
fespace Vh(Th, P0);
Vh n = nuTriangle;
```

4.2.21 P
Current point.

```plaintext
real cx = P.x;
real cy = P.y;
real cz = P.z;
```
4.2.22 pi

Pi = 3.14159.

```c
real Pi = pi;
```

This is a real value.

4.2.23 region

Region number of the current point. If the point is outside, then `region == notaregion` where `notaregion` is a FreeFEM integer constant.

```c
int R = Th(xR, yR).region;
```

4.2.24 sparsesolver

Sparse matrix solver.

4.2.25 true

True boolean value.

```c
bool b = true;
```

4.2.26 verbosity

Verbosity level.

```c
int Verbosity = verbosity;
verbosity = 0;
```

0 = nothing, 1 = little information, 10 = a lot of information, ...

This is an integer value.

4.2.27 version

FreeFEM version.

```c
cout << version << endl;
```

4.2.28 volume

Volume of the current tetrahedra.

```c
fespace Vh0(Th, P0);
Vh0 V = volume;
```
4.2.29 \( x \)

The \( x \) coordinate at the current point.

\[
\text{real } \text{CurrentX} = x;
\]

This is a real value.

4.2.30 \( y \)

The \( y \) coordinate at the current point.

\[
\text{real } \text{CurrentY} = y;
\]

This is a real value.

4.2.31 \( z \)

The \( z \) coordinate at the current point.

\[
\text{real } \text{CurrentZ} = z;
\]

This is a real value.

4.3 Quadrature formulae

The quadrature formula is like the following:

\[
\int_{D} f(x) \approx \sum_{\ell=1}^{L} \omega_\ell f(\xi_\ell)
\]

4.3.1 \texttt{int1d}

Quadrature formula on an edge.

\textbf{Notations}

\(|D|\) is the measure of the edge \( D \).

For a shake of simplicity, we denote:

\[
f(x) = g(t)
\]

with \( 0 \leq t \leq 1; x = (1-t)x_0 + tx_1.\)
This quadrature formula is exact on $P_1$. 

$$
\int_D f(x) \approx |D| g \left( \frac{1}{2} \right)
$$

This quadrature formula is exact on $P_3$. 

$$
\int_D f(x) \approx \frac{|D|}{2} \left( g \left( \frac{1 + \sqrt{1/3}}{2} \right) + g \left( \frac{1 - \sqrt{1/3}}{2} \right) \right)
$$

This quadrature formula is exact on $P_5$. 

$$
\int_D f(x) \approx \frac{|D|}{18} \left( 5g \left( \frac{1 + \sqrt{3/5}}{2} \right) + 8g \left( \frac{1}{2} \right) + 5g \left( \frac{1 - \sqrt{3/5}}{2} \right) \right)
$$

This quadrature formula is exact on $P_7$. 

$$
\int_D f(x) \approx \frac{|D|}{72} \left( (18 - \sqrt{30})g \left( \frac{1 - \sqrt{525+70\sqrt{35}}}{2} \right) + (18 + \sqrt{30})g \left( \frac{1 + \sqrt{525+70\sqrt{35}}}{2} \right) + (18 - \sqrt{30})g \left( \frac{1 - \sqrt{525-70\sqrt{35}}}{2} \right) + (18 + \sqrt{30})g \left( \frac{1 + \sqrt{525-70\sqrt{35}}}{2} \right) \right)
$$
qf5pE

\[ \int_{D} f(x) \approx |D| \left( \frac{332 - 13\sqrt{70}}{1800} g \left( \frac{1 - \frac{\sqrt{245+14\sqrt{70}}}{21}}{2} \right) + \frac{332 - 13\sqrt{70}}{1800} g \left( \frac{1 + \frac{\sqrt{245+14\sqrt{70}}}{21}}{2} \right) + \frac{64}{225} g \left( \frac{1}{2} \right) + \frac{332 + 13\sqrt{70}}{1800} g \left( \frac{1}{2} \right) \right) \]

This quadrature formula is exact on \( P_9 \).

qf1pElump

\[ \int_{D} f(x) \approx |D| \left( g(0) + g(1) \right) \]

This quadrature formula is exact on \( P_2 \).

4.3.2 int2d

Note: Complete formulas are no longer detailed

qf1pT

\[ \int_{D} f(x) \approx |D| \left( \frac{g(0) + g(1)}{2} \right) \]

This quadrature formula is exact on \( P_1 \).

qf2pT

\[ \int_{D} f(x) \approx |D| \left( \frac{g(0) + g(1)}{2} \right) \]

This quadrature formula is exact on \( P_2 \).
FreeFEM Documentation, Release 4.6

$qf5pT$

```latex
int2d(Th, qfe=qf5pT)( ... )
```

or

```latex
int2d(Th, qforder=6)( ... )
```

This quadrature formula is exact on $P_5$.

$qf1pTlump$

```latex
int2d(Th, qfe=qf1pTlump)( ... )
```

This quadrature formula is exact on $P_1$.

$qf2pT4P1$

```latex
int2d(Th, qfe=qf2pT4P1)( ... )
```

This quadrature formula is exact on $P_1$.

$qf7pT$

```latex
int2d(Th, qfe=qf7pT)( ... )
```

or

```latex
int2d(Th, qforder=8)( ... )
```

This quadrature formula is exact on $P_7$.

$qf9pT$

```latex
int2d(Th, qfe=qf9pT)( ... )
```

or

```latex
int2d(Th, qforder=10)( ... )
```

This quadrature formula is exact on $P_9$.

4.3.3 int3d

$qfV1$

```latex
int3d(Th, qfe=qfV1)( ... )
```

or

```latex
int3d(Th, qforder=7)( ... )
```

This quadrature formula is exact on $P_7$. 

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\[ \int_{3d}(Th, \ \text{qforder}=2)( \ldots ) \]

This quadrature formula is exact on \( P_1 \).

\textbf{qfV2}

\[ \int_{3d}(Th, \ \text{qfe=qfV2})( \ldots ) \]

or

\[ \int_{3d}(Th, \ \text{qforder}=3)( \ldots ) \]

This quadrature formula is exact on \( P_2 \).

\textbf{qfV5}

\[ \int_{3d}(Th, \ \text{qfe=qfV5})( \ldots ) \]

or

\[ \int_{3d}(Th, \ \text{qforder}=6)( \ldots ) \]

This quadrature formula is exact on \( P_5 \).

\textbf{qfV1lump}

\[ \int_{3d}(Th, \ \text{qfe=qfV1lump})( \ldots ) \]

This quadrature formula is exact on \( P_1 \).

\section*{4.4 Operators}

\subsection*{4.4.1 Addition operator +}

\[ \text{real} \ a = 1. + 2.; \]

Works for \text{int, real, complex, string, mesh, mesh3, array}.

\subsection*{4.4.2 Increment operator ++}

Pre-increment:

\[ \text{int} \ i = 0; \]
\[ ++i; \]

Post-increment:

\[ \text{int} \ i = 0; \]
\[ i++; \]
4.4.3 Substraction operator -

```plaintext
real a = 1. - 2.;
```

Works for `int, real, complex, array`.

4.4.4 Decrement operator –

Pre-decrement:

```plaintext
int i = 0;
--i;
```

Post-decrement:

```plaintext
int i = 0;
i--;
```

4.4.5 Multiplication operator *

```plaintext
real[int] b;
matrix A
real[int] x = A^-1*b;
```

Works for `int, real, complex, array, matrix`.

4.4.6 Equal operator =

```plaintext
real a = 1.;
```

4.4.7 Comparison operator ==

```plaintext
real a = 1.;
real b = 1.;
cout << (a == b) << endl;
```

4.4.8 Comparison operator !=

```plaintext
real a = 1.;
real b = 2.;
cout << (a != b) << endl;
```
4.4.9 Comparison operator <, <=

```plaintext
real a = 1.;
real b = 2.;
cout << (a < b) << endl;
cout << (a <= b) << endl;
```

4.4.10 Comparison operator >, >=

```plaintext
real a = 3.;
real b = 2.;
cout << (a > b) << endl;
cout << (a >= b) << endl;
```

4.4.11 Compound operator +=, -=, *=, /=

```plaintext
real a = 1;
a += 2.;
a -= 1.;
a *= 3.;
a /= 2.;
```

4.4.12 Term by term multiplication .*

```plaintext
matrix A = B .* C;
```

4.4.13 Division operator /

```plaintext
real a = 1. / 2.;
```

Works for int, real, complex.

4.4.14 Term by term division ./

```plaintext
matrix A = B ./ C;
```

4.4.15 Remainder from the division %

```plaintext
int a = 1 % 2;
```

Works for int, real.
4.4.16 Power operator ^

```plaintext
real a = 2.^2;
```

Works for `int, real, complex, matrix`.

4.4.17 Inverse of a matrix ^-1

```plaintext
real[int] Res = A^-1 * b;
```

**Warning**: This operator can not be used to directly create a matrix, see *Matrix inversion*.

4.4.18 Transpose operator '

```plaintext
real[int] a = b';
```

Works for array and `matrix`.

**Note**: For `matrix<complex>`, the `::freefem'''` operator return the Hermitian transpose.

4.4.19 Tensor scalar product :

\[ A : B = \sum_{i,j} A_{ij} B_{ij} \]

4.4.20 C++ arithmetical if expression ? :

\[ a \ ? \ b \ : \ c \] is equal to \( b \) if the \( a \) is true, \( c \) otherwise.

**Tip**: Example with `int`

```plaintext
int a = 12; int b = 5;
```

```plaintext
cout << a << " + " << b << " = " << a + b << endl;
cout << a << " - " << b << " = " << a - b << endl;
cout << a << " * " << b << " = " << a * b << endl;
cout << a << " / " << b << " = " << a / b << endl;
cout << a << " % " << b << " = " << a % b << endl;
cout << a << " ^ " << b << " = " << a ^ b << endl;
cout << ( " << a << " < " << b << " ? " << a << " : " << b << " ) = " << (a < b ? a : b) << endl;
```

The output of this script is:
Tip: Example with real

```cpp
real a = sqrt(2.); real b = pi;

cout << a << " + " << b << " = " << a + b << endl;
cout << a << " - " << b << " = " << a - b << endl;
cout << a << " * " << b << " = " << a * b << endl;
cout << a << " / " << b << " = " << a / b << endl;
cout << a << " % " << b << " = " << a % b << endl;
cout << "( " << a << " ? " << b << " ? " << a << " : " << b << ") = " << (a < b ? a : b) << endl;
```

The output of this script is:

```
1.41421 + 3.14159 = 4.55581
1.41421 - 3.14159 = -1.72738
1.41421 * 3.14159 = 4.44288
1.41421 / 3.14159 = 0.450158
1.41421 % 3.14159 = 1
1.41421 ^ 3.14159 = 2.97069
```
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4.5.3 else

See if.

4.5.4 while

While loop.

```plaintext
while (condition)
  ...
```

4.5.5 continue

Continue a loop.

```plaintext
for (int i = 0; i < N; ++i){
  ...
  if (condition) continue;
  ...
}
```

4.5.6 break

Break a loop.

```plaintext
while (condition){
  ...
  if (condition) break;
  ...
}
```

4.5.7 try

Try a part of code.

```plaintext
try{
  ...
} catch(...){
  ...
}
```

See Basic error handling example and Error handling example.
4.5.8 catch

Catch an error, see try

4.5.9 Implicit loop

Array with one index:

```cpp
for [i, ai : a]
```

If real[int] a(10), then i=0:9 and ai is a reference to a[i].

Array with two indices or matrix:

```cpp
for [i, j, aij : a]
```

If real[int] a(10, 11), then i=0:9, j=1:10 and aij is a reference to a(i, j).

See Implicit loop example.

4.6 I/O

See I/O example

See File stream example.

4.6.1 cout

Standard C++ output device (default: console).

```cpp
cout << "Some text" << endl;
```

4.6.2 cin

Standard C++ input device (default: keyboard).

```cpp
cin >> var;
```

4.6.3 endl

End of line.

```cpp
cout << "Some text" << endl;
```

4.6.4 ifstream

Open a file in read mode.
ifstream file("file.txt");

Note: A file is closed at the end of a block.

4.6.5 ofstream

Open a file in write mode.

ofstream file("file.txt");

Note: A file is closed at the end of a block.

4.6.6 append

Append data to an existing file.

ofstream file("file.txt", append);

4.6.7 binary

Write a file in binary.

ofstream file("file.btxt", binary);

4.6.8 seekg

Set the file position.

file.seekg(Pos);

4.6.9 tellg

Get the file position.

int Pos = file.tellg();

4.6.10 flush

Flush the buffer of the file.

file.flush
4.6.11 getline

Get the current line.

```cpp
string s;
getline(file, s);
```

4.6.12 Output format

In the descriptions below, `f` is an output stream, for example `cout` or an `ofstream`.

All this methods, excepted the first, return a stream, so they can be chained:

```cpp
cout.scientific.showpos << 3 << endl;
```

**precision**

Set the number of digits printed to the right of the decimal point. This applies to all subsequent floating point numbers written to that output stream. However, this won’t make floating-point “integers” print with a decimal point. It’s necessary to use `fixed` for that effect.

```cpp
int np = f.precision(n)
```

**scientific**

Formats floating-point numbers in scientific notation

```cpp
f.scientific
```

**fixed**

Used fixed point notation for floating-point numbers. Opposite of scientific.

```cpp
f.fixed
```

**showbase**

Converts insertions to an external form that can be read according to the C++ lexical conventions for integral constants. By default, showbase is not set.

```cpp
f.showbase
```

**noshowbase**

Unset showbase flags.

```cpp
f.noshowbase
```
showpos

Inserts a plus sign (+) into a decimal conversion of a positive integral value.

```cpp
f.showpos
```

noshowpos

Unset showpos flags.

```cpp
f.noshowpos
```

default

Reset all the previous flags to the default expect precision.

```cpp
f.default
```

setw

Behaves as if member width were called with \( n \) as argument on the stream on which it is inserted as a manipulator (it can be inserted on output streams).

```cpp
f.setw(n)
```

4.7 Functions

4.7.1 abs

Return the absolute value.

```cpp
real a = abs(b);
```

Parameters:

- \( b(\text{int, real, complex, fespace function, real}[\text{int}] \text{or} \text{real}[\text{int, int}]) \)

Output:

- \( a(\text{int, real, real}[\text{int}] \text{or} \text{real}[\text{int, int}]) \)

4.7.2 acos

arccos function.

```cpp
real theta = acos(x);
```

Parameter:

- \( x(\text{real, real}[\text{int}] \text{or} \text{real}[\text{int, int}]) \)
Output:

- \( \theta(\text{real, real[int]} \text{ or real[int, int]}) \)

\[ \text{Output:} \]

- \( \theta(\text{real}) \)

**4.7.3 acosh**

```plaintext
real \theta = \text{acosh}(x);
```

\[ \text{arccosh}(x) = \ln \left( x + \sqrt{x^2 - 1} \right) \]

Parameter:

- \( x(\text{real}) \)

Output:

- \( \theta(\text{real}) \)
4.7.4 adaptmesh

Mesh adaptation function.

```c
mesh Thnew = adaptmesh(Th, [fx, fy], hmin=HMin, hmax=HMax, err=Err, errg=ErrG, ...
  nbvx=NbVx, nbsmooth= NbSmooth, nbjacoby= NbJacoby, ratio= Ratio, omega= Omega, iso= Iso, ...
  abserror= AbsError, cutoff= Cutoff, verbosity= Verosity, inquire= Inquire, ...
  splitpbedge= SplitPbEdge, maxsubdiv= MaxSubdiv, rescaling= Rescaling, ...
  keepbackvertices= KeepBackVertices, IsMetric= isMetric, power= Power, ...
  thetamax= ThetaMax, splitin2= SplitIn2, metric= Metric, ...
  nomeshgeneration= NoMeshGeneration, periodic= Periodic);
```

Parameters:
- `Th` (mesh) Mesh to refine
- `[fx, fy]` (func or fespace function), scalar or vectorial Function to follow for the mesh adaptation
- `hmin` (real) Minimum edge size
- `hmax` (real) Maximum edge size
- `err` (real) Error level (P1 interpolation)
- `errg` (real) Relative geometrical error
- `nbvx` (int) Maximum number of vertices
- `nbsmooth` (int) Number of smoothing iterations
- `nbjacoby` (int) Number of iterations for the smoothing procedure
- `ratio` (real) Ratio of the triangles
- `omega` (real) Relaxation parameter for the smoothing procedure
- `iso` (bool) Isotropic adaptation (if true)
- `abserror` (bool) Error (if true) - Relative error (if false)
- `cutoff` (real) Lower limit of the relative error evaluation
- `verbosity` (real) Verbosity level
- `inquire` (bool) If true, inquire graphically
- `splitpbedge` (bool) If true, split all internal edges in half
- `maxsubdiv` (int) Bound the maximum subdivisions
- `rescaling` (bool) Rescale the function in [0, 1]
- `keepbackvertices` (bool) If true, try to keep vertices of the original mesh
- `IsMetric` (bool) If true, the metric is defined explicitly
- `power` (int) Exponent of the Hessian
- `thetamax` (int) Minimum corner angle (in degree)
- `splitin2` (bool) Split all triangles into 4 sub-triangles if true
- `metric` ([real[int], real[int], real[int]]) Array of 3 real arrays defining the metric
- `nomeshgeneration` (bool) If true, the mesh is not generated
- `periodic` ([real[int], int]) Build an adapted periodic mesh

Output:
• Thnew (mesh or mesh3)

### 4.7.5 adj

Adjacent triangle of the triangle $k$ by the edge $e$

```c
int T = Th[k].adj(e);
```

Parameter:
- $e$ (int) Edge number

Output:
- $T$ (int) Triangle number

### 4.7.6 AffineCG

Affine conjugate gradient solver

Used to solve a problem like $A x = b$

```c
int Conv = AffineCG(A, x, precon=Precon, nbiter=NbIter, eps=Eps, veps=VEps, ...
                      stop=Stop);
```

Parameters:
- $A$ (matrix) Matrix of the problem $A x = b$
- $x$ (real[int]) Solution vector
- $precon$ (real[int]) Preconditionning function
- $nbiter$ (int) Maximum number of iterations
- $eps$ (real) Convergence criterion
  - If $\varepsilon > 0$: test $||A(x)||_p \leq \varepsilon ||A(x_0)||_p$
  - If $\varepsilon < 0$: test $||A(x)||_p^2 \leq |\varepsilon|$
- $veps$ (real) Same as $eps$, but return $-eps$
- $stop$ (func) Convergence criterion as a function

Prototype is func bool StopFunc (int Iter, real[int] U, real[int] g)
  - $u$: current solution, $g$: current gradient (not preconditionned)

Output:
- $Conv$ (int) 0: converged - !0: not converged

### 4.7.7 AffineGMRES

Affine GMRES solver

Parameters and output are the same as AffineCG
4.7.8 arg

Return the argument of a complex number.

```c
real a = arg(c);
```

Parameters:
- `c` (complex)

Output:
- `r` (real)

4.7.9 asin

Arcsine function.

```c
real theta = asin(x);
```

Parameter:
- `x` (real, real[int] or real[int, int])

Output:
- `theta` (real, real[int] or real[int, int])

![Fig. 4.3: arcsin function](image)

4.7.10 asinh

```c
real theta = asinh(x);
```

Parameter:
- `x` (real)

\[
\text{asinh}(x) = \ln \left( x + \sqrt{x^2 + 1} \right)
\]
4.7.11 assert

Verify if a condition is true (same as C), if not the program stops.

\[
\text{assert}(x==0)
\]

Parameter:
- Boolean condition

Output:
- None

4.7.12 atan

arctan function.

\[
\text{real } \theta = \text{atan}(x);
\]

Parameter:
- \(x\) (real)

Output:
- \(\theta\) (real)

4.7.13 atan2

\(\text{arctan} \left( \frac{y}{x} \right)\) function, returning the correct sign for \(\theta\).
real theta = atan2(y, x)

Parameter:
  • x(real)

Output:
  • theta(real)

4.7.14 atanh

real theta = atanh(x);

Parameter:
  • x(real)

Output:
  • theta(real)

4.7.15 atoi

Convert a string to an integer.

int a = atoi(s);

Parameter:
  • s(string)

Output:
  • a(int)
4.7.16 atof

Convert a string to a real.

```c
real a = atof(s);
```

Parameter:
- `s` (string)

Output:
- `a` (real)

4.7.17 BFGS

Todo: todo

4.7.18 buildmesh

Build a 2D mesh using border elements.

```c
mesh Th = buildmesh(b1(nn) + b2(nn) + b3(nn) + b4(nn), [points=Points], [nbvx=Nbvx],
                   [fixedborder=FixedBorder]);
```

Parameters:
- `b1,b2,b3,b4` (border)
  - Geometry border, `b1(nn)` means `b1` border discretized by `nn` vertices
- `points(real[int, int]) [Optional]`
  - Specify a set of points
  - The size of `Points` array is `(nbp, 2)`, containing a set of `nbp` points with `x` and `y` coordinates

4.7. Functions
• `nbvx=(int) [Optional]`
  Maximum number of vertices Default: 9000
• `fixedborder=(bool) [Optional]`
  If true, mesh generator cannot change the boundary mesh
  Default: false

Output:
  • `Th(mesh)` Resulting mesh

### 4.7.19 ceil

Round fractions up of \( x \).

```bash
int c = ceil(x);
```

Parameter:
  • `x(real)`

Output:
  • `c(int)`

### 4.7.20 change

Change a property of a mesh.

```bash
int[int] L = [0, 1];
TheNew = change(Th, label=L);
```

Parameters:
  • `Th(mesh)` Original mesh
  • `label=L(int[int])` Pair of old and new label
  • `region=R(int[int])` Pair of old and new region
  • `flabel=l(func int)` Function of int given the new label
  • `fregion=r(func int)` Function of int given the new region

Output:
  • `Thnew(mesh)` Mesh with changed parameters

### 4.7.21 checkmovemesh

Check a `movemesh` without mesh generation.

```bash
real minT = checkmovemesh(Th, [Dx, Dy]);
```
Parameters:
Same as *movemesh*

Output:
- \( \text{minT}(\text{real}) \) Minimum triangle area

### 4.7.22 chi

Characteristic function of a mesh.

```cpp
int IsInMesh = chi(Th)(x, y);
```

Parameters:
- \( \text{Th} \) (mesh or mesh3)
- \( x \) (real) Position \( x \)
- \( y \) (real) Position \( y \)

Output:
- \( \text{IsInMesh}(\text{int}) \) \( 1 \) if \((x, y) \in \text{Th}\) \( 0 \) if \((x, y) \notin \text{Th}\)

### 4.7.23 clock

Get the clock in second.

```cpp
real t = clock();
```

Parameter:
- None

Output:
- \( t \) (real) Current CPU time

### 4.7.24 complexEigenValue

Same as *EigenValue* for complex problems.

### 4.7.25 conj

Calculate the conjugate of a complex number.

```cpp
complex C1 = 1 + 1i;
complex C2 = conj(C1);
```

Parameter:
- \( C1 \) (complex) Complex number

Output:
- \( C2 \) (complex) Conjugate of \( C1 \)
### 4.7.26 convect

Characteristics Galerkin method.

```cpp
real cgm = convect([Ux, Uy], dt, c);
real cgm = convect([Ux, Uy, Uz], dt, c);
```

Compute $c \circ X$ with $X(x) = x_\tau$ and $x_\tau$ is the solution of:

\[
\begin{align*}
\dot{x}_\tau &= u(x_\tau) \\
x_\tau &= x
\end{align*}
\]

Parameters:

- $u_x$ (fespace function) Velocity: $x$ component
- $u_y$ (fespace function) Velocity: $y$ component
- $u_z$ (fespace function) **3D only** Velocity: $z$ component
- $dt$ (real) Time step
- $c$ (fespace function) Function to convect

Output:

- $cgm$ (real) Result

### 4.7.27 copysign

C++ `copysign` function.

```cpp
real s = copysign(a, b);
```

### 4.7.28 cos

`cos` function.

```cpp
real x = cos(theta);
```

Parameters:

- $\theta$ (real or complex)

Output:

- $x$ (real or complex)

### 4.7.29 cosh

`cosh` function.

```cpp
real x = cosh(theta);
```

Parameters:

\[ \cosh(x) = \frac{e^x + e^{-x}}{2} \]
Fig. 4.7: cos function

• \texttt{theta(real)}
Output:
• \texttt{x(real)}

4.7.30 \texttt{diffnp}
Arithmetic useful function.

\begin{verbatim}
1 diffnp(a, b) = (a<0) && (0<b) ? (b-a) : 0;
\end{verbatim}

4.7.31 \texttt{diffpos}
Arithmetic useful function.

\begin{verbatim}
1 diffpos(a, b) = \texttt{max}(b-a, 0);
\end{verbatim}

4.7.32 \texttt{dist}
Arithmetic useful function.

\begin{verbatim}
1 dist(a, b) = \texttt{sqrt}(a^2 + b^2);
2 dist(a, b, c) = \texttt{sqrt}(a^2 + b^2 + c^2);
\end{verbatim}

4.7.33 \texttt{dumptable}
Show all types, operators and functions in \texttt{FreeFEM}.

\begin{verbatim}
1 dumptable(out);
\end{verbatim}

Parameters:
• `out (ostream) cout of ofstream file.

Output:
• None

### 4.7.34 dx

$x$ derivative.

```cpp
Uh up = dx(u);
```

\[
\frac{\partial u}{\partial x}
\]

Parameters:
• `u (fespace function)`

Output:
• `up (fespace function)`

### 4.7.35 dxx

$x$ double derivative.

```cpp
Uh upp = dxx(u);
```

\[
\frac{\partial^2 u}{\partial x^2}
\]

Parameters:
• `u (fespace function)`

Output:
• `upp (fespace function)`

### 4.7.36 dxy

$x y$ derivative.

```cpp
Uh upp = dxy(u);
```

\[
\frac{\partial^2 u}{\partial x \partial y}
\]

Parameters:
• `u (fespace function)`

Output:
• `upp (fespace function)`
4.7.37 dxz

$xz$ derivative.

```plaintext
Uh upp = dxz(u);
```

$\frac{\partial^2 u}{\partial x \partial z}$

Parameters:
- $u$ (fespace function)

Output:
- $upp$ (fespace function)

4.7.38 dy

$y$ derivative.

```plaintext
Uh up = dy(u);
```

$\frac{\partial u}{\partial y}$

Parameters:
- $u$ (fespace function)

Output:
- $upp$ (fespace function)

4.7.39 dyx

$yx$ derivative.

```plaintext
Uh upp = dyx(u);
```

$\frac{\partial^2 u}{\partial y \partial x}$

Parameters:
- $u$ (fespace function)

Output:
- $upp$ (fespace function)

4.7.40 dyy

$y$ double derivative.

```plaintext
Uh upp = dyy(u);
```

$\frac{\partial^2 u}{\partial x^2}$

Parameters:
• \( u \) (fespace function)

Output:
• \( u \) (fespace function)

### 4.7.41 dyz

**yz derivative.**

\[
\frac{\partial^2 u}{\partial y \partial z}
\]

\[\text{Uh upp = dyz(u);}\]

Parameters:
• \( u \) (fespace function)

Output:
• \( u \) (fespace function)

### 4.7.42 dz

**z derivative.**

\[
\frac{\partial u}{\partial z}
\]

\[\text{Uh up = dz(u);}\]

Parameters:
• \( u \) (fespace function)

Output:
• \( u \) (fespace function)

### 4.7.43 dzx

**zx derivative.**

\[
\frac{\partial^2 u}{\partial z \partial x}
\]

\[\text{Uh upp = dzx(u);}\]

Parameters:
• \( u \) (fespace function)

Output:
• \( u \) (fespace function)
### 4.7.44 dzy

$zy$ derivative.

```
Uh upp = dzy(u);
```

$\frac{\partial^2 u}{\partial z \partial y}$

Parameters:
- $u$ (fespace function)

Output:
- $upp$ (fespace function)

### 4.7.45 dzz

$z$ double derivative.

```
Uh upp = dzz(u);
```

$\frac{\partial^2 u}{\partial z^2}$

Parameters:
- $u$ (fespace function)

Output:
- $upp$ (fespace function)

### 4.7.46 EigenValue

Compute the generalized eigenvalue of $Au = \lambda Bu$. The shifted-inverse method is used by default with $\text{sigma}=\sigma$, the shift of the method. The function $\text{EigenValue}$ can be used for either matrices or functions returning a matrix vector product. The use of the matrix version is shown below.

```
int k = EigenValue(A, B, nev=, sigma=);
```

Parameters:
- $A, B$: matrices of same size
- $nev=n$: number of desired eigenvalues given by an integer $n$
- $sym=:$ the problem is symmetric or not
- $tol=:$ the relative accuracy to which eigenvalues are to be determined
- $value=:$ an array to store the real part of the eigenvalues
- $ivalue=:$ an array to store the imaginary part of the eigenvalues
- $vector=:$ a Finite Element function array to store the eigenvectors
- $sigma=:$ the shift value
- Other parameters are available for more advanced use: see the ARPACK documentation.
Output: The output is the number of converged eigenvalues, which can be different than the number of requested eigenvalues given by \texttt{nev=}. Note that the eigenvalues and the eigenvectors are stored for further purposes using the parameters \texttt{value=} and \texttt{vector=}.

### 4.7.47 emptymesh

Build an empty mesh.

Useful to handle Lagrange multipliers in mixed and Mortar methods.

```plaintext
mesh eTh = emptymesh(Th, ssd);
```

Parameters:
- \texttt{Th (mesh)} Mesh to empty
- \texttt{ssd (int [int])} Pseudo subregion label

Output:
- \texttt{eTh (mesh)} Empty mesh

### 4.7.48 erf

The error function:

\[ erf(x) = \frac{2}{\sqrt{\pi}} \int_0^x \exp(-t^2)dt \]

```plaintext
real err = erf(x);
```

Parameters:
- \texttt{x (real)}

Output:
- \texttt{err (real)}

### 4.7.49 erfc

Complementary of the error function:

\[ erfc(x) = 1 - erf(x) \]

```plaintext
real erfc = erfc(x);
```

Parameters:
- \texttt{x (real)}

Output:
- \texttt{err (real)}
4.7.50 exec

Execute an external command.

```cpp
int v = exec(command);
```

Parameters:
- `command` (string) Command to execute

Output:
- `v` (int) Value returned by the command

4.7.51 exit

Exit function, equivalent to `return`.

```cpp
exit(N);
```

Parameters:
- `N` (int) Return value

Output:
- None

4.7.52 exp

Exponential function.

```cpp
real a = exp(b);
```

Parameters:
- `b` (real or complex)

Output:
- `a` (real or complex)

4.7.53 fdim

Positive difference (`cmath` function).

```cpp
real fd = fdim(a, b);
```

Parameters:
- `a` (real)
- `b` (real)

Output:
- `fd` (real) If $x > y$, return $x - y$; if $x \leq y$, return 0
4.7.54 floor

Floor function.

```plaintext
real a = floor(b);
```

Return the largest integer value not greater than \( b \).

Parameters:

- \( b \) (real)

Output:

- \( a \) (real)

4.7.55 fmax

Maximum (cmath function).

```plaintext
real Max = fmax(a, b);
```

Parameters:

- \( a \) (real)
- \( b \) (real)

Output:

- \( Max \) (real)

4.7.56 fmin

Minimum (cmath function).

```plaintext
real Min = fmin(a, b);
```

Parameters:

- \( a \) (real)
- \( b \) (real)

Output:

- \( Min \) (real)

4.7.57 fmod

Remainder of \( a/b \) (cmath function).

```plaintext
real Mod = fmod(a, b);
```

Parameters:

- \( a \) (real)
- \( b \) (real)
Output:

- \textit{\textbf{Mod(real)}}

### 4.7.58 imag

Imaginary part of a complex number.

```c
complex c = 1. + 1i;
real Im = imag(c);
```

### 4.7.59 int1d

1D integral.

```c
int1d(Th, [Label], [qfe=Qfe], [qforder=Qforder])(
    ...
)
```

Used in \textit{problem, solve} or \textit{varf} definition to impose a boundary condition only (\textbf{FreeFEM} does not support 1D simulation), or outside to calculate a quantity.

**Parameters:**

- \textit{Th} (mesh) Mesh where the integral is calculated
- \textit{Label} (int) \textit{Optional}
  - Label of the 1D border Default: all borders of the mesh
- \textit{qfe= (quadrature formula)} \textit{Optional}
  - Quadrature formula, see \textit{quadrature formulae}
- \textit{qforder= (quadrature formula)} \textit{Optional}
  - Quadrature order, see \textit{quadrature formulae}

**Output:**

- Depending on the situation: In a \textit{problem, solve} or \textit{varf} definition: Non relevant.
- Outside: \texttt{real (example: real l = int1d(Th, 1)(1.));}.

**Warning:** In a \textit{problem, solve} or \textit{varf} definition, the content of \texttt{int1d} must be a linear or bilinear form.

### 4.7.60 int2d

2D integral.

```c
int2d(Th, [Region], [qfe=Qfe], [qforder=Qforder])(
    ...
)
```

Or

4.7. Functions
int2d(Th, [Label], [qfe=Qfe], [qforder=Qforder])(
    ...
)

Used in `problem`, `solve` or `varf` definition to:
- Calculate integral in 2D simulation
- Impose a boundary condition in 3D simulation
Or outside to calculate a quantity.

Parameters:

- `Th` (mesh or mesh3) Mesh where the integral is calculated
- `Region (int)` [Optional] Label of the 2D region (2D simulation) Default: all regions of the mesh
- `Label (int)` [Optional] Label of the 2D border (3D simulation) Default: all borders of the mesh
- `qfe= (quadrature formula)` [Optional]
  Quadrature formula, see `quadrature formulae`
- `qforder= (quadrature formula)` [Optional]
  Quadrature order, see `quadrature formulae`

Output:

- Depending on the situation: In a `problem`, `solve` or `varf` definition: Non relevant. Outside: real (example: real `s = int2d(Th, 1)(1.)`).

**Warning:** In a `problem`, `solve` or `varf` definition, the content of the `int2d` must be a linear or bilinear form.

4.7.61 int3d

3D integral.

int3d(Th, [Region], [qfe=Qfe], [qforder=Qforder])(
    ...
)

Used in `problem`, `solve` or `varf` definition to calculate integral in 3D simulation, or outside to calculate a quantity.

Parameters:

- `Th` (mesh3) Mesh where the integral is calculated
- `Region (int)` [Optional]
  Label of the 3D region
  Default: all regions of the mesh
- `qfe= (quadrature formula)` [Optional]
  Quadrature formula, see `quadrature formulae`
- `qforder= (quadrature formula)` [Optional]
  Quadrature order, see `quadrature formulae`

Output:

- Depending on the situation: In a `problem`, `solve` or `varf` definition: Non relevant. Outside: real (example: real `v = int3d(Th, 1)(1.)`).
Warning: In a problem, solve or varf definition, the content of the int3d must be a linear or bilinear form.

4.7.62 intalledges

Integral on all edges.

```plaintext
intalledges(Th, [Region])(
    ...
)
```

Parameters:
- Th (mesh) Mesh where the integral is calculated
- Region (int) [Optional]
  Label of the region
  Default: all regions of the mesh

Output:
- Non relevant

4.7.63 intallfaces

Integral on all faces.

Same as intalledges for mesh3.

4.7.64 interpolate

Interpolation operator from a finite element space to another.

```plaintext
matrix I = interpolate(Wh, Vh, [inside=Inside], [t=T], [op=Op], [U2Vc=U2VC]);
```

Parameters:
- Wh (fespace) Target finite element space
- Vh (fespace) Original finite element space
- inside= (bool) If true, create a zero extension outside the Vh domain
- t= (bool) If true, return the transposed matrix
- op= (int) 0: interpolate the function (default value) 1: interpolate \( \partial_x \) 2: interpolate \( \partial_y \) 3: interpolate \( \partial_z \)
- U2Vc= (int [int]) Array of the same size of Wh describing which component of Vh is interpolated in Wh

Output:
- I (matrix) Interpolation matrix operator
4.7.65 invdiff

Arithmetic useful function.

```
invdiff(a, b) = \(\text{abs}(a-b) < 10^{-30}\) ? (a-b) : 1/(a-b)
```

```
invdiff(a, b, e) = \(\text{abs}(a-b) < e\) ? (a-b) : 1/(a-b)
```

4.7.66 invdiffnp

Arithmetic useful function.

```
invdiffnp(a, b) = (a<0) \& (0<b) ? 1/(b-a) : 0
```

4.7.67 invdiffpos

Arithmetic useful function.

```
invdiffpos(a, b) = (a<b) ? 1./(b-a) : 0
```

4.7.68 isInf

The C++ isInf function.

```
bool b = isInf(a);
```

4.7.69 isNaN

The C++ isNan function.

```
bool b = isNaN(a);
```

4.7.70 isNormal

The C++ isNormal function.

4.7.71 j0

Bessel function of first kind, order 0.

```
real b = j0(x);
```

Parameters:

- \(x\) (real)

Output:

- \(b\) (real)
4.7.72 j1

Bessel function of first kind, order 1.

```plaintext
real b = j1(x);
```

Parameters:
- `x` (real)

Output:
- `b` (real)

4.7.73 jn

Bessel function of first kind, order n.

```plaintext
real b = jn(n, x);
```

\[
J_n(x) = \sum_{p=0}^{\infty} \frac{(1)^p}{p! (n + p)!} \left(\frac{x}{2}\right)^{2p+n}
\]

Parameters:
- `n` (int)
- `x` (real)

Output:
- `b` (real)

4.7.74 jump

Jump function across an edge.

```plaintext
intalledges ( ... jump(c) ... )
```

Parameters:
- `c` (fespace function) Discontinuous function

Output:
- Non relevant

4.7.75 LinearCG

Linear CG solver

Parameters and output are the same as `AffineCG`
4.7.76 LinearGMRES

Linear GMRES solver
Parameters and output are the same as AffineCG

4.7.77 lgamma

Natural logarithm of the absolute value of the \( \Gamma \) function of \( x \).

```cpp
real lg = lgamma(x);
```

Parameters:
- \( x \) (real)

Output:
- \( lg \) (real)

4.7.78 log

Natural logarithm.

```cpp
real l = log(x);
```

Parameters:
- \( x \) (real or complex)

Output:
- \( l \) (real or complex)

**Note:** Complex value

For complex value, the \( \log \) function is defined as:

\[
\log(z) = \log(|z|) + i \arg(z)
\]

4.7.79 log10

Common logarithm.

```cpp
real l = log10(x);
```

Parameters:
- \( x \) (real)

Output:
- \( l \) (real)
4.7.80 lrint

Integer value nearest to $x$.

```c
int l = lrint(a);
```

Parameters:
- $a$ (real)

Output:
- $l$ (int)

4.7.81 lround

Round a value, and return an integer value.

```c
int l = lround(a);
```

Parameters:
- $a$ (real)

Output:
- $l$ (int)

4.7.82 ltime

Return the current time since the Epoch.

```c
int t = ltime();
```

Parameter:
- None

Output:
- $t$ (int)

4.7.83 max

Maximum value of two, three or four values.

```c
real m = max(a, b);
real m = max(a, b, c);
real m = max(a, b, c, d);
```

Parameters:
- $a$ (int or real)
- $b$ (int or real)
- $c$ (int or real) [Optional]
- $d$ (int or real) [Optional]
Output:
• \( b(\text{int or real}) \)

### 4.7.84 min

Minimum value of two, three or four values.

\[
\begin{align*}
\text{real } m &= \min(a, b); \\
\text{real } m &= \min(a, b, c); \\
\text{real } m &= \min(a, b, c, d);
\end{align*}
\]

Parameters:
• \( a(\text{int or real}) \)
• \( b(\text{int or real}) \)
• \( c(\text{int or real}) [\text{Optional}] \)
• \( d(\text{int or real}) [\text{Optional}] \)

Output:
• \( b(\text{int or real}) \)

### 4.7.85 movemesh

Move a mesh.

\[
\begin{align*}
\text{mesh } \text{MovedTh} &= \text{movemesh}(\text{Th}, [\text{Dx}, \text{Dy}]); \\
\text{mesh3 } \text{MovedTh} &= \text{movemesh}(\text{Th}, [\text{Dx}, \text{Dy}, \text{Dz}], [\text{region}=>\text{Region}], [\text{label}=>\text{Label}], \ldots [\text{facemerge}=>\text{FaceMerge}], [\text{ptmerge}=>\text{PtMerge}], [\text{orientation}=>\text{Orientation}]);
\end{align*}
\]

Parameters:
• \( \text{Th} (\text{mesh of mesh3}) \) Mesh to move
• \( \text{Dx} (\text{fespace function}) \) Displacement along \( x \)
• \( \text{Dy} (\text{fespace function}) \) Displacement along \( y \)
• \( \text{Dz} (\text{fespace function}) \) \textbf{3D only} \\
  Displacement along \( z \)
• \( \text{region}=(\text{int}) [\text{Optional}] \) \textbf{3D only} \\
  Set label to tetrahedra
• \( \text{label}=(\text{int} [\text{int}]) [\text{Optional}] \) \textbf{3D only} \\
  Set label of faces (see \text{change} for more information)
• \( \text{facemerge}=(\text{int}) [\text{Optional}] \) \textbf{3D only} \\
  If equal to 1, some faces can be merged during the mesh moving Default: 1
• \( \text{ptmerge}=(\text{real}) [\text{Optional}] \) \textbf{3D only} \\
  Criteria to define when two points merge
• orientation=(int) [Optional] 3D only
  If equal to 1, allow orientation reverse if tetrahedra is not positive Default: 1

Output:
  • MovedTh (mesh or mesh3) Moved mesh

4.7.86 NaN

C++ nan function.

```cpp
real n = NaN([String]);
```

Parameters:
  • String (string) Default: ""

4.7.87 NLCG

Non-linear conjugate gradient.
Parameters and output are the same as AffineCG

4.7.88 on

Dirichlet condition function.

```cpp
problem (u, v)
  ...
  + on(Label, u=uD)
  ...
```

Warning: Used only in problem, solve and varf

Parameters:
  • Label (int or border in 2D)
  Boundary reference where to impose the Dirichlet condition
  • uD(fespace function, func or real or int)
  Dirichlet condition (u is an unknown of the problem)

Output:
  • Non relevant

4.7.89 plot

Plot meshes and results.

```cpp
plot([Th, [u], [[Ux, Uy, Uz]], [wait=Wait], [ps=PS], [coef=Coef], [fill=Fill], ...
    cmm=Cmm], [value=Value], [aspectratio=AspectRatio], [bb=Bb], [nbiso=NbIso], ...
    nbarrow=NbArrow], [viso=Viso], [arrow=Arrow], [bw=Bw], [grey=Grey], ...
    boundary=Boundary], [dim=Dim], [prev=Prev], [WindowTitle=WI]);
```

4.7. Functions
Note: Only one of $\text{Th}$, $u$ or $[\text{Ux, Uy}] / [\text{Ux, Uy, Uz}]$ is needed for the plot command.

Parameters:

- $\text{Th}$ (mesh or mesh3) Mesh to display
- $u$ (fespace function) Scalar fespace function to display
- $[\text{Ux, Uy}] / [\text{Ux, Uy, Uz}]$ (fespace function array) Vectorial fespace function to display
- $[\text{Ux, Uy}([\text{real[int]}, \text{real[int]}])]$ Couple a real array to display a curve
- $\text{wait}=$ (bool) If true, wait before continue
- $\text{ps}=$ (string) Name of the file to save the plot (.ps or .eps format)
- $\text{coef}=$ (real) Arrow size
- $\text{fill}=$ (bool) If true, fill color between isovalue (usable with scalar fespace function only)
- $\text{cmm}=$ (string) Text comment in the graphic window
- $\text{value}=$ (bool) If true, show the value scale
- $\text{asr}=$ (bool) If true, preserve the aspect ratio
- $\text{bb}=$ ([real[int], real[int]]) Specify a bounding box using two corner points
- $\text{nbiso}=$ (int) Number of isovalues
- $\text{nbarrow}=$ (int) Number of colors of arrows values
- $\text{viso}=$ (real[int]) Specify an array of isovalues
- $\text{varrow}=$ (real[int]) Specify an array of arrows values color
- $\text{bw}=$ (bool) If true, the plot is in black and white
- $\text{grey}=$ (bool) If true, the plot is in grey scale
- $\text{hsv}=$ (real[int]) Array of $3 \times n$ values defining HSV color model $[h_1, s_1, v_1, ..., h_n, s_n, v_n]$
- $\text{boundary}=$ (bool) If true, display the boundary of the domain
- $\text{dim}=$ (int) Set the dimension of the plot: 2 or 3
- $\text{prev}=$ (bool) Use the graphic state of the previous state
- $\text{WindowIndex}=$ (int) Specify window index for multiple windows graphics

Output:

- None

See the plot section for in-graphic commands.

### 4.7.90 polar

Polar coordinates.

```cpp
complex p = polar(a, b);
```
Parameters:
- $a$ (real)
- $b$ (real)

Output:
- $p$ (complex)

### 4.7.91 pow

Power function.

```cpp
real p = pow(a, b);
```

$p = a^b$

Parameters:
- $a$ (real)
- $b$ (real)

Output:
- $p$ (real)

### 4.7.92 projection

Arithmetic useful function.

```cpp
real p = projection(a, b, x);
```

Projection is equivalent to:

```cpp
projection(a, b, x) = min(max(a, x), b)*a < b + min(max(b, x), a)*(1-a < b);
```

Parameters:
- $a$ (real)
- $b$ (real)
- $x$ (real)

Output:
- $p$ (real)

### 4.7.93 randinit

Initialize the state vector by using a seed.

```cpp
randinit(seed);
```

Parameters:
- $seed$ (int)
Output:
• None

### 4.7.94 randint31

Generate unsigned int (31 bits) random number.

```plaintext
int r = randint31();
```

Parameters:
• None

Output:
• \( r \) (int)

### 4.7.95 randint32

Generate unsigned int (32 bits) random number.

```plaintext
int r = randint32();
```

Parameters:
• None

Output:
• \( r \) (int)

### 4.7.96 randreal1

Generate uniform real in \([0, 1]\) (32 bits).

```plaintext
real r = randreal1();
```

Parameters:
• None

Output:
• \( r \) (real)

### 4.7.97 randreal2

Generate uniform real in \([0, 1)\) (32 bits).

```plaintext
real r = randreal2();
```

Parameters:
• None

Output:
4.7.98 randreal3

Generate uniform real in (0, 1) (32 bits).

```cpp
real r = randreal3();
```

Parameters:
- None

Output:
- r(real)

4.7.99 randres53

Generate uniform real in [0, 1) (53 bits).

```cpp
real r = randres53();
```

Parameters:
- None

Output:
- r(real)

4.7.100 readmesh

Read a 2D mesh file at different formats (see Mesh Generation).

```cpp
mesh Th = readmesh(MeshFileName);
```

Parameters:
- MeshFileName (string)

Output:
- Th(mesh)

4.7.101 readmesh3

Read a 3D mesh file at different formats (see Mesh Generation).

```cpp
mesh3 Th = readmesh3(MeshFileName);
```

Parameters:
- MeshFileName (string)

Output:
- Th(mesh3)
4.7.102 real

Return the real part of a complex number.

```
real r = real(c);
```

Parameters:
- c (complex)

Output:
- r (real)

4.7.103 rint

Integer value nearest to \( x \) (real value).

```
real r = rint(a);
```

Parameters:
- a (real)

Output:
- r (real)

4.7.104 round

Round a value (real value).

```
real r = round(a);
```

Parameters:
- a (real)

Output:
- r (real)

4.7.105 savemesh

Save a 2D or 3D mesh in different formats (see Mesh Generation 2D and Mesh Generation 3D).

```
savemesh(Th, MeshFileName);
```

Parameters:
- Th (mesh or mesh3)
- MeshFileName (string)

Output:
- None
4.7.106 set

Set a property to a matrix. See \textit{matrix}.

4.7.107 sign

Sign of a value.

\begin{verbatim}
int s = sign(a);
\end{verbatim}

Parameters:
- \texttt{a} (real or int)

Output:
- \texttt{s} (int)

4.7.108 signbit

C++ \texttt{signbit} function

\begin{verbatim}
int s = signbit(a);
\end{verbatim}

4.7.109 sin

\texttt{sin} function.

\begin{verbatim}
real x = sin(theta);
\end{verbatim}

Parameter:
- \texttt{theta} (real or complex)

Output:
- \texttt{x} (real or complex)

4.7.110 sinh

\texttt{sinh} function.

\begin{verbatim}
real x = sinh(theta);
\end{verbatim}

\begin{equation}
\text{sinh}(x) = \frac{e^x - e^{-x}}{2}
\end{equation}

Parameter:
- \texttt{theta} (real)

Output:
- \texttt{x} (real)
Fig. 4.8: sin function

Fig. 4.9: sinh function
### 4.7.111 sort

Sort two array in parallel

```c
sort(A, B);
```

**Parameters:**
- `A` (real[int])
- `B` (int[int])

**Output:**
- None

`A` is sorted in ascending order, `B` is sorted as `A`.

### 4.7.112 splitmesh

Split mesh triangles according to a function.

```c
Th = splitmesh(Th0, f);
```

**Parameters:**
- `Th0` (mesh)
- `f` (func or fespace function)

**Output:**
- `Th` (mesh)

### 4.7.113 sqrt

Square root

```c
real s = sqrt(a);
```

**Parameter:**
- `a` (real)

**Output:**
- `s` (real)

### 4.7.114 square

1. Square of a number.

```c
real S = square(a);
```

**Parameter:**
- `a` (real)

**Output:**
2. Build a structured square mesh.

```c
mesh Th = square(nnX, nnY, [[L*x, H*y]], [flags=Flags], [label=Labels],,
    [region=Region]);
```

Parameters:
- `nnX(int)` Discretization along \( x \)
- `nnY(int)` Discretization along \( y \)
- `L(real)` [Optional] Length along \( x \)
- `H(real)` [Optional] Height along \( y \)
- `flags=(int)` [Optional]
- `label=(int[int])` [Optional]
- `region=(int)` [Optional]

Structured mesh type, see Mesh Generation chapter for more information

Output:
- `Th(mesh)`

### 4.7.115 storagetotal

```c
int total = storagetotal();
```

### 4.7.116 storageused

```c
int used = storageused();
```

### 4.7.117 strtok

C++ `strtok` function

```c
string text = "10.5";
real number = strtok(text);
```

Parameter:
- `text(string)`

Output:
- `number(real)`

### 4.7.118 strtol

C++ `strtol` function
```c
string text = "10";
int number = strtol(text);
int base = 16;
int number = strtol(text, base);
```

Parameter:
- `text` (string)
- `base` (int) Base [Optional]

Output:
- `number` (int)

### 4.7.119 swap

**Swap values.**

```c
swap(a, b);
```

Parameters:
- `a` (real)
- `b` (real)

Output:
- None

### 4.7.120 system

**Execute a system command.**

```c
int Res = system(Command);
```

Parameter:
- `Command` (string) System command

Output:
- `Res` (int) Value returned by the system command

**Note:** On Windows, the full path of the command is needed. For example, to execute `ls.exe`:

```c
int Res = exec("C:\cygwin\bin\ls.exe");
```

### 4.7.121 tan

**tan function.**

```c
real x = tan(theta);
```
4.7.122 tanh

tanh function.

```cpp
real x = tanh(theta);
```

Parameter:
- `theta` (real)

Output:
- `x` (real)

4.7.123 tgamma

Calculate the Γ function of `x`.

```cpp
real tg = tgamma(x);
```

Parameter:
- `x` (real)

Output:
- `tg` (real)
4.7.124 time

Return the current time (C++ function).

```cpp
real t = time();
```

Parameter:
- None

Output:
- t(real)

4.7.125 trace

Matrix trace

```cpp
real tr = trace([[1, 2], [3, 4]]);
```

Parameters:
- Matrix

Output:
- Trace of the matrix (real)

4.7.126 trunc

Split triangle of a mesh.

```cpp
mesh Th = trunc(Th0, R, [split=Split], [label=Label]);
```

Parameters:
- Th0 (mesh)
• \( R \) (bool or int) Split triangles where \( R \) is true or different from 0
• \( \text{split}=(\text{int}) \) [Optional]
  Level of splitting Default: 1
• \( \text{label}=(\text{int}) \) [Optional]
  Label number of new boundary item Default: 1

Output:
  • \( \text{Th} \) (mesh)

4.7.127 \( y_0 \)

Bessel function of second kind, order 0.

```plaintext
real B = y0(x);
```

Parameters:
  • \( x \) (real)

Output:
  • \( b \) (real)

4.7.128 \( y_1 \)

Bessel function of second kind, order 1.

```plaintext
real B = y1(x);
```

Parameters:
  • \( x \) (real)

Output:
  • \( b \) (real)

4.7.129 \( y_n \)

Bessel function of second kind, order \( n \).

```plaintext
real B = yn(n, x);
```

\[
Y_n(x) = \lim_{\lambda \to n} \frac{J_\lambda(x) \cos(\lambda \pi) - J_{-\lambda}(x)}{\sin(\lambda \pi)}
\]

Parameters:
  • \( n \) (int)
  • \( x \) (real)

Output:
  • \( b \) (real)
4.8 External libraries

4.8.1 aniso

boundaniso

Todo: todo

4.8.2 BEC

BECtrap

Todo: todo

GPvortex

Todo: todo

dxGPVortex

Todo: todo

dyGPVortex

Todo: todo

4.8.3 Binary I/O

LoadVec

Todo: todo
LoadFlag

TODO: todo

SaveVec

TODO: todo

flag

TODO: todo

4.8.4 buildlayer

buildlayers

TODO: todo

4.8.5 ClosePoints

radiusSearch

TODO: todo

Voisinage

TODO: todo

neighborhood

TODO: todo
4.8.6 Curvature

extractborder

Extract a border of a mesh.

```c
int Res = extractborder(Th, Label, Points);
```

Parameters:
- \( \text{Th} (\text{mesh or mesh3}) \)
- \( \text{Label} (\text{int}) \) Label of the border to extract
- \( \text{Points} (\text{real}[\text{int}, \text{int}]) \) Extracted points Must be allocated as real[int, int] Points(3, 1);

Output:
- \( \text{Res} (\text{real}) \) Length of the extracted border

curvature

Todo: todo

rxicurvature

Todo: todo

4.8. External libraries
curves

Todo: todo

setecurveabcisse

Todo: todo

equiparameter

Todo: todo

Tresca

Todo: todo

VonMises

Todo: todo

4.8.7 dfft

Refer to the FFTW documentation for more informations.

plandfft

Todo: todo

execute

Todo: todo
Todo: todo

dfft
Todo: todo

map
Todo: todo

4.8.8 distance

Need

```cpp
load "distance"
```

distance

```cpp
distance(Th, d, dist, [distmax=DistMax]);
```

Parameters:
- `Th` (mesh)
- `d`
- `dist` (real[int])

Output:
- 

Todo: todo

checkdist
Todo: todo
4.8.9 DxWriter

Dxaddmesh

Todo: todo

Dxaddtimeseries

Todo: todo

Dxaddsol2ts

Todo: todo

4.8.10 Element_P1bl

expert

Todo: todo

4.8.11 exactpartition

exactpartition

Todo: todo

4.8.12 ff-AiryBiry

airy

Todo: todo

biry

Todo: todo
4.8.13 ff-cmaes

cmaes

Todo: todo

4.8.14 ff_gsl_awk

Refer to the GSL documentation for more informations

gslcdfugaussianP

Link to:
```
gsl_cdf_ugaussian_P(a)
```

gslcdfugaussianQ

Link to:
```
gsl_cdf_ugaussian_Q(a)
```

gslcdfugaussianPinv

Link to:
```
gsl_cdf_ugaussian_Pinv(a)
```

gslcdfugaussianQinv

Link to:
```
gsl_cdf_ugaussian_Qinv(a)
```

gslcdfgaussianP

Link to:
```
gsl_cdf_gaussian_P(a, b)
```

gslcdfgaussianQ

Link to:
```
gsl_cdf_gaussian_Q(a, b)
```
gslcdfgaussianPinv

Link to:

\[ \text{gsl\_cdf\_gaussian\_Pinv}(a, b) \]


gslcdfgaussianQinv

Link to:

\[ \text{gsl\_cdf\_gaussian\_Qinv}(a, b) \]


gslcdfgammaP

Link to:

\[ \text{gsl\_cdf\_gamma\_P}(a, b, c) \]


gslcdfgammaQ

Link to:

\[ \text{gsl\_cdf\_gamma\_Q}(a, b, c) \]


gslcdfgammaPinv

Link to:

\[ \text{gsl\_cdf\_gamma\_Pinv}(a, b, c) \]


gslcdfgammaQinv

Link to:

\[ \text{gsl\_cdf\_gamma\_Pinv}(a, b, c) \]


gslcdfcauchyP

Link to:

\[ \text{gsl\_cdf\_cauchy\_P}(a, b) \]


gslcdfcauchyQ

Link to:

\[ \text{gsl\_cdf\_cauchy\_Q}(a, b) \]
gslcdfcauchyPinv

Link to:

```
gsl_cdf_cauchy_Pinv(a, b)
```

gslcdfcauchyQinv

Link to:

```
gsl_cdf_cauchy_Qinv(a, b)
```

gslcdflaplaceP

Link to:

```
gsl_cdf_laplace_P(a, b)
```

gslcdflaplaceQ

Link to:

```
gsl_cdf_laplace_Q(a, b)
```

gslcdflaplacePinv

Link to:

```
gsl_cdf_laplace_Pinv(a, b)
```

gslcdflaplaceQinv

Link to:

```
gsl_cdf_laplace_Qinv(a, b)
```

gslcdfrayleighP

Link to:

```
gsl_cdf_rayleigh_P(a, b)
```

gslcdfrayleighQ

Link to:

```
gsl_cdf_rayleigh_Q(a, b)
```

4.8. External libraries 423
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**gslcdfrayleighPinv**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_rayleigh_Pinv(double a, double b);
```

**gslcdfrayleighQinv**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_rayleigh_Qinv(double a, double b);
```

**gslcdfchisqP**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_chisq_P(double a, double b);
```

**gslcdfchisqQ**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_chisq_Q(double a, double b);
```

**gslcdfchisqPinv**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_chisq_Pinv(double a, double b);
```

**gslcdfchisqQinv**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_chisq_Qinv(double a, double b);
```

**gslcdfexponentialP**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_exponential_P(double a, double b);
```

**gslcdfexponentialQ**

Link to:

```c
#include <gsl/gsl_cdf.h>

double gsl_cdf_exponential_Q(double a, double b);
```
**gslcdfexponentialPinv**

Link to:

```c
gsl_cdf_exponential_Pinv(a, b)
```

**gslcdfexponentialQinv**

Link to:

```c
gsl_cdf_exponential_Qinv(a, b)
```

**gslcdfexppowP**

Link to:

```c
gsl_cdf_exppow_P(a, b, c)
```

**gslcdfexppowQ**

Link to:

```c
gsl_cdf_exppow_Q(a, b, c)
```

**gslcdftdistP**

Link to:

```c
gsl_cdf_t_dist_P(a, b)
```

**gslcdftdistQ**

Link to:

```c
gsl_cdf_t_dist_Q(a, b)
```

**gslcdftdistPinv**

Link to:

```c
gsl_cdf_t_dist_Pinv(a, b)
```

**gslcdftdistQinv**

Link to:

```c
gsl_cdf_t_dist_Qinv(a, b)
```
### gslcdfdistP

Link to:

| gsl_cdf_fdist_P(a, b, c) |

### gslcdfdistQ

Link to:

| gsl_cdf_fdist_Q(a, b, c) |

### gslcdfdistPinv

Link to:

| gsl_cdf_fdist_Pinv(a, b, c) |

### gslcdfdistQinv

Link to:

| gsl_cdf_fdist_Qinv(a, b, c) |

### gslcdfbetaP

Link to:

| gsl_cdf_beta_P(a, b, c) |

### gslcdfbetaQ

Link to:

| gsl_cdf_beta_Q(a, b, c) |

### gslcdfbetaPinv

Link to:

| gsl_cdf_beta_Pinv(a, b, c) |

### gslcdfbetaQinv

Link to:

| gsl_cdf_beta_Qinv(a, b, c) |
gslcdfflatP

Link to:

gsl_cdf_flat_P(a, b, c)

gslcdfflatQ

Link to:

gsl_cdf_flat_Q(a, b, c)

gslcdfflatPinv

Link to:

gsl_cdf_flat_Pinv(a, b, c)

gslcdfflatQinv

Link to:

gsl_cdf_flat_Qinv(a, b, c)

gslcdflognormalP

Link to:

gsl_cdf_lognormal_P(a, b, c)

gslcdflognormalQ

Link to:

gsl_cdf_lognormal_Q(a, b, c)

gslcdflognormalPinv

Link to:

gsl_cdf_lognormal_Pinv(a, b, c)

gslcdflognormalQinv

Link to:

gsl_cdf_lognormal_Qinv(a, b, c)

4.8. External libraries


**gslcdfgumbel1P**

Link to:

`gsl_cdf_gumbel1_P(a, b, c)`

**gslcdfgumbel1Q**

Link to:

`gsl_cdf_gumbel1_Q(a, b, c)`

**gslcdfgumbel1Pinv**

Link to:

`gsl_cdf_gumbel1_Pinv(a, b, c)`

**gslcdfgumbel1Qinv**

Link to:

`gsl_cdf_gumbel1_Qinv(a, b, c)`

**gslcdfgumbel2P**

Link to:

`gsl_cdf_gumbel2_P(a, b, c)`

**gslcdfgumbel2Q**

Link to:

`gsl_cdf_gumbel2_Q(a, b, c)`

**gslcdfgumbel2Pinv**

Link to:

`gsl_cdf_gumbel2_Pinv(a, b, c)`

**gslcdfgumbel2Qinv**

Link to:

`gsl_cdf_gumbel2_Qinv(a, b, c)`
**gslcdfweibullP**

Link to:

```c
gsl_cdf_weibull_P(a, b, c)
```

**gslcdfweibullQ**

Link to:

```c
gsl_cdf_weibull_Q(a, b, c)
```

**gslcdfweibullPinv**

Link to:

```c
gsl_cdf_weibull_Pinv(a, b, c)
```

**gslcdfweibullQinv**

Link to:

```c
gsl_cdf_weibull_Qinv(a, b, c)
```

**gslcdfparetoP**

Link to:

```c
gsl_cdf_pareto_P(a, b, c)
```

**gslcdfparetoQ**

Link to:

```c
gsl_cdf_pareto_Q(a, b, c)
```

**gslcdfparetoPinv**

Link to:

```c
gsl_cdf_pareto_Pinv(a, b, c)
```

**gslcdfparetoQinv**

Link to:

```c
gsl_cdf_pareto_Qinv(a, b, c)
```
**gslcdflogisticP**

Link to:

```
gsl_cdf_logistic_P(a, b)
```

**gslcdflogisticQ**

Link to:

```
gsl_cdf_logistic_Q(a, b)
```

**gslcdflogisticPinv**

Link to:

```
gsl_cdf_logistic_Pinv(a, b)
```

**gslcdflogisticQinv**

Link to:

```
gsl_cdf_logistic_Qinv(a, b)
```

**gslcdfbinomialP**

Link to:

```
gsl_cdf_binomial_P(a, b, c)
```

**gslcdfbinomialQ**

Link to:

```
gsl_cdf_binomial_Q(a, b, c)
```

**gslcdfpoissonP**

Link to:

```
gsl_cdf_poisson_P(a, b)
```

**gslcdfpoissonQ**

Link to:

```
gsl_cdf_poisson_Q(a, b)
```
gslcdfgeometricP

Link to:
```
gsl_cdf_geometric_P(a, b)
```


gslcdfgeometricQ

Link to:
```
gsl_cdf_geometric_Q(a, b)
```


gslcdfnegativebinomialP

Link to:
```
gsl_cdf_negative_binomial_P(a, b, c)
```


gslcdfnegativebinomialQ

Link to:
```
gsl_cdf_negative_binomial_Q(a, b, c)
```


gslcdfpascalP

Link to:
```
gsl_cdf_pascal_P(a, b, c)
```


gslcdfpascalQ

Link to:
```
gsl_cdf_pascal_Q(a, b, c)
```


gslranbernoullipdf

Link to:
```
gsl_ran_bernoulli_pdf(a, b)
```


gslranbeta

Link to:
```
gsl_ran_beta(a, b, c)
```

4.8. External libraries
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**gslranbetapdf**

Link to:

```
gsl_ran_beta_pdf(a, b, c)
```

**gslranbinomialpdf**

Link to:

```
gsl_ran_binomial_pdf(a, b, c)
```

**gslranexponential**

Link to:

```
gsl_ran_exponential(a, b)
```

**gslranexponentialpdf**

Link to:

```
gsl_ran_exponential_pdf(a, b)
```

**gslranexppow**

Link to:

```
gsl_ran_exppow(a, b, c)
```

**gslranexppowpdf**

Link to:

```
gsl_ran_exppow_pdf(a, b, c)
```

**gslrancauchy**

Link to:

```
gsl_ran_cauchy(a, b)
```

**gslrancauchypdf**

Link to:

```
gsl_ran_cauchy_pdf(a, b)
```
### gslranchisq

**Link to:**

```c
#include <gsl/gsl-randist>

// Random numbers from the chi-squared distribution.

double gsl_ran_chisq (const gsl_rng * r, double a)
```

### gslranchisqpdf

**Link to:**

```c
#include <gsl/gsl-randist>

// Probability density function.

double gsl_ran_chisq_pdf (const gsl_rng * r, double a, double x)
```

### gslranerlang

**Link to:**

```c
#include <gsl/gsl-randist>

// Random numbers from the Erlang distribution.

double gsl_ran_erlang (const gsl_rng * r, double a, double b, double c)
```

### gslranerlangpdf

**Link to:**

```c
#include <gsl/gsl-randist>

// Probability density function.

double gsl_ran_erlang_pdf (const gsl_rng * r, double a, double b, double c)
```

### gslranfdist

**Link to:**

```c
#include <gsl/gsl-randist>

// Random numbers from the F-distribution.

double gsl_ran_fdist (const gsl_rng * r, double a, double b, double c)
```

### gslranfdistpdf

**Link to:**

```c
#include <gsl/gsl-randist>

// Probability density function.

double gsl_ran_fdist_pdf (const gsl_rng * r, double a, double b, double c)
```

### gslranflat

**Link to:**

```c
#include <gsl/gsl-randist>

// Random numbers from the flat distribution.

double gsl_ran_flat (const gsl_rng * r, double a, double b, double c)
```

### gslranflatpdf

**Link to:**

```c
#include <gsl/gsl-randist>

// Probability density function.

double gsl_ran_flat_pdf (const gsl_rng * r, double a, double b, double c)
```
gslrangamma

Link to:

```
gsl_ran_gamma(a, b, c)
```

gslrangammaint

Link to:

```
gsl_ran_gamma_int(a, b, c)
```

gslrangammapdf

Link to:

```
gsl_ran_gamma_pdf(a, b, c)
```

gslrangammamt

Link to:

```
gsl_ran_gamma_mt(a, b, c)
```

gslrangammaknuth

Link to:

```
gsl_ran_gamma_knuth(a, b, c)
```

gslrangaussian

Link to:

```
gsl_ran_gaussian(a, b)
```

gslrangaussianratiomethod

Link to:

```
gsl_ran_gaussian_ratio_method(a, b)
```

gslrangaussianziggurat

Link to:

```
gsl_ran_gaussian_ziggurat(a, b)
```
**gslrangaussianpdf**

Link to:

```c
int gsl_ran_gaussian_pdf(double x, double mean, double sigma);
```

**gslranugaussian**

Link to:

```c
int gsl_ran_ugaussian(double *x);
```

**gslranugaussianratiomethod**

Link to:

```c
int gsl_ran_ugaussian_ratio_method(double *x);
```

**gslranugaussianpdf**

Link to:

```c
int gsl_ran_ugaussian_pdf(double *x);
```

**gslrangaussiantail**

Link to:

```c
int gsl_ran_gaussian_tail(double x, double mean, double sigma);
```

**gslrangaussiantailpdf**

Link to:

```c
int gsl_ran_gaussian_tail_pdf(double x, double mean, double sigma);
```

**gslranugaussiantail**

Link to:

```c
int gsl_ran_ugaussian_tail(double *x);
```

**gslranugaussiantailpdf**

Link to:

```c
int gsl_ran_ugaussian_tail_pdf(double *x);
```
gslranlandau

Link to:

   gsl_ran_landau(a)

gslranlandaupdf

Link to:

   gsl_ran_landau_pdf(a)

gslrangeometricpdf

Link to:

   gsl_ran_geometric_pdf(a, b)

gslrangumbel1

Link to:

   gsl_ran_gumbell(a, b, c)

gslrangumbel1pdf

Link to:

   gsl_ran_gumbell_pdf(a, b, c)

gslrangumbel2

Link to:

   gsl_ran_gumbel2(a, b, c)

gslrangumbel2pdf

Link to:

   gsl_ran_gumbel2_pdf(a, b, c)

gslranlogistic

Link to:

   gsl_ran_logistic(a, b)
**gslranlogisticpdf**

Link to:

```c
int gsl_ran_logistic_pdf(a, b);
```

**gslranlognormal**

Link to:

```c
int gsl_ran_lognormal(a, b, c);
```

**gslranlognormalpdf**

Link to:

```c
int gsl_ran_lognormal_pdf(a, b, c);
```

**gslranlogarithmicpdf**

Link to:

```c
int gsl_ran_logarithmic_pdf(a, b);
```

**gslrnegativebinomialpdf**

Link to:

```c
int gsl_ran_negative_binomial_pdf(a, b, c);
```

**gslranpascalpdf**

Link to:

```c
int gsl_ran_pascal_pdf(a, b, c);
```

**gslranpareto**

Link to:

```c
int gsl_ran_pareto(a, b, c);
```

**gslranparetopdf**

Link to:

```c
int gsl_ran_pareto_pdf(a, b, c);
```
<table>
<thead>
<tr>
<th>Function</th>
<th>Description</th>
<th>Link to</th>
</tr>
</thead>
<tbody>
<tr>
<td>gslranpoissonpdf</td>
<td></td>
<td>gsl_ran_poisson_pdf(a, b)</td>
</tr>
<tr>
<td>gslranrayleigh</td>
<td></td>
<td>gsl_ran_rayleigh(a, b)</td>
</tr>
<tr>
<td>gslranrayleighpdf</td>
<td></td>
<td>gsl_ran_rayleigh_pdf(a, b)</td>
</tr>
<tr>
<td>gslranrayleightail</td>
<td></td>
<td>gsl_ran_rayleigh_tail(a, b, c)</td>
</tr>
<tr>
<td>gslranrayleightailpdf</td>
<td></td>
<td>gsl_ran_rayleigh_tail_pdf(a, b, c)</td>
</tr>
<tr>
<td>gslrantdist</td>
<td></td>
<td>gsl_ran_tdsit(a, b)</td>
</tr>
<tr>
<td>gslrantdistpdf</td>
<td></td>
<td>gsl_ran_tdsit_pdf(a, b)</td>
</tr>
<tr>
<td>gslranlaplace</td>
<td></td>
<td>gsl_ran_laplace(a, b)</td>
</tr>
</tbody>
</table>
**gslranlaplacepdf**

Link to:

```c
int gsl_ran_laplace_pdf(double x, double a, double b)
```

**gslranlevy**

Link to:

```c
int gsl_ran_levy(double x, double a, double b, double c)
```

**gslranweibull**

Link to:

```c
int gsl_ran_weibull(double x, double a, double b, double c)
```

**gslranweibullpdf**

Link to:

```c
int gsl_ran_weibull_pdf(double x, double a, double b, double c)
```

**gslsfairyAi**

Link to:

```c
int gsl_sf_airy_Ai(double x, double b)
```

**gslsfairyBi**

Link to:

```c
int gsl_sf_airy_Bi(double x, double b)
```

**gslsfairyAiscaled**

Link to:

```c
int gsl_sf_airy_Ai_scaled(double x, double b)
```

**gslsfairyBiscaled**

Link to:

```c
int gsl_sf_airy_Bi_scaled(double x, double b)
```
glsfairyAideriv
Link to:

```
gsl_sf_airy_Ai_deriv(a, b)
```

glsfairyBideriv
Link to:

```
gsl_sf_airy_Bi_deriv(a, b)
```

glsfairyAiderivscaled
Link to:

```
gsl_sf_airy_Ai_deriv_scaled(a, b)
```

glsfairyBiderivscaled
Link to:

```
gsl_sf_airy_Bi_deriv_scaled(a, b)
```

glsfairyzeroAi
Link to:

```
gsl_sf_airy_Ai(a, b)
```

glsfairyzeroBi
Link to:

```
gsl_sf_airy_aero_Bi(a)
```

glsfairyzeroAideriv
Link to:

```
gsl_sf_airy_aero_Ai_deriv(a)
```

glsfairyzeroBideriv
Link to:

```
gsl_sf_airy_aero_Bi_deriv(a)
```
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:\texttt{gslsfbesselJ0}
Link to:
\begin{verbatim}
gsl_sf_bessel_J0(a)
\end{verbatim}

:\texttt{gslsfbesselJ1}
Link to:
\begin{verbatim}
gsl_sf_bessel_J1(a)
\end{verbatim}

:\texttt{gslsfbesselJn}
Link to:
\begin{verbatim}
gsl_sf_bessel_Jn(a, b)
\end{verbatim}

:\texttt{gslsfbesselY0}
Link to:
\begin{verbatim}
gsl_sf_bessel_Y0(a)
\end{verbatim}

:\texttt{gslsfbesselY1}
Link to:
\begin{verbatim}
gsl_sf_bessel_Y1(a)
\end{verbatim}

:\texttt{gslsfbesselYn}
Link to:
\begin{verbatim}
gsl_sf_bessel_Yn(a, b)
\end{verbatim}

:\texttt{gslsfbessell0}
Link to:
\begin{verbatim}
gsl_sf_bessel_I0(a)
\end{verbatim}

:\texttt{gslsfbessell1}
Link to:
\begin{verbatim}
gsl_sf_bessel_I1(a)
\end{verbatim}
glsfbselln

Link to:

```
gsl_sf_bessel_In(a, b)
```

glsfbsell0scaled

Link to:

```
gsl_sf_bessel_I0_scaled(a)
```

glsfbsell1scaled

Link to:

```
gsl_sf_bessel_I1_scaled(a)
```

glsfbsellnscaled

Link to:

```
gsl_sf_bessel_In_scaled(a, b)
```

glsfbsellK0

Link to:

```
gsl_sf_bessel_K0(a)
```

glsfbsellK1

Link to:

```
gsl_sf_bessel_K1(a)
```

glsfbsellKn

Link to:

```
gsl_sf_bessel_Kn(a, b)
```

glsfbsellK0scaled

Link to:

```
gsl_sf_bessel_K0_scaled(a)
```
gslsfbesselK1scaled

Link to:

\[
gsl\_sf\_bessel\_K1\_scaled(a)
\]

gslsfbesselKnScaled

Link to:

\[
gsl\_sf\_bessel\_Kn\_scaled(a, b)
\]

gslsfbesselj0

Link to:

\[
gsl\_sf\_bessel\_j0(a)
\]

gslsfbesselj1

Link to:

\[
gsl\_sf\_bessel\_j1(a)
\]

gslsfbesselj2

Link to:

\[
gsl\_sf\_bessel\_j2(a)
\]

gslsfbesseljl

Link to:

\[
gsl\_sf\_bessel\_jl(a, b)
\]

gslsfbessely0

Link to:

\[
gsl\_sf\_bessel\_y0(a)
\]

gslsfbessely1

Link to:

\[
gsl\_sf\_bessel\_y0(a)
\]
**gslsfbessely2**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_y0(a)
```

**gslsfbesselyl**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_jl(a, b)
```

**gslsfbesseli0scaled**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_i0_scaled(a)
```

**gslsfbesseli1scaled**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_i1_scaled(a)
```

**gslsfbesseli2scaled**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_i2_scaled(a)
```

**gslsfbesseliiscaled**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_il_scaled(a, b)
```

**gslsfbesselk0scaled**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_k0_scaled(a)
```

**gslsfbesselk1scaled**

Link to:

```c
#include <gsl/gsl_sf_bessel.h>

gsl_sf_bessel_k1_scaled(a)
```
glsfbsesselk2scaled

Link to:

```
gsl_sf_bessel_k2_scaled(a)
```

glsfbsesselklsscaled

Link to:

```
gsl_sf_bessel_kl_scaled(a, b)
```

glsfbsesselJnu

Link to:

```
gsl_sf_bessel_Jnu(a, b)
```

glsfbsesselYnu

Link to:

```
gsl_sf_bessel_Ynu(a, b)
```

glsfbsesselNuscaled

Link to:

```
gsl_sf_bessel_Inu_scaled(a, b)
```

glsfbsesselInu

Link to:

```
gsl_sf_bessel_Inu(a, b)
```

glsfbsesselKnuscaled

Link to:

```
gsl_sf_bessel_Knu_scaled(a, b)
```

glsfbsesselKnu

Link to:

```
gsl_sf_bessel_Knu(a, b)
```
gslsfbessellnKnu

Link to:

```c
gsl_sf_bessel_lnKnu(a, b)
```

glsfbesselzeroJ0

Link to:

```c
gsl_sf_bessel_zero_J0(a)
```

glsfbesselzeroJ1

Link to:

```c
gsl_sf_bessel_zero_J1(a)
```

glsfbesselzeroJnu

Link to:

```c
gsl_sf_bessel_zero_Jnu(a, b)
```

glsfclausen

Link to:

```c
gsl_sf_clausen(a)
```

glsfhydrogenicR1

Link to:

```c
gsl_sf_hydrogenicR_1(a, b)
```

glsfdawson

Link to:

```c
gsl_sf_dawnson(a)
```

glsfdebye1

Link to:

```c
gsl_sf_debye_1(a)
```
**gslsfdebye2**

Link to:

```c
gsl_sf_debye_2(a)
```

**glsfdebye3**

Link to:

```c
gsl_sf_debye_3(a)
```

**glsfdebye4**

Link to:

```c
gsl_sf_debye_4(a)
```

**glsfdebye5**

Link to:

```c
gsl_sf_debye_5(a)
```

**glsfdebye6**

Link to:

```c
gsl_sf_debye_6(a)
```

**glsfdilog**

Link to:

```c
gsl_sf_dilog(a)
```

**glsf_multiply**

Link to:

```c
gsl_sf_multiply(a, b)
```

**glsfellintKcomp**

Link to:

```c
gsl_sf_ellint_Kcomp(a, b)
```
glsfelltintEcomp

Link to:

\begin{verbatim}
gsl_sf_ellint_Ecomp(a, b)
\end{verbatim}

glsfelltintPcomp

Link to:

\begin{verbatim}
gsl_sf_ellint_Pcomp(a, b, c)
\end{verbatim}

glsfelltintDcomp

Link to:

\begin{verbatim}
gsl_sf_ellint_Dcomp(a, b)
\end{verbatim}

glsfelltintF

Link to:

\begin{verbatim}
gsl_sf_ellint_F(a, b, c)
\end{verbatim}

glsfelltintE

Link to:

\begin{verbatim}
gsl_sf_ellint_E(a, b, c)
\end{verbatim}

glsfelltintRC

Link to:

\begin{verbatim}
gsl_sf_ellint_RC(a, b, c)
\end{verbatim}

glsferfc

Link to:

\begin{verbatim}
gsl_sf_erfc(a)
\end{verbatim}

glsflogerfc

Link to:

\begin{verbatim}
gsl_sf_log_erfc(a)
\end{verbatim}
### gslsferf

Link to:

```
gsl_sf_erf(a)
```

### gslsferfZ

Link to:

```
gsl_sf_erf_Z(a)
```

### gslsferfQ

Link to:

```
gsl_sf_erf_Q(a)
```

### gslsfhazard

Link to:

```
gsl_sf_hazard(a)
```

### gslsfexp

Link to:

```
gsl_sf_exp(a)
```

### gslsfexpmult

Link to:

```
gsl_sf_exp_mult(a, b)
```

### gslsfexpm1

Link to:

```
gsl_sf_expm1(a)
```

### gslsfexprel

Link to:

```
gsl_sf_exprel(a)
```
gslsfexprel2

Link to:

\begin{verbatim}
gsl_sf_exprel_2(a)
\end{verbatim}

gslsfexpreln

Link to:

\begin{verbatim}
gsl_sf_exprel_n(a, b)
\end{verbatim}

gslsfexpintE1

Link to:

\begin{verbatim}
gsl_sf_expint_E1(a)
\end{verbatim}

gslsfexpintE2

Link to:

\begin{verbatim}
gsl_sf_expint_E2(a)
\end{verbatim}

gslsfexpintEn

Link to:

\begin{verbatim}
gsl_sf_expint_En(a, b)
\end{verbatim}

gslsfexpintE1scaled

Link to:

\begin{verbatim}
gsl_sf_expint_E1_scaled(a)
\end{verbatim}

gslsfexpintE2scaled

Link to:

\begin{verbatim}
gsl_sf_expint_E1_scaled(a)
\end{verbatim}

gslsfexpintEnscaled

Link to:

\begin{verbatim}
gsl_sf_expint_En_scaled(a, b)
\end{verbatim}
gslsfexpintEi

Link to:

```
gsl_sf_expint_Ei(a)
```

gslsfexpintEiscaled

Link to:

```
gsl_sf_expint_Ei_scaled(a)
```

gslsfShi

Link to:

```
gsl_sf_Shi(a)
```

gslsfChi

Link to:

```
gsl_sf_Chi(a)
```

gslsfexpint3

Link to:

```
gsl_sf_expint_3(a)
```

gslsfSi

Link to:

```
gsl_sf_Si(a)
```

gslsfCi

Link to:

```
gsl_sf_Ci(a)
```

gslsfatanint

Link to:

```
gsl_sf_atanint(a)
```

4.8. External libraries
gsl_sf_fermi_dirac_m1

Link to:
```
gsl_sf_fermi_dirac_m1(a)
```

gsl_sf_fermi_dirac_0

Link to:
```
gsl_sf_fermi_dirac_0(a)
```

gsl_sf_fermi_dirac_1

Link to:
```
gsl_sf_fermi_dirac_1(a)
```

gsl_sf_fermi_dirac_2

Link to:
```
gsl_sf_fermi_dirac_2(a)
```

gsl_sf_fermi_dirac_int

Link to:
```
gsl_sf_fermi_dirac_int(a, b)
```

gsl_sf_fermi_dirac_mhalf

Link to:
```
gsl_sf_fermi_dirac_mhalf(a)
```

gsl_sf_fermi_dirac_half

Link to:
```
gsl_sf_fermi_dirac_half(a)
```

gsl_sf_fermi_dirac_3half

Link to:
```
gsl_sf_fermi_dirac_3half(a)
```
\texttt{gslsffermidiracin0}

Link to:

\begin{verbatim}
gsl sf_fermi_dirac_inc_0(a, b)
\end{verbatim}

\texttt{glsflngamma}

Link to:

\begin{verbatim}
gsl sf_lngamma(a)
\end{verbatim}

\texttt{glsfgamma}

Link to:

\begin{verbatim}
gsl sf_gamma(a)
\end{verbatim}

\texttt{glsfgammastar}

Link to:

\begin{verbatim}
gsl sf_gammastar(a)
\end{verbatim}

\texttt{glsfgammainv}

Link to:

\begin{verbatim}
gsl sf_gammainv(a)
\end{verbatim}

\texttt{glsftaylorcoeff}

Link to:

\begin{verbatim}
gsl sf_taylorcoeff(a, b)
\end{verbatim}

\texttt{glsffact}

Link to:

\begin{verbatim}
gsl sf_fact(a)
\end{verbatim}

\texttt{glsfdoublefact}

Link to:

\begin{verbatim}
gsl sf_doublefact(a)
\end{verbatim}
gslsflnfact

Link to:

\[ \text{gsl}_\text{sf}_\text{lnfact}(a) \]

gslsflndoublefact

Link to:

\[ \text{gsl}_\text{sf}_\text{lndoublefact}(a) \]

glsflnchoose

Link to:

\[ \text{gsl}_\text{sf}_\text{lnchoose}(a, b) \]

glsfschoose

Link to:

\[ \text{gsl}_\text{sf}_\text{choose}(a, b) \]

glsflnpoch

Link to:

\[ \text{gsl}_\text{sf}_\text{lnpoch}(a, b) \]

glsfpoch

Link to:

\[ \text{gsl}_\text{sf}_\text{poch}(a, b) \]

glsfpochrel

Link to:

\[ \text{gsl}_\text{sf}_\text{pochrel}(a, b) \]

glsfgammaincQ

Link to:

\[ \text{gsl}_\text{sf}_\text{gamma_inc}_\text{Q}(a, b) \]
gslsfgammaincP

Link to:

```
<code>gsl_sf_gamma_inc_P(a, b)</code>
```

glsfgamma\text{inc}

Link to:

```
<code>gsl_sf_gamma_inc(a, b)</code>
```

gsflnbeta

Link to:

```
<code>gsl_sf_lnbeta(a, b)</code>
```

gsfbeta

Link to:

```
<code>gsl_sf_beta(a, b)</code>
```

gsfbetainc

Link to:

```
<code>gsl_sf_betainc(a, b, c)</code>
```

gsfgegenpoly1

Link to:

```
<code>gsl_sf_gegenpoly_1(a, b)</code>
```

gsfgegenpoly2

Link to:

```
<code>gsl_sf_gegenpoly_2(a, b)</code>
```

gsfgegenpoly3

Link to:

```
<code>gsl_sf_gegenpoly_3(a, b)</code>
```
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\textbf{gslsfgegenpolyn}

Link to:

\begin{verbatim}
gsl_sf_gegenpoly_n(a, b, c)
\end{verbatim}

\textbf{glsfhyperg0F1}

Link to:

\begin{verbatim}
gsl_sf_hyperg_0F1(a, b)
\end{verbatim}

\textbf{glsfhyperg1F1int}

Link to:

\begin{verbatim}
gsl_sf_hyperg_1F1_inc(a, b, c)
\end{verbatim}

\textbf{glsfhyperg1F1}

Link to:

\begin{verbatim}
gsl_sf_hyperg_1F1(a, b, c)
\end{verbatim}

\textbf{glsfhypergUint}

Link to:

\begin{verbatim}
gsl_sf_hyperg_U_inc(a, b, c)
\end{verbatim}

\textbf{glsfhypergU}

Link to:

\begin{verbatim}
gsl_sf_hyperg_U(a, b, c)
\end{verbatim}

\textbf{glsfhyperg2F0}

Link to:

\begin{verbatim}
gsl_sf_hyperg_U_2F0(a, b, c)
\end{verbatim}

\textbf{glsflaguerre1}

Link to:

\begin{verbatim}
gsl_sf_laguerre_1(a, b)
\end{verbatim}
**gslflaguerr2**

Link to:

```c
gsl_sf_laguerre_2(a, b)
```

**gslflaguerr3**

Link to:

```c
gsl_sf_laguerre_3(a, b)
```

**gslflaguerrnen**

Link to:

```c
gsl_sf_laguerre_n(a, b, c)
```

**gslflambertW0**

Link to:

```c
gsl_sf_lambert_W0(a)
```

**gslflambertWm1**

Link to:

```c
gsl_sf_lambert_Wm1(a)
```

**gslsflegendrePl**

Link to:

```c
gsl_sf_legendre_Pl(a, b)
```

**gslsflegendreP1**

Link to:

```c
gsl_sf_legendre_P1(a, b)
```

**gslsflegendreP2**

Link to:

```c
gsl_sf_legendre_P2(a)
```
**gslsflegendreP3**

Link to:

```
gsl_sf_legendre_P3(a)
```

**gslsflegendreQ0**

Link to:

```
gsl_sf_legendre_Q0(a)
```

**gslsflegendreQ1**

Link to:

```
gsl_sf_legendre_Q1(a)
```

**gslsflegendreQl**

Link to:

```
gsl_sf_legendre_Ql(a, b)
```

**gslsflegendrePlm**

Link to:

```
gsl_sf_legendre_Plm(a, b, c)
```

**gslsflegendresphPlm**

Link to:

```
gsl_sf_legendre_sphPlm(a, b, c)
```

**gslsflegendrearraysize**

Link to:

```
gsl_sf_legendre_array_size(a, b)
```

**glsfconicalPhalf**

Link to:

```
gsl_sf_conicalP_half(a, b)
```

gslsfconicalPmhalf

Link to:

\[ \text{gsl\_sf\_conicalP\_mhalf}(a, b) \]

gslsfconicalP0

Link to:

\[ \text{gsl\_sf\_conicalP\_0}(a, b) \]

gslsfconicalP1

Link to:

\[ \text{gsl\_sf\_conicalP\_1}(a, b) \]

gslsfconicalPsphreg

Link to:

\[ \text{gsl\_sf\_conicalP\_sph\_reg}(a, b, c) \]

gslsfconicalPcylreg

Link to:

\[ \text{gsl\_sf\_conicalP\_cyl\_reg}(a, b, c) \]

gslsflegendreH3d0

Link to:

\[ \text{gsl\_sf\_legendre\_H3d\_0}(a, b) \]

gslsflegendreH3d1

Link to:

\[ \text{gsl\_sf\_legendre\_H3d\_1}(a, b) \]

gslsflegendreH3d

Link to:

\[ \text{gsl\_sf\_legendre\_H3d}(a, b, c) \]

4.8. External libraries
gslsflog

Link to:

```c
gsl_sf_log(a)
```

glsflogabs

Link to:

```c
gsl_sf_log_abs(a)
```

glsflog1plusx

Link to:

```c
gsl_sf_log_1plusx(a)
```

glsflog1plusxmx

Link to:

```c
gsl_sf_log_1plusx_mx(a)
```

glsfpowint

Link to:

```c
gsl_sf_pow_int(a, b)
```

glsfpsiint

Link to:

```c
gsl_sf_psi_int(a)
```

glsfpsi

Link to:

```c
gsl_sf_psi(a)
```

glsfpsi1piy

Link to:

```c
gsl_sf_psi_1piy(a)
```
**gslfpsi1int**

Link to:
```
gsl_sf_psi_1_int(a)
```

**gslfpsi1**

Link to:
```
gsl_sf_psi_1(a)
```

**gslfpsin**

Link to:
```
gsl_sf_psi_n(a, b)
```

**glsfsynchrotron1**

Link to:
```
gsl_sf_synchrotron_1(a)
```

**glsfsynchrotron2**

Link to:
```
gsl_sf_synchrotron_2(a)
```

**glsftransport2**

Link to:
```
gsl_sf_transport_2(a)
```

**glsftransport3**

Link to:
```
gsl_sf_transport_3(a)
```

**glsftransport4**

Link to:
```
gsl_sf_transport_4(a)
```

---

4.8. External libraries 461
gslsftransport5

Link to:

```
gsl_sf_transport_5(a)
```

gslsfsin

Link to:

```
gsl_sf_sin(a)
```

gslsfcos

Link to:

```
gsl_sf_cos(a)
```

gslsfhypot

Link to:

```
gsl_sf_hypot(a, b)
```

gslsfinc

Link to:

```
gsl_sf_sinc(a)
```

gslsflnsinh

Link to:

```
gsl_sf_lnsinh(a)
```

gslsflncosh

Link to:

```
gsl_sf_lncosh(a)
```

gslsfangledrestrictsymm

Link to:

```
gsl_sf_andle_restrict_symm(a)
```
**gslsfanglerestrictpos**

Link to:

```
gsl_sf_angle_restrict_pos(a)
```

**glsfzetaint**

Link to:

```
gsl_sf_zeta_int(a)
```

**glsfzeta**

Link to:

```
gsl_sf_zeta(a)
```

**glsfzetam1**

Link to:

```
gsl_sf_zetam1(a)
```

**glsfzetam1int**

Link to:

```
gsl_sf_zetam1_int(a)
```

**glsfhzeta**

Link to:

```
gsl_sf_hzeta(a, b)
```

**glsfetaint**

Link to:

```
gsl_sf_eta_int(a)
```

**glsfeta**

Link to:

```
gsl_sf_eta(a)
```

4.8. External libraries
4.8.15  ff-Ipopt

Refer to the Ipopt documentation for more informations.

Todo: todo

4.8.16  fflapack

Refer to the LAPACK documentation for more informations.

inv

Todo: todo

dgeev

Todo: todo

zgeev

Todo: todo

geev

Todo: todo

geev

Todo: todo

dggev

Todo: todo
zggev

Todo: todo

dsygvd

Todo: todo

dgesdd

Todo: todo

zhegv

Todo: todo

dsyev

Todo: todo

zheev

Todo: todo

4.8.17 ff-mmap-semaphore

Wait

Todo: todo

trywait

Todo: todo

4.8. External libraries
Post

Todo: todo

msync

Todo: todo

Read

Todo: todo

Write

Todo: todo

4.8.18 ffnewuoa

newuoa

Todo: todo

4.8.19 ff-NLopt

Refer to the NLOPT documentation for more informations.

nloptDIRECT

Todo: todo

nloptDIRECTL

Todo: todo
nloptDIRECTLRand

Todo: todo

nloptDIRECTScal

Todo: todo

nloptDIRECTNoScal

Todo: todo

nloptDIRECTLNoScal

Todo: todo

nloptDIRECTLRandNoScal

Todo: todo

nloptOrigDIRECT

Todo: todo

nloptOrigDIRECTL

Todo: todo

nloptStoGO

Todo: todo
nloptStoGORand

Todo: todo

nloptLBFGS

Todo: todo

nloptPRAXIS

Todo: todo

nloptVar1

Todo: todo

nloptVar2

Todo: todo

nloptTNewton

Todo: todo

nloptTNewtonRestart

Todo: todo

nloptTNewtonPrecond

Todo: todo
nlOptimizerNewtonPrecondRestart

Todo: todo

nlOptimizerCRS2

Todo: todo

nlOptimizerMMA

Todo: todo

nlOptimizerCOBYLA

Todo: todo

nlOptimizerNEWUOA

Todo: todo

nlOptimizerNEWUOABound

Todo: todo

nlOptimizerNelderMead

Todo: todo

nlOptimizerSbplx

Todo: todo

4.8. External libraries
nloptBOBYQA

Todo: todo

nloptISRES

Todo: todo

nloptSLSQP

Todo: todo

nloptMLSL

Todo: todo

nloptMLSLDS

Todo: todo

nloptAUGLAG

Todo: todo

nloptAUGLAGEQ

Todo: todo

4.8.20 ffrandom

srandomdev

Todo: todo
FreeFEM Documentation, Release 4.6

srandom

Todo: todo

random

Todo: todo

4.8.21 FreeFemQA

MeshGenQA

Todo: todo

4.8.22 freeyams

freeyams

Todo: todo

4.8.23 gmsh

Need

```
load "gmsh"
```

The gmsh software is available [here](#)

gmshload

Load a 2D mesh build with Gmsh.

```
mesh Th = gmshload(MeshFile, [reftri=RefTri], [renum=Renum]);
```

Parameters:

- `MeshFile` (string) Mesh file name
- `reftri` (.. todo:: todo)
- `renum` (.. todo:: todo)

Output:

- `Th(mesh)`

4.8. External libraries 471
gmshload3

Load a 3D mesh build with Gmsh.

```cpp
mesh3 Th = gmshload(MeshFile, [reftet=RefTet], [renum=Renum]);
```

Parameters:

- `MeshFile` (string) Mesh file name
- `reftet` (.. todo:: todo)
- `renum` (.. todo:: todo)

Output:

- `Th` (mesh3)

savegmsh

Todo: todo

4.8.24 gsl

gslpolysolvequadratic

Todo: todo

gslpolysolvecubic

Todo: todo

gslpolycomplexsolve

Todo: todo

gsirnguniform

Todo: todo
gslrnguniformpos

Todo: todo

gslname

Todo: todo

gslrngget

Todo: todo

gslrngmin

Todo: todo

gslrngmax

Todo: todo

gslrngset

Todo: todo

gslrngtype

Todo: todo

4.8.25 ilut

applyilutPrecond

Todo: todo
makellutPrecond

Todo: todo

4.8.26 iohdf5

savehdf5sol

Todo: todo

4.8.27 iovtk

savevtk

Save mesh or solution in vtk/vtu format.

\begin{verbatim}
savevtk(FileName, Th, \{Ux, Uy, Uz\}, p, \{dataname=DataName\},
       \{withsurfacemesh=WithSurfaceMesh\}, \{order=Order\}, \{floatmesh=FloatMesh\},
       \{floatsol=FloatSol\}, \{bin=Bin\}, \{swap=Swap\});
\end{verbatim}

Parameters:

- FileName (string) File name: *.vtk or *.vtu
- Th (mesh or mesh3)
- Ux, Uy, Uz, p (fespace function of vector of fespace functions) Solutions to save, as much as wanted
- dataname= (string) Name of solutions, sepreated by a space
- withsurfacemesh= (bool) .. todo:: todo
- order= (int [int]) Order of solutions.
  Available: 0 or 1
  floatmesh= (bool) .. todo:: todo
  floatsol= (bool) .. todo:: todo
  bin= (bool) If true, save file in binary format
  swap (bool) .. todo:: todo

Output:

- None

vtkload

Todo: todo
vtkload3

Todo: todo

4.8.28 isoline

Need

```cpp
todo
```

isoline

```cpp
int N = isoline(Th, u, xy, iso=Iso, close=Close, smoothing=Smoothing, ratio=Ratio, eps=Eps, beginend=BeginEnd, file=File);
```

Todo: todo

Curve

Todo: todo

Area

Todo: todo

findalloocalmin

Todo: todo

4.8.29 lapack

inv

Todo: todo
dgee

Todo: todo

zgee

Todo: todo

gee

Todo: todo

dgge

Todo: todo

zgge

Todo: todo

dsygvd

Todo: todo

dgesdd

Todo: todo

zhegv

Todo: todo
dsyev

Todo: todo

zheev

Todo: todo

dgelsy

Todo: todo

4.8.30 lgbmo

bmo

Todo: todo

4.8.31 mat_dervieux

MatUpWind1

Todo: todo

4.8.32 mat_psi

MatUpWind0

Todo: todo

4.8.33 medit

medit

Todo: todo
savesol

Todo: todo

readsol

Todo: todo

4.8.34metis

metisnodal

Todo: todo

metisdual

Todo: todo

4.8.35 MetricKuate

MetricKuate

Todo: todo

4.8.36 MetricPk

MetricPk

Todo: todo

4.8.37mmg3d

mmg3d

Todo: todo
4.8.38 mmg3d-v4.0

mmg3d

Todo: todo

4.8.39 msh3

change

Todo: todo

movemesh23

Todo: todo

movemesh2D3Dsurf

Todo: todo

movemesh3

Todo: todo

movemesh

Todo: todo

movemesh3D

Todo: todo
deplacement

Todo: todo

checkbemesh

Todo: todo

buildlayers

Todo: todo

bcube

Todo: todo

cube

Construct a cubic mesh.

```cpp
mesh3 Th = cube(nnX, nnY, nnZ, [X(x), Y(y), Z(z)], [label=Label], [flags=Flags], ...
    [region=Region]);
```

Parameters:

- `nnX(int)` Number of discretization point along \( x \)
- `nnY(int)` Number of discretization point along \( y \)
- `nnZ(int)` Number of discretization point along \( z \)
- `X(x)` (func) \([Optional]\) Affine function of \( x \) to define the length Default: \( x \)
- `Y(y)` (func) \([Optional]\) Affine function of \( y \) to define the width Default: \( y \)
- `Z(z)` (func) \([Optional]\) Affine function of \( z \) to define the height Default: \( z \)
- `label=(int[int])` \([Optional]\)
  - List of surface labels Default: \([1, 2, 3, 4, 5, 6]\)
- `flags=(int)` \([Optional]\)
  - Refer to `square`
- `region=(int)` \([Optional]\)
  - Region number of the cube volume Default: 0

Output:
• Th(mesh3) Cube mesh

trunc

Todo: todo

gluemesh

Todo: todo

extract

Todo: todo

showborder

Todo: todo

getborder

Todo: todo

AddLayers

Todo: todo

4.8.40 mshmet

mshmet

Todo: todo
4.8.41 MUMPS

default to MUMPS seq

Todo: todo

4.8.42 MUMPS_seq

default to MUMPS seq

Todo: todo

4.8.43 netgen

netg

Todo: todo

netgsl

Todo: todo

netgload

Todo: todo

4.8.44 NewSolver

default to UMFPACK

Todo: todo
4.8.45 PARDISO

defaulttoPARDISO

Todo: todo

ompsetnumthreads

Todo: todo

ompgetnumthreads

Todo: todo

ompgetmaxthreads

Todo: todo

4.8.46 pcm2rnm

readpcm

Todo: todo

4.8.47 pipe

flush

Todo: todo

sleep

Todo: todo
usleep

Todo: todo

4.8.48 qf11to25

QF1d

Todo: todo

QF2d

Todo: todo

QF3d

Todo: todo

tripleQF

4.8.49 scotch

scotch

Todo: todo

4.8.50 shell

readdir

Todo: todo

unlink

Todo: todo
rmdir

Todo: todo

cddir

Todo: todo

chdir

Todo: todo

basename

Todo: todo

dirname

Todo: todo

mkdir

Todo: todo

chmod

Todo: todo

cpfile

Todo: todo
FreeFEM Documentation, Release 4.6

stat

Todo: todo

isdir

Todo: todo

getsenv

Todo: todo

setenv

Todo: todo

unsetenv

Todo: todo

4.8.51 splitedges

SplitedgeMesh

Todo: todo

4.8.52 splitmesh12

splitmesh12

Todo: todo
4.8.53 splitmesh3

Todo: todo

4.8.54 splitmesh4

Todo: todo

4.8.55 splitmesh6

Todo: todo

4.8.56 SuperLu

defaulttoSuperLu

Todo: todo

4.8.57 symmetrizeCSR

Todo: todo

4.8.58 tetgen

Refer to the Tetgen documentation for more informations.
tetgconvexhull

Todo: todo

tetgtransfo

Todo: todo

tetg

Build a 3D mesh from a surface.

```c
mesh3 Th = tetg(Th0, [reftet=RefTet], [label=Label], [switch=Switch],
               [nbofholes=NbOfHoles], [holelist=HoleList], [nbofregions=NbOfRegions],
               [regionlist=RegionList], [nboffacetcl=NbOfFaceTcl], [facetcl=FaceTcl])
```

Todo: todo

tetgreconstruction

Todo: todo

4.8.59 UMFPACK64
defaulyttouMFPACK64

Todo: todo

4.8.60 VTK_writer_3d

Vtkaddmesh

Todo: todo
Vtkaddscalar

Todo: todo

4.8.61 VTK_writer

Vtkaddmesh

Todo: todo

Vtkaddscalar
CHAPTER FIVE

MATHEMATICAL MODELS

Summary:
This chapter goes deeper into a number of problems that FreeFEM can solve. It is a complement to the Tutorial part which was only an introduction.

Users are invited to contribute to make this models database grow.

5.1 Static problems

5.1.1 Soap Film

Our starting point here will be the mathematical model to find the shape of soap film which is glued to the ring on the \( xy \)-plane:

\[
C = \{(x, y); \; x = \cos t, \; y = \sin t, \; 0 \leq t \leq 2\pi\}
\]

We assume the shape of the film is described by the graph \((x, y, u(x, y))\) of the vertical displacement \(u(x, y) (x^2 + y^2 < 1)\) under a vertical pressure \(p\) in terms of force per unit area and an initial tension \(\mu\) in terms of force per unit length.

Consider the “small plane” ABCD, \(A: (x, y, u(x, y))\), \(B: (x, y, u(x + \delta x, y))\), \(C: (x, y, u(x + \delta x, y + \delta y))\) and \(D: (x, y, u(x, y + \delta y))\).

Denote by \(\vec{n}(x, y) = (n_x(x, y), n_y(x, y), n_z(x, y))\) the normal vector of the surface \(z = u(x, y)\). We see that the vertical force due to the tension \(\mu\) acting along the edge AD is \(-\mu n_x(x, y)\delta y\) and the the vertical force acting along the edge AD is:

\[
\mu n_x(x + \delta x, y)\delta y \simeq \mu \left( n_x(x, y) + \frac{\partial n_x}{\partial x}(x, y)\right) \delta y
\]
Similarly, for the edges AB and DC we have:

\[-\mu n_y(x, y) \delta x, \quad \mu (n_y(x, y) + \partial n_y/\partial y)(x, y) \delta x\]

The force in the vertical direction on the surface ABCD due to the tension \(\mu\) is given by:

\[\mu (\partial n_x/\partial x) \delta x \delta y + T (\partial n_y/\partial y) \delta y \delta x\]

Assuming small displacements, we have:

\[v_x = (\partial u/\partial x)/\sqrt{1 + (\partial u/\partial x)^2 + (\partial u/\partial y)^2} \approx \partial u/\partial x,\]

\[v_y = (\partial u/\partial y)/\sqrt{1 + (\partial u/\partial x)^2 + (\partial u/\partial y)^2} \approx \partial u/\partial y\]

Letting \(\delta x \to dx, \delta y \to dy\), we have the equilibrium of the vertical displacement of soap film on ABCD by \(p\):

\[\mu dx dy \partial^2 u/\partial x^2 + \mu dx dy \partial^2 u/\partial y^2 + p dx dy = 0\]

Using the Laplace operator \(\Delta = \partial^2/\partial x^2 + \partial^2/\partial y^2\), we can find the virtual displacement write the following:

\[-\Delta u = f \quad \text{in } \Omega\]

where \(f = p/\mu, \Omega = \{(x, y): x^2 + y^2 < 1\}\).

Poisson’s equation appears also in electrostatics taking the form of \(f = \rho/\epsilon\) where \(\rho\) is the charge density, \(\epsilon\) the dielectric constant and \(u\) is named as electrostatic potential.

The soap film is glued to the ring \(\partial \Omega = C\), then we have the boundary condition:

\[u = 0 \quad \text{on } \partial \Omega\]

If the force is gravity, for simplify, we assume that \(f = -1\).

```
// Parameters
int nn = 50;
func f = -1;
func ue = (x^2+y^2-1)/4; //ue: exact solution

// Mesh
border a(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
mesh disk = buildmesh(a(nn));
plot(disk);

// Fespace
fespace fem1(disk, P1);
fem1 u, v;

// Problem
problem laplace (u, v)
    = int2d(disk)( //bilinear form
        dx(u) * dx(v) + dy(u) * dy(v)
    ) - int2d(disk)( //linear form
        f * v
    ) + on(1, u=0); //boundary condition
```

(continues on next page)
// Solve
laplace;

// Plot
plot (u, value=true, wait=true);

// Error
fem1 err = u - ue;
plot(err, value=true, wait=true);

cout << "error L2 = " << sqrt(int2d(disk)(err^2)) << endl;
cout << "error H10 = " << sqrt(int2d(disk)((dx(u)-x/2)^2) + int2d(disk)((dy(u)-y/2)^2)) << endl;

/// Re-run with a mesh adaptation ///

// Mesh adaptation
disk = adaptmesh(disk, u, err=0.01);
plot(disk, wait=true);

// Solve
laplace;
plot (u, value=true, wait=true);

// Error
err = u - ue; //become FE-function on adapted mesh
plot(err, value=true, wait=true);

cout << "error L2 = " << sqrt(int2d(disk)(err^2)) << endl;
cout << "error H10 = " << sqrt(int2d(disk)((dx(u)-x/2)^2) + int2d(disk)((dy(u)-y/2)^2)) << endl;

Fig. 5.1: Isovalue of u
In the 37th line, the $L^2$-error estimation between the exact solution $u_e$,

$$\|u_h - u_e\|_{0, \Omega} = \left( \int_{\Omega} |u_h - u_e|^2 \, dxdy \right)^{1/2}$$

and in the following line, the $H^1$-error seminorm estimation:

$$|u_h - u_e|_{1, \Omega} = \left( \int_{\Omega} |\nabla u_h - \nabla u_e|^2 \, dxdy \right)^{1/2}$$

are done on the initial mesh. The results are $\|u_h - u_e\|_{0, \Omega} = 0.000384045$, $|u_h - u_e|_{1, \Omega} = 0.0375506$.

After the adaptation, we have $\|u_h - u_e\|_{0, \Omega} = 0.000109043$, $|u_h - u_e|_{1, \Omega} = 0.0188411$. So the numerical solution is improved by adaptation of mesh.

### 5.1.2 Electrostatics

We assume that there is no current and a time independent charge distribution. Then the electric field $E$ satisfies:

\[
\begin{align*}
\text{div} E &= \rho / \epsilon \\
\text{curl} E &= 0
\end{align*}
\]

where $\rho$ is the charge density and $\epsilon$ is called the permittivity of free space.

From the equation (5.1) we can introduce the electrostatic potential such that $E = -\nabla \phi$. Then we have Poisson’s equation $-\Delta \phi = f$, $f = -\rho / \epsilon$.

We now obtain the equipotential line which is the level curve of $\phi$, when there are no charges except conductors $\{C_i\}_{1 \ldots N}$. Let us assume $K$ conductors $C_1, \ldots, C_K$ within an enclosure $C_0$.

Each one is held at an electrostatic potential $\phi_i$. We assume that the enclosure $C_0$ is held at potential 0. In order to know $\phi(x)$ at any point $x$ of the domain $\Omega$, we must solve:

\[-\Delta \phi = 0 \quad \text{in } \Omega\]

where $\Omega$ is the interior of $C_0$ minus the conductors $C_i$, and $\Gamma$ is the boundary of $\Omega$, that is $\bigcup_{i=0}^{N} C_i$.

Here $g$ is any function of $x$ equal to $\phi_i$ on $C_i$ and to 0 on $C_0$. The boundary equation is a reduced form for:

\[\phi = \phi_i \text{ on } C_i, \quad i = 1 \ldots N, \phi = 0 \text{ on } C_0.\]

First we give the geometrical informations: $C_0 = \{(x, y); \quad x^2 + y^2 = 5^2\}$, $C_1 = \{(x, y); \quad \frac{1}{2} x^2 (x - 2)^2 + \frac{1}{2} y^2 = 1\}$, $C_2 = \{(x, y); \quad \frac{1}{2} (x^2 + 2)^2 + \frac{1}{2} y^2 = 1\}$.

Let $\Omega$ be the disk enclosed by $C_0$ with the elliptical holes enclosed by $C_1$ and $C_2$. Note that $C_0$ is described counterclockwise, whereas the elliptical holes are described clockwise, because the boundary must be oriented so that the computational domain is to its left.
5.1.3 Aerodynamics

Let us consider a wing profile $S$ in a uniform flow. Infinity will be represented by a large circle $\Gamma_{\infty}$. As previously, we must solve:

$$\Delta \varphi = 0 \quad \text{in } \Omega, \quad \varphi|_S = c, \quad \varphi|_{\Gamma_{\infty}} = u_{\infty 1x} - u_{\infty 2x}$$  \hspace{1cm} (5.2)

where $\Omega$ is the area occupied by the fluid, $u_{\infty}$ is the air speed at infinity, $c$ is a constant to be determined so that $\partial_n \varphi$ is continuous at the trailing edge $P$ of $S$ (so-called Kutta-Joukowski condition). Lift is proportional to $c$. 

5.1. Static problems 495
To find $c$ we use a superposition method. As all equations in (5.2) are linear, the solution $\varphi_c$ is a linear function of $c$

$$\varphi_c = \varphi_0 + c\varphi_1$$

where $\varphi_0$ is a solution of (5.2) with $c = 0$ and $\varphi_1$ is a solution with $c = 1$ and zero speed at infinity.

With these two fields computed, we shall determine $c$ by requiring the continuity of $\partial \varphi / \partial n$ at the trailing edge. An equation for the upper surface of a NACA0012 (this is a classical wing profile in aerodynamics; the rear of the wing is called the trailing edge) is:

$$y = 0.17735\sqrt{x} - 0.075597x - 0.212836x^2 + 0.17363x^3 - 0.06254x^4$$

Taking an incidence angle $\alpha$ such that $\tan \alpha = 0.1$, we must solve:

$$-\Delta \varphi = 0 \quad \text{in } \Omega, \quad \varphi|_{\Gamma_1} = y - 0.1x, \quad \varphi|_{\Gamma_2} = c$$

where $\Gamma_2$ is the wing profile and $\Gamma_1$ is an approximation of infinity. One finds $c$ by solving:

$$-\Delta \varphi_0 = 0 \quad \text{in } \Omega, \quad \varphi_0|_{\Gamma_1} = y - 0.1x, \quad \varphi_0|_{\Gamma_2} = 0,$$

$$-\Delta \varphi_1 = 0 \quad \text{in } \Omega, \quad \varphi_1|_{\Gamma_1} = 0, \quad \varphi_1|_{\Gamma_2} = 1$$

The solution $\varphi = \varphi_0 + c\varphi_1$ allows us to find $c$ by writing that $\partial \varphi / \partial n$ has no jump at the trailing edge $P = (1, 0)$.

We have $\partial_n \varphi - (\varphi(P^+) - \varphi(P))/\delta$ where $P^+$ is the point just above $P$ in the direction normal to the profile at a distance $\delta$. Thus the jump of $\partial_n \varphi$ is $(\varphi_0|_{P^+} + c(\varphi_1|_{P^+} - 1)) + (\varphi_0|_{P^-} + c(\varphi_1|_{P^-} - 1))$ divided by $\delta$ because the normal changes sign between the lower and upper surfaces. Thus

$$c = -\frac{\varphi_0|_{P^+} + \varphi_0|_{P^-}}{(\varphi_1|_{P^+} + \varphi_1|_{P^-} - 2)},$$

which can be programmed as:

$$c = -\frac{\varphi_0(0.99, 0.01) + \varphi_0(0.99, -0.01)}{(\varphi_1(0.99, 0.01) + \varphi_1(0.99, -0.01) - 2)}.$$
Isovalue of $c_p = -(\partial_x \psi)^2 - (\partial_y \psi)^2$

Zooming of $c_p$

```plaintext
// Problem
solve Joukowski0(psi0, vh) = int2d(Th)(
    dx(psi0) * dx(vh)
    + dy(psi0) * dy(vh)
) + on(a, psi0=y-0.1*x)
    + on(upper, lower, psi0=0);
plot(psi0);

solve Joukowski1(psi1, vh) = int2d(Th)(
    dx(psi1) * dx(vh)
    + dy(psi1) * dy(vh)
) + on(a, psi1=0)
    + on(upper, lower, psi1=1);
plot(psi1);

// continuity of pressure at trailing edge
real beta = psi0(0.99,0.01) + psi0(0.99,-0.01);
beta = -beta / (psi1(0.99,0.01) + psi1(0.99,-0.01)-2);
Vh psi = beta*psi1 + psi0;
plot(psi);

ZVh Zpsi = psi;
plot(Zpsi, bw=true);

ZVh cp = -dx(psi)^2 - dy(psi)^2;
plot(cp);

ZVh Zcp = cp;
plot(Zcp, nbiso=40);
```

5.1. Static problems
5.1.4 Error estimation

There are famous estimation between the numerical result $u_h$ and the exact solution $u$ of the Poisson’s problem:

If triangulations $\{T_h\}_{h \downarrow 0}$ is regular (see Regular Triangulation), then we have the estimates:

$$
|\nabla u - \nabla u_h|_{0, \Omega} \leq C_1 h
$$

$$
\|u - u_h\|_{0, \Omega} \leq C_2 h^2
$$

with constants $C_1, C_2$ independent of $h$, if $u$ is in $H^2(\Omega)$. It is known that $u \in H^2(\Omega)$ if $\Omega$ is convex.

In this section we check (5.3). We will pick up numericall error if we use the numerical derivative, so we will use the following for (5.3).

$$
\int_{\Omega} |\nabla u - \nabla u_h|^2 \, dx \, dy = \int_{\Omega} \nabla u \cdot (u - 2u_h) \, dx \, dy + \int_{\Omega} \nabla u_h \cdot \nabla u_h \, dx \, dy
$$

The constants $C_1, C_2$ are depend on $T_h$ and $f$, so we will find them by FreeFEM.

In general, we cannot get the solution $u$ as a elementary functions even if spetical functions are added. Instead of the exact solution, here we use the approximate solution $u_0$ in $V_h(T_h, P_2)$, $h \sim 0$.

```
1 // Parameters
2 func f = x*y;
3
4 //Mesh
5 mesh Th0 = square(100, 100);
6
7 // Fespace
8 fespace V0h(Th0, P2);
9 V0h u0, v0;
10
11 // Problem
12 solve Poisson0 (u0, v0)
13 = int2d(Th0) (dx(u0)*dx(v0) + dy(u0)*dy(v0))
14 - int2d(Th0) (f*v0)
15 + on(1, 2, 3, 4, u0=0)
16 ;
17 plot(u0);
18
19 // Error loop
20 real[int] errL2(10), errH1(10);
21 for (int i = 1; i <= 10; i++) {
22    // Mesh
23    mesh Th = square(5+i*3,5+i*3);
24
25    // Fespace
26    fespace Vh(Th, P1);
27    Vh u, v;
28    fespace Ph(Th, P0);
29    Ph h = hTriangle;  //get the size of all triangles
30
31    // Problem
32    solve Poisson (u, v)
```

(continues on next page)
38 \begin{verbatim}
= int2d(Th)(
   dx(u)\times dx(v)
   + dy(u)\times dy(v)
 )
- int2d(Th)(
   f\times v
)
+ on(1, 2, 3, 4, u=0)
;
// Error
V0h uu = u; //interpolate solution on first mesh
errL2[i-1] = sqrt( int2d(Th0)((uu - u0)^2) )/h[].max^2;
errH1[i-1] = sqrt( int2d(Th0)(f*(u0 - 2*uu + uu)) )/h[].max;
}

// Display
cout << "C1 = " << errL2.max << "("<<errL2.min<<")" << endl;
cout << "C2 = " << errH1.max << "("<<errH1.min<<")" << endl;
\end{verbatim}

We can guess that $C_1 = 0.0179253(0.0173266)$ and $C_2 = 0.0729566(0.0707543)$, where the numbers inside the parentheses are minimum in calculation.

### 5.1.5 Periodic Boundary Conditions

We now solve the Poisson equation:

\[-\Delta u = \sin(x + \pi/4) \times \cos(y + \pi/4)\]

on the square $]0, 2\pi[^2$ under bi-periodic boundary condition $u(0, y) = u(2\pi, y)$ for all $y$ and $u(x, 0) = u(x, 2\pi)$ for all $x$.

These boundary conditions are achieved from the definition of the periodic finite element space.

\begin{verbatim}
// Parameters
func f = sin(x*pi/4.)\times cos(y*pi/4.); //right hand side

// Mesh
mesh Th = square(10, 10, [2*x*pi, 2*y*pi]);

// Fespace
//defined the fespace with periodic condition
//label: 2 and 4 are left and right side with y the curve abscissa
// 1 and 2 are bottom and upper side with x the curve abscissa
fespace Vh(Th, P2, periodic=[[2, y], [4, y], [1, x], [3, x]]);
Vh uh, vh;

// Problem
problem laplace (uh, vh)
   = int2d(Th)(
   dx(uh)\times dx(vh)
   + dy(uh)\times dy(vh)
 )
   + int2d(Th)(
   - f\times vh
\end{verbatim}
The periodic condition does not necessarily require parallel boundaries. The following example give such example.

**Tip:** Periodic boundary conditions - non-parallel boundaries

```plaintext
// Parameters
int n = 10;
real r = 0.25;
real r2 = 1.732;
func f = (y+x+1)*(y+x-1)*(y-x+1)*(y-x-1);

// Mesh
border a(t=0, 1){x=-t+1; y=t; label=1;};
border b(t=0, 1){x=-t; y=1-t; label=2;};
border c(t=0, 1){x=t-1; y=-t; label=3;};
border d(t=0, 1){x=t; y=-1+t; label=4;};
border e(t=0, 2*pi){x=r*cos(t); y=-r*sin(t); label=0;};
mesh Th = buildmesh(a(n) + b(n) + c(n) + d(n) + e(n));
plot(Th, wait=true);

// Fespace
//warning for periodic condition:
//side a and c
```

Fig. 5.5: The isovalue of solution \( u \) with periodic boundary condition
// on side a (label 1) $ x \in [0,1] $ or $ x-y\in [-1,1] $ 
// on side c (label 3) $ x \in [-1,0]$ or $ x-y\in[-1,1] $ 
//so the common abscissa can be respectively $x$ and $x+1$ 
//or you can can try curviline abscissa $x-y$ and $x+y$

// 1 first way
fespace Vh(Th, P2, periodic=[[2, 1+x], [4, x], [1, x], [3, 1+x]]);

Vh uh, vh;

// Problem
real intf = int2d(Th)(f);
real mTh = int2d(Th)(1);
real k = intf / mTh;
problem laplace (uh, vh)
= int2d(Th) ( dx(uh)*dx(vh) + dy(uh)*dy(vh) )
  + int2d(Th) ( (k-f)*vh )
;

// Solve
laplace;

// Plot
plot(uh, wait=true);

Fig. 5.6: The isovalue of solution $u$ for $\Delta u = ((y+x)^2 + 1)((y-x)^2 + 1) - k$, in $\Omega$ and $\partial_\nu u = 0$ on hole, and with two periodic boundary condition on external border

5.1. Static problems
Tip: Periodic boundary conditions - non-equal border

```c
// Macro
//irregular boundary condition to build border AB
macro LINEBORDER(A, B, lab)
  border A#B(t=0,1){ real t1=1.-t;
  x=A#x+t1*B#x*t;
  y=A#y+t1*B#y*t;
  label=lab; } //EOM
// compute \||AB|| A=(ax,ay) et B=(bx,by)
macro dist(ax, ay, bx, by) sqrt(square((ax)-(bx)) + square((ay)-(by))) //EOM
macro Grad(u) [dx(u), dy(u)] //EOM

// Parameters
int n = 10;
real Ax = 0.9, Ay = 1;
real Bx = 2, By = 1;
real Cx = 2.5, Cy = 2.5;
real Dx = 1, Dy = 2;
real gx = (Ax+Bx+Cx+Dx)/4.;
real gy = (Ay+By+Cy+Dy)/4.;

// Mesh
LINEBORDER(A, B, 1)
LINEBORDER(B, C, 2)
LINEBORDER(C, D, 3)
LINEBORDER(D, A, 4)
mesh Th=buildmesh(AB(n)+BC(n)+CD(n)+DA(n), fixedborder=1);

// Fespace
real l1 = dist(Ax,Ay,Bx,By);
real l2 = dist(Bx,By,Cx,Cy);
real l3 = dist(Cx,Cy,Dx,Dy);
real l4 = dist(Dx,Dy,Ax,Ay);
func s1 = dist(Ax,Ay,x,y)/l1; //absisse on AB = ||AX||/||AB||
func s2 = dist(Bx,By,x,y)/l2; //absisse on BC = ||BX||/||BC||
func s3 = dist(Cx,Cy,x,y)/l3; //absisse on CD = ||CX||/||CD||
func s4 = dist(Dx,Dy,x,y)/l4; //absisse on DA = ||DX||/||DA||
verbosity = 6; //to see the abscissee value of the periodic condition
fespace Vh(Th, P1, periodic=[[1, s1], [3, s3], [2, s2], [4, s4]]);
verbosity = 1; //reset verbosity
Vh u, v;
real cc = 0;
cc = int2d(Th)((x-gx)*(y-gy)-cc)/Th.area;
cout << "compatibility = " << int2d(Th)((x-gx)*(y-gy)-cc) <<endl;

// Problem
solve Poisson(u, v)
  = int2d(Th)(
    Grad(u)'*Grad(v)
    + 1e-10*u*v
  )
  - int2d(Th)(
      ...
  )
(continues on next page)
```
10*v*((x-gx)*(y-gy)-cc)
)
;
// Plot
plot(u, value=true);

Tip: Periodic boundry conditions - Poisson cube-balloon

load "msh3" load "tetgen" load "medit"

// Parameters
real hs = 0.1; //mesh size on sphere
int[int] N = [20, 20, 20];
real [int,int] B = [[-1, 1], [-1, 1], [-1, 1]];
int [int,int] L = [[1, 2], [3, 4], [5, 6]];
real x0 = 0.3, y0 = 0.4, z0 = 0.6;
func f = sin(x*2*pi+x0)*sin(y*2*pi+y0)*sin(z*2*pi+z0);

// Mesh
bool buildTh = 0;
mesh3 Th;
try { //a way to build one time the mesh or read it if the file exist
   Th = readmesh3("Th-hex-sph.mesh");
}
catch (...){
   buildTh = 1;
}

if (buildTh){
   include "MeshSurface.idp"

   // Surface Mesh
   mesh3 ThH = SurfaceHex(N, B, L, 1);
   mesh3 ThS = Sphere(0.5, hs, 7, 1);
   mesh3 ThHS = ThH + ThS;
   real voltet = (hs^3)/6.;
   real[int] domain = [0, 0, 0, 1, voltet, 0, 0, 0.7, 2, voltet];
   Th = tetg(ThHS, switch="pqaAAYYQ", nbofregions=2, regionlist=domain);
   savemesh(Th, "Th-hex-sph.mesh");
}

// Fespace
fespace Ph(Th, P0);
Ph reg = region;
cout << " centre = " << reg(0,0,0) << endl;
cout << " exterier = " << reg(0,0,0.7) << endl;
verbosity = 50;
fespace Vh(Th, P1, periodic=[[3, x, z], [4, x, z], [1, y, z], [2, y, z], [5, x, y], ...
   →[5, x, y]]);

5.1. Static problems
5.1.6 Poisson Problems with mixed boundary condition

Here we consider the Poisson equation with mixed boundary conditions:

For given functions $f$ and $g$, find $u$ such that:

$$
-\Delta u = f \quad \text{in } \Omega
$$

$$
u = g \quad \text{on } \Gamma_D
$$

$$
\frac{\partial u}{\partial n} = 0 \quad \text{on } \Gamma_N
$$
where $\Gamma_D$ is a part of the boundary $\Gamma$ and $\Gamma_N = \Gamma \setminus \Gamma_D$.

The solution $u$ has the singularity at the points $\{\gamma_1, \gamma_2\} = \Gamma_D \cap \Gamma_N$.

When $\Omega = \{(x, y); -1 < x < 1, 0 < y < 1\}$, $\Gamma_N = \{(x, y); -1 \leq x < 0, y = 0\}$, $\Gamma_D = \partial \Omega \setminus \Gamma_N$, the singularity will appear at $\gamma_1 = (0, 0)$, $\gamma_2 = (-1, 0)$, and $u$ has the expression:

$$u = K_i u_S + u_R, \ u_R \in H^2(\text{near } \gamma_i), \ i = 1, 2$$

with a constants $K_i$.

Here $u_S = r_j^{1/2} \sin(\theta_j/2)$ by the local polar coordinate $(r_j, \theta_j)$ at $\gamma_j$ such that $(r_1, \theta_1) = (r, \theta)$.

Instead of polar coordinate system $(r, \theta)$, we use that $r = \sqrt{x^2 + y^2}$ and $\theta = \text{atan2}(y, x)$ in FreeFEM.

Assume that $f = -2 \times 30(x^2 + y^2)$ and $g = u_e = 10(x^2 + y^2)^{1/4} \sin\left(\frac{\tan^{-1}(y/x)}{2}\right) + 30(x^2y^2)$, where $u_e S$ is the exact solution.

```plaintext
1 // Parameters
2 func f = -2*30*(x^2+y^2); //given function
3 //the singular term of the solution is K*us (K: constant)
4 func us = sin(atan2(y,x)/2)*sqrt( sqrt(x^2+y^2) );
5 real K = 10.;
6 func ue = K*us + 30*(x^2*y^2);
7
8 // Mesh
9 border N(t=0, 1){x=-1+t; y=0; label=1;};
10 border D1(t=0, 1){x=t; y=0; label=2;};
11 border D2(t=0, 1){x=1; y=t; label=2;};
12 border D3(t=0, 2){x=1-t; y=1; label=2;};
13 border D4(t=0, 1){x=-1; y=1-t; label=2;};
14 mesh T0h = buildmesh(N(10) + D1(10) + D2(10) + D3(20) + D4(10));
15 plot(T0h, wait=true);
16
17 // Fespace
18 fespace V0h(T0h, P1);
19 V0h u0, v0;
20
21 //Problem
22 solve Poisson0 (u0, v0)
23 = int2d(T0h) ( dx(u0)*dx(v0) + dy(u0)*dy(v0) )
24 - int2d(T0h) ( f*v0 )
25 + on(2, u0=ue);
26
27 // Mesh adaptation by the singular term
28 mesh Th = adaptmesh(T0h, us);
29 for (int i = 0; i < 5; i++)
30 mesh Th = adaptmesh(Th, us);
31
32 // Fespace
33 fespace Vh(Th, P1);
34 Vh u, v;
35```

(continues on next page)
From line 35 to 37, mesh adaptations are done using the base of singular term.

In line 61, $H1e = |u_e|_{1,\Omega}$ is calculated.

In lines 64 and 65, the relative errors are calculated, that is:

\[
\frac{||u^0_h - u_e||_{1,\Omega}}{H1e} = 0.120421 \\
\frac{||u^a_h - u_e||_{1,\Omega}}{H1e} = 0.0150581
\]

where $u^0_h$ is the numerical solution in $T0h$ and $u^a_h$ is $u$ in this program.

### 5.1.7 Poisson with mixed finite element

Here we consider the Poisson equation with mixed boundary value problems:

For given functions $f, g_d, g_n$, find $p$ such that

\[
\begin{align*}
-\Delta p &= 1 \quad \text{in } \Omega \\
p &= g_d \quad \text{on } \Gamma_D \\
\partial p / \partial n &= g_n \quad \text{on } \Gamma_N
\end{align*}
\]

where $\Gamma_D$ is a part of the boundary $\Gamma$ and $\Gamma_N = \Gamma \setminus \Gamma_D$.

The mixed formulation is: find $p$ and $u$ such that:

\[
\begin{align*}
\nabla p + u &= 0 \quad \text{in } \Omega \\
\nabla . u &= f \quad \text{in } \Omega \\
p &= g_d \quad \text{on } \Gamma_D \\
\partial u . n &= g_n . n \quad \text{on } \Gamma_N
\end{align*}
\]

where $g_n$ is a vector such that $g_n . n = g_n$. 

From line 35 to 37, mesh adaptations are done using the base of singular term.

In line 61, $H1e = |u_e|_{1,\Omega}$ is calculated.

In lines 64 and 65, the relative errors are calculated, that is:

\[
\frac{||u^0_h - u_e||_{1,\Omega}}{H1e} = 0.120421 \\
\frac{||u^a_h - u_e||_{1,\Omega}}{H1e} = 0.0150581
\]

where $u^0_h$ is the numerical solution in $T0h$ and $u^a_h$ is $u$ in this program.
The variational formulation is:

\[ \forall v \in V_0 : \int_\Omega p \nabla \cdot v + vv = \int_\Gamma g_d v_n \]
\[ \forall q \in P : \int_\Omega q \nabla \cdot u = \int_\Omega q f \]
\[ \frac{\partial u}{\partial n} = g_n \text{ on } \Gamma_N \]

where the functional spaces are:

\[ P = L^2(\Omega), \quad V = H(div) = \{ v \in L^2(\Omega)^2, \nabla \cdot v \in L^2(\Omega) \} \]

and:

\[ V_0 = \{ v \in V; \ v.n = 0 \text{ on } \Gamma_N \} \]

To write the **FreeFEM** example, we have just to choose the finite elements spaces.

Here \( V \) space is discretized with Raviart-Thomas finite element \( RT_0 \) and \( P \) is discretized by constant finite element \( P_0 \).

**Example 9.10** LaplaceRT.edp

```plaintext
// Parameters
func gd = 1.;
func g1n = 1.;
func g2n = 1.;

// Mesh
mesh Th = square(10, 10);

// Fespace
fespace Vh(Th, RT0);
Vh [u1, u2];
Vh [v1, v2];

fespace Ph(Th, P0);
Ph p, q;

// Problem
problem laplaceMixte ([u1, u2, p], [v1, v2, q], solver=GMRES, eps=1.0e-10, tgv=1e30, ...
  dimKrylov=150)
  = int2d(Th)(
      p*q*1e-15 //this term is here to be sure
      // that all sub matrix are inversible (LU requirement)
      + u1*v1
      + u2*v2
      + p*(dx(v1)+dy(v2))
      + (dx(u1)+dy(u2))*q
      )
  + int2d(Th) (q
    )
  - int1d(Th, 1, 2, 3)(
      gd*(vl*N.x +v2*N.y)
    )
  + on(4, u1=g1n, u2=g2n)
;

// Solve
laplaceMixte;
```

(continues on next page)
5.1.8 Metric Adaptation and residual error indicator

We do metric mesh adaption and compute the classical residual error indicator $\eta_T$ on the element $T$ for the Poisson problem.

First, we solve the same problem as in a previous example.

```cpp
// Parameters
real[int] viso(21);
for (int i = 0; i < viso.n; i++)
viso[i] = 10.^(+(i-16.)/2.);
real error = 0.01;
func f = (x-y);

// Mesh
border ba(t=0, 1.0){x=t; y=0; label=1;}
border bb(t=0, 0.5){x=1; y=t; label=2;}
border bc(t=0, 0.5){x=1-t; y=0.5; label=3;}
border bd(t=0.5, 1){x=0.5; y=t; label=4;}
border be(t=0.5, 1){x=1-t; y=1; label=5;}
border bf(t=0.0, 1){x=0; y=1-t; label=6;}
mesh Th = buildmesh(ba(6) + bb(4) + bc(4) + bd(4) + be(4) + bf(6));

// Fespace
fespace Vh(Th, P2);
Vh u, v;

fespace Nh(Th, P0);
Nh rho;

// Problem
problem Problem1 (u, v, solver=CG, eps=1.0e-6)
= int2d(Th, qforder=5)
  u*v+1.0e-10
  + dx(u)*dx(v)
  + dy(u)*dy(v)
  + int2d(Th, qforder=5)
  - f*v
;
```

Now, the local error indicator $\eta_T$ is:

$$\eta_T = \left( h_T^2 \| f + \Delta u_h \|^2_{L^2(T)} + \sum_{e \in \mathcal{E}_T} h_e \left\| \frac{\partial u_h}{\partial n_e} \right\|^2_{L^2(e)} \right)^{\frac{1}{2}}$$

where $h_T$ is the longest edge of $T$, $\mathcal{E}_T$ is the set of $T$ edge not on $\Gamma = \partial \Omega$, $n_T$ is the outside unit normal to $K$, $h_e$ is the length of edge $e$, $[g]$ is the jump of the function $g$ across edge (left value minus right value).

Of course, we can use a variational form to compute $\eta_T^2$, with test function constant function in each triangle.
5.1.9 Adaptation using residual error indicator

In the previous example we compute the error indicator, now we use it, to adapt the mesh. The new mesh size is given by the following formulae:

\[
h_{n+1}(x) = \frac{h_n(x)}{f_n(\eta_K(x))}
\]

If the method is correct, we expect to look the graphics by an almost constant function \( \eta \) on your computer as in Fig. 5.8a and Fig. 5.8b.
where $\eta_n(x)$ is the level of error at point $x$ given by the local error indicator, $h_n$ is the previous “mesh size” field, and $f_n$ is a user function defined by $f_n = \min(3, \max(1/3, \eta_n/\eta_n^*))$ where $\eta_n^* = \text{mean}(\eta_n)c$, and $c$ is a user coefficient generally close to one.

First a macro `MeshSizeComputation` is defined to get a $P_1$ mesh size as the average of edge length.

```plaintext
// macro the get the current mesh size parameter
// in:
// Th the mesh
// Vh P1 fespace on Th
// out :
// h: the Vh finite element finite set to the current mesh size
macro MeshSizeComputation (Th, Vh, h)
{
    real[int] count(Th.nv);
    /*mesh size (lenEdge = integral(e) 1 ds)*/
    varf vmeshsizen (u, v) = intalledges(Th, qfnbpE=1)(v);
    /*number of edges per vertex*/
    varf vedgecount (u, v) = intalledges(Th, qfnbpE=1)(v/lenEdge);
    /*mesh size*/
    count = vedgecount(0, Vh);
    h[] = 0.;
    h[] = vmeshsizen(0, Vh);
    cout << "count min = " << count.min << " max = " << count.max << endl;
    h[] = h[].count;
    cout << "-- bound meshsize = " << h[].min << " " << h[].max << endl;
}
```

A second macro to re-mesh according to the new mesh size.

```plaintext
// macro to remesh according the de residual indicator
// in:
// Th the mesh
// Ph P0 fespace on Th
// Vh P1 fespace on Th
// vindicator the varf to evaluate the indicator
// coef on etameam
macro ReMeshIndicator (Th, Ph, Vh, vindicator, coef)
{
    Vh h=0;
    //evaluate the mesh size*/
    MeshSizeComputation(Th, Vh, h);
    Ph etak;
    etak[] = vindicator(0, Ph);
    etak[] = sqrt(etak[]);
    real etastar= coef*(etak[].sum/etak[].n);
    cout << "etastar = " << etastar << " sum = " << etak[].sum << " " << endl;
    /*etaK is discontinuous*/
    /*we use P1 L2 projection with mass lumping*/
    Vh fn, sigma;
    varf veta(unused, v) = int2d(Th)(etak*v);
    varf vun(unused, v) = int2d(Th)(1*v);
    fn[] = veta(0, Vh);
    sigma[] = vun(0, Vh);
    fn[] = fn[]./ sigma[];
    fn = max(min(fn/etastar,3.),0.3333);
}
```

(continues on next page)
```csharp
/*new mesh size*/
  h = h / fn;
/*build the mesh*/
  Th = adaptmesh(Th, IsMetric=1, h, splitpbedge=1, nbvx=10000);
}

// Parameters
real hinit = 0.2; // initial mesh size
func f=(x-y);

// Mesh
border ba(t=0, 1.0){x=t; y=0; label=1;}
border bb(t=0, 0.5){x=1-t; y=t; label=2;}
border bc(t=0, 0.5){x=1-t; y=0.5; label=3;}
border bd(t=0.5, 1){x=0.5; y=t; label=4;}
border be(t=0.5, 1){x=1-t; y=1; label=5;}
border bf(t=0.0, 1){x=0; y=1-t; label=6;}
mesh Th = buildmesh(ba(6) + bb(4) + bc(4) + bd(4) + be(4) + bf(6));

// Fespace
fespace Vh(Th, P1); // for the mesh size and solution
  Vh h = hinit; // the FE function for the mesh size
fespace Ph(Th, P0); // for the error indicator

// Build a mesh with the given mesh size hinit
  Th = adaptmesh(Th, h, IsMetric=1, splitpbedge=1, nbvx=10000);
  plot(Th, wait=1);

// Problem
problem Poisson (u, v)
  = int2d(Th, qforder=5)(
          u*v+1.0e-10
          + dx(u)*dx(v)
          + dy(u)*dy(v)
        )
    - int2d(Th, qforder=5)(
          f*v
        )
  ;

varf indicator2 (unused, chiK)
  = intalledges(Th){
      chiK*lenEdge*square(jump(N.x*dx(u) + N.y*dy(u)))
    }
  + int2d(Th){
      chiK*square(hTriangle*(f + dxx(u) + dyy(u)))
    }
  ;

// Mesh adaptation loop
for (int i = 0; i < 10; i++){
  u = u;
}

// Solve
```

5.1. Static problems
5.2 Elasticity

Consider an elastic plate with undeformed shape $\Omega \times [0, h]$ in $\mathbb{R}^3$, $\Omega \subset \mathbb{R}^2$.

By the deformation of the plate, we assume that a point $P(x_1, x_2, x_3)$ moves to $P(\xi_1, \xi_2, \xi_3)$. The vector $u = (u_1, u_2, u_3) = (\xi_1 - x_1, \xi_2 - x_2, \xi_3 - x_3)$ is called the displacement vector.

By the deformation, the line segment $x, x + \Delta x$ moves approximately to $x + u(x), x + \tau \Delta x + u(x + \tau \Delta x)$ for small $\tau$, where $x = (x_1, x_2, x_3), \Delta x = (\Delta x_1, \Delta x_2, \Delta x_3)$.

We now calculate the ratio between two segments:

$$\eta(\tau) = \frac{\tau^{-1}|\Delta x|^{-1}}{|(u(x + \tau \Delta x) - u(x) + \tau \Delta x) - \tau |\Delta x|}$$

then we have (see e.g. [NECAS2017], p.32)

$$\lim_{\tau \to 0} \eta(\tau) = (1 + 2\epsilon_{ij}\nu_i \nu_j)^{1/2} - 1, \quad 2\epsilon_{ij} = \frac{\partial u_k}{\partial x_i} \frac{\partial u_k}{\partial x_j} + \left( \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right)$$

where $\nu_i = \Delta x_i |\Delta x|^{-1}$. If the deformation is small, then we may consider that:

$$(\partial u_k / \partial x_i)(\partial u_k / \partial x_i) \approx 0$$

and the following is called small strain tensor:

$$\varepsilon_{ij}(u) = \frac{1}{2} \left( \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right)$$
The tensor $e_{ij}$ is called finite strain tensor. Consider the small plane $\Delta \Pi(x)$ centered at $x$ with the unit normal direction $n = (n_1, n_2, n_3)$, then the surface on $\Delta \Pi(x)$ at $x$ is:

$$\left(\sigma_{ij}(x) n_j, \sigma_{2j}(x) n_j, \sigma_{3j}(x) n_j\right)$$

where $\sigma_{ij}(x)$ is called stress tensor at $x$. Hooke’s law is the assumption of a linear relation between $\sigma_{ij}$ and $\epsilon_{ij}$ such as:

$$\sigma_{ij}(x) = c_{ijkl}(x) \epsilon_{ij}(x)$$

with the symmetry $c_{ijkl} = c_{jikl}, c_{ijkl} = c_{ijlk}, c_{ijkl} = c_{klij}$. If Hooke’s tensor $c_{ijkl}(x)$ do not depend on the choice of coordinate system, the material is called isotropic at $x$. If $c_{ijkl}$ is constant, the material is called homogeneous. In homogeneous isotropic case, there is Lamé constants $\lambda, \mu$ (see e.g. [NECAS2017], p.43) satisfying

$$\sigma_{ij} = \lambda \delta_{ij} \text{div} + 2\mu \epsilon_{ij}$$

where $\delta_{ij}$ is Kronecker’s delta.

We assume that the elastic plate is fixed on $\Gamma_D \times ]-h, h[$, $\Gamma_D \subset \partial \Omega$. If the body force $f = (f_1, f_2, f_3)$ is given in $\Omega \times ]-h, h[\text{ and surface force } g \text{ is given in } \Omega_N \times ]-h, h[, \partial \Omega \setminus \Gamma_D\text{, then the equation of equilibrium is given as follows:}$

$$\begin{align*}
-\partial_i \sigma_{ij} &= f_i \text{ in } \Omega \times ]-h, h[,
\sigma_{ij} n_j &= g_i \text{ on } \Omega_N \times ]-h, h[, u_i = 0 \text{ on } \Gamma_D \times ]-h, h[, i = 1, 2, 3
\end{align*}$$

(5.4)

We now explain the plain elasticity. 

- **Plain strain:**

  On the end of plate, the contact condition $u_3 = 0, g_3 = 0$ is satisfied. In this case, we can suppose that $f_3 = g_3 = u_3 = 0$ and $u(x_1, x_2, x_3) = \pi(x_1, x_2)$ for all $-h < x_3 < h$.

- **Plain stress:**

  The cylinder is assumed to be very thin and subjected to no load on the ends $x_3 = \pm h$, that is,

  $$\sigma_{3i} = 0, \quad x_3 = \pm h, \quad i = 1, 2, 3$$

  The assumption leads that $\sigma_{3i} = 0$ in $\Omega \times ]-h, h[$ and $u(x_1, x_2, x_3) = \pi(x_1, x_2)$ for all $-h < x_3 < h$.

- **Generalized plain stress:**

  The cylinder is subjected to no load at $x_3 = \pm h$. Introducing the mean values with respect to thickness,

  $$\overline{\pi}(x_1, x_2) = \frac{1}{2h} \int_{-h}^{h} u(x_1, x_2, x_3)dx_3$$

  and we derive $\overline{u}_3 \equiv 0$. Similarly we define the mean values $\overline{f}, \overline{g}$ of the body force and surface force as well as the mean values $\overline{\pi}_{ij}$ and $\overline{\sigma}_{ij}$ of the components of stress and strain, respectively.

In what follows we omit the overlines of $\overline{\pi}, \overline{f}, \overline{g}, \overline{\pi}_{ij}$ and $\overline{\sigma}_{ij}$. Then we obtain similar equation of equilibrium given in (5.4) replacing $\Omega \times ]-h, h[$ with $\Omega$ and changing $i = 1, 2$. In the case of plane stress, $\sigma_{ij} = \lambda^* \delta_{ij} \text{div} + 2\mu \epsilon_{ij}, \lambda^* = (2\mu)/(\lambda + \mu)$.

The equations of elasticity are naturally written in variational form for the displacement vector $u(x) \in V$ as:

$$\int_{\Omega} [2\mu \epsilon_{ij}(u) \epsilon_{ij}(v) + \lambda \epsilon_{ii}(u) \epsilon_{jj}(v)] = \int_{\Omega} f \cdot v + \int_{\Gamma} g \cdot v, \forall v \in V$$
where $V$ is the linear closed subspace of $H^1(\Omega)^2$.

**Tip**: Beam

Consider an elastic plate with the undeformed rectangle shape $[0, 10] \times [0, 2]$. The body force is the gravity force $f$ and the boundary force $g$ is zero on lower and upper side. On the two vertical sides of the beam are fixed.

```cpp
// Parameters
real E = 21.5;
real sigma = 0.29;
real gravity = -0.05;

// Mesh
border a(t=2, 0) {x=0; y=t; label=1;}
border b(t=0, 10) {x=t; y=0; label=2;}
border c(t=0, 2) {x=10; y=t; label=1;}
border d(t=0, 10) {x=10-t; y=2; label=3;}
mesh th = buildmesh(b(20) + c(5) + d(20) + a(5));

// Fespace
fespace Vh(th, [P1, P1]);
Vh [uu, vv];
Vh [w, s];

// Macro
real sqrt2 = sqrt(2.);
macro epsilon(u1, u2) [dx(u1), dy(u2), (dy(u1)+dx(u2))/sqrt2] //
macro div(u,v) (dx(u) + dy(v)) //

// Problem
real mu = E/(2*(1+sigma));
real lambda = E*sigma/((1+sigma)*(1-2*sigma));
solve Elasticity ([uu, vv], [w, s])
    = int2d(th)(
        lambda*div(w,s)*div(uu,vv)
        + 2.*mu*( epsilon(w,s)'*epsilon(uu,vv) )
    )
    + int2d(th)(
        - gravity*s
    )
    + on(1, uu=0, vv=0);

// Plot
plot([uu, vv], wait=true);
plot([uu,vv], wait=true, bb=[[-0.5, 2.5], [2.5, -0.5]]);

// Movemesh
mesh th1 = movemesh(th, [x+uu, y+vv]);
plot(th1, wait=true);
```

**Tip**: Beam 3D

Consider elastic box with the undeformed parallelepiped shape $[0, 5] \times [0, 1] \times [0, 1]$. The body force is the gravity force $f$ and the boundary force $g$ is zero on all face except one the one vertical left face where the beam is fixed.
include "cube.idp"

// Parameters
int [int] Nxyz = [20, 5, 5];
real [int, int] Bxyz = [[0., 5.], [0., 1.], [0., 1.]];
int [int, int] Lxyz = [[1, 2], [2, 2], [2, 2]];

real E = 21.5e4;
real sigma = 0.29;
real gravity = -0.05;

// Mesh
mesh3 Th = Cube(Nxyz, Bxyz, Lxyz);

// Fespace
fespace Vh(Th, [P1, P1, P1]);
Vh [u1, u2, u3], [v1, v2, v3];

// Macro
real sqrt2 = sqrt(2.);
macro epsilon(u1, u2, u3) [
  dx(u1), dy(u2), dz(u3),
  (dz(u2) + dy(u3))/sqrt2,
  (dz(u1) + dx(u3))/sqrt2,
  (dy(u1) + dx(u2))/sqrt2]
macro div(u1, u2, u3) (dx(u1) + dy(u2) + dz(u3))

// Problem
real mu = E/(2*(1+sigma));
real lambda = E*sigma/((1+sigma)*(1-2*sigma));
solve Lame ([u1, u2, u3], [v1, v2, v3])
  = int3d(Th)(
    lambda*div(u1, u2, u3)*div(v1, v2, v3)
    + 2.*mu*( epsilon(u1, u2, u3)'*epsilon(v1, v2, v3) )
  )
  - int3d(Th)(
    gravity*v3
  )
  + on(1, u1=0, u2=0, u3=0);

// Display
real dmax = u1[].max;
cout << "max displacement = " << dmax << endl;

// Movemesh
real coef = 0.1/dmax;
int[int] ref2 = [1, 0, 2, 0];
mesh3 Thm = movemesh3(Th, transfo=[x+u1*coef, y+u2*coef, z+u3*coef], label=ref2);
Thm = change(Thm, label=ref2);

// Plot
plot(Th, Thm, wait=true, cmm="coef amplification = "+coef);
5.2.1 Fracture Mechanics

Consider the plate with the crack whose undeformed shape is a curve $\Sigma$ with the two edges $\gamma_1, \gamma_2$.

We assume the stress tensor $\sigma_{ij}$ is the state of plate stress regarding $(x, y) \in \Omega_{\Sigma} = \Omega \setminus \Sigma$. Here $\Omega$ stands for the undeformed shape of elastic plate without crack.

If the part $\Gamma_N$ of the boundary $\partial \Omega$ is fixed and a load $\mathcal{L} = (f, g) \in L^2(\Omega)^2 \times L^2(\Gamma_N)^2$ is given, then the displacement $u$ is the minimizer of the potential energy functional:

$$
\mathcal{E}(v; \mathcal{L}, \Omega_{\Sigma}) = \int_{\Omega_{\Sigma}} \{w(x, v) - f \cdot v\} - \int_{\Gamma_N} g \cdot v
$$

over the functional space $V(\Omega_{\Sigma})$,

$$
V(\Omega_{\Sigma}) = \{v \in H^1(\Omega_{\Sigma})^2; \; v = 0 \text{ on } \Gamma_D = \partial \Omega \setminus \Gamma_N\},
$$

where $w(x, v) = \sigma_{ij}(v)\varepsilon_{ij}(v)/2$,

$$
\sigma_{ij}(v) = C_{ijkl}(x)\varepsilon_{kl}(v), \quad \varepsilon_{ij}(v) = (\partial v_i/\partial x_j + \partial v_j/\partial x_i)/2, \quad (C_{ijkl} : \text{ Hooke’s tensor}).
$$

If the elasticity is homogeneous isotropic, then the displacement $u(x)$ is decomposed in an open neighborhood $U_k$ of $\gamma_k$ as in (see e.g. [OHTSUH1980])

$$
u(x) = \sum_{l=1}^{2} K_l(\gamma_k) r_k^{1/2} S_{kl}^C(\theta_k) + u_{k,R}(x) \quad \text{for } x \in \Omega_{\Sigma} \cap U_k, \; k = 1, 2 \tag{5.5}
$$

with $u_{k,R} \in H^2(\Omega_{\Sigma} \cap U_k)^2$, where $U_k, \; k = 1, 2$ are open neighborhoods of $\gamma_k$ such that $\partial L_1 \cap U_1 = \gamma_1, \; \partial L_m \cap U_2 = \gamma_2$, and

$$
S_{k1}^C(\theta_k) = \frac{1}{4\mu} \frac{1}{(2\pi)^{1/2}} \begin{bmatrix}
2\kappa - 1 \cos(\theta_k/2) - \cos(3\theta_k/2) \\
-2\kappa + 1 \sin(\theta_k/2) + \sin(3\theta_k/2)
\end{bmatrix},
$$

$$
S_{k2}^C(\theta_k) = \frac{1}{4\mu} \frac{1}{(2\pi)^{1/2}} \begin{bmatrix}
-2\kappa - 1 \sin(\theta_k/2) + 3 \sin(3\theta_k/2) \\
[2\kappa - 1] \cos(\theta_k/2) + \cos(3\theta_k/2)
\end{bmatrix},
$$

where $\mu$ is the shear modulus of elasticity, $\kappa = 3 - 4\nu$ ($\nu$ is the Poisson’s ratio) for plane strain and $\kappa = \frac{3-\nu}{1+\nu}$ for plane stress.

The coefficients $K_1(\gamma_i)$ and $K_2(\gamma_i)$, which are important parameters in fracture mechanics, are called stress intensity factors of the opening mode (mode I) and the sliding mode (mode II), respectively.
For simplicity, we consider the following simple crack
\[ \Omega = \{(x, y) : -1 < x < 1, -1 < y < 1\}, \quad \Sigma = \{(x, y) : -1 \leq x \leq 0, y = 0\} \]
with only one crack tip \( \gamma = (0, 0) \). Unfortunately, FreeFEM cannot treat crack, so we use the modification of the domain with U-shape channel (see U-shape example, Fig. 3.19) with \( d = 0.0001 \). The undeformed crack \( \Sigma \) is approximated by
\[ \Sigma_d = \{(x, y) : -1 \leq x \leq -10 \times d, -d \leq y \leq d\} \cup \{(x, y) : -10 \times d \leq x \leq 0, -d = 0.1 \times x \leq y \leq d - 0.1 \times x\} \]
and \( \Gamma_D = \text{R} \) in U-shape example, Fig. 3.19.

In this example, we use three technique:

- Fast Finite Element Interpolator from the mesh Th to Zoom for the scale-up of near \( \gamma \).
- After obtaining the displacement vector \( \mathbf{u} = (u, v) \), we shall watch the deformation of the crack near \( \gamma \) as follows,
  \[ \text{mesh Plate} = \text{movemesh}(\text{Zoom, \{x+u, y+v\}}); \]
  \[ \text{plot(Plate);} \]
- Adaptivity is an important technique here, because a large singularity occurs at \( \gamma \) as shown in (5.5).

The first example creates mode I deformation by the opposed surface force on \( B \) and \( T \) in the vertical direction of \( \Sigma \), and the displacement is fixed on \( R \).

In a laboratory, fracture engineers use photoelasticity to make stress field visible, which shows the principal stress difference
\[ \sigma_1 - \sigma_2 = \sqrt{(\sigma_{11} - \sigma_{22})^2 + 4\sigma_{12}^2} \]
where \( \sigma_1 \) and \( \sigma_2 \) are the principal stresses.

In opening mode, the photoelasticity make symmetric pattern concentrated at \( \gamma \).

**Tip:** Crack Opening, \( K_2(\gamma) = 0 \)

```
//Parameters
real d = 0.0001; int n = 5; real cb = 1, ca = 1, tip = 0.0;
real E = 21.5;
real sigma = 0.29;

// Mesh
border L1(t=0, ca-d){x=-cb; y=-d-t;}
border L2(t=0, ca-d){x=-cb; y=ca-t;}
border B(t=0, 2){x=cb*(t-1); y=-ca;}
border C1(t=0, 1){x=-ca*(1-t)+(tip-10*d)*t; y=d;}
border C21(t=0, 1){x=(tip-10*d)*(1-t)+tip*t; y=d*(1-t);}
border C22(t=0, 1){x=(tip-10*d)*t+tip*(1-t); y=-d+t;}
border C3(t=0, 1){x=(tip-10*d)*(1-t)-ca+t; y=-d;}
border C4(t=0, 2*d){x=-ca; y=-d+t;}
border R(t=0, 2){x=cb; y=cb*(t-1);}
border T(t=0, 2){x=cb*(1-t); y=ca;}

mesh Th = buildmesh(L1(n/2) + L2(n/2) + B(n)
    + C1(n) + C21(3) + C22(3) + C3(n) + R(n) + T(n));
plot(Th, wait=true);
```
(continues on next page)
cb=0.1; ca=0.1;

mesh Zoom = buildmesh(L1(n/2) + L2(n/2) + B(n) + C1(n) + C21(3) + C22(3) + C3(n) + R(n) + T(n));
plot(Zoom, wait=true);

// Fespace
fespace Vh(Th, [P2, P2]);
Vh [u, v];
Vh [w, s];

fespace zVh(Zoom, P2);
zVh Sx, Sy, Sxy, N;

// Problem
real mu = E/(2*(1+sigma));
real lambda = E*sigma/((1+sigma)*(1-2*sigma));
solve Problem ([u, v], [w, s])
  = int2d(Th)(
    2*mu*(dx(u)*dx(w) + (dx(v)+dy(u))*(dx(s)+dy(w)))/4
    + lambda*(dx(u) + dy(v))*(dx(w) + dy(s))/2
  )
- int1d(Th, T) (0.1*(1-x)*s)
+ int1d(Th, B) (0.1*(1-x)*s)
+on(R, u=0, v=0)
;

// Loop
for (int i = 1; i <= 5; i++){
  mesh Plate = movemesh(Zoom, [x+u, y+v]); //deformation near gamma
  Sx = lambda*dx(u) + 2*mu*dx(u);
  Sy = lambda*dy(v) + 2*mu*dy(v);
  Sxy = mu*(dy(u) + dx(v));
  N = 0.1*1*sqrt((Sx-Sy)^2 + 4*Sxy^2); //principal stress difference
  if (i == 1){
    plot(Plate, bw=1);
    plot(N, bw=1);
  }
  else if (i == 5){
    plot(Plate, bw=1);
    plot(N, bw=1);
    break;
  }
}

// Adaptmesh
Th = adaptmesh(Th, [u, v]);

// Solve
Problem;
}

It is difficult to create mode II deformation by the opposed shear force on B and T that is observed in a laboratory. So
(a) Crack open displacement (COD) on the first mesh

(b) Principal stress difference on the first mesh

(c) COD on the last adaptive mesh

(d) Principal stress difference on the last adaptive mesh
we use the body shear force along $\Sigma$, that is, the $x$-component $f_1$ of the body force $f$ is given by

$$f_1(x, y) = H(y - 0.001) * H(0.1 - y) - H(-y - 0.001) * H(y + 0.1)$$

where $H(t) = 1$ if $t > 0; = 0$ if $t < 0$.

**Tip:** Crack Sliding, $K_2(\gamma) = 0$

```plaintext
// Parameters
real d = 0.0001; int n = 5; real cb = 1, ca = 1, tip = 0.0;
real E = 21.5;
real sigma = 0.29;

// Mesh
border L1(t=0, ca-d){x=-cb; y=-d-t;}
border L2(t=0, ca-d){x=-cb; y=ca-t;}
border B(t=0, 2){x=cb*(t-1); y=-ca;}
border C1(t=0, 1){x=-ca*(1-t)+(tip-10*d)*t; y=d;}
border C21(t=0, 1){x=(tip-10*d)*t; y=-d+t;}
border C22(t=0, 1){x=(tip-10*d)*t; y=-d-t;}
border C3(t=0, 1){x=(tip-10*d)*t+ca*t; y=-d;}
border C4(t=0, 2*d){x=-ca; y=-d+t;}
border R(t=0, 2){x=cb; y=cb*(t-1);}
border T(t=0, 2){x=cb*(1-t); y=ca;}
mesh Th = buildmesh(L1(n/2) + L2(n/2) + B(n) + C1(n) + C21(3) + C22(3) + C3(n) + R(n) + T(n));
plot(Th, wait=true);

cb=0.1; ca=0.1;
mesh Zoom = buildmesh(L1(n/2) + L2(n/2) + B(n) + C1(n) + C21(3) + C22(3) + C3(n) + R(n) + T(n));
plot(Zoom, wait=true);

// Fespace
fespace Vh(Th, [P2, P2]);
Vh [u, v];
Vh [w, s];

fespace zVh(Zoom, P2);
zVh Sx, Sy, Sxy, N;

fespace Vh1(Th, P1);
Vh1 fx = (((y>0.001)*(y<0.1)) - ((y<0.001)*(y>=0.1)));

// Problem
real mu = E/(2*(1+sigma));
real lambda = E*sigma/(((1+sigma)*(1-2*sigma)));
solve Problem ([u, v], [w, s]) {
  = int2d(Th)(
    2*mu*(dx(u)*dx(w) + ((dx(v) + dy(u))*(dx(s)+ dy(w))))/4
    + lambda*(dx(u) + dy(v))*(dx(w) + dy(s))/2
  )
  -int2d(Th)(
    fx*w
  )
  +on(R, u=0, v=0)
}
```

(continues on next page)
5.3 Non-linear static problems

Here we propose to solve the following non-linear academic problem of minimization of a functional:

\[ J(u) = \int_{\Omega} \frac{1}{2} f(|\nabla u|^2) - u \cdot b \]

where \( u \) is function of \( H^1_0(\Omega) \) and \( f \) defined by:

\[ f(x) = a \cdot x + x - \ln(1 + x), \quad f'(x) = a + \frac{x}{1 + x}, \quad f''(x) = \frac{1}{(1 + x)^2} \]

5.3.1 Newton-Raphson algorithm

Now, we solve the Euler problem \( \nabla J(u) = 0 \) with Newton-Raphson algorithm, that is:

\[ u^{n+1} = u^n - (\nabla^2 J(u^n))^{-1} \cdot \nabla J(u^n) \]
(a) COD on the first mesh

(b) Principal stress difference in the first mesh

(c) COD on the last adaptive mesh

(d) Principal stress difference on the last adaptive mesh
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5.3. Non-linear static problems

```plaintext
plot(Th, wait=true);

// Fespace
fespace Vh(Th, P1);
Vh u=0;
Vh v, w;

fespace Ph(Th, P1dc);
Ph alpha; // to store |nabla u|^2
Ph dalpha; // to store 2f''(|nabla u|^2)

// Function
func real f (real u){
    return u*a + u - log(1.+u);
}
func real df (real u){
    return a +u/(1.+u);
}
func real ddf (real u){
    return 1. / ((1.+u)*(1.+u));
}

// Problem
//the variational form of evaluate dJ = nabla J
//dJ = f'()*(dx(u)*dx(vh) + dy(u)*dy(vh))
varf vJ (uh, vh) = int2d(Th)(
    alpha*(dx(u)*dx(vh) + dy(u)*dy(vh))
    - b*vh
) + on(1, 2, 3, 4, uh=0);

//the variational form of evaluate d^2J = nabla^2 J
//hJ(uh,vh) = f'()*(dx(uh)*dx(vh) + dy(uh)*dy(vh))
// + 2*f''() (dx(u)*dx(uh) + dy(u)*dy(uh)) * (dx(u)*dx(vh) + dy(u)*dy(vh))
varf vhJ (uh, vh) = int2d(Th)(
    alpha*(dx(uh)*dx(vh) + dy(uh)*dy(vh))
    + dalpha*(dx(u)*dx(vh) + dy(u)*dy(vh))* (dx(u)*dx(uh) + dy(u)*dy(uh))
) + on(1, 2, 3, 4, uh=0);

// Newton algorithm
for (int i = 0; i < 100; i++){
    // Compute f' and f''
    alpha = df(dx(u)*dx(u) + dy(u)*dy(u));
dalpha = 2*ddf(dx(u)*dx(u) + dy(u)*dy(u));

    // nabla J
    v[] = vJ(0, Vh);

    // Residual
    real res = v[].'*v[];
    cout << i << " residu^2 = " << res << endl;
    if( res < 1e-12) break;
```

(continues on next page)
5.4 Eigen value problems

This section depends on your installation of FreeFEM; you need to have compiled ARPACK. This tool is available in FreeFEM if the word eigenvalue appears in line Load:, like:

```
-- FreeFem++ v*.* (date *** *** ** **:**:** CET ****)
file : ***.edp
Load: lg_fem lg_mesh eigenvalue
```

This tool is based on arpack++, the object-oriented version of ARPACK eigenvalue package [LEHOUCQ1998].

The function EigenValue computes the generalized eigenvalue of \( A u = \lambda B u \). The Shift-invert method is used by default, with \( \sigma = \sigma \) the shift of the method.

The matrix \( OP \) is defined with \( A - \sigma B \).

The return value is the number of converged eigenvalues (can be greater than the number of requested eigenvalues nev=)

```
int k = EigenValue(OP, B, nev=Nev, sigma=Sigma);
```

where the matrix \( OP = A - \sigma B \) with a solver and boundary condition, and the matrix \( B \).

There is also a functional interface:

```
int k = EigenValue(n, FOP1, FB, nev=Nev, sigma=Sigma);
```

where \( n \) is the size of the problem, and the operators are now defined through functions, defining respectively the matrix product of \( OP^{-1} \) and \( B \), as in

```
int n = OP1.n;
func real[int] FOP1(real[int] & u){ real[int] Au = OP^-1*u; return Au; }
func real[int] FB(real[int] & u){ real[int] Au = B*u; return Au; }
```

If you want finer control over the method employed in ARPACK, you can specify which mode ARPACK will work with (mode=, see ARPACK documentation [LEHOUCQ1998]). The operators necessary for the chosen mode can be passed through the optional parameters \( A=, A1=, B=, B1=, \) (see below).

- \( \text{mode}=1 \): Regular mode for solving \( A u = \lambda u \)

```
int k = EigenValue(n, A=FOP, mode=1, nev=Nev);
```

where the function \( FOP \) defines the matrix product of \( A \)
• mode=2: Regular inverse mode for solving $Au = \lambda Bu$

```cpp
int k = EigenValue(n, A=FOP, B=FB, B1=FB1, mode=2, nev=Nev);
```

where the functions FOP, FB and FB1 define respectively the matrix product of $A$, $B$ and $B^{-1}$

• mode=3: Shift-invert mode for solving $Au = \lambda Bu$

```cpp
int k = EigenValue(n, A1=FOP1, B=FB, mode=3, sigma=Sigma, nev=Nev);
```

where the functions FOP1 and FB define respectively the matrix product of $OP^{-1} = (A - \sigma B)^{-1}$ and $B$

You can also specify which subset of eigenvalues you want to compute (which=). The default value is which="LM", for eigenvalues with largest magnitude. "SM" is for smallest magnitude, "LA" for largest algebraic value, "SA" for smallest algebraic value, and "BE" for both ends of the spectrum.

Remark: For complex problems, you need to use the keyword `complexEigenValue` instead of `EigenValue` when passing operators through functions.

**Note:** Boundary condition and Eigenvalue Problems

The locking (Dirichlet) boundary condition is make with exact penalization so we put $1e30=tgv$ on the diagonal term of the locked degree of freedom (see [Finite element chapter](#)). So take Dirichlet boundary condition just on $A$ and not on $B$ because we solve $w = OP^{-1} * B * v$.

If you put locking (Dirichlet) boundary condition on $B$ matrix (with key work on) you get small spurious modes ($10^{-30}$), due to boundary condition, but if you forget the locking boundary condition on $B$ matrix (no keywork on) you get huge spurious ($10^{30}$) modes associated to these boundary conditons. We compute only small mode, so we get the good one in this case.

• sym= The problem is symmetric (all the eigen value are real)

• nev= The number desired eigenvalues (nev) close to the shift.

• value= The array to store the real part of the eigenvalues

• ivalue= The array to store the imaginary part of the eigenvalues

• vector= The FE function array to store the eigenvectors

• rawvector= An array of type real[int,int] to store eigenvectors by column.

For real non symmetric problems, complex eigenvectors are given as two consecutive vectors, so if eigenvalue $k$ and $k + 1$ are complex conjugate eigenvalues, the $k$th vector will contain the real part and the $k + 1$th vector the imaginary part of the corresponding complex conjugate eigenvectors.

• tol= The relative accuracy to which eigenvalues are to be determined;

• sigma= The shift value;

• maxit= The maximum number of iterations allowed;

• ncv= The number of Arnoldi vectors generated at each iteration of ARPACK;

• mode= The computational mode used by ARPACK (see above);

• which= The requested subset of eigenvalues (see above).

**Tip:** Laplace eigenvalue

In the first example, we compute the eigenvalues and the eigenvectors of the Dirichlet problem on square $\Omega = [0, \pi]^2$. 

5.4. Eigen value problems
The problem is to find: $\lambda$, and $\nabla u_\lambda$ in $\mathbb{R} \times H^1_0(\Omega)$

$$\int_\Omega \nabla u_\lambda \nabla v = \lambda \int_\Omega uv \quad \forall v \in H^1_0(\Omega)$$

The exact eigenvalues are $\lambda_{n,m} = (n^2 + m^2)$, $(n, m) \in \mathbb{N}^2$ with the associated eigenvectors are $u_{m,n} = \sin(nx) \ast \sin(my)$.

We use the generalized inverse shift mode of the arpack++ library, to find 20 eigenvalues and eigenvectors close to the shift value $\sigma = 20$.

```cpp
// Parameters
verbosity=0;
real sigma = 20; //value of the shift
int nev = 20; //number of computed eigen value close to sigma

// Mesh
mesh Th = square(20, 20, [pi*x, pi*y]);

// Fespace
fespace Vh(Th, P2);
Vh u1, u2;

// Problem
// OP = A - sigma B ; // the shifted matrix
varf op (u1, u2)
  = int2d(Th)(
      dx(u1)*dx(u2)
    + dy(u1)*dy(u2)
    - sigma* u1*u2
  )
  + on(1, 2, 3, 4, u1=0);

varf b ([u1], [u2]) = int2d(Th)(u1*u2); //no boundary condition

matrix OP = op(Vh, Vh, solver=Crout, factorize=1); //crout solver because the matrix... in not positive

matrix B = b(Vh, Vh, solver=CG, eps=1e-20);

// important remark:
// the boundary condition is make with exact penalization:
// we put le30=tgv on the diagonal term of the lock degree of freedom.
// So take Dirichlet boundary condition just on $a$ variational form
// and not on $b$ variational form.
// because we solve $ w=OP^-1*B*v $ 

// Solve
real[int] ev(nev); //to store the nev eigenvalue
Vh[int] eV(nev); //to store the nev eigenvector

int k = EigenValue(OP, B, sym=true, sigma=sigma, value=ev, vector=eV,
    tol=1e-10, maxit=0, ncv=0);

// Display & Plot
for (int i = 0; i < k; i++){
  u1 = eV[i];
  real qq = int2d(Th)(dx(u1)*dx(u1) + dy(u1)*dy(u1));
  real mm = int2d(Th)(u1*u1);```
The output of this example is:

\begin{verbatim}
lambda[0] = 5.0002, err= -1.46519e-11
lambda[1] = 8.00074, err= -4.05158e-11
lambda[7] = 17.0048, err= 1.03883e-10
lambda[8] = 18.0083, err= -4.05497e-11
lambda[10] = 20.0096, err= -4.16212e-14
lambda[12] = 25.0283, err= 6.77444e-10
lambda[16] = 29.0273, err= 1.38242e-10
lambda[17] = 32.0449, err= 1.2522e-10
lambda[18] = 34.049, err= 3.40213e-11
lambda[19] = 34.0492, err= 2.41751e-10
\end{verbatim}

5.4. Eigen value problems

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5.5 Evolution problems

**FreeFEM** also solves evolution problems such as the heat equation:

\[
\frac{\partial u}{\partial t} - \mu \Delta u = f \quad \text{in } \Omega \times [0,T[
\]
\[
u(x,0) = \nu_0(x) \quad \text{in } \Omega
\]
\[
(\partial u / \partial n)(x, t) = 0 \quad \text{on } \partial \Omega \times [0,T[
\]

(5.6)

with a positive viscosity coefficient \( \mu \) and homogeneous Neumann boundary conditions.

We solve (5.6) by FEM in space and finite differences in time.

We use the definition of the partial derivative of the solution in the time derivative:

\[
\frac{\partial u}{\partial t}(x, y, t) = \lim_{\tau \to 0} \frac{u(x, y, t) - u(x, y, t - \tau)}{\tau}
\]

which indicates that \( u^m(x, y) = u(x, y, m\tau) \) will satisfy approximatively:

\[
\frac{\partial u}{\partial t}(x, y, m\tau) \approx \frac{u^m(x, y) - u^{m-1}(x, y)}{\tau}
\]

The time discretization of heat equation (5.6) is as follows, \( \forall m = 0, \ldots, [T/\tau] \):

\[
\frac{u^{m+1} - u^m}{\tau} - \mu \Delta u^{m+1} = f^{m+1} \quad \text{in } \Omega
\]
\[
u^0(x) = \nu_0(x) \quad \text{in } \Omega
\]
\[
\frac{\partial u^{m+1}}{\partial n}(x) = 0 \quad \text{on } \partial \Omega
\]

which is so-called *backward Euler method* for (5.6).

To obtain the variational formulation, multiply with the test function \( v \) both sides of the equation:

\[
\int_{\Omega} \{ u^{m+1} v - \tau \Delta u^{m+1} v \} = \int_{\Omega} \{ u^m + \tau f^{m+1} \} v
\]

By the divergence theorem, we have:

\[
\int_{\Omega} \{ u^{m+1} v + \tau \nabla u^{m+1} \cdot \nabla v \} - \int_{\partial \Omega} \tau (\partial u^{m+1} / \partial n) v = \int_{\Omega} \{ u^m v + \tau f^{m+1} v \}
\]

By the boundary condition \( \partial u^{m+1} / \partial n = 0 \), it follows that:

\[
\int_{\Omega} \{ u^{m+1} v + \tau \nabla u^{m+1} \cdot \nabla v \} - \int_{\Omega} \{ u^m v + \tau f^{m+1} v \} = 0
\]

(5.7)

Using the identity just above, we can calculate the finite element approximation \( u^m_h \) of \( u^m \) in a step-by-step manner with respect to \( t \).

**Tip:** Example

We now solve the following example with the exact solution \( u(x, y, t) = tx^4 \), \( \Omega = ]0, 1[^2 \).

\[
\frac{\partial u}{\partial t} - \mu \Delta u = x^4 - \mu 12tx^2 \quad \text{in } \Omega \times ]0, 3[ \]
\[
u(x, y, 0) = 0 \quad \text{in } \Omega \]
\[
u|_{\partial \Omega} = t * x^4 \quad \text{on } \Omega
\]
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// Parameters
real dt = 0.1;
real mu = 0.01;

// Mesh
mesh Th = square(16, 16);

// Fespace
fespace Vh(Th, P1);
Vh u, v, uu, f, g;

// Problem
problem dHeat (u, v) = int2d(Th)(
    u*v
    + dt*mu*(dx(u)*dx(v) + dy(u)*dy(v))
) + int2d(Th)(
    - uu*v
    - dt*f*v
) + on(1, 2, 3, 4, u=g);

// Time loop
real t = 0;
uu = 0;
for (int m = 0; m <= 3/dt; m++){
    // Update
    t = t + dt;
f = x^4 - mu*t*12*x^2;
g = t*x^4;
uu = u;

    // Solve
    dHeat;

    // Plot
    plot(u, wait=true);
    cout << "t=" << t << " - L^2-Error=" << sqrt(int2d(Th)((u-t*x^4)^2)) << endl;
}

In the last statement, the $L^2$-error $\left( \int_\Omega |u - tx^4|^2 \right)^{1/2}$ is calculated at $t = m\tau, \tau = 0.1$. At $t = 0.1$, the error is 0.000213269. The errors increase with $m$ and 0.00628589 at $t = 3$.
The iteration of the backward Euler (5.7) is made by for loop.

**Note:** The stiffness matrix in the loop is used over and over again. FreeFEM support reuses of stiffness matrix.

### 5.5.1 Mathematical Theory on Time Difference Approximations.

In this section, we show the advantage of implicit schemes. Let $V, H$ be separable Hilbert space and $V$ is dense in $H$. Let $a$ be a continuous bilinear form over $V \times V$ with coercivity and symmetry.
Then $\sqrt{a(v, v)}$ become equivalent to the norm $\|v\|$ of $V$.

**Problem Ev(f, Omega):** For a given $f \in L^2(0, T; V')$, $u^0 \in H$

$$
\frac{d}{dt}(u(t), v) + a(u(t), v) = (f(t), v) \quad \forall v \in V, \quad a.e. \ t \in [0, T]
$$

where $V'$ is the dual space of $V$.

Then, there is an unique solution $u \in L^\infty(0, T; H) \cap L^2(0, T; V)$.

Let us denote the time step by $\tau > 0$, $N_T = [T/\tau]$. For the discretization, we put $u^n = u(n\tau)$ and consider the time difference for each $\theta \in [0, 1]$

$$
\frac{1}{\tau} (u_h^{n+1} - u_h^n, \phi_i) + a(u_h^{n+\theta}, \phi_i) = (f^{n+\theta}, \phi_i) \\
M = (m_{ij}), \quad m_{ij} = (\phi_j, \phi_i), \quad A = (a_{ij}), \quad a_{ij} = a(\phi_j, \phi_i)
$$

Formula (5.8) is the **forward Euler scheme** if $\theta = 0$, Crank-Nicolson scheme if $\theta = 1/2$, the **backward Euler scheme** if $\theta = 1$.

Unknown vectors $u^n = (u_h^1, \ldots, u_h^M)^T$ in

$$
u^n_h(x) = u^n_1 \phi_1(x) + \cdots + u^n_m \phi_m(x), \quad u^n_1, \ldots, u^n_m \in \mathbb{R}
$$

are obtained from solving the matrix

$$(M + \theta \tau A)u^{n+1} = \{M - (1 - \theta)\tau A\}u^n + \tau \{\theta f^{n+1} + (1 - \theta)f^n\}
$$

Refer [TABATA1994], pp.70–75 for solvability of (5.8). The stability of (5.8) is in [TABATA1994], Theorem 2.13:

Let $\{T_h\}_{h \downarrow 0}$ be regular triangulations (see **Regular Triangulation**). Then there is a number $c_0 > 0$ independent of $h$ such that,

$$
|u_h^n|^2 \leq \begin{cases} 
\frac{1}{\theta} \left\{ |u_h^0|^2 + \tau \sum_{k=0}^{n-1} \|f^{k+\theta}\|_{V^*_h}^2 \right\} & \theta \in [0, 1/2) \\
|u_h^0|^2 + \tau \sum_{k=0}^{n-1} \|f^{k+\theta}\|_{V^*_h}^2 & \theta \in [1/2, 1]
\end{cases}
$$

if the following are satisfied:

1. When $\theta \in [0, 1/2)$, then we can take a time step $\tau$ in such a way that

$$
\tau < \frac{2(1-\delta)}{(1-2\theta)c_0^2} h^2
$$

for arbitrary $\delta \in (0, 1)$.

2. When $1/2 \leq \theta \leq 1$, we can take $\tau$ arbitrary.

**Tip:** Example

```c
// Parameters
def real tau = 0.1; real theta = 0.;

// Mesh
```

(continues on next page)
mesh Th = square(12, 12);

// Fespace
fespace Vh(Th, P1);
Vh u, v, oldU;
Vh f1, f0;

fespace Ph(Th, P0);
Ph h = hTriangle; // mesh sizes for each triangle

// Function
func real f (real t){
    return x^2*(x-1)^2 + t*(-2 + 12*x - 11*x^2 - 2*x^3 + x^4);
}

// File
ofstream out("err02.csv"); // file to store calculations
out << "mesh size = " << h[].max << ", time step = " << tau << endl;
for (int n = 0; n < 5/tau; n++)
    out << n*tau << ",";
out << endl;

// Problem
problem aTau (u, v)
    = int2d(Th)(
        u*v
        + theta*tau*(dx(u)*dx(v) + dy(u)*dy(v) + u*v)
    )
    - int2d(Th)(
        oldU*v
        - (1-theta)*tau*(dx(oldU)*dx(v) + dy(oldU)*dy(v) + oldU*v)
    )
    - int2d(Th)(
        tau*(theta*f1 + (1-theta)*f0)*v
    );

// Theta loop
while (theta <= 1.0){
    real t = 0;
    real T = 3;
    oldU = 0;
    out << theta << ",";
    for (int n = 0; n < T/tau; n++)
    // Update
        t = t + tau;
    f0 = f(n*tau);
    f1 = f((n+1)*tau);
    // Solve
    aTau;
    oldU = u;
    // Plot
    plot(u);
    // Error
}
We can see in Fig. 5.14 that \( u^\alpha_n(\theta) \) become unstable at \( \theta = 0.4 \), and figures are omitted in the case \( \theta < 0.4 \).

### 5.5.2 Convection

The hyperbolic equation

\[
\partial_t u + \alpha \cdot \nabla u = f;
\]

for a vector-valued function \( \alpha \)

appears frequently in scientific problems, for example in the Navier-Stokes equations, in the Convection-Diffusion equation, etc.

In the case of 1-dimensional space, we can easily find the general solution \( (x,t) \mapsto u(x,t) = u^0(x-\alpha t) \) of the
From (5.11), it follows that

\[ \partial_t u + \alpha \partial_x u = 0 \]

\[ u(x, 0) = u^0(x), \quad (5.9) \]

because \( \partial_t u + \alpha \partial_x u = -\alpha \dot{u}^0 + a \dot{u}^0 = 0 \), where \( \dot{u}^0 = du^0(x)/dx \).

Even if \( \alpha \) is not constant, the construction works on similar principles. One begins with the ordinary differential equation (with the convention that \( \alpha \) is prolonged by zero apart from \((0, L) \times (0, T)) \):

\[ \dot{X}(\tau) = +\alpha(X(\tau), \tau), \quad \tau \in (0, t) \quad X(t) = x \]

In this equation \( \tau \) is the variable and \( x, t \) are parameters, and we denote the solution by \( X_{x,t}(\tau) \). Then it is noticed that \( (x, t) \rightarrow v(X(\tau), \tau) \) in \( \tau = t \) satisfies the equation

\[ \partial_t v + \alpha \partial_x v = \partial_t X \dot{v} + a \partial_x X \dot{v} = 0 \]

and by the definition \( \partial_t X = \dot{X} = +\alpha \) and \( \partial_x X = \partial_x x \) in \( \tau = t \), because if \( \tau = t \) we have \( X(\tau) = x \).

The general solution of (5.9) is thus the value of the boundary condition in \( X_{x,t}(0) \), that is to say \( u(x, t) = u^0(X_{x,t}(0)) \) where \( X_{x,t}(0) \) is on the \( x \) axis, \( u(x, t) = u^0(X_{x,t}(0)) \) if \( X_{x,t}(0) \) is on the axis of \( t \).

In higher dimension \( \Omega \subset \mathbb{R}^d \), \( d = 2, 3 \), the equation for the convection is written

\[ \partial_t u + \alpha \cdot \nabla u = 0 \quad \text{in} \quad \Omega \times (0, T) \]

where \( a(x, t) \in \mathbb{R}^d \).

**FreeFEM** implements the Characteristic-Galerkin method for convection operators. Recall that the equation (5.8) can be discretized as

\[ \frac{D u}{D t} = f \quad \text{i.e.} \quad \frac{d u}{d t}(X(t), t) = f(X(t), t) \quad \text{where} \quad \frac{d X}{d t}(t) = \alpha(X(t), t) \]

where \( D \) is the total derivative operator. So a good scheme is one step of backward convection by the method of Characteristics-Galerkin

\[ \frac{1}{\tau} \left( u^{m+1}(x) - u^m(X^m(x)) \right) = f^m(x) \quad (5.10) \]

where \( X^m(x) \) is an approximation of the solution at \( t = m \tau \) of the ordinary differential equation

\[ \frac{d X}{d t}(t) = \alpha^m(X(t)), X((m + 1)\tau) = x. \]

where \( \alpha^m(x) = (\alpha_1(x, m\tau), \alpha_2(x, m\tau)) \). Because, by Taylor's expansion, we have

\[ u^m(X(m\tau)) = u^m(X((m + 1)\tau)) - \tau \sum_{i=1}^{d} \frac{\partial u^m}{\partial x_i}(X((m + 1)\tau)) \frac{\partial X_i}{\partial t}((m + 1)\tau) + o(\tau) \]

\[ = u^m(x) - \tau \alpha^m(x) \cdot \nabla u^m(x) + o(\tau) \]

where \( X_i(t) \) are the \( i \)-th component of \( X(t) \), \( u^m(x) = u(x, m\tau) \) and we used the chain rule and \( x = X((m + 1)\tau) \). From (5.11), it follows that

\[ u^m(X^m(x)) = u^m(x) - \tau \alpha^m(x) \cdot \nabla u^m(x) + o(\tau) \]

Also we apply Taylor's expansion for \( t \rightarrow u^m(x - \alpha^m(x)t), 0 \leq t \leq \tau \), then

\[ u^m(x - \alpha t) = u^m(x) - \tau \alpha^m(x) \cdot \nabla u^m(x) + o(\tau). \]

Putting

5.5. Evolution problems
convect($\alpha, -\tau, u^m$) $\approx u^m(x - \alpha^m\tau)$

we can get the approximation

$$u^m(X^m(x)) \approx \text{convect}([a_1^m, a_2^m], -\tau, u^m) \text{ by } X^m \approx x \mapsto x - \tau[a_1^m(x), a_2^m(x)]$$

A classical convection problem is that of the “rotating bell” (quoted from [LUCQUIN1998], p.16).

Let $\Omega$ be the unit disk centered at 0, with its center rotating with speed $\alpha_1 = y$, $\alpha_2 = -x$. We consider the problem (5.8) with $f = 0$ and the initial condition $u(x, 0) = u^0(x)$, that is, from (5.10)

$$u^{m+1}(x) = u^m(X^m(x)) \approx \text{convect}(\alpha, -\tau, u^m)$$

The exact solution is $u(x, t) = u(X(t))$ where $X$ equals $x$ rotated around the origin by an angle $\theta = -t$ (rotate in clockwise). So, if $u^0$ in a 3D perspective looks like a bell, then $u$ will have exactly the same shape, but rotated by the same amount. The program consists in solving the equation until $T = 2\pi$, that is for a full revolution and to compare the final solution with the initial one; they should be equal.

Tip: Convect

```cpp
// Parameters
real dt = 0.17;

// Mesh
border C(t=0, 2*pi){x=cos(t); y=sin(t);}
mesh Th = buildmesh(C(70));

// Fespace
fespace Vh(Th, P1);
Vh u0;
Vh a1 = -y, a2 = x; //rotation velocity
Vh u;

// Initialization
u = exp(-10*((x-0.3)^2+(y-0.3)^2));

// Time loop
real t = 0.;
for (int m = 0; m < 2*pi/dt; m++)
{
    // Update
    t += dt;
    u0 = u;

    // Convect
    u = convect([a1, a2], -dt, u0); //u^{m+1}=u^m(X^m(x))

    // Plot
    plot(u, cmm= " t="+t", min=":u[].min", max=":u[].max");
}
```

Note: The scheme \texttt{convect} is unconditionally stable, then the bell become lower and lower (the maximum of $u^{37}$ is 0.406 as shown in Fig. 5.15a.)
5.5.3 2D Black-Scholes equation for an European Put option

In mathematical finance, an option on two assets is modeled by a Black-Scholes equation in two space variables, (see for example [WILMOTT1995] or [ACHDOU2005]).

\[
\begin{align*}
\partial_t u & \quad + \frac{\sigma_1 x^2}{2} \frac{\partial^2 u}{\partial x^2} + \frac{\sigma_2 y^2}{2} \frac{\partial^2 u}{\partial y^2} \\
& \quad + \rho xy \frac{\partial^2 u}{\partial x \partial y} + r S_1 \frac{\partial u}{\partial x} + r S_2 \frac{\partial u}{\partial y} - r P = 0
\end{align*}
\]

which is to be integrated in \((0,T) \times \mathbb{R}^+ \times \mathbb{R}^+\) subject to, in the case of a put

\[
u(x,y,T) = (K - \max(x,y))^+
\]

Boundary conditions for this problem may not be so easy to device. As in the one dimensional case the PDE contains boundary conditions on the axis \(x_1 = 0\) and on the axis \(x_2 = 0\), namely two one dimensional Black-Scholes equations driven respectively by the data \(u(0,+,T)\) and \(u(+,0,T)\). These will be automatically accounted for because they are embedded in the PDE. So if we do nothing in the variational form (i.e. if we take a Neumann boundary condition at these two axis in the strong form) there will be no disturbance to these. At infinity in one of the variable, as in 1D, it makes sense to impose \(u = 0\). We take

\[
\sigma_1 = 0.3, \quad \sigma_2 = 0.3, \quad \rho = 0.3, \quad r = 0.05, \quad K = 40, \quad T = 0.5
\]

An implicit Euler scheme is used and a mesh adaptation is done every 10 time steps. To have an unconditionally stable scheme, the first order terms are treated by the Characteristic Galerkin method, which, roughly, approximates

\[
\frac{\partial u}{\partial t} + a_1 \frac{\partial u}{\partial x} + a_2 \frac{\partial u}{\partial y} \approx \frac{1}{\tau} \left( u^{n+1}(x) - u^n(x - \alpha \tau) \right)
\]

Tip: Black-Scholes

```c
// Parameters
int m = 30; int L = 80; int LL = 80; int j = 100; real sigx = 0.3; real sigy = 0.3;
real rho = 0.3; real r = 0.05; real K = 40; real dt = 0.01;

// Mesh
mesh th = square(m, m, [L*x, LL*y]);

// Fespace
fespace Vh(th, P1);
```

(continues on next page)
Vh u = max(K - max(x, y), 0.);
Vh xveloc, yveloc, v, uold;

// Time loop
for (int n = 0; n*dt <= 1.0; n++) {
  // Mesh adaptation
  if (j > 20) {
    th = adaptmesh(th, u, verbosity=1, abserror=1, nbjacoby=2, err=0.001, nbvx=5000, omega=1.8, ratio=1.8, nbsmooth=3, splitpbedge=1, maxsubdiv=5, rescaling=1);
    j = 0;
    xveloc = -x*r + x*sigx^2 + x*rho*sigx*sigy/2;
    yveloc = -y*r + y*sigy^2 + y*rho*sigx*sigy/2;
    u = u;
  }
  // Update
  uold = u;
  // Solve
  solve eq1(u, v, init=j, solver=LU) = int2d(th)(
    u*v*(r+1/dt) + dx(u)*dx(v)*x*sigx^2/2 + dy(u)*dy(v)*y*sigy^2/2 + (dy(u)*dx(v) + dx(u)*dy(v))*rho*sigx*sigy*x*y/2
  - int2d(th)(v*convect([xveloc, yveloc], dt, uold)/dt
  ) + on(2, 3, u=0);
  // Update
  j = j+1;
}
// Plot
plot(u, wait=true, value=true);

Results are shown on Fig. 5.16a and Fig. 5.16b.

5.6 Navier-Stokes equations

The Stokes equations are: for a given f ∈ L^2(Ω)^2:

\[
\begin{align*}
-\Delta u + \nabla p &= f \\
\nabla \cdot u &= 0
\end{align*}
\]

in Ω \ (5.11)

where u = (u_1, u_2) is the velocity vector and p the pressure. For simplicity, let us choose Dirichlet boundary conditions on the velocity, u = u_G on Γ.

In [TEMAM1977], Theorem 2.2, there is a weak form of (5.11):
Find \( \mathbf{v} = (v_1, v_2) \in \mathbf{V}(\Omega) \):

\[
\mathbf{V}(\Omega) = \{ \mathbf{w} \in H_0^1(\Omega)^2 \mid \text{div} \mathbf{w} = 0 \}
\]

which satisfy:

\[
\sum_{i=1}^{2} \int_{\Omega} \nabla u_i \cdot \nabla v_i = \int_{\Omega} \mathbf{f} \cdot \mathbf{w} \quad \text{for all } \mathbf{v} \in \mathbf{V}
\]

Here it is used the existence \( p \in H^1(\Omega) \) such that \( \mathbf{u} = \nabla p \), if:

\[
\int_{\Omega} \mathbf{u} \cdot \mathbf{v} = 0 \quad \text{for all } \mathbf{v} \in \mathbf{V}
\]

Another weak form is derived as follows: We put:

\[
\mathbf{V} = H_0^1(\Omega)^2; \quad \mathbf{W} = \left\{ q \in L^2(\Omega) \mid \int_{\Omega} q = 0 \right\}
\]

By multiplying the first equation in (5.11) with \( \mathbf{v} \in \mathbf{V} \) and the second with \( q \in \mathbf{W} \), subsequent integration over \( \Omega \), and an application of Green’s formula, we have:

\[
\int_{\Omega} \nabla \mathbf{u} \cdot \nabla \mathbf{v} - \int_{\Omega} \text{div} \mathbf{u} \cdot \text{div} p = \int_{\Omega} \mathbf{f} \cdot \mathbf{v}
\]

\[
\int_{\Omega} \text{div} \mathbf{u} \cdot q = 0
\]

This yields the weak form of (5.11):

Find \( (\mathbf{u}, p) \in \mathbf{V} \times \mathbf{W} \) such that:

\[
a(\mathbf{u}, \mathbf{v}) + b(\mathbf{v}, p) = (\mathbf{f}, \mathbf{v})
\]

\[
b(\mathbf{u}, q) = 0
\]

for all \( (\mathbf{v}, q) \in \mathbf{V} \times \mathbf{W} \), where:

\[
a(\mathbf{u}, \mathbf{v}) = \int_{\Omega} \nabla \mathbf{u} \cdot \nabla \mathbf{v} = \sum_{i=1}^{2} \int_{\Omega} \nabla u_i \cdot \nabla v_i
\]

\[
b(\mathbf{u}, q) = -\int_{\Omega} \text{div} \mathbf{u} q
\]
Now, we consider finite element spaces \( V_h \subset V \) and \( W_h \subset W \), and we assume the following basis functions:

\[
V_h = V_h \times V_h, \quad V_h = \{ v_h \mid v_h = v_1 \phi_1 + \cdots + v_M \phi_M \}, \\
W_h = \{ q_h \mid q_h = q_1 \varphi_1 + \cdots + q_M \varphi_M \}
\]

The discrete weak form is: Find \((u_h, p_h) \in \mathbb{V}_h \times W_h\) such that:

\[
a(u_h, v_h) + b(v_h, p) = (f, v_h), \quad \forall v_h \in \mathbb{V}_h \\
b(u_h, q_h) = 0, \quad \forall q_h \in W_h
\]  \hspace{1cm} (5.12)

Note: Assume that:

1. There is a constant \( \alpha_h > 0 \) such that:

\[
a(v_h, v_h) \geq \alpha \|v_h\|^2_{1,\Omega} \quad \text{for all} \quad v_h \in \mathbb{V}_h
\]

where:

\[
\mathbb{V}_h = \{ v_h \in \mathbb{V}_h \mid b(w_h, q_h) = 0 \quad \text{for all} \quad q_h \in W_h \}
\]

2. There is a constant \( \beta_h > 0 \) such that:

\[
\sup_{v_h \in \mathbb{V}_h} \frac{b(v_h, q_h)}{\|v_h\|_{1,\Omega}} \geq \beta_h \|q_h\|_{0,\Omega} \quad \text{for all} \quad q_h \in W_h
\]

Then we have an unique solution \((u_h, p_h)\) of (5.12) satisfying:

\[
\|u - u_h\|_{1,\Omega} + \|p - p_h\|_{0,\Omega} \leq C \left( \inf_{v_h \in \mathbb{V}_h} \|u - v_h\|_{1,\Omega} + \inf_{q_h \in W_h} \|p - q_h\|_{0,\Omega} \right)
\]

with a constant \( C > 0 \) (see e.g. [ROBERTS1993], Theorem 10.4).

Let us denote that:

\[
A = (A_{ij}), A_{ij} = \int_{\Omega} \nabla \phi_i \cdot \nabla \phi_j, \quad i, j = 1, \ldots, M_v \\
B = (B_{xij}, B_{yij}), B_{xij} = -\int_{\Omega} \partial \phi_j / \partial x \phi_i, \quad B_{yij} = -\int_{\Omega} \partial \phi_j / \partial y \phi_i \\
i = 1, \ldots, M_W; j = 1, \ldots, M_V
\]

then (5.12) is written by:

\[
\begin{pmatrix}
A & B^* \\
B & 0
\end{pmatrix}
\begin{pmatrix}
U_h \\
p_h
\end{pmatrix}
= 
\begin{pmatrix}
F_h \\
0
\end{pmatrix}
\]

where:

\[
A = \begin{pmatrix} A & 0 \\ 0 & A \end{pmatrix}, \quad B^* = \begin{pmatrix} B_{x^T} & B_{y^T} \end{pmatrix}, \quad U_h = \begin{cases} \{ u_{1,h} \} \\
\{ u_{2,h} \} \end{cases}, \quad F_h = \begin{cases} \{ \int_{\Omega} f_1 \phi_1 \} \\
\{ \int_{\Omega} f_2 \phi_1 \} \end{cases}
\]

**Penalty method:** This method consists of replacing (5.12) by a more regular problem:

Find \((v_h, p_h') \in \mathbb{V}_h \times \tilde{W}_h\) satisfying:

\[
a(u_h', v_h) + b(v_h, p_h') = (f, v_h), \quad \forall v_h \in \mathbb{V}_h \\
b(u_h', q_h) - \epsilon (p_h', q_h) = 0, \quad \forall q_h \in \tilde{W}_h
\]  \hspace{1cm} (5.13)
where $\hat{W}_h \subset L^2(\Omega)$. Formally, we have:

$$\text{div} u_h^\epsilon = \epsilon p_h^\epsilon$$

and the corresponding algebraic problem:

$$\begin{pmatrix} A & B^* \\ B & -\epsilon I \end{pmatrix} \begin{pmatrix} U_h^\epsilon \\ \{p_h^\epsilon\} \end{pmatrix} = \begin{pmatrix} F_h \\ 0 \end{pmatrix}$$

**Note:** We can eliminate $p_h^\epsilon = (1/\epsilon)B^*U_h^\epsilon$ to obtain:

$$(A + (1/\epsilon)B^*B)U_h^\epsilon = F_h$$

(5.14)

Since the matrix $A + (1/\epsilon)B^*B$ is symmetric, positive-definite, and sparse, (5.14) can be solved by known technique. There is a constant $C > 0$ independent of $\epsilon$ such that:

$$\|u_h - u_h^\epsilon\|_{1,\Omega} + \|p_h - p_h^\epsilon\|_{0,\Omega} \leq C\epsilon$$

(see e.g. [ROBERTS1993], 17.2)

---

**Tip:** Cavity

The driven cavity flow problem is solved first at zero Reynolds number (Stokes flow) and then at Reynolds 100. The velocity pressure formulation is used first and then the calculation is repeated with the stream function vorticity formulation.

We solve the driven cavity problem by the penalty method (5.13) where $u_{\Gamma} \cdot n = 0$ and $u_{\Gamma} \cdot s = 1$ on the top boundary and zero elsewhere ($n$ is the unit normal to $\Gamma$, and $s$ the unit tangent to $\Gamma$).

The mesh is constructed by:

```plaintext
mesh Th = square(8, 8);
```

We use a classical Taylor-Hood element technique to solve the problem:

The velocity is approximated with the $P_2$ FE ($X_h$ space), and the pressure is approximated with the $P_1$ FE ($M_h$ space), where:

$$X_h = \{ v \in H^1([0,1]^2) \mid \forall K \in T_h \  v|_K \in P_2 \}$$

and:

$$M_h = \{ v \in H^1([0,1]^2) \mid \forall K \in T_h \  v|_K \in P_1 \}$$

The FE spaces and functions are constructed by:

```plaintext
fespace Xh(Th, P2); //definition of the velocity component space
fespace Mh(Th, P1); //definition of the pressure space
Xh u2, v2;
Xh u1, v1;
Mh p, q;
```

The Stokes operator is implemented as a system-solve for the velocity $(u_1, u_2)$ and the pressure $p$. The test function for the velocity is $(v_1, v_2)$ and $q$ for the pressure, so the variational form (5.12) in freefem language is:

---

### 5.6. Navier-Stokes equations
Each unknown has its own boundary conditions.

If the streamlines are required, they can be computed by finding $\psi$ such that $\text{rot}\psi = u$ or better:

$$-\Delta \psi = \nabla \times u$$

Now the Navier-Stokes equations are solved:

$$\frac{\partial u}{\partial t} + u \cdot \nabla u - \nu \Delta u + \nabla p = 0, \nabla \cdot u = 0$$

with the same boundary conditions and with initial conditions $u = 0$.

This is implemented by using the convection operator $\text{convect}$ for the term $\frac{\partial u}{\partial t} + u \cdot \nabla u$, giving a discretization in time

$$\frac{1}{\tau} (u^{n+1} - u^n \circ X^n) - \nu \Delta u^{n+1} + \nabla p^{n+1} = 0,$$

$$\nabla \cdot u^{n+1} = 0$$

The term $u^n \circ X^n(x) \approx u^n(x - u^n(x)\tau)$ will be computed by the operator $\text{convect}$, so we obtain:
9 )
  - p*q*(0.000001)
  - p*dx(v1) - p*dy(v2)
  - dx(u1)*q - dy(u2)*q
)
+ int2d(Th)(
  - alpha*convect([up1,up2],-dt,up1)*v1
  - alpha*convect([up1,up2],-dt,up2)*v2
)
+ on(3, u1=1, u2=0)
+ on(1, 2, 4,u1=0, u2=0)
;
22 // Time loop
for (i = 0; i <= 10; i++){
  // Update
  up1 = u1;
  up2 = u2;

  // Solve
  NS;

  // Plot
  if (!(i % 10))
    plot(coef=0.2, cmm="[u1,u2] and p", p, [u1, u2]);
}

Notice that the stiffness matrices are reused (keyword init=i)
The complete script is available in cavity example.

5.6.1 Uzawa Algorithm and Conjugate Gradients

We solve Stokes problem without penalty. The classical iterative method of Uzawa is described by the algorithm (see e.g. [ROBERTS1993], 17.3, [GLOWINSKI1979], 13 or [GLOWINSKI1985], 13):

- **Initialize:** Let $p_h^0$ be an arbitrary chosen element of $L^2(\Omega)$.
- **Calculate :math:`\mathbf{u}_h`:** Once $p_h^n$ is known, $v_h^n$ is the solution of:

$$ u_h^n = A^{-1}(f_h - B^*p_h^n) $$

- **Advance :math:`p_h`:** Let $p_h^{n+1}$ be defined by:

$$ p_h^{n+1} = p_h^n + \rho_n B u_h^n $$

There is a constant $\alpha > 0$ such that $\alpha \leq \rho_n \leq 2$ for each $n$, then $u_h^n$ converges to the solution $u_h$, and then $B v_h^n \to 0$ as $n \to \infty$ from the *Advance p_h*. This method in general converges quite slowly.

First we define mesh, and the Taylor-Hood approximation. So $X_h$ is the velocity space, and $M_h$ is the pressure space.

**Tip:** Stokes Uzawa

5.6. Navier-Stokes equations
// Mesh
mesh Th = square(10, 10);

// Fespace
fespace Xh(Th, P2);
Xh u1, u2;
Xh bc1, bc2;
Xh b;

fespace Mh(Th, P1);
Mh p;
Mh ppp; // ppp is a working pressure

// Problem
varf bx (u1, q) = int2d(Th)(-(dx(u1)*q));
varf by (u1, q) = int2d(Th)(-(dy(u1)*q));
varf a (u1, u2)
= int2d(Th)(
  dx(u1)*dx(u2)
  + dy(u1)*dy(u2)
) + on(3, u1=1)
  + on(1, 2, 4, u1=0) ;
// remark: put the on(3,u1=1) before on(1,2,4,u1=0)
// because we want zero on intersection

matrix A = a(Xh, Xh, solver=CG);
matrix Bx = bx(Xh, Mh); // B=(Bx, By)
matrix By = by(Xh, Mh);

bc1[] = a(0,Xh); // boundary condition contribution on u1
bc2 = 0; // no boundary condition contribution on u2

// p_h^n + B A^-1 - B^* p_h^n = -div u_h
// is realized as the function divup
func real[int] divup (real[int] & pp)
{
  // compute u1(pp)
  b[] = Bx'*pp;
  b[] *= -1;
  b[] += bc1[];
  u1[] = A^-1*b[];
  // compute u2(pp)
  b[] = By'*pp;
  b[] *= -1;
  b[] += bc2[];
  u2[] = A^-1*b[];
  // u^n = (A^-1 Bx^T p_h^n, By^T p_h^n)^T
  ppp[] = Bx*u1[]; // ppp = Bx u_1
  ppp[] += By*u2[]; //+ By u_2
  return ppp[] ;
}

// Initialization
p=0; //p_h^0 = 0
LinearCG(divup, p[], eps=1.e-6, nbiter=50); // p_h^(n+1) = p_h^n + B u_h^n
// if n> 50 or |p_h^(n+1) - p_h^n| <= 10^-6, then the loop end
5.6.2 NSUzawaCahouetChabart.edp

In this example we solve the Navier-Stokes equation past a cylinder with the Uzawa algorithm preconditioned by the Cahouet-Chabart method (see [GLOWINSKI2003] for all the details).

The idea of the preconditioner is that in a periodic domain, all differential operators commute and the Uzawa algorithm comes to solving the linear operator ∇.((αId + νΔ)^(-1)∇, where Id is the identity operator. So the preconditioner suggested is αΔ^{-1} + νId.

To implement this, we do:

```
Tip: NS Uzawa Cahouet Chabart
```

```
// Parameters
verbosity = 0;
real D = 0.1;
real H = 0.41;
real cx0 = 0.2;
real cy0 = 0.2; //center of cylinder
real xa = 0.15;
real ya = 0.2;
real xe = 0.25;
real ye = 0.2;
int nn = 15;

//TODO
real Um = 1.5; //max velocity (Rey 100)
real nu = 1e-3;

func U1 = 4.*Um*y*(H-y)/(H*H); //Boundary condition
func U2 = 0.;
real T=2;
real dt = D/nn/Um; //CFL = 1
real epspq = 1e-10;
real eps = 1e-6;

// Variables
func Ub = Um*2./3.;
real alpha = 1/dt;
real Rey = Ub*D/nu;
real t = 0.;

// Mesh
border fr1(t=0, 2.2){x=t; y=0; label=1;}
border fr2(t=0, H){x=2.2; y=t; label=2;}
border fr3(t=2.2, 0){x=t; y=H; label=1;}
border fr4(t=H, 0){x=0; y=t; label=1;}
border fr5(t=2*pi, 0){x=cx0*D*sin(t)/2; y=cy0*D*cos(t)/2; label=3;}

mesh Th = buildmesh(fr1(5*nn) + fr2(nn) + fr3(5*nn) + fr4(nn) + fr5(-nn*3));

(continues on next page)
```
// Fespace
fespace Mh(Th, [P1]);
Mh p;

fespace Xh(Th, [P2]);
Xh u1, u2;

fespace Wh(Th, [P1dc]);
Wh w; // vorticity

// Macro
macro grad(u) [dx(u), dy(u)] //
macro div(u1, u2) (dx(u1) + dy(u2)) //

// Problem
varf von1 ([u1, u2, p], [v1, v2, q])
  = on(3, u1=0, u2=0)
  + on(1, u1=U1, u2=U2)
;

// remark : the value 100 in next varf is manually fitted, because free outlet.
varf vA (p, q) =
  int2d(Th) (grad(p)' * grad(q)
  + int1d(Th, 2) (100*p*q
  )
;

varf vM (p, q)
  = int2d(Th, qft=qf2pT) (p*q
  + on(2, p=0)
  )
;

varf vu ([u1], [v1])
  = int2d(Th) (alpha*(u1*v1)
  + nu*(grad(u1)' * grad(v1))
  )
  + on(1, 3, u1=0)

;

varf vu1 ([p], [v1]) = int2d(Th) (p*dx(v1));
varf vu2 ([p], [v1]) = int2d(Th) (p*dy(v1));

varf vonu1 ([u1], [v1]) = on(1, u1=U1) + on(3, u1=0);
varf vonu2 ([u1], [v1]) = on(1, u1=U2) + on(3, u1=0);

matrix pAM = vM(Mh, Mh, solver=UMFPACK);
matrix pAA = vA(Mh, Mh, solver=UMFPACK);
matrix AU = vu(Xh, Xh, solver=UMFPACK);
matrix B1 = vu1(Mh, Xh);
matrix B2 = vu2(Mh, Xh);
real[int] brhs1 = vonu1(0, Xh);
real[int] brhs2 = vonu2(0, Xh);

varf vrhs1(uu, vv) = int2d(Th)(convect([u1, u2], -dt, u1)*vv*alpha) + vonu1;
varf vrhs2(v2, v1) = int2d(Th)(convect([u1, u2], -dt, u2)*v1*alpha) + vonu2;

// Uzawa function
func real[int] JUzawa (real[int] & pp){
    real[int] b1 = brhs1; b1 += B1*pp;
    real[int] b2 = brhs2; b2 += B2*pp;
    u1[] = AU^-1 * b1;
    u2[] = AU^-1 * b2;
    pp = B1'*u1[];
    pp += B2'*u2[];
    pp = -pp;
    return pp;
}

// Preconditioner function
func real[int] Precon (real[int] & p){
    real[int] pa = pAA^-1*p;
    real[int] pm = pAM^-1*p;
    real[int] pp = alpha*pa + nu*pm;
    return pp;
}

// Initialization
p = 0;

// Time loop
int ndt = T/dt;
for(int i = 0; i < ndt; ++i){
    // Update
    brhs1 = vrhs1(0, Xh);
    brhs2 = vrhs2(0, Xh);

    // Solve
    int res = LinearCG(JUzawa, p[], precon=Precon, nbiter=100, verbosity=10, veps=eps);
    assert(res==1);
    eps = -abs(eps);

    // Vorticity
    w = -dy(u1) + dx(u2);
    plot(w, fill=true, wait=0, nbiso=40);

    // Update
    dt = min(dt, T-t);
    t += dt;
    if(dt < 1e-10*T) break;
}

// Plot
plot(w, fill=true, nbiso=40);

// Display
cout << "u1 max = " << u1[].linfty

5.6. Navier-Stokes equations
Warning: Stop test of the conjugate gradient
Because we start from the previous solution and the end the previous solution is close to the final solution, don’t take a relative stop test to the first residual, take an absolute stop test (negative here).

Fig. 5.17: The vorticity at Reynolds number 100 a time 2s with the Cahouet-Chabart method.

5.7 Variational Inequality

We present, a classical example of variational inequality.

Let us denote $C = \{ u \in H^1_0(\Omega), u \leq g \}$

The problem is:

$$ u = \arg \min_{u \in C} J(u) = \frac{1}{2} \int_{\Omega} \nabla u \cdot \nabla u - \int_{\Omega} f u $$

where $f$ and $g$ are given function.

The solution is a projection on the convex $C$ of $f^*$ for the scalar product $((v, w)) = \int_{\Omega} \nabla v \cdot \nabla w$ of $H^1_0(\Omega)$ where $f^*$ is solution of:

$$ (f^*, v) = \int_{\Omega} f v, \forall v \in H^1_0(\Omega) $$

The projection on a convex satisfy clearly $\forall v \in C, \ ((u - v, u - f)) \leq 0$, and after expanding, we get the classical inequality:

$$ \forall v \in C, \ \int_{\Omega} \nabla (u - v) \nabla u \leq \int_{\Omega} (u - v) f $$

We can also rewrite the problem as a saddle point problem:

Find $\lambda, u$ such that:

$$ \max_{\lambda \in L^2(\Omega), \lambda \geq 0} \min_{u \in H^1_0(\Omega)} L(u, \lambda) = \frac{1}{2} \int_{\Omega} \nabla u \cdot \nabla u - \int_{\Omega} f u + \int_{\Omega} \lambda (u - g)^+ $$

where $((u - g)^+) = \max(0, u - g)$.
This saddle point problem is equivalent to find \( u, \lambda \) such that:

\[
\begin{align*}
\int_{\Omega} \nabla u \cdot \nabla v + \lambda v^+ & = \int_{\Omega} f u, \quad \forall v \in H^1_0(\Omega) \\
\int_{\Omega} (u - g)^+ & = 0, \quad \forall \mu \in L^2(\Omega), \mu \geq 0, \lambda \geq 0,
\end{align*}
\]

An algorithm to solve the previous problem is:

1. \( k=0 \), and choose \( \lambda_0 \) belong \( H^{-1}(\Omega) \)
2. Loop on \( k = 0 \), ..... 
   - set \( I_k = \{ x \in \Omega / \lambda_k + c (u_{k+1} - g) \leq 0 \} \)
   - \( V_{g,k+1} = \{ v \in H^1_0(\Omega) / v = g \text{ on } I_k \} \)
   - \( V_{0,k+1} = \{ v \in H^1_0(\Omega) / v = 0 \text{ on } I_k \} \)
   - Find \( u_{k+1} \in V_{g,k+1} \) and \( \lambda_{k+1} \in H^{-1}(\Omega) \) such that
     \[
     \begin{align*}
     \int_{\Omega} \nabla u_{k+1} \cdot \nabla v_{k+1} d\omega & = \int_{\Omega} f v_{k+1}, \quad \forall v_{k+1} \in V_{0,k+1} \\
     < \lambda_{k+1}, v > & = \int_{\Omega} \nabla u_{k+1} \cdot \nabla v - f v d\omega
     \end{align*}
     \]

   where \( <, > \) is the duality bracket between \( H^1_0(\Omega) \) and \( H^{-1}(\Omega) \), and \( c \) is a penalty constant (large enough).

You can find all the mathematics about this algorithm in [ITO2003] [HINTERMULLER2002].

Now how to do that in FreeFEM? The full example is:

```plaintext
Tip: Variational inequality

1. load "medit"
2. // Parameters
3. real eps = 1e-5;
4. real c = 1000; //penalty parameter of the algorithm
5. real tgv = 1e30; //a huge value for exact penalization
6. func f = 1; //right hand side function
7. func fd = 0; //Dirichlet boundary condition function
8. // Mesh
9. mesh Th = square(20, 20);
10. // Fespace
11. fespace Vh(Th, P1);
12. int n = Vh.ndof; //number of degree of freedom
13. Vh uh, uhp; //u^n+1 and u^n
14. Vh Ik; //to define the set where the contain is reached.
15. Vh g = 0.05; //discret function g
16. Vh lambda = 0;
17. // Problem
18. varf a (uh, vh) = int2d(Th) (dx(uh)*dx(vh)
19. + dy(uh)*dy(vh)
20. )

(continues on next page)
```

5.7. Variational Inequality 547
- \begin{verbatim}
27 - int2d(Th) (f*vh
28 ) + on(1, 2, 3, 4, uh=fv)
29 ;
30 //the mass Matrix construction
31 varf vM (uh, vh) = int2d(Th) (uh*vh);
32 //two versions of the matrix of the problem
33 matrix A = a(Vh, Vh, tgv-tgv, solver=CG); //one changing
34 matrix AA = a(Vh, Vh, solver=CG); //one for computing residual
35 matrix M = vM(Vh, Vh); //to do a fast computing of L^2 norm : sqrt(u'*(w=M*u))
36 real[int] Aiin(n);
37 real[int] Aii = A.diag; //get the diagonal of the matrix
38 real[int] rhs = a(0, Vh, tgv-tgv);
39 // Initialization
40 Ik = 0;
41 uh\[0\] = -tgv;
42 // Loop
43 for(int iter = 0; iter < 100; ++iter) {
44 // Update
45 real[int] b = rhs; //get a copy of the Right hand side
46 real[int] Ak(n); //the complementary of Ik ( !Ik = (Ik-1) )
47 Ak = 1.; Ak -= Ik\[\]\[\];
48 //adding new locking condition on b and on the diagonal if (Ik ==1 )
49 b = Ik\[\]\[\]* g\[\]; b *= tgv; b -= Ak .\[\]* rhs;
50 Aiin = Ik\[\]\[\]* tgv; Aiin += Ak .\[\]* Aii; //set Aii= tgv i in Ik
51 A.diag = Aiin; //set the matrix diagonal
52 set(A, solver=CG); //important to change preconditioning for solving
53 // Solve
54 uh\[\]\[\] = A^-1\[\]* b; //solve the problem with more locking condition
55 // Residual
56 lambda\[\]\[\] = AA \[\]* uh\[\]\[\]; //compute the residual (fast with matrix)
57 lambda\[\]\[\] += rhs; //remark rhs = -\int f v
58 Ik = (lambda + c*( g- uh)) < 0.; //the new locking value
59 // Plot
60 plot(Ik, wait=true, cmm=" lock set ", value=true, fill=true);
61 plot(uh, wait=true, cmm="uh");
62 // Error
63 //trick to compute L^2 norm of the variation (fast method)
64 real[int] diff(n), Mdiff(n);
65 diff = uh\[\]\[\] - uh\[\]\[\];
66 Mdiff = M*diff;
67 real err = sqrt(Mdiff'*diff);
68 cout << "|| u_{k=1} - u_{k} ||_2 = " << err << endl;
69 // Stop test
\end{verbatim}
Note: As you can see on this example, some vector or matrix operator are not implemented so a way is to skip the expression and we use operator `+=, -=` to merge the result.

5.8 Domain decomposition

We present three classic examples of domain decomposition technique: first, Schwarz algorithm with overlapping, second Schwarz algorithm without overlapping (also call Shur complement), and last we show to use the conjugate gradient to solve the boundary problem of the Shur complement.

5.8.1 Schwarz overlapping

To solve:

\[-\Delta u = f, \text{ in } \Omega = \Omega_1 \cup \Omega_2 \quad u|_{\Gamma} = 0\]

the Schwarz algorithm runs like this:

\[-\Delta u_1^{n+1} = f \text{ in } \Omega_1 \quad u_1^{n+1}|_{\Gamma_1} = u_2^n\]
\[-\Delta u_2^{n+1} = f \text{ in } \Omega_2 \quad u_2^{n+1}|_{\Gamma_2} = u_1^n\]

where \(\Gamma_i\) is the boundary of \(\Omega_i\) and on the condition that \(\Omega_1 \cap \Omega_2 \neq \emptyset\) and that \(u_i\) are zero at iteration 1.

Here we take \(\Omega_1\) to be a quadrangle, \(\Omega_2\) a disk and we apply the algorithm starting from zero.

Tip: Schwarz overlapping

```cpp
if(err < eps) break;

// Update
uhp[] = uh[];
}

// Plot
medit("uh", Th, uh);
```

(continued on next page)
Fig. 5.18: The 2 overlapping mesh TH and th

```plaintext
plot(th, TH, wait=true); // to see the 2 meshes

// Fespace
fespace vh(th, P1);
vh u=0, v;

fespace VH(TH, P1);
VH U, V;

// Problem
int i = 0;
problem PB (U, V, init=i, solver=Cholesky)
  = int2d(TH) (dx(U)*dx(V) + dy(U)*dy(V))
  + int2d(TH) (-V)
  + on(inside, U=u)
  + on(outside, U=0);

problem pb (u, v, init=i, solver=Cholesky)
  = int2d(th) (dx(u)*dx(v) + dy(u)*dy(v))
  + int2d(th) (-v)
  + on(inside, U=u)
  + on(outside, U=0);
```

(continues on next page)
5.8.2 Schwarz non overlapping Scheme

To solve:

\[-\Delta u = f \text{ in } \Omega = \Omega_1 \cup \Omega_2 \quad u|_{\Gamma} = 0\]

the Schwarz algorithm for domain decomposition without overlapping runs like this

Let introduce \(\Gamma_i\) is common the boundary of \(\Omega_1\) and \(\Omega_2\) and \(\Gamma_i^e = \partial \Omega_i \setminus \Gamma_i\).

The problem find \(\lambda\) such that \((u_1|_{\Gamma_i} = u_2|_{\Gamma_i})\) where \(u_i\) is solution of the following Laplace problem:

\[-\Delta u_i = f \text{ in } \Omega_i \quad u_i|_{\Gamma_i} = \lambda \quad u_i|_{\Gamma_i^e} = 0\]

To solve this problem we just make a loop with upgrading \(\lambda\) with

\[\lambda = \lambda \pm \frac{(u_1 - u_2)}{2}\]
where the sign $+$ or $-$ of $\pm$ is choose to have convergence.

**Tip:** Schwarz non-overlapping

```plaintext
// Parameters
int inside = 2; int outside = 1; int n = 4;

// Mesh
border a(t=1, 2){x=t; y=0; label=outside;};
border b(t=0, 1){x=2; y=t; label=outside;};
border c(t=2, 0){x-t; y=1; label=outside;};
border d(t=1, 0){x=1-t; y=t; label=inside;};
border e(t=0, 1){x=1-t; y=t; label=inside;};
border e1(t=pi/2, 2*pi){x=cos(t); y=sin(t); label=outside;};
mesh th = buildmesh(a(5*n) + b(5*n) + c(10*n) + d(5*n));
mesh TH = buildmesh(e(5*n) + e1(25*n));
plot(th, TH, wait=true);

// Fespace
fespace vh(th, P1);
vh u=0, v;
vh lambda=0;

fespace VH(TH, P1);
VH U, V;

// Problem
int i = 0;
problem PB (U, V, init=i, solver=Cholesky)
  = int2d(TH)(
    dx(U)*dx(V)
  + dy(U)*dy(V)
```

(continues on next page)
5.8.3 Schwarz conjugate gradient

To solve \(-\Delta u = f\) in \(\Omega = \Omega_1 \cup \Omega_2\) \(u|_{\Gamma_1} = 0\) the Schwarz algorithm for domain decomposition without overlapping runs like this.

Let introduce \(\Gamma_i\) is common the boundary of \(\Omega_1\) and \(\Omega_2\) and \(\Gamma_i^c = \partial \Omega_i \setminus \Gamma_i\).

The problem find \(\lambda\) such that \((u_1|_{\Gamma_i} = u_2|_{\Gamma_i})\) where \(u_i\) is solution of the following Laplace problem:

\[-\Delta u_i = f \text{ in } \Omega_i \quad u_i|_{\Gamma_i} = \lambda \quad u_i|_{\Gamma_i^c} = 0\]

The version of this example uses the Shur complement. The problem on the border is solved by a conjugate gradient method.

**Tip:** Schwarz conjugate gradient

First, we construct the two domains:
FreeFEM Documentation, Release 4.6

(a) Isovalues of the solution at iteration 0 without overlapping

(b) Isovalues of the solution at iteration 9 without overlapping

1 // Parameters
2 int inside = 2; int outside = 1; int n = 4;
3
4 // Mesh
5 border Gamma1(t=1, 2){x=t; y=0; label=outside;}
6 border Gamma2(t=0, 1){x=2; y=t; label=outside;}
7 border Gamma3(t=2, 0){x=t; y=1; label=outside;}
8 border GammaInside(t=1, 0){x=1-t; y=t; label=inside;}
9 border GammaArc(t=pi/2, 2*pi){x=cos(t); y=sin(t); label=outside;}
10 mesh Th1 = buildmesh(Gamma1(5*n) + Gamma2(5*n) + GammaInside(5*n) + Gamma3(5*n));
11 mesh Th2 = buildmesh(GammaInside(-5*n) + GammaArc(25*n));
12 plot(Th1, Th2);

Now, define the finite element spaces:

1 // Fespace
2 fespace Vh1(Th1, P1);
3 Vh1 u1, v1;
4 Vh1 lambda;
5 Vh1 p=0;
6
7 fespace Vh2(Th2, P1);
8 Vh2 u2, v2;

Note: It is impossible to define a function just on a part of boundary, so the λ function must be defined on the all domain Ω such as:

Vh1 lambda;

The two Poisson’s problems:
And, we define a border matrix, because the \( \lambda \) function is none zero inside the domain \( \Omega_1 \):

```plaintext
varf b(u2, v2, solver=CG) = int1d(Th1, inside)(u2*v2);
matrix B = b(Vh1, Vh1, solver=CG);
```

The boundary problem function,

\[
\lambda \mapsto \int_{\Gamma_1} (u_1 - u_2)v_1
\]

```plaintext
func real[int] BoundaryProblem (real[int] &l){
lambda[] = l; //make FE function form l
Pb1;
Pb2;
i++; //no refactorization i != 0
v1 = -(u1-u2);
lambda[] = B*v1[];
return lambda[];
}
```

**Note:** The difference between the two notations \( v_1 \) and \( v_1[] \) is: \( v_1 \) is the finite element function and \( v_1[] \) is the vector in the canonical basis of the finite element function \( v_1 \).
5.9 Fluid-structure coupled problem

This problem involves the Lamé system of elasticity and the Stokes system for viscous fluids with velocity \( \mathbf{u} \) and pressure \( p \):

\[
\begin{align*}
-\Delta \mathbf{u} + \nabla p &= 0 \quad \text{in } \Omega \\
\nabla \cdot \mathbf{u} &= 0 \quad \text{in } \Omega \\
\mathbf{u} &= \mathbf{u}_\Gamma \quad \text{on } \Gamma = \partial \Omega
\end{align*}
\]

where \( \mathbf{u}_\Gamma \) is the velocity of the boundaries. The force that the fluid applies to the boundaries is the normal stress

\[ h = (\nabla \mathbf{u} + \nabla \mathbf{u}^T) \mathbf{n} - pn \]

Elastic solids subject to forces deform: a point in the solid at \((x,y)\) goes to \((X,Y)\) after. When the displacement vector \( \mathbf{v} = (v_1, v_2) = (X - x, Y - y) \) is small, Hooke’s law relates the stress tensor \( \sigma \) inside the solid to the deformation tensor \( \epsilon \):

\[
\sigma_{ij} = \lambda \delta_{ij} \nabla \cdot \mathbf{v} + 2\mu \epsilon_{ij}, \quad \epsilon_{ij} = \frac{1}{2} \left( \frac{\partial v_i}{\partial x_j} + \frac{\partial v_j}{\partial x_i} \right)
\]

where \( \delta \) is the Kronecker symbol and where \( \lambda, \mu \) are two constants describing the material mechanical properties in terms of the modulus of elasticity, and Young’s modulus.

The equations of elasticity are naturally written in variational form for the displacement vector \( \mathbf{v}(x) \in V \) as:

\[
\int_{\Omega} \left[ 2\mu \epsilon_{ij}(\mathbf{v}) \epsilon_{ij}(\mathbf{w}) + \lambda \epsilon_{ii}(\mathbf{v}) \epsilon_{jj}(\mathbf{w}) \right] = \int_{\Omega} \mathbf{g} \cdot \mathbf{w} + \int_{\Gamma} h \cdot \mathbf{w}, \forall \mathbf{w} \in V
\]

The data are the gravity force \( \mathbf{g} \) and the boundary stress \( h \).

**Tip:** Fluid-structure In our example, the Lamé system and the Stokes system are coupled by a common boundary on which the fluid stress creates a displacement of the boundary and hence changes the shape of the domain where the Stokes problem is integrated. The geometry is that of a vertical driven cavity with an elastic lid. The lid is a beam with weight so it will be deformed by its own weight and by the normal stress due to the fluid reaction. The cavity is the \( 10 \times 10 \) square and the lid is a rectangle of height \( l = 2 \).

A beam sits on a box full of fluid rotating because the left vertical side has velocity one. The beam is bent by its own weight, but the pressure of the fluid modifies the bending.

The bending displacement of the beam is given by \((uu, vv)\) whose solution is given as follows.
real gravity = -0.05;
real coef = 0.2;

// Mesh (solid)
border a(t=2, 0){x=0; y=t; label=1;}
border b(t=0, 10){x=t; y=0; label=bottombeam;}
border c(t=0, 2){x=10; y=t; label=1;}
border d(t=0, 10){x=10-t; y=2; label=3;}
mesh th = buildmesh(b(20) + c(5) + d(20) + a(5));

// Fespace (solid)
fespace Vh(th, P1);
Vh uu, w, vv, s;

// Macro
real sqrt2 = sqrt(2.);
macro epsilon(u1, u2) [dx(u1), dy(u2), (dy(u1)+dx(u2))/sqrt2] //
macro div(u1, u2) (dx(u1) + dy(u2)) //

// Problem (solid)
real mu = E/(2*(1+sigma));
real lambda = E*sigma/({1+sigma}*(1-2*sigma));
solve Elasticity([uu, vv], [w, s])
  = int2d(th){
    lambda*div(w,s)*div(uu,vv)
    + 2.*mu*(epsilon(w,s)'*epsilon(uu,vv))
  }
  + int2d(th){
    - gravity*s
  }
  + on(1, uu=0, vv=0)
;
plot([uu, vv], wait=true);

mesh th1 = movemesh(th, [x+uu, y+vv]);
plot(th1, wait=true);

Then Stokes equation for fluids at low speed are solved in the box below the beam, but the beam has deformed the box (see border h):

// Mesh (fluid)
border e(t=0, 10){x=t; y=-10; label=1;}
border f(t=0, 10){x=10; y=-10+t; label=1;}
border g(t=0, 10){x=0; y=t; label=2;}
border h(t=0, 10){x=t; y=vv(t,0)*(t>=0.001)*(t <= 9.999); label=3;}
mesh sh = buildmesh(h(-20) + f(10) + e(10) + g(10));
plot(sh, wait=true);

We use the Uzawa conjugate gradient to solve the Stokes problem like in Navier-Stokes equations.
fespace Mh(sh, P1);
Mh p, ppp;

// Problem (fluid)
varf bx (u1, q) = int2d(sh)(-(dx(u1)*q));
varf by (u1, q) = int2d(sh)(-(dy(u1)*q));
varf Lap (u1, u2) = int2d(sh)(
  dx(u1)*dx(u2) + dy(u1)*dy(u2)
  + on(2, u1=1) + on(1, 3, u1=0));
bcl[] = Lap(0, Xh);

matrix A = Lap(Xh, Xh, solver=CG);
matrix Bx = bx(Xh, Mh);
matrix By = by(Xh, Mh);

func real[int] divup (real[int] & pp){
  int verb = verbosity;
  verbosity = 0;
  brhs[] = Bx'*pp;
  brhs[] += bcl[] .*bcx[];
  u1[] = A^-1*brhs[];
  brhs[] = By'*pp;
  brhs[] += bcl[] .*bcy[];
  u2[] = A^-1*brhs[];
  ppp[] = Bx*u1[];
  ppp[] += By*u2[];
  verbosity = verb;
  return ppp[];
}

do a loop on the two problems

// Coupling loop
for(int step = 0; step < 10; ++step){
  // Solve (fluid)
  LinearCG(divup, p[], eps=1.e-3, nbiter=50);
  divup(p[]);
}

Now the beam will feel the stress constraint from the fluid:

// Forces
Vh sigmain1, sigma22, sigma12;
Vh uu1=uu, vv1=vv;

sigmain1([x+uu, y+vv]) = (2*dx(u1) - p);
sigma22([x+uu, y+vv]) = (2*dy(u2) - p);
sigma12([x+uu, y+vv]) = (dx(u1) + dy(u2));

which comes as a boundary condition to the PDE of the beam:
// Solve (solid)
solve Elasticity2 ([uu, vv], [w, s], init=step)
  = int2d(th)(
    lambda*div(w,s)*div(uu,vv)
    + 2.*mu*(epsilon(w,s)'*epsilon(uu,vv))
  )
  + int2d(th)(
    - gravity*s
  )
  + int1d(th, bottombeam)(
    - coef*(sigma11*N.x*w + sigma22*N.y*s + sigma12*(N.y*w*N.x*s))
  )
  + on(1, uu=0, vv=0)
;

// Plot
plot([uu, vv], wait=1);

// Error
real err = sqrt(int2d(th)((uu-uu1)^2 + (vv-vv1)^2));
cout << "Erreur L2 = " << err << endl;

Notice that the matrix generated by Elasticity2 is reused (see init=i). Finally we deform the beam:

// Movemesh
th1 = movemesh(th, [x+0.2*uu, y+0.2*vv]);
plot(th1, wait=true);

Fluid velocity and pressure, displacement vector of the structure and displaced geometry in the fluid-structure interaction of a soft side and a driven cavity are shown Fig. 5.22, Fig. 5.23a and Fig. 5.23b

![Fig. 5.22: Velocity and pressure](image-url)
5.10 Transmission problem

Consider an elastic plate whose displacement change vertically, which is made up of three plates of different materials, welded on each other.

Let \( \Omega_i, i = 1, 2, 3 \) be the domain occupied by \( i \)-th material with tension \( \mu_i \) (see Soap film).

The computational domain \( \Omega \) is the interior of \( \Omega_1 \cup \Omega_2 \cup \Omega_3 \). The vertical displacement \( u(x, y) \) is obtained from:

\[
-\mu_i \Delta u = f \quad \text{in} \quad \Omega_i \\
\mu_i \partial_n u |_{\Gamma_i} = -\mu_j \partial_n u |_{\Gamma_j} \quad \text{on} \quad \Omega_i \cap \Omega_j \quad \text{if} \quad 1 \leq i < j \leq 3
\]

(5.15)

where \( \partial_n u |_{\Gamma_i} \) denotes the value of the normal derivative \( \partial_n u \) on the boundary \( \Gamma_i \) of the domain \( \Omega_i \).

By introducing the characteristic function \( \chi_i \) of \( \Omega_i \), that is:

\[
\chi_i(x) = 1 \quad \text{if} \quad x \in \Omega_i; \quad \chi_i(x) = 0 \quad \text{if} \quad x \notin \Omega_i
\]

we can easily rewrite (5.15) to the weak form. Here we assume that \( u = 0 \) on \( \Gamma = \partial \Omega \).

Transmission problem: For a given function \( f \), find \( u \) such that:

\[
\begin{align*}
    a(u, v) &= \ell(f, v) \quad \text{for all} \quad v \in H^1_0(\Omega) \\
    a(u, v) &= \int_\Omega \mu \nabla u \cdot \nabla v \\
    \ell(f, v) &= \int_\Omega f v
\end{align*}
\]

where \( \mu = \mu_1 \chi_1 + \mu_2 \chi_2 + \mu_3 \chi_3 \). Here we notice that \( \mu \) become the discontinuous function.

This example explains the definition and manipulation of region, i.e. sub-domains of the whole domain. Consider this L-shaped domain with 3 diagonals as internal boundaries, defining 4 sub-domains:

```plaintext
// Mesh
border a(t=0, 1)(x=t; y=0);
border b(t=0, 0.5)(x=1; y=t);
border c(t=0, 0.5)(x=1-t; y=0.5);
border d(t=0.5, 1)(x=0.5; y=t);
border e(t=0.5, 1)(x=1-t; y=1);
border f(t=0, 1)(x=0; y=1-t);
border i1(t=0, 0.5)(x=t; y=1-t);
border i2(t=0, 0.5)(x=t; y=t);
border i3(t=0, 0.5)(x=1-t; y=t);

mesh th = buildmesh(a(6) + b(4) + c(4) + d(4) + e(4) + f(6) + i1(6) + i2(6) + i3(6));

// Fespace
fespace Ph(th, P0); //constant discontinuous functions / element
Ph reg=region; //defined the P0 function associated to region number

// Plot
plot(reg, fill=true, wait=true, value=true);
```
Fig. 5.24: The function \texttt{reg}

\texttt{region} is a keyword of \texttt{FreeFEM} which is in fact a variable depending on the current position (is not a function today, use \texttt{Ph \ reg=region;} to set a function). This variable value returned is the number of the sub-domain of the current position. This number is defined by \texttt{buildmesh} which scans while building the mesh all its connected component.

So to get the number of a region containing a particular point one does:

```
1  // Characteristic function
2  int nupper = reg(0.4, 0.9); //get the region number of point (0.4,0.9)
3  int nlower = reg(0.9, 0.1); //get the region number of point (0.4,0.1)
4  cout << "nlower = " << nlower << " , nupper = " << nupper<< endl;
5  Ph nu = 1 + 5*(region==nlower) + 10*(region==nupper);
6
7  // Plot
8  plot(nu, fill=true, wait=true);
```

This is particularly useful to define discontinuous functions such as might occur when one part of the domain is copper and the other one is iron, for example.

We this in mind we proceed to solve a Laplace equation with discontinuous coefficients ($\nu$ is 1, 6 and 11 below).

```
1  // Problem
2  solve lap (u, v)
3      = int2d(th)(
4          nu*(dx(u)*dx(v) + dy(u)*dy(v))
5      )
6     + int2d(th)(
7        - 1*v
8      )
9     + on(a, b, c, d, e, f, u=0)
10  ;
11
12  // Plot
13  plot(u);
```

5.10. Transmission problem
Fig. 5.25: The function \( nu \)

Fig. 5.26: The isovalue of the solution \( u \)
5.11 Free boundary problems

The domain $\Omega$ is defined with:

```plaintext
// Parameters
real L = 10; //length
real hl = 2.1; //left height
real hr = 0.35; //right height
int n = 4;

// Mesh
border a(t=0, L){x=t; y=0;} //bottom: Gamma_a
border b(t=0, hr){x=L; y=t;} //right: Gamma_b
border f(t=L, 0){x=t; y=t*(hr-hl)/L+hl;} //free surface: Gamma_f
border d(t=hl, 0){x=0; y=t;} //left: Gamma_d
mesh Th = buildmesh(a(10*n) + b(6*n) + f(8*n) + d(3*n));
plot(Th);
```

![Fig. 5.27: The mesh of the domain $\Omega$.](image)

The free boundary problem is:

Find $u$ and $\Omega$ such that:

$$
\begin{cases}
    -\Delta u &= 0 & \text{in } \Omega \\
    u &= y & \text{on } \Gamma_b \\
    \frac{\partial u}{\partial n} &= 0 & \text{on } \Gamma_d \cup \Gamma_a \\
    \frac{\partial u}{\partial n} &= \frac{q}{\pi} n_x & \text{on } \Gamma_f \\
    u &= y & \text{on } \Gamma_f
\end{cases}
$$

We use a fixed point method;

$\Omega^0 = \Omega$

In two step, fist we solve the classical following problem:

$$
\begin{cases}
    -\Delta u &= 0 & \text{in } \Omega^n \\
    u &= y & \text{on } \Gamma_b^n \\
    \frac{\partial u}{\partial n} &= 0 & \text{on } \Gamma_d^n \cup \Gamma_a^n \\
    u &= y & \text{on } \Gamma_f^n
\end{cases}
$$

The variational formulation is:

Find $u$ on $V = H^1(\Omega^n)$, such that $u = y$ on $\Gamma_b^n$ and $\Gamma_f^n$

$$
\int_{\Omega^n} \nabla u \nabla u' = 0, \forall u' \in V \text{ with } u' = 0 \text{ on } \Gamma_b^n \cup \Gamma_f^n
$$
And secondly to construct a domain deformation $\mathcal{F}(x, y) = [x, y - v(x, y)]$ where $v$ is solution of the following problem:

$$
\begin{align*}
-\Delta v &= 0 & \text{in } \Omega^n \\
v &= 0 & \text{on } \Gamma^n_a \\
\frac{\partial v}{\partial n} &= 0 & \text{on } \Gamma^n_b \cup \Gamma^n_d \\
\frac{\partial v}{\partial n} &= \frac{q}{K} n_x & \text{on } \Gamma^n_f
\end{align*}
$$

The variational formulation is:

Find $v$ on $V$, such than $v = 0$ on $\Gamma^n_a$:

$$
\int_{\Omega^n} \nabla v \nabla v' = \int_{\Gamma^n_f} \left( \frac{\partial u}{\partial n} - \frac{q}{K} n_x \right) v', \quad \forall v' \in V \text{ with } v' = 0 \text{ on } \Gamma^n_a
$$

Finally the new domain $\Omega^{n+1} = \mathcal{F}(\Omega^n)$

**Tip:** Free boundary

The FreeFEM implementation is:

```plaintext
// Parameters
real L = 10; //length
real hr = 2.1; //left height
real hl = 0.35; //right height
int n = 4;
real q = 0.02; //incoming flow
real K = 0.5; //permeability

// Mesh
border a(t=0, L){x=t; y=0;}; //bottom: Gamma_a
border b(t=0, hr){x=L; y=t;}; //right: Gamma_b
border f(t=L, 0){x=t; y=t*(hr-hl)/L+hl;}; //free surface: Gamma_f
border d(t=hl, 0){x=0; y=t;}; // left: Gamma_d
mesh Th = buildmesh(a(10*n) + b(6*n) + f(8*n) + d(3*n));
plot(Th);

// Fespace
fespace Vh(Th, P1);
Vh u, v;
Vh uu, vv;

// Problem
problem Pu (u, uu, solver=CG) = int2d(Th)(
    dx(u)*dx(uu) + dy(u)*dy(uu)
) + on(b, f, u=y) ;

problem Pv (v, vv, solver=CG) = int2d(Th)(
    dx(v)*dx(vv) + dy(v)*dy(vv)
) + on(a, v=0)
```

(continues on next page)
+ int1d(Th, f)(
    vv*((q/K)*N.y - (dx(u)*N.x + dy(u)*N.y))
  )
  ;

// Loop
int j = 0;
real errv = 1.;
real erradap = 0.001;
while (errv > 1e-6) {
  // Update
  j++;
  // Solve
  Pu;
  Pv;
  // Plot
  plot(Th, u, v);
  // Error
  errv = int1d(Th, f){v*v};
  // Movemesh
  real coef = 1.;
  real mintcc = checkmovemesh(Th, [x, y])/5.;
  real mint = checkmovemesh(Th, [x, y-v*coef]);
  if (mint < mintcc || j%10==0) { //mesh too bad => remeshing
    Th = adaptmesh(Th, u, err=erradap);
    mintcc = checkmovemesh(Th, [x, y])/5.;
  }
  while (1) {
    real mint = checkmovemesh(Th, [x, y-v*coef]);
    if (mint > mintcc) break;
    cout << "min |T| = " << mint << endl;
    coef /= 1.5;
  }
  Th=movemesh(Th, [x, y-coef*v]);
  // Display
  cout << endl << j << " - errv = " << errv << endl;
}
// Plot
plot(Th);
plot(u, wait=true);

5.11. Free boundary problems
5.12 Non-linear elasticity

The non-linear elasticity problem is: find the displacement \((u_1, u_2)\) minimizing \(J\):

\[
\min J(u_1, u_2) = \int_{\Omega} f(F) - \int_{\Gamma_p} P_a u_2
\]

where \(F(u_1, u_2) = A(E[u_1, u_2], E[u_1, u_2])\) and \(A(X, Y)\) is bilinear symmetric positive form with respect two matrix \(X, Y\).

where \(f\) is a given \(C^2\) function, and \(E[u_1, u_2] = (E_{ij})_{i=1,2, j=1,2}\) is the Green-Saint Venant deformation tensor defined with:

\[
E_{ij} = 0.5((\partial_i u_j + \partial_j u_i) + \sum_k \partial_i u_k \times \partial_j u_k)
\]

Denote \(u = (u_1, u_2), v = (v_1, v_2), w = (w_1, w_2)\). So, the differential of \(J\) is:

\[
DJ(u)(v) = \int DF2(u)(v) f'(F2(u))) - \int P_a v_2
\]

where \(DF2(u)(v) = 2A(DE[u](v), E[u])\) and \(DE\) is the first differential of \(E\).

The second order differential is:

\[
D^2 J(u)((v), (w)) = \int DF2(u)(v) DF2(u)(w) f''(F2(u)))
+ \int D^2 F2(u)(v, w) f'(F2(u)))
\]

where:

\[
D^2 F2(u)(v, w) = 2A( D^2 E[u](v, w), E[u] ) + 2A( DE[u](v), DE[u](w) ) .
\]

and \(D^2 E\) is the second differential of \(E\).

So all notations can be define with \texttt{macro}:

```plaintext
macro EL(u, v) [dx(u), (dx(v)+dy(u)), dy(v)] //is [epsilon_11, 2epsilon_12, epsilon_22]

macro ENL(u, v) [4(dx(u)*dx(u) + dx(v)*dx(v))*0.5, 5(dx(u)*dy(u) + dx(v)*dy(v)), 6(dy(u)*dy(u) + dy(v)*dy(v))*0.5]

//

macro dENL(u, v, uu, vv) [dx(u)*dx(uu) + dx(v)*dx(vv)), dx(u)*dy(uu) + dx(v)*dy(vv) + dy(u)*dy(uu) + dy(v)*dy(vv) + dx(uu)*dy(u) + dx(vv)*dy(v),
```

(continues on next page)
The Newton Method is:
choose $n = 0$, and $u_0, v_0$ the initial displacement

- loop:
  - find $(du, dv)$: solution of
    $D^2J(u_n, v_n)((w, s), (du, dv)) = DJ(u_n, v_n)(w, s), \forall w, s$
  - $u_{n+1} = u_n - du$, $v_{n+1} = v_n - dv$
  - until $(du, dv)$ small is enough

The way to implement this algorithm in FreeFEM is use a macro tool to implement $A$ and $F_2, f, f', f''$.

A macro is like in ccp preprocessor of C++, but this begin by macro and the end of the macro definition is before the comment // . In this case the macro is very useful because the type of parameter can be change. And it is easy to make automatic differentiation.
(continued from previous page)

```cpp
    (dy(u)*dy(u) + dy(v)*dy(v))*0.5
             ]  //
    
macro dENL(u, v, uu, vv) [
    (dx(u)*dx(uu) + dx(v)*dx(vv)),
    (dx(u)*dy(uu) + dx(v)*dy(vv) + dx(uu)*dy(u) + dx(vv)*dy(v)),
    (dy(u)*dy(uu) + dy(v)*dy(vv))
             ]  //

macro E(u, v) (EL(u, v) + ENL(u, v)) // is \[ E_{11}, 2E_{12}, E_{22} \]
macro dE(u, v, uu, vv) (EL(uu, vv) + dENL(u, v, uu, vv)) //
macro ddE(u, v, uu, vv, uuu, vvv) (dENL(uuu, vvv, uu, vv)) //

macro F2(u, v) (E(u, v)'*A*E(u, v)) //
macro dF2(u, v, uu, vv) (E(u, v)'*A*dE(u, v, uu, vv)*2.) //
macro ddF2(u, v, uu, vv, uuu, vvv) ( (dE(u, v, uu, vv)'*A*dE(u, v, uu, vvv)) + (E(u, v)'*A*ddE(u, v, uu, vv, uu, vvv)) ) //

macro f(u) (u)*(u)*0.25) //
macro df(u) (u)*0.5 //
macro ddf(u) (0.5) //

// Parameters
real mu = 0.012e5; // kg/cm^2
real lambda = 0.4e5; // kg/cm^2
real Pa = 1e2;

// sigma = 2*mu*E + lambda*tr(E)*Id
// A(u,v) = sigma(u):E(v)
//
// ( a b c )
// ( b c d )
//
// tr*Id : (a,b,c) -> (a+c,0,a+c)
// so the associated matrix is:
// ( 1 0 1 )
// ( 0 0 0 )
// ( 1 0 1 )
real a11 = 2*mu + lambda;
real a22 = mu; // because [0, 2*t12, 0]' A [0, 2*s12, 0] = 2*mu*(t12*s12 + t21*s21) =
            4*mu*t12*s12
real a33 = 2*mu + lambda;
real a12 = 0;
real a13 = lambda;
real a23 = 0;
// symmetric part
real a21 = a12;
real a31 = a13;
real a32 = a23;

// the matrix A
func A = [[a11, a12, a13], [a21, a22, a23], [a31, a32, a33]];
// Mesh
int n = 30;
```

(continues on next page)
int m = 10;
mesh Th = square(n, m, [x, .3*y]); // label: 1 bottom, 2 right, 3 up, 4 left;
int bottom = 1, right = 2, upper = 3, left = 4;
plot(Th);

// Fespace
fespace Wh(Th, P1dc);
Wh e2, fe2, dfe2, ddfe2;
fespace Vh(Th, [P1, P1]);
Vh [uu, vv] = [0, 0], [w, s], [un, vn] = [0, 0];
fespace Sh(Th, P1);
Sh u1, v1;

// Problem
varf vmass ([uu, vv], [w, s], solver=CG) = int2d(Th)(uu*w + vv*s);
matrix M = vmass(Vh, Vh);
problem NonLin([uu, vv], [w, s], solver=LU)
  = int2d(Th, qforder=1) // (D^2 J(un))
    dF2(un, vn, uu, vv)*dF2(un, vn, w, s)*ddfe2
    + ddF2(un, vn, uu, vv, w, s)*ddfe2
  - int1d(Th, upper){
    Pa*s
  }
  - int2d(Th, qforder=1) // (D J(un))
    dF2(un, vn, w, s)*dfe2
  + on(right, left, uu=0, vv=0);

// Newton's method
for (int i = 0; i < 10; i++){
  cout << "Loop " << i << endl;
  // Update
e2 = F2(un, vn);
dfe2 = df(e2);
  ddfe2 = ddf(e2);
  cout << "e2 max = " << e2[].max << ", min = " << e2[].min << endl;
  cout << "dfe2 max = " << dfe2[].max << ", min = " << dfe2[].min << endl;
  cout << "ddfe2 max = " << ddfe2[].max << ", min = " << ddfe2[].min << endl;
  // Solve
  NonLin;
  w[] = M*uu[];
  // Residual
  real res = sqrt(w[].* uu[]); // L^2 norm of [uu, vv]
  cout << " L^2 residual = " << res << endl;
  // Update
  v1 = vv;
  ul = uu;
  cout << "ul min = " << ul[].min << ", ul max = " << ul[].max << endl;
  cout << "v1 min = " << v1[].min << ", v1 max = " << v1[].max << endl;

(continues on next page)
// Plot
plot([uu, vv], wait=true, cmm="uu, vv");

// Update
un[] -= uu[];
plot([un, vn], wait=true, cmm="displacement");

if (res < 1e-5) break;
}

// Plot
plot([un, vn], wait=true);

// Movemesh
mesh th1 = movemesh(Th, [x+un, y+vn]);

// Plot
plot(th1, wait=true);

5.13 Compressible Neo-Hookean materials

Author: Alex Sadovsky

5.13.1 Notation

In what follows, the symbols $u$, $F$, $B$, $C$, $\sigma$ denote, respectively, the displacement field, the deformation gradient, the left Cauchy-Green strain tensor $B = FF^T$, the right Cauchy-Green strain tensor $C = F^TF$, and the Cauchy stress tensor.

We also introduce the symbols $I_1 := \text{tr} C$ and $J := \det F$. Use will be made of the identity:

$$\frac{\partial J}{\partial C} = JC^{-1}$$

The symbol $I$ denotes the identity tensor. The symbol $\Omega_0$ denotes the reference configuration of the body to be deformed. The unit volume in the reference (resp., deformed) configuration is denoted $dV$ (resp., $dV_0$); these two are related by:

$$dV = JdV_0,$$

which allows an integral over $\Omega$ involving the Cauchy stress $T$ to be rewritten as an integral of the Kirchhoff stress $\kappa = JT$ over $\Omega_0$.

5.13.2 Recommended References

For an exposition of nonlinear elasticity and of the underlying linear and tensor algebra, see [OGDEN1984]. For an advanced mathematical analysis of the Finite Element Method, see [RAVIART1998].
5.13.3 A Neo-Hookean Compressible Material

Constitutive Theory and Tangent Stress Measures

The strain energy density function is given by:

\[ W = \frac{\mu}{2} (I_1 - \text{tr} \, I - 2 \ln J) \]

(see [HORGAN2004], formula (12)).

The corresponding 2nd Piola-Kirchoff stress tensor is given by:

\[ S_n := \frac{\partial W}{\partial E}(F_n) = \mu (I - C^{-1}) \]

The Kirchhoff stress, then, is:

\[ \kappa = FSF^T = \mu (B - I) \]

The tangent Kirchhoff stress tensor at \( F_n \) acting on \( \delta F_{n+1} \) is, consequently:

\[ \frac{\partial \kappa}{\partial F}(F_n) \delta F_{n+1} = \mu \left[ F_n (\delta F_{n+1})^T + \delta F_{n+1} (F_n)^T \right] \]

The Weak Form of the BVP in the Absence of Body (External) Forces

The \( \Omega_0 \) we are considering is an elliptical annulus, whose boundary consists of two concentric ellipses (each allowed to be a circle as a special case), with the major axes parallel. Let \( P \) denote the dead stress load (traction) on a portion \( \partial \Omega_0 \) (= the inner ellipse) of the boundary \( \partial \Omega_0 \). On the rest of the boundary, we prescribe zero displacement.

The weak formulation of the boundary value problem is:

\[ 0 = \int_{\Omega_0} \kappa [F_n] : \left\{ (\nabla \otimes w)(F_n + \delta F_{n+1})^{-1} \right\} \]

\[ - \int_{\partial \Omega_0} P \cdot \hat{N}_0 \]

For brevity, in the rest of this section we assume \( P = 0 \). The provided FreeFEM code, however, does not rely on this assumption and allows for a general value and direction of \( P \).

Given a Newton approximation \( u_n \) of the displacement field \( u \) satisfying the BVP, we seek the correction \( \delta u_{n+1} \) to obtain a better approximation:

\[ u_{n+1} = u_n + \delta u_{n+1} \]

by solving the weak formulation:

\[ 0 = \int_{\Omega_0} \kappa [F_n + \delta F_{n+1}] : \left\{ (\nabla \otimes w)(F_n + \delta F_{n+1})^{-1} \right\} - \int_{\partial \Omega_0} P \cdot \hat{N}_0 \]

\[ = \int_{\Omega_0} \kappa [F_n] + \frac{\partial \kappa}{\partial F}[F_n] \delta F_{n+1} \cdot \left\{ (\nabla \otimes w)(F_n + \delta F_{n+1})^{-1} \right\} \]

\[ = \int_{\Omega_0} \kappa [F_n] + \frac{\partial \kappa}{\partial F}[F_n] \delta F_{n+1} \cdot \left\{ (\nabla \otimes w)(F_n^{-1} + F_n^{-2} \delta F_{n+1}) \right\} \]

\[ = \int_{\Omega_0} \kappa [F_n] : \left\{ (\nabla \otimes w)F_n^{-1} \right\} \]

\[ - \int_{\Omega_0} \kappa [F_n] : \left\{ (\nabla \otimes w)(F_n^{-2} \delta F_{n+1}) \right\} \]

\[ + \int_{\Omega_0} \left\{ \frac{\partial \kappa}{\partial F}[F_n] \delta F_{n+1} \right\} : \left\{ (\nabla \otimes w)F_n^{-1} \right\} \]

for all test functions \( w \),

where we have taken:

\[ \delta F_{n+1} = \nabla \otimes \delta u_{n+1} \]

Note: Contrary to standard notational use, the symbol \( \delta \) here bears no variational context. By \( \delta \) we mean simply an increment in the sense of Newton’s Method. The role of a variational virtual displacement here is played by \( w \).
5.13.4 An Approach to Implementation in FreeFEM

Introducing the code-like notation, where a string in < >’s is to be read as one symbol, the individual components of the tensor:

\[ < TanK > := \frac{\partial K}{\partial F} [F_n] \delta F_{n+1} \]

will be implemented as the macros \(< TanK11 >, < TanK12 >, \ldots \). The individual components of the tensor quantities:

\[ D_1 := F_n (\delta F_{n+1})^T + \delta F_{n+1} (F_n)^T, \]
\[ D_2 := F_n^T \delta F_{n+1}, \]
\[ D_3 := (\nabla \otimes w) F_n^{-2} \delta F_{n+1}, \]

and

\[ D_4 := (\nabla \otimes w) F_n^{-1}, \]

will be implemented as the macros:

\[ < d1Aux11 >, < d1Aux12 >, \ldots , < d1Aux22 >, \]
\[ < d2Aux11 >, < d2Aux12 >, \ldots , < d2Aux22 >, \]
\[ < d3Aux11 >, < d3Aux12 >, \ldots , < d3Aux22 >, \]
\[ < d4Aux11 >, < d4Aux12 >, \ldots , < d4Aux22 > \}

respectively.

In the above notation, the tangent Kirchhoff stress term becomes

\[ \frac{\partial K}{\partial F} (F_n) \delta F_{n+1} = \mu D_1 \]

while the weak BVP formulation acquires the form:

\[ 0 = \int_{\Omega} \kappa [F_n] : D_4 \]
\[ - \int_{\Omega} \kappa [F_n] : D_3 \]
\[ + \int_{\Omega} \{ \frac{\partial K}{\partial F} [F_n] \delta F_{n+1} \} : D_4 \}

for all test functions \( w \)

---

1 // Macro
2 //Macros for the gradient of a vector field (u1, u2)
3 macro grad11(u1, u2) (dx(u1)) //
4 macro grad21(u1, u2) (dy(u1)) //
5 macro grad12(u1, u2) (dx(u2)) //
6 macro grad22(u1, u2) (dy(u2)) //
7 //Macros for the deformation gradient
8 macro F11(u1, u2) (1.0 + grad11(u1, u2)) //
9 macro F12(u1, u2) (0.0 + grad12(u1, u2)) //
10 macro F21(u1, u2) (0.0 + grad21(u1, u2)) //
11 macro F22(u1, u2) (1.0 + grad22(u1, u2)) //
12 //Macros for the incremental deformation gradient
13 macro dF11(varu1, varu2) (grad11(varu1, varu2)) //
14 macro dF12(varu1, varu2) (grad12(varu1, varu2)) //
15 (continues on next page)
macro dF21(varu1, varu2) (grad21(varu1, varu2)) //
macro dF22(varu1, varu2) (grad22(varu1, varu2)) //

//Macro for the determinant of the deformation gradient
macro J(u1, u2) {
    F11(u1, u2)*F22(u1, u2) - F12(u1, u2)*F21(u1, u2)
} //

//Macros for the inverse of the deformation gradient
macro Finv11 (u1, u2) {
    F22(u1, u2) / J(u1, u2)
} //
macro Finv22 (u1, u2) {
    F11(u1, u2) / J(u1, u2)
} //
macro Finv12 (u1, u2) {
    - F12(u1, u2) / J(u1, u2)
} //
macro Finv21 (u1, u2) {
    - F21(u1, u2) / J(u1, u2)
} //

//Macros for the square of the inverse of the deformation gradient
macro FFinv11 (u1, u2) {
    Finv11(u1, u2)^2 + Finv12(u1, u2)*Finv21(u1, u2)
} //
macro FFinv12 (u1, u2) {
    Finv12(u1, u2)*(Finv11(u1, u2) + Finv22(u1, u2))
} //
macro FFinv21 (u1, u2) {
    Finv21(u1, u2)*(Finv11(u1, u2) + Finv22(u1, u2))
} //
macro FFinv22 (u1, u2) {
    Finv12(u1, u2)*Finv21(u1, u2) + Finv22(u1, u2)^2
} //

//Macros for the inverse of the transpose of the deformation gradient
macro FinvT11(u1, u2) (Finv11(u1, u2)) //
macro FinvT12(u1, u2) (Finv21(u1, u2)) //
macro FinvT21(u1, u2) (Finv12(u1, u2)) //
macro FinvT22(u1, u2) (Finv22(u1, u2)) //

//The left Cauchy-Green strain tensor
macro B11(u1, u2) {
    F11(u1, u2)^2 + F12(u1, u2)^2
} //
macro B12(u1, u2) {
    F11(u1, u2)*F21(u1, u2)
} //

(continues on next page)
+ F11(u1, u2)\cdot F22(u1, u2)
} //

macro B21(u1, u2) {
  F11(u1, u2)\cdot F21(u1, u2)
  + F12(u1, u2)\cdot F22(u1, u2)
} //

macro B22(u1, u2) {
  F21(u1, u2)^2 + F22(u1, u2)^2
} //

//The macros for the auxiliary tensors (D0, D1, D2, ...): Begin

//The tensor quantity D0 = F(n) (delta F(n+1))^T
macro d0Aux11 (u1, u2, varu1, varu2) {
  dF11(varu1, varu2) * F11(u1, u2)
  + dF12(varu1, varu2) * F12(u1, u2)
} //

macro d0Aux12 (u1, u2, varu1, varu2) {
  dF21(varu1, varu2) * F11(u1, u2)
  + dF22(varu1, varu2) * F12(u1, u2)
} //

macro d0Aux21 (u1, u2, varu1, varu2) {
  dF11(varu1, varu2) * F21(u1, u2)
  + dF12(varu1, varu2) * F22(u1, u2)
} //

macro d0Aux22 (u1, u2, varu1, varu2) {
  dF21(varu1, varu2) * F21(u1, u2)
  + dF22(varu1, varu2) * F22(u1, u2)
} //

///The tensor quantity D1 = D0 + D0^T
macro d1Aux11 (u1, u2, varu1, varu2) {
  2.0 \cdot d0Aux11 (u1, u2, varu1, varu2)
} //

macro d1Aux12 (u1, u2, varu1, varu2) {
  d0Aux11 (u1, u2, varu1, varu2)
  + d0Aux12 (u1, u2, varu1, varu2)
} //

macro d1Aux21 (u1, u2, varu1, varu2) {
  d1Aux12 (u1, u2, varu1, varu2)
} //

macro d1Aux22 (u1, u2, varu1, varu2) {
  2.0 \cdot d0Aux22 (u1, u2, varu1, varu2)
} //

///The tensor quantity D2 = F^{-T}_{n} \cdot dF_{n+1}
macro d2Aux11 (u1, u2, varu1, varu2) {
  dF11(varu1, varu2) * FinvT11(u1, u2)
  + dF21(varu1, varu2) * FinvT12(u1, u2)
} //

(continues on next page)
macro d2Aux12 (u1, u2, varu1, varu2) (  
    dF12(varu1, varu2) * FinvT11(u1, u2)  
    + dF22(varu1, varu2) * FinvT12(u1, u2)  
) //

macro d2Aux21 (u1, u2, varu1, varu2) (  
    dF11(varu1, varu2) * FinvT21(u1, u2)  
    + dF21(varu1, varu2) * FinvT22(u1, u2)  
) //

macro d2Aux22 (u1, u2, varu1, varu2) (  
    dF12(varu1, varu2) * FinvT21(u1, u2)  
    + dF22(varu1, varu2) * FinvT22(u1, u2)  
) //

///The tensor quantity \( D_3 = F^{-2}_{n} dF_{n+1} \)
macro d3Aux11 (u1, u2, varu1, varu2, w1, w2) (  
    dF11(varu1, varu2) * FFinv11(u1, u2) * grad11(w1, w2)  
    + dF21(varu1, varu2) * FFinv12(u1, u2) * grad11(w1, w2)  
    + dF11(varu1, varu2) * FFinv21(u1, u2) * grad12(w1, w2)  
    + dF21(varu1, varu2) * FFinv22(u1, u2) * grad12(w1, w2)  
) //

macro d3Aux12 (u1, u2, varu1, varu2, w1, w2) (  
    dF12(varu1, varu2) * FFinv11(u1, u2) * grad11(w1, w2)  
    + dF22(varu1, varu2) * FFinv12(u1, u2) * grad11(w1, w2)  
    + dF12(varu1, varu2) * FFinv21(u1, u2) * grad12(w1, w2)  
    + dF22(varu1, varu2) * FFinv22(u1, u2) * grad12(w1, w2)  
) //

macro d3Aux21 (u1, u2, varu1, varu2, w1, w2) (  
    dF11(varu1, varu2) * FFinv11(u1, u2) * grad21(w1, w2)  
    + dF21(varu1, varu2) * FFinv12(u1, u2) * grad21(w1, w2)  
    + dF11(varu1, varu2) * FFinv21(u1, u2) * grad22(w1, w2)  
    + dF21(varu1, varu2) * FFinv22(u1, u2) * grad22(w1, w2)  
) //

macro d3Aux22 (u1, u2, varu1, varu2, w1, w2) (  
    dF12(varu1, varu2) * FFinv11(u1, u2) * grad21(w1, w2)  
    + dF22(varu1, varu2) * FFinv12(u1, u2) * grad21(w1, w2)  
    + dF12(varu1, varu2) * FFinv21(u1, u2) * grad22(w1, w2)  
    + dF22(varu1, varu2) * FFinv22(u1, u2) * grad22(w1, w2)  
) //

///The tensor quantity \( D_4 = (grad w) * Finv \)
macro d4Aux11 (w1, w2, u1, u2) (  
    Finv11(u1, u2)*grad11(w1, w2)  
    + Finv21(u1, u2)*grad12(w1, w2)  
) //

macro d4Aux12 (w1, w2, u1, u2) (  
    Finv12(u1, u2)*grad11(w1, w2)  
    + Finv22(u1, u2)*grad12(w1, w2)  
) //

macro d4Aux21 (w1, w2, u1, u2) (
Finv11(u1, u2)*grad21(w1, w2) + Finv21(u1, u2)*grad22(w1, w2) }

macro d4Aux22 (w1, w2, u1, u2) {
    Finv12(u1, u2)*grad21(w1, w2) + Finv22(u1, u2)*grad22(w1, w2) }

//The macros for the auxiliary tensors (D0, D1, D2, ...): End

//The macros for the various stress measures: BEGIN
//The Kirchhoff stress tensor
macro StressK11(u1, u2) {
    mu * (B11(u1, u2) - 1.0) }

//The Kirchhoff stress tensor
macro StressK12(u1, u2) {
    mu * B12(u1, u2) }

//The Kirchhoff stress tensor
macro StressK21(u1, u2) {
    mu * B21(u1, u2) }

//The Kirchhoff stress tensor
macro StressK22(u1, u2) {
    mu * (B22(u1, u2) - 1.0) }

//The tangent Kirchhoff stress tensor
macro TanK11(u1, u2, varu1, varu2) {
    mu * d1Aux11(u1, u2, varu1, varu2) }

macro TanK12(u1, u2, varu1, varu2) {
    mu * d1Aux12(u1, u2, varu1, varu2) }

macro TanK21(u1, u2, varu1, varu2) {
    mu * d1Aux21(u1, u2, varu1, varu2) }

macro TanK22(u1, u2, varu1, varu2) {
    mu * d1Aux22(u1, u2, varu1, varu2) }

//The macros for the stress tensor components: END

// Parameters
real mu = 5.e2; //Elastic coefficients (kg/cm^2)
real D = 1.e3; // (1 / compressibility)
real Pa = -3.e2; // Stress loads
real InnerRadius = 1.e0; // The wound radius
real OuterRadius = 4.e0; // The outer (truncated) radius
real tol = 1.e-4; // Tolerance (L^2)
real InnerEllipseExtension = 1.e0; //Extension of the inner ellipse ((major axis) - (minor axis))
int m = 40, n = 20;

// Mesh
border InnerEdge(t=0, 2.*pi){x=(1.0 + InnerEllipseExtension)*InnerRadius*cos(t);
y=InnerRadius*sin(t); label=1;}
border OuterEdge(t=0, 2.*pi){x=(1.0 + 0.0*InnerEllipseExtension)*OuterRadius*cos(t);
y=OuterRadius*sin(t); label=2;}
mesh Th = buildmesh(InnerEdge(-m) + OuterEdge(n));
int bottom = 1, right = 2, upper = 3, left = 4;
plot(Th);

// Fespace
fespace Wh(Th, P1dc);
fespace Vh(Th, [P1, P1]);
Vh [w1, w2], [u1n, u2n];
Vh [ehat1x, ehat1y], [ehat2x, ehat2y];
Vh [auxVec1, auxVec2]; //The individual elements of the total 1st Piola-Kirchoff stress
Vh [ef1, ef2];

fespace Sh(Th, P1);
Sh p, ppp;
Sh StrK11, StrK12, StrK21, StrK22;
Sh u1, u2;

// Problem
varf vfMass1D(p, q) = int2d(Th)(p*q);
matrix Mass1D = vfMass1D(Sh, Sh, solver=CG);
p[] = 1;
ppp[] = Mass1D * p[];
real DomainMass = ppp[].sum;
cout << "DomainMass = " << DomainMass << endl;

varf vmass ([u1, u2], [v1, v2], solver=CG)
  = int2d(Th)( (u1*v1 + u2*v2) / DomainMass );
matrix Mass = vmass(Vh, Vh);
matrix Id = vmass(Vh, Vh);

//Define the standard Euclidean basis functions
[ehat1x, ehat1y] = [1.0, 0.0];
[ehat2x, ehat2y] = [0.0, 1.0];

real ContParam, dContParam;

problem neoHookeanInc ([varu1, varu2], [w1, w2], solver=LU)
  = int2d(Th, qforder=1)(
    - ( StressK11 (u1n, u2n) * d3Aux11(u1n, u2n, varu1, varu2, w1, w2) 
      + StressK12 (u1n, u2n) * d3Aux12(u1n, u2n, varu1, varu2, w1, w2) 
      + StressK21 (u1n, u2n) * d3Aux21(u1n, u2n, varu1, varu2, w1, w2) 
      + StressK22 (u1n, u2n) * d3Aux22(u1n, u2n, varu1, varu2, w1, w2) ) 
  + StressK11 (u1n, u2n) * d3Aux11(u1n, u2n, varu1, varu2, w1, w2)
  + StressK12 (u1n, u2n) * d3Aux12(u1n, u2n, varu1, varu2, w1, w2)
  + StressK21 (u1n, u2n) * d3Aux21(u1n, u2n, varu1, varu2, w1, w2)
  + StressK22 (u1n, u2n) * d3Aux22(u1n, u2n, varu1, varu2, w1, w2) 
  + StressK11 (u1n, u2n) * d3Aux11(u1n, u2n, varu1, varu2, w1, w2)
  + StressK12 (u1n, u2n) * d3Aux12(u1n, u2n, varu1, varu2, w1, w2)
  + StressK21 (u1n, u2n) * d3Aux21(u1n, u2n, varu1, varu2, w1, w2)
  + StressK22 (u1n, u2n) * d3Aux22(u1n, u2n, varu1, varu2, w1, w2) 
  + StressK11 (u1n, u2n) * d3Aux11(u1n, u2n, varu1, varu2, w1, w2)
  + StressK12 (u1n, u2n) * d3Aux12(u1n, u2n, varu1, varu2, w1, w2)
  + StressK21 (u1n, u2n) * d3Aux21(u1n, u2n, varu1, varu2, w1, w2)
  + StressK22 (u1n, u2n) * d3Aux22(u1n, u2n, varu1, varu2, w1, w2) 
  (continues on next page)
+ StressK22 (u1n, u2n) * d3Aux22(u1n, u2n, varu1, varu2, w1, w2)
)
+ TanK11 (u1n, u2n, varu1, varu2) * d4Aux11(w1, w2, u1n, u2n)
+ TanK12 (u1n, u2n, varu1, varu2) * d4Aux12(w1, w2, u1n, u2n)
+ TanK21 (u1n, u2n, varu1, varu2) * d4Aux21(w1, w2, u1n, u2n)
+ TanK22 (u1n, u2n, varu1, varu2) * d4Aux22(w1, w2, u1n, u2n)
)
+ int2d(Th, qforder=1)(
  StressK11 (u1n, u2n) * d4Aux11(w1, w2, u1n, u2n)
+ StressK12 (u1n, u2n) * d4Aux12(w1, w2, u1n, u2n)
+ StressK21 (u1n, u2n) * d4Aux21(w1, w2, u1n, u2n)
+ StressK22 (u1n, u2n) * d4Aux22(w1, w2, u1n, u2n)
)
//Choose one of the following two boundary conditions involving Pa:
// Load vectors normal to the boundary:
  - int1d(Th, 1)(
    Pa * (w1*N.x + w2*N.y)
  )
//Load vectors tangential to the boundary:
//  - int1d(Th, 1)(
//    Pa * (w1*N.y - w2*N.x)
//  )
+ on(2, varu1=0, varu2=0)
;

//Auxiliary variables
matrix auxMat;

// Newton's method
ContParam = 0.;
dContParam = 0.01;

//Initialization:
[varu1, varu2] = [0., 0.];
[u1n, u2n] = [0., 0.];
real res = 2.*tol;
real eforceres;
int loopcount = 0;
int loopmax = 45;

// Iterations
while (loopcount <= loopmax && res >= tol){
  loopcount ++;
  cout << "Loop " << loopcount << endl;

  // Solve
  neoHookeanInc;

  // Update
  u1 = varu1;
u2 = varu2;

  // Residual
  w1[] = Mass*varu1[];
  res = sqrt(w1[]’ * varu1[]); //L^2 norm of [varu1, varu2]
  cout << " L^2 residual = " << res << endl;
}
5.14 Whispering gallery modes

Author: I. S. Grudinin

In whispering gallery mode (WGM) resonators, which are typically spheres or disks, electromagnetic field is trapped by total internal reflections from the boundary. Modes of such resonators are distinguished by compact volume and record high quality factors (Q) in a broad range of frequencies.

Modern applications of such resonators include microwave and optical cavities for atomic clocks, cavity optomechanics, nonlinear and quantum optics. Analytical solutions for WG modes are only available for a limited number of idealized geometries, such as sphere or ellipsoid. Since resonator dimensions are typically much larger than optical wavelength, direct application of numerical 3D finite difference time domain (FDTD) or finite element methods (FEM) is not practical. It’s possible to solve the vectorial wave equation by reducing it to a two dimensional case by taking axial symmetry into account.

Such reduction leads to a system of 3 equations to be solved in a 2D “\(\rho - z\)” section of a resonator. Please refer to [OXBORROW2007] for a detailed derivation and to [GRUDININ2012] for an example of using FreeFEM to compute WGMs.

5.14.1 Wave equation for the WGMs

Since electric field is discontinuous on the surface of a dielectric and magnetic field is typically not, we derive our equations for the magnetic field. The electric field can be easily derived at a later stage from \(\vec{E} = \frac{i \omega}{\varepsilon_0} \hat{\epsilon}^{-1} \nabla \times \vec{H}\).

Following a standard procedure starting with Maxwell equations we derive a wave equation in a single-axis anisotropic medium such as an optical crystal:

\[
\nabla \times \left( \hat{\epsilon}^{-1} \nabla \times \vec{H} \right) - k_0^2 \vec{H} - \alpha \nabla \left( \nabla \cdot \vec{H} \right) = 0
\]

(5.16)

Here \(k_0 = \omega/c\) is the wavenumber, \(\alpha\) is the penalty term added to fight spurious FEM solutions. For anisotropic single-axis medium with \(\partial \hat{\epsilon} / \partial \phi = 0\) in cylindrical system of coordinates we have:

\[
\hat{\epsilon} = \begin{pmatrix}
\epsilon_{\rho} & 0 & 0 \\
0 & \epsilon_{\rho} & 0 \\
0 & 0 & \epsilon_z
\end{pmatrix}.
\]
We now assume axial symmetry of our electromagnetic fields and insert an imaginary unity in front of the $H_{\phi}$ to allow all field components to be real numbers and also to account for the phase shift of this component $\vec{H}(\rho, \phi, z) = \{H_{\rho}(\rho, z), iH_{\phi}(\rho, z), H_{z}(\rho, z)\} \times e^{im\phi}$.

We write the wave equation (5.16) explicitly in cylindrical coordinates, thus obtaining a set of three differential equations for the domain $\Omega$ given by the resonator’s cross section and some space outside:

\[
\begin{align*}
A_1 \{H^t_{\rho}, H^t_{\phi}, H^t_{z}\} &= 0 \\
A_2 \{H^t_{\rho}, H^t_{\phi}, H^t_{z}\} &= 0 \\
A_3 \{H^t_{\rho}, H^t_{\phi}, H^t_{z}\} &= 0
\end{align*}
\]

The numerical solutions of these equations and boundary conditions can be found with FreeFEM if we write the system in the weak, or integral form.

5.14.2 Weak formulation

In general, to obtain the integral or “weak” statements equivalent to system (5.17) and boundary conditions we form a scalar dot product between an arbitrary magnetic field test function $H^t = \{H^t_{\rho}, H^t_{\phi}, H^t_{z}\}$ and the components of our vectorial equation $A_1, A_2, A_3$, and integrate over the resonator’s cross section domain $\Omega$ (and its boundary for the boundary conditions):

\[
\int_{\Omega} (H^t_{\rho}A_1 + H^t_{\phi}A_2 + H^t_{z}A_3) d\Omega
\]

We can reduce the order of partial derivatives in this integral by using the Green’s formula for integration by parts. For example:

\[
\int_{\Omega} H^t_{z} \frac{\partial^2 H_{z}}{\partial \rho^2} d\Omega = -\int_{\Omega} \frac{\partial H^t_{z}}{\partial \rho} \frac{\partial H_{z}}{\partial \rho} d\Omega + \oint_{\Gamma} H^t_z \frac{\partial H_{z}}{\partial \rho} n_\rho d\Gamma
\]

Thus converting equations (5.17) we obtain a large expression for the weak form.

5.14.3 A dielectric sphere example with FreeFEM

We now compute the fundamental mode frequency for a fused silica sphere. The sphere is 36 micrometer in diameter, the refractive index is 1.46, the boundary condition is the magnetic wall (which can actually be omitted as it holds automatically).

```plaintext
1 // Parameters
2 real radius = 36;  // approximate radius of the cavity
3 real yb = -10, yt = -yb;  // window yb=bottom and yt=top coordinates
4 real xl = radius-5, xr = radius+3;  // window xl=left and xr=right coordinates
5 real angle = asin((yt)/radius);  // angle of the sphere segment to model in radians
6 int Nm = 60;  // number of mesh vertices per border
7 real ne = 1.46;  // n_e-extraordinary refractive index (root of permittivity parallel to z-axis, epara)
8 real no = 1.46;  // n_o-ordinary refractive index (root of permittivity orthogonal to z-axis, eorto)
9 real nm = 1;  // refractive index of surrounding medium (air)
10 int nev = 4;  // number of eigen values to find
11 int M = 213;  // azimuthal mode order ~ 2pi*n*R/lambda
12 real alpha = 1;  // penalty term
```

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// Mesh
border W1l(t=0, 1){x=xl+(radius*cos(angle)-xl)*(1-t); y=yt; label=1;}
border W1r(t=0, 1){x=xr-(xr-radius*cos(angle))*t; y=yt; label=1;}
border W2(t=0, 1){x=xr-(xr-radius*cos(angle))*t; y=yt; label=1;}
border W3l(t=0, 1){x=xl+(radius*cos(angle)-xl)*(t); y=yb; label=1;}
border W3r(t=0, 1){x=xr-(xr-radius*cos(angle))*(1-t); y=yb; label=1;}
border W4(t=0, 1){x=xl+(radius*cos(angle)-xl)*(t); y=yb; label=1;}
border W4(t=0, 1){x=xl+(radius*cos(angle)-xl)*(t); y=yb; label=1;}

mesh Th = buildmesh(W1r(Nm/4) + W1l(Nm/4) + W4(Nm) + W3l(Nm/4) + W3r(Nm/4) + W2(Nm) + S(Nm));
plot(Th, WindowIndex=0);

// Fespace
fespace Ph(Th, P0);
Ph reg = region;

int ncav = reg(xl+1, 0); // cavity
int nair = reg(xr-1, 0); //air
Ph eorto = no^2*(region==ncav) + nm^2*(region==nair); //subdomains for epsilon values inside and outside the resonators
Ph epara = ne^2*(region==ncav) + nm^2*(region==nair); //subdomains for epsilon values inside and outside the resonators

//supplementary variables to store eigenvectors, defined on mesh Th with P2 elements - Largange quadratic.
fespace Supp(Th, P2);
Supp eHsqr;

//3d vector FE space
fespace Vh(Th, [P2, P2, P2]);
Vh [Hr, Hphi, Hz], [vHr, vHphi, vHz]; //magnetic field components on Vh space and test functions vH

// Macro
//boundary condition macros
macro EWall(Hr, Hphi, Hz) (dy(Hr) - dx(Hz) + Hr*N.x + Hz*N.y - eorto*(Hphi - Hr*M+dx(Hphi)*x)*N.y)
macro MWall(Hr, Hphi, Hz) (Hphi + Hz*N.x - Hr*N.y + epara*(Hz+M - dy(Hphi)*x)*N.x + eorto*(Hphi - Hr*M+dx(Hphi)*x)*N.y )

// Problem
real sigma = (M/(ne*radius))^2+2; // value of the shift (k^2), where the modes will be found
varf b ([Hr, Hphi, Hz], [vHr, vHphi, vHz]) = int2d(Th)(x*(Hr*vHr+Hphi*vHphi+Hz*vHz))
)

; // OP = A - sigma B ; // the shifted matrix
varf op ([Hr, Hphi, Hz], [vHr, vHphi, vHz]) = int2d(Th)(

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\[
\begin{align*}
& \text{-}2vHzHHz/x/A/r \\
& + eorto*(dx(vHphi)*(Hphi - MvHr) + dx(Hphi)*(vHphi - MvHr)) //B \\
& + x*(eorto*dx(vHphi)*dx(Hphi) + epara*((dx(vHz) - dy(vHr))*(dx(Hz) - dy(Hr)) + dy(vHphi)*dy(Hphi))) //C \\
& + alpha*(vHrHr - M(vHphiHr + HphiHvHr) + M^2vHphiHphi)/x/D/r \\
& + (dx(vHr) + dy(vHz))*(Hr - M*Hphi) + (vHr - M*vHphi)*(dx(Hr) + dy(Hz)) //E \\
& - sigma*x*(vHrHr + vHphiHvHphi + vHzHHz) \\
& \text{electric wall boundary condition on the boundary of computation domain} \\
& + int1d(Th, 1)(EWall(Hr, Hphi, Hz)*EWall(vHr, vHphi, vHz) \\
& ) \\
& \text{setting sparse matrices and assigning the solver UMFPACK to solve eigenvalue problem} \\
& \text{matrix } B = b(Vh, Vh, \text{ solver=UMFPACK}); \\
& \text{matrix } OP = op(Vh, Vh, \text{ solver=UMFPACK}); \\
& \text{// Solve} \\
& \text{real [int] ev(nev); //to store the nev eigenvalue} \\
& \text{Vh [int] [eHr, eHphi, eHz] (nev); //to store the nev eigenvector} \\
& \text{// calling ARPACK on sparse matrices with the assigned solver UMFPACK:} \\
& \text{int } k = \text{EigenValue (OP, B, sym=true, sigma=sigma, value=ev, vector=eHr, tol=1e-10,} \\
& \text{ maxit=0, ncv=0);} \\
& \text{k = min(k, nev); //sometimes the number of converged eigen values} \\
& \text{//can be greater than nev} \\
& \text{//file to output mode values} \\
& \text{ofstream f("modes.txt");} \\
& \text{//setting number of digits in the file output} \\
& \text{int nold = f.precisio(11);} \\
& \text{// Plot & Save} \\
& \text{for (int i = 0; i < k; i++)} \\
& \quad \text{real lambda = 2*pi/sqrt(ev[i]);} \\
& \quad \text{eHsqr = (sqrt(eHr[i]^2 + eHphi[i]^2 + eHz[i]^2)); //intensity from magnetic field components} \\
& \quad \text{plot (eHsqr, WindowIndex=i, value=i, nbiso=20, LabelColors=1, aspectratio=1, cmm="Mode "+i", lambda="+lambda"", F="+(299792.458/lambda)");} \\
& \quad f << "Mode "+i << "", ka=" << sqrt(ev[i])*radius << endl; \\
\end{align*}
\]
6.1 Misc

6.1.1 Poisson's Equation

```cpp
// Parameters
int nn = 20;
real L = 1.;
real H = 1.;
real l = 0.5;
real h = 0.5;
func f = 1.;
func g = 0.;
int NAdapt = 10;

// Mesh
border b1(t=0, L){x=t; y=0;};
border b2(t=0, h){x=L; y=t;};
border b3(t=L, l){x=t; y=h;};
border b4(t=h, H){x=l; y=t;};
border b5(t=1, 0){x=t; y=H};
border b6(t=H, 0){x=0; y=t};

mesh Th = buildmesh(b1(nn*L) + b2(nn*h) + b3(nn*(L-l)) + b4(nn*(H-h)) + b5(nn*l) + b6(nn*H));

// Fespace
fespace Vh(Th, P1); // Change P1 to P2 to test P2 finite element
Vh u, v;

// Macro
macro grad(u) [dx(u), dy(u)]; //

// Problem
problem Poisson (u, v, solver=CG, eps=-1.e-6) = int2d(Th) (grad(u)' * grad(v) ) + int2d(Th) ( f * v )
```

(continues on next page)
6.1.2 Poisson’s equation 3D

```c
load "tetgen"

// Parameters
real hh = 0.1;
func ue = 2. * x * x + 3. * y * y + 4. * z * z + 5. * x * y + 6. * x * z + 1.;
func f = -18.;

// Mesh
mesh Th = square(10, 20, [x*pi/2, 2*y*pi]); // ]-pi/2, pi/2[X]0,2pi[
func f1 = cos(x)*cos(y);
func f2 = cos(x)*sin(y);
```
func f3 = sin(x);
func f1x = sin(x)*cos(y);
func f1y = -cos(x)*sin(y);
func f2x = -sin(x)*sin(y);
func f2y = cos(x)*cos(y);
func f3x = cos(x);
func f3y = 0;
func m11 = f1x^2 + f2x^2 + f3x^2;
func m21 = f1x*f1y + f2x*f2y + f3x*f3y;
func m22 = f1y^2 + f2y^2 + f3y^2;
func perio = [[4, y], [2, y], [1, x], [3, x]];
real vv = 1/square(hh);
Th = adaptmesh(Th, m11*vv, m21*vv, m22*vv, IsMetric=1, periodic=perio);
Th = adaptmesh(Th, m11*vv, m21*vv, m22*vv, IsMetric=1, periodic=perio);
plot(Th);
border cc(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
mesh Th2 = buildmesh(cc(50));
// Fespace
fespace Vh(Th3, P23d);
Vh u, v;
Vh uhe = ue;
cout << "uhe min: " << uhe[].min << " - max: " << uhe[].max << endl;
cout << uhe(0.,0.,0.) << endl;
fespace Vh2(Th2, P2); Vh2 u2, u2e;
// Macro
macro Grad3(u) [dx(u), dy(u), dz(u)] //
// Problem
problem Lap3d (u, v, solver=CG) = int3d(Th3)(
    Grad3(v)' * Grad3(u) ) - int3d(Th3)( f * v ) + on(0, 1, u=ue);
// Solve
Lap3d;
cout << "u min: " << u[].min << " - max: " << u[].max << endl;
// Error
real err = int3d(Th3)(square(u-ue));
cout << int3d(Th3)(1.) << " = " << Th3.measure << endl;
Vh d = ue - u;
cout << " err = " << err << " - diff l^intfy = " << d[].linfty << endl;

// Plot
u2 = u;
ue = u2;
plot(u2, wait=true);
plot(u2, u2e, wait=true);

Fig. 6.2: Iso-surfaces of the solution

6.1.3 Stokes Equation on a cube

load "msh3"
load "medit" // Dynamically loaded tools for 3D

// Parameters
int nn = 8;

// Mesh
mesh Th0 = square(nn, nn);
int[int] rup = [0, 2];
int[int] rdown = [0, 1];
int[int] rmid = [1, 1, 2, 1, 3, 1, 4, 1];
real zmin = 0, zmax = 1;

mesh3 Th = buildlayers(Th0, nn, zbound=[zmin, zmax],
                        reffacemid=rmid, reffaceup=rup, reffacelow=rdown);

medit("c8x8x8", Th); // 3D mesh visualization with medit

// Fespaces
fespace Vh2(Th0, P2);

fespace VVh(Th, [P2, P2, P2, P1]);

fespace VVh(Th, [P2, P2, P2, P1]);

Vh2 [ux, uz, p2];
VVh [u1, u2, u3, p];
VVh [v1, v2, v3, q];

// Macro
macro Grad(u) [dx(u), dy(u), dz(u)] //

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(a) Solution

(b) Associated mesh

Fig. 6.3: Stokes

(continued from previous page)

```cpp
macro div(u1,u2,u3) (dx(u1) + dy(u2) + dz(u3)) //

// Problem (directly solved)
solve vStokes ([u1, u2, u3, p], [v1, v2, v3, q]) = int3d(Th, qforder=3) {
  Grad(u1)' * Grad(v1) + Grad(u2)' * Grad(v2) + Grad(u3)' * Grad(v3)
  - div(u1, u2, u3) * q - div(v1, v2, v3) * p + 1e-10 * q * p
  + on(2, u1=1., u2=0, u3=0) + on(1, u1=0, u2=0, u3=0);
}

// Plot
plot(p, wait=1, nbiso=5); // 3D visualization of pressure isolines

// See 10 plan of the velocity in 2D
for(int i = 1; i < 10; i++){
  // Cut plane
  real yy = i/10.;
  // 3D to 2D interpolation
  ux = u1(x,yy,y);
  uz = u3(x,yy,y);
  p2 = p(x,yy,y);
  // Plot
  plot([ux, uz], p2, cmm="cut y = "+yy, wait=1);
}
```

6.1. Misc
6.1.4 Cavity

```cpp
//Parameters
int m = 300;
real L = 1;
real rho = 500.;
real mu = 0.1;
real uin = 1;
func fx = 0;
func fy = 0;
int[int] noslip = [1, 2, 4];
int[int] inflow = [3];
real dt = 0.1;
real T = 50;
real eps = 1e-3;

//Macros
macro div(u) (dx(u#x) + dy(u#y)) //
macro grad(u) [dx(u), dy(u)] //
macro Grad(u) [grad(u#x), grad(u#y)] //

//Time
real cpu;
real tabcpu;

//mesh
border C1(t = 0, L){ x = t; y = 0; label = 1; }
border C2(t = 0, L){ x = L; y = t; label = 2; }
border C3(t = 0, L){ x = L-t; y = L; label = 3; }
border C4(t = 0, L){ x = 0; y = L-t; label = 4; }
mesh th = buildmesh( C1(m) + C2(m) + C3(m) + C4(m) );

fespace UPh(th, [P2,P2,P1]);
UPh [ux, uy, p];
UPh [uhx, uhy, ph];
UPh [upx, upy, pp];

//Solve
varf navierstokes([ux, uy, p], [uhx, uhy, ph])
  = int2d(th){
    rho/dt*[ux, uy]'*[uhx, uhy]
    + mu*(Grad(u):Grad(uh))
    - p* div(uh)
    - ph* div(u)
    - 1e-10 *p*ph
  }

  + int2d(th) {
    [fx, fy]'*[uhx, uhy]
    + rho/dt*[convect([upx, upy], -dt, upx), convect([upx, upy], -dt, upy)]'*[uhx, -uhy]
  }

  + on(noslip, ux=0, uy=0)
```

(continues on next page)
55 + on(inflow, ux=uin, uy=0)
56 ;
57
//Initialization
58 [ux, uy, p]=[0, 0, 0];
59
matrix<real> NS = navierstokes(UPh, UPh, solver=sparsesolver);
60 real[int] NSrhs = navierstokes(0, UPh);
61
//Time loop
62 for(int j = 0; j < T/dt; j++){
63    [upx, upy, pp]=[ux, uy, p];
64    NSrhs = navierstokes(0, UPh);
65    ux[] = NS^-1 * NSrhs;
66    plot( [ux,uy], p, wait=0, cmm=j);
67 }
68
//CPU
69 cout << " CPU = " << clock()-cpu << endl;
70 tabcpu = clock()-cpu;

6.2 Mesh Generation

6.2.1 Square mesh

mesh Th0 = square(10 ,10);
mesh Th1 = square(4, 5);
real x0 = 1.2;
real x1 = 1.8;
real y0 = 0;
real y1 = 1;
int n = 5;
real m = 20;
mesh Th2 = square(n, m, [x0+(x1-x0)*x, y0+(y1-y0)*y]);
for (int i = 0; i < 5; ++i){
    int[int] labs = [11, 12, 13, 14];
    mesh Th3 = square(3, 3, flags=i, label=labs, region=10);
    plot(Th3, wait=1, cmm="square flags = "i");
}

6.2.2 Mesh adaptation

// Parameters
real eps = 0.0001;
real h = 1;
real hmin = 0.05;
6.2.3 Mesh adaptation for the Poisson’s problem

/// Parameters
real error = 0.1;

/// Mesh
border ba(t=0, 1){x=t; y=0; label=1;}
border bb(t=0, 0.5){x=1; y=t; label=1;}
border bc(t=0, 0.5){x=1-t; y=0.5; label=1;}
border bd(t=0.5, 1){x=0.5; y=t; label=1;}
border be(t=0.5, 1){x=1-t; y=1; label=1;}
border bf(t=0, 1){x=0; y=1-t; label=1;}
mesh Th = buildmesh(ba(6) + bb(4) + bc(4) + bd(4) + be(4) + bf(6));

/// Fespace
fespace Vh(Th, P1);
Vh u, v;
// Function
func f = 1;

// Problem
problem Poisson(u, v, solver=CG, eps=1.e-6)
  = int2d(Th)(
    dx(u)*dx(v) + dy(u)*dy(v)
  ) - int2d(Th)(
    f*v + on(1, u=0);

// Adaptmesh loop
for (int i = 0; i < 4; i++){
  Poisson;
  Th = adaptmesh(Th, u, err=error);
  error = error/2;
}

// Plot
plot(u);

6.2.4 Uniform mesh adaptation

mesh Th = square(2, 2); // The initial mesh
plot(Th, wait=true);

Th = adaptmesh(Th, 1./30., IsMetric=1, nbvx=10000);
plot(Th, wait=true);

Th = adaptmesh(Th, 1./30., IsMetric=1, nbvx=10000); // More than one time due to the
Th = adaptmesh(Th, 1./30., IsMetric=1, nbvx=10000); // adaptation bound `maxsubdiv`=
plot(Th, wait=true);

6.2.5 Borders

{
  int upper = 1;
  int others = 2;
  int inner = 3;

  border C01(t=0, 1){x=0; y=-1+t; label=upper;}
  border C02(t=0, 1){x=1.5-1.5*t; y=-1; label=upper;}
  border C03(t=0, 1){x=1.5; y=-t; label=upper;}
  border C04(t=0, 1){x=1+0.5*t; y=0; label=others;}
  border C05(t=0, 1){x=0.5+0.5*t; y=0; label=others;}
  border C06(t=0, 1){x=0.5*t; y=0; label=others;}
  border C11(t=0, 1){x=0.5; y=-0.5*t; label=inner;}
  border C12(t=0, 1){x=0.5+0.5*t; y=-0.5; label=inner;}
  border C13(t=0, 1){x=1; y=-0.5+0.5*t; label=inner;}
}
Fig. 6.5: Mesh adaptation (Poisson)

Fig. 6.6: Uniform mesh adaptation
```
int n = 10;
plot(C01(-n) + C02(-n) + C03(-n) + C04(-n) + C05(-n)
    + C06(-n) + C11(n) + C12(n) + C13(n), wait=true);
mesh Th = buildmesh(C01(-n) + C02(-n) + C03(-n) + C04(-n) + C05(-n)
    + C06(-n) + C11(n) + C12(n) + C13(n));
plot(Th, wait=true);

cout << "Part 1 has region number " << Th(0.75, -0.25).region << endl;
cout << "Part 2 has region number " << Th(0.25, -0.25).region << endl;
```

```}
```
```
```
```
border a(t=0, 2*pi){x=cos(t); y=sin(t); label=1;}
border b(t=0, 2*pi){x=0.3+0.3*cos(t); y=0.3*sin(t); label=2;}
plot(a(50) + b(30)); //to see a plot of the border mesh
mesh Thwithouthole = buildmesh(a(50) + b(-30));
plot(Thwithouthole);

real r=1;
border a(t=0, 2*pi){x=r*cos(t); y=r*sin(t); label=1;}
r=0.3;
border b(t=0, 2*pi){x=r*cos(t); y=r*sin(t); label=1;}
// mesh Thwithouthole = buildmesh(a(50) + b(-30)); // do not do this because the two
// circles have the same radius = 0.3$
```

### 6.2.6 Change

```verbatim
verbosity=3;
// Mesh
mesh Th1 = square(10, 10);
mesh Th2 = square(20, 10, [x+1, y]);
int[int] r1=[2, 0];
plot(Th1, wait=true);
Th1 = change(Th1, label=r1); // Change edges' label from 2 to 0
plot(Th1, wait=true);
int[int] r2=[4, 0];
Th2 = change(Th2, label=r2); // Change edges' label from 4 to 0
plot(Th2, wait=true);

mesh Th = Th1 + Th2; // 'gluing together' Th1 and Th2 meshes
cout << "nb lab = " << int1d(Th1,1,3) (1./lenEdge)+int1d(Th2,1,2,3) (1./lenEdge)
<< " == " << int1d(Th,1,2,3,4) (1./lenEdge) " == " ((10+20)+10)*2 << endl;
plot(Th, wait=true);
```
Fig. 6.7: Borders
fespace Vh(Th, P1);
Vh u, v;

macro Grad(u) [dx(u),dy(u)] // Definition of a macro

solve P(u, v) = int2d(Th)(Grad(u)'*Grad(v)) - int2d(Th)(v) + on(1, 3, u=0);

plot(u, wait=1);

6.2.7 Cube

load "msh3"

int[int] 16 = [37, 42, 45, 40, 25, 57];
int r11 = 11;
mesh3 Th = cube(4, 5, 6, [x*2-1, y*2-1, z*2-1], label=16, flags = 3, region=r11);
cout << "Volume = " << Th.measure << ", border area = " << Th.bordermeasure << endl;

int err = 0;
for(int i = 0; i < 100; ++i){
    real s = int2d(Th,i){1.};
    real sx = int2d(Th,i){x};
    real sy = int2d(Th,i){y};
    real sz = int2d(Th,i){z};
}
if(s){
    int ix = (sx/s+1.5);
    int iy = (sy/s+1.5);
    int iz = (sz/s+1.5);
    int ii = (ix + 4*(iy+1) + 16*(iz+1) );
    //value of ix,iy,iz => face min 0, face max 2, no face 1
    cout "Label = " << i << " s = " << s << " ix = " << iy << " iy = " << iz << " ii = " << ii << endl;
    if( i != ii ) err++;
}

real volr11 = int3d(Th,r11)(1.);
cout "Volume region = " << 11 << " volr11 = " << volr11 << endl;
if((volr11 - Th.measure )>1e-8) err++;
plot(Th, fill=false);
cout "Nb err = " << err << endl;
assert(err==0);

---

### 6.2.8 Empty mesh

```cpp
[网通2.t=0, 2*pi]{x=cos(t); y=sin(t); label=1;}
```

mesh Th = buildmesh(a(20));
Th = emptymesh(Th);
plot(Th);

mesh Th = square(10, 10);
int[int] ssd(Th.nt);
// Builds the pseudo region numbering
for(int i = 0; i < ssd.n; i++){
    int iq = i/2; // Because we have 2 triangles per quad
    int ix = iq%10;
    int iy = iq/10;
    ssd[i] = 1 + (ix>=5) + (iy>=5)*2;
```
6.2.9 3 points

// Square for Three-Point Bend Specimens fixed on Fix1, Fix2
// It will be loaded on Load
real a = 1, b = 5, c = 0.1;
int n = 5, m = b*n;
border Left(t=0, 2*a){x=-b; y=a-t;}
border Bot1(t=0, b/2-c){x=-b+t; y=a;}
border Fix1(t=0, 2*c){x=-b/2-c+t; y=a;}
border Bot2(t=0, b-2*c){x=b/2+c+t; y=a;}
border Fix2(t=0, 2*c){x=b/2-c+t; y=a;}
border Bot3(t=0, b/2-c){x=b/2+c+t; y=a;}
border Right(t=0, 2*a){x=b; y=-a+t;}
border Top1(t=0, b-c){x=b-t; y=a;}
border Load(t=0, 2*c){x=c-t; y=a;}
border Top2(t=0, b-c){x=c-t; y=a;}

mesh Th = buildmesh(Left(n) + Bot1(m/4) + Fix1(5) + Bot2(m/2) + Fix2(5) + Bot3(m/4) + Right(n) + Top1(m/2) + Load(10) + Top2(m/2));
plot(Th, bw=true);

6.2.10 Bezier

// A cubic Bezier curve connecting two points with two control points
func real bzi(real p0, real p1, real q1, real q2, real t){

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```cpp
return p0*(1-t)^3 + q1*3*(1-t)^2*t + q2*3*(1-t)*t^2 + p1*t^3;
```

```cpp
real[int] p00 = [0, 1], p01 = [0, -1], q00 = [-2, 0.1], q01 = [-2, -0.5];
real[int] p11 = [1, -0.9], q10 = [0.1, -0.95], q11 = [0.5, -1];
real[int] p21 = [2, 0.7], q20 = [3, -0.4], q21 = [4, 0.5];
real[int] q30 = [0.5, 1.1], q31 = [1.5, 1.2];
```

```cpp
// Parameters
int C1 = 99;
```

### 6.2.11 Build layer mesh

```cpp
load "msh3"
load "tetgen"
load "medit"
```

```cpp
// Parameters
int m = 5;
```

```cpp
mesh Th = buildmesh(G1(t=0, 1) + G2(t=0, 1) + G3(t=0, 1) + G4(t=0, 1));
plot(Th, bw=true);
```
\begin{verbatim}
7    int C2 = 98;
8
9    // 2D mesh
10   border C01(t=0, pi){ x=t; y=0; label=1;}
11   border C02(t=0, 2*pi){ x=pi; y=t; label=1;}
12   border C03(t=0, pi){ x=pi-t; y=2*pi; label=1;}
13   border C04(t=0, 2*pi){ x=0; y=2*pi-t; label=1;}
14
15   border C11(t=0, 0.7){ x=0.5+t; y=2.5; label=C1;}
16   border C12(t=0, 2){ x=1.2; y=2.5+t; label=C1;}
17   border C13(t=0, 0.7){ x=1.2-t; y=4.5; label=C1;}
18   border C14(t=0, 2){ x=0.5; y=4.5-t; label=C1;}
19
20   border C21(t=0, 0.7){ x=2.3+t; y=2.5; label=C2;}
21   border C22(t=0, 2){ x=3; y=2.5+t; label=C2;}
22   border C23(t=0, 0.7){ x=3-t; y=4.5; label=C2;}
23   border C24(t=0, 2){ x=2.3; y=4.5-t; label=C2;}
24
25   mesh Th = buildmesh(C01(10) + C02(10) + C03(10) + C04(10)
27       + C21(-5) + C22(-5) + C23(-5) + C24(-5));
28
29   mesh Ths = buildmesh(C01(10) + C02(10) + C03(10) + C04(10)
31
32   // Construction of a box with one hole and two regions
33   func zmin = 0.;
34   func zmax = 1.;
35   int MaxLayer = 10;
36
37   func XX = x*cos(y);
38   func YY = x*sin(y);
39   func ZZ = z;
40
41   int[int] r1 = [0, 41], r2 = [98, 98, 99, 99, 1, 56];
42   int[int] r3 = [4, 12]; // Change upper surface mesh's triangles labels
43   // generated by the 2D mesh's triangles Th
\end{verbatim}
```c++
44 // from label 4 to label 12
45 int[int] r4 = [4, 45]; // Change lower surface mesh's triangles labels
46 // generated by the 2D mesh's triangles Th
47 // from label 4 to label 45
48 mesh3 Th3 = buildlayers(Th, MaxLayer, 
49   zbound=[zmin, zmax], region=r1,
50   labelmid=r2, labelup=r3, labeldown=r4);
51 medit("box 2 regions 1 hole", Th3);
52
53 // Construction of a sphere with TetGen
54 func XX1 = cos(y)*sin(x);
55 func YY1 = sin(y)*sin(x);
56 func ZZ1 = cos(x);
57
58 real[int] domain = [0., 0., 0., 0., 0.001];
59 string test = "paACQ";
60 cout << "test = " << test << endl;
61 mesh3 Th3sph = tetgtransfo(Ths, 
62   transf=[XX1, YY1, ZZ1],
63   switch=test, nbofregions=1, regionlist=domain);
64 medit("sphere 2 regions", Th3sph);
```

### 6.2.12 Sphere

```c++
// Parameter
real hh = 0.1;

// Mesh 2D
mesh Th = square(10, 20, [x*pi/2, 2*y*pi]); // ]-pi/2, pi/2[X]0, 2pi[
// A parametrization of a sphere
func f1 = cos(x)*cos(y);
func f2 = cos(x)*sin(y);
func f3 = sin(x);
// Partial derivative of the parametrization DF
func f1x = sin(x)*cos(y);
func f1y = -cos(x)*sin(y);
```

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### 6.3 Finite Element

#### 6.3.1 Periodic 3D

```plaintext
load "msh3"
load "medit"

// Parameters
searchMethod = 1; // More safe search algo
real a = 1, d = 0.5, h = 0.5;
int nnb = 7, nni = 10;
```

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```c
int nz = 3;
func zmin = 0;
func zmax = h;

// Mesh 2D
border b1(t=0.5, -0.5){x=a*t; y=-a/2; label=1;}
border b2(t=0.5, -0.5){x=a/2; y=a*t; label=2;}
border b3(t=0.5, -0.5){x=a*t; y=a/2; label=3;}
border b4(t=0.5, -0.5){x=-a/2; y=a*t; label=4;}
border il(t=0.5, -0.5){x=d/2*cos(t); y=-d/2*sin(t); label=7;}
mesh Th = buildmesh(b1(-nnb) + b3(nnb) + b2(-nnb) + b4(nnb) + il(nni));

{ // Cleaning the memory correctly
  int[int] old2new(0:Th.nv-1);
  fespace Vh2(Th, P1);
  Vh2 sorder = x + y;
  sort(sorder[], old2new);
  int[int] new2old = old2new^-1; // Inverse permutation
  Th = change(Th, renumv=new2old);
  sorder[] = 0:Th.nv-1;
}

{ // Mesh 3D
  int[int] rup = [0, 5], rlow = [0, 6], rmid = [1, 1, 2, 2, 3, 3, 4, 4, 7, 7], rtet = [0, 41];
  mesh3 Th3 = buildlayers(Th, nz, zbound=[zmin, zmax],
                          reftet=rtet, reffacemid=rmid, reffaceup=rup, reffacelow=rlow);
  for(int i = 1; i <= 6; ++i)
    cout << " int " << i << " : " << int2d(Th3,i)(1.) << " " << int2d(Th3,i)(1./area)
         << endl;
  plot(Th3, wait=true);
  medit("Th3", Th3);
  fespace Vh(Th3, P2, periodic=[[1, x, z], [3, x, z], [2, y, z], [4, y, z], [5, x, y],
                             [6, x, y]]);
}
```

### 6.3.2 Lagrange multipliers

```c
// Parameters
func f = 1 + x - y;

// Mesh
mesh Th = square(10, 10);

// Fespace
fespace Vh(Th, P1);
int n = Vh.ndof;
int nl = n+1;
```

(continues on next page)
Vh uh, vh;

// Problem
varf va (uh, vh)
  = int2d(Th)(
    dx(uh)*dx(vh)
    + dy(uh)*dy(vh)
  );

varf vL (uh, vh) = int2d(Th)(f*vh);
varf vb (uh, vh) = int2d(Th)(l*vh);

matrix A = va(Vh, Vh);
real[int] b = vL(0, Vh);
real[int] B = vb(0, Vh);

// Block matrix
matrix AA = [
  [ A, B ],
  [ B', 0 ]
];
set(AA, solver=sparsesolver);

real[int] bb(n+1), xx(n+1), l1(l1);
bl = 0;
// Builds the right hand side block
bb = [b, bl];

// Solve
xx = AA^-1 * bb;

// Set values
[uh[], l1] = xx;

// Display
cout << " l = " << l(0) << " b(u, l) = " B'*uh[] << endl;

// Plot
plot(uh);
6.4 Visualization

6.4.1 Plot

```plaintext
mesh Th = square(5,5);
fespace Vh(Th, P1);

// Plot scalar and vectorial FE function
Vh uh=x*x+y*y, vh=-y^2+x^2;
plot(Th, uh, [uh, vh], value=true, wait=true);

// Zoom on box defined by the two corner points [0.1,0.2] and [0.5,0.6]
plot(uh, [uh, vh], bb=[[0.1, 0.2], [0.5, 0.6]],
     wait=true, grey=true, fill=true, value=true);

// Compute a cut
int n = 10;
real[int] xx(10), yy(10);
for (int i = 0; i < n; i++)
  x = i/real(n);
  y = i/real(n);
  xx[i] = i;
  yy[i] = uh; // Value of uh at point (i/10., i/10.)
plot([xx, yy], wait=true);

// File for gnuplot
ofstream gnu("plot.gp");
for (int i = 0; i < n; i++)
  gnu << xx[i] << " " << yy[i] << endl;

// Calls the gnuplot command, waits 5 seconds and generates a postscript plot (UNIX ONLY)
exec("echo 'plot "plot.gp" w l \n pause 5 \n set term postscript \n set output \n "gnuplot.eps" \n replot \n quit' | gnuplot");
```
Fig. 6.17: Plot
6.4.2 HSV

// From: http://en.wikipedia.org/wikiHSV_color_space
// The HSV (Hue, Saturation, Value) model defines a color space
// in terms of three constituent components:
// HSV color space as a color wheel
// Hue, the color type (such as red, blue, or yellow):
// Ranges from 0-360 (but normalized to 0-100% in some applications like here)
// Saturation, the "vibrancy" of the color: Ranges from 0-100%
// The lower the saturation of a color, the more "grayness" is present
// and the more faded the color will appear.
// Value, the brightness of the color: Ranges from 0-100%

mesh Th = square(10, 10, [2*x-1, 2*y-1]);
fespace Vh(Th, P1);
Vh uh=2-x^2-y^2;

real[int] colorhsv=[  // Color hsv model
        4./6., 1 , 0.5, // Dark blue
        4./6., 1 , 1,  // Blue
        5./6., 1 , 1, // Magenta
        1, 1 , 1, // Red
        1, 0.5 , 1 // Light red
];
real[int] viso(31);
for (int i = 0; i < viso.n; i++)
    viso[i] = i*0.1;
plot(uh, viso=viso(0:viso.n-1), value=true, fill=true, wait=true, hsv=colorhsv);

Fig. 6.18: Result

6.4.3 Medit

load "medit"

mesh Th = square(10, 10, [2*x-1, 2*y-1]);
FreeFEM Documentation, Release 4.6

**6.4.4 Paraview**

```cpp
load "iovtk"

mesh Th = square(10, 10, [2*x-1, 2*y-1]);

fespace Vh(Th, P1);
Vh u=2-x*x-y*y;

int[int] Order = [1];
string DataName = "u";
savevtk("u.vtu", Th, u, dataname=DataName, order=Order);
```

Fig. 6.19: Medit

(continued from previous page)
6.5 Algorithms & Optimizations

6.5.1 Algorithms

```c++
// Parameters
int nerr = 0;
int debugJ = 0;
int debugdJ = 0;
real umax = 0;

// Algorithms tests
{  
    func bool stop (int iter, real[int] u, real[int] g){
        cout << " stop = " << iter << " " << u.linfty << " " << g.linfty << endl;
        return g.linfty < 1e-5 || iter > 15;
    }
    // minimization of J(u) = 1./2 * sum (i+1) u_i^2 - b_i
    real[int] b(10), u(10);
    
    //J
    func real J (real[int] & u){
        real s = 0;
        for (int i = 0; i < u.n; i++)
            s += (i+1)*u[i]*u[i]*0.5 - b[i]*u[i];
        if (debugJ)
            cout << "J = " << s << " u = " << u[0] << " u[1] << endl;
        return s;
    }
    
    //the gradient of J (this is a affine version (the RHS is in)
    func real[int] DJ (real[int] &u){
        for (int i = 0; i < u.n; i++)
            u[i] = (i+1)*u[i];
        if (debugdJ)
        u -= b;
        if (debugdJ)
            cout << "
    }
    
    // minimization of J(u)
    stop(0, u, J);  
    
    // gradient of J
    DJ();
}
```

(continues on next page)
    return u; //return of global variable ok
}

//the gradient of the bilinear part of J (the RHS is remove)
func real[int] DJ0 (real[int] &u){
    for (int i = 0 ; i < u.n; i++)
        u[i] = (i+1)*u[i];
    if(debugdJ)
        cout << "dJ0: u =" << u[0] << " " << u[1] << " " << u[2] << endl;
    return u; //return of global variable ok
}

//erro calculation
func real error (real[int] & u, real[int] & b){
    real s = 0;
    for (int i = 0; i < u.n; i++)
        s += abs((i+1)*u[i] - b[i]);
    return s;
}

func real[int] matId (real[int] &u){ return u; }

int verb=5; //verbosity
b = 1.; //set right hand side
u = 0.; //set initial guess

LinearCG(DJ, u, eps=1.e-6, nbiter=20, precon=matId, verbosity=verb);
    cout << "LinearCG (Affine) : J(u) = " << J(u) << ", err = " << error(u, b) << endl;
    nerr += !(error(u,b) < 1e-5);
    if(nerr) cout << "sol: u = " << u[0] << " " << u[1] << " " << u[2] << endl;

b = 1;
    u = 0;
LinearCG(DJ, u, eps=1.e-15, nbiter=20, precon=matId, verbosity=verb, stop=stop);
    cout << "LinearCG (Affine with stop) : J(u) = " << J(u) << ", err = " << error(u, b) << endl;
    nerr += !(error(u,b) < 1e-5);
    if(nerr) cout << "sol: u = " << u[0] << " " << u[1] << " " << u[2] << endl;

b = 1;
    u = 0;
LinearCG(DJ0, u, b, eps=1.e-6, nbiter=20, precon=matId, verbosity=verb);
    cout << "LinearCG (Linear) : J(u) = " << J(u) << ", err = " << error(u, b) << endl;
    nerr += !(error(u,b) < 1e-5);
    if(nerr) cout << "sol: u = " << u[0] << " " << u[1] << " " << u[2] << endl;

b = 1;
    u = 0;
AffineGMRES(DJ, u, eps=1.e-6, nbiter=20, precon=matId, verbosity=verb);
    cout << "AffineGMRES (Affine) : J(u) = " << J(u) << ", err = " << error(u, b) << endl;
    nerr += !(error(u,b) < 1e-5);
    if(nerr) cout << "sol: u = " << u[0] << " " << u[1] << " " << u[2] << endl;
87 b=1;
88 u=0;
89 LinearGMRES(DJ0, u, b, eps=1.e-6, nbiter=20, precon=matId, verbosity=verb);
90 cout << "LinearGMRES (Linear) : J(u) = " << J(u) << " , err = " << error(u, b) << endl;
91 nerr += !(error(u,b) < 1e-5);
92 if(nerr) cout << "sol: u = " << u[0] << " " << u[1] << " " << u[2] << endl;
93 b=1;
94 u=0;
95 NLCG(DJ, u, eps=1.e-6, nbiter=20, precon=matId, verbosity=verb);
96 cout << "NLCG: J(u) = " << J(u) << " , err = " << error(u, b) << endl;
97 nerr += !(error(u,b) < 1e-5);
99 //warning: BFGS use a full matrix of size nxn (where n=u.n)
100 b=1;
101 u=2;
102 BFGS(J, DJ, u, eps=1.e-6, nbiter=20, nbiterline=20);
103 cout << "BFGS: J(u) = " << J(u) << " , err = " << error(u, b) << endl;
104 is(error(u,b) < 1e-5);
105 if(nerr) cout << "sol: u = " << u[0] << " " << u[1] << " " << u[2] << endl;
106 assert(nerr==0);
107 }
108 // A real non linear test
109 // Parameters
110 real a = 0.001;
111 real eps = 1e-6;
112 //f(u) = a*u + u-ln(1+u), f'(u) = a+ u/(1+u), f''(u) = 1/(1+u)^2
113 func real f(real u) { return u*a+u-log(1+u); }
114 func real df(real u) { return a+u/(1+u); }
115 func real ddf(real u) { return 1/(1+u)*(1+u); }
116 // Mesh
117 mesh Th = square(20, 20);
118 // Fespace
119 fespace Vh(Th, P1);
120 Vh b = 1;
121 Vh u = 0;
122 fespace Ph(Th, P0);
123 Ph alpha; //store df(|nabla u|^2)
124 // The functionnal J
125 //J(u) = 1/2 int_Omega f(|nabla u|^2) - int_Omega u b
126 func real J (real[int] & u){
127 Vh w;
128 w[] = u;
129 real r = int2d(Th)(0.5*f(dx(w)*dx(w) + dy(w)*dy(w)) - b*w);
130 cout << "J(u) = " << r << " " << u.min << " " << u.max << endl;
131 return r;
// The gradient of J
func real[int] dJ(real[int] & u) {
    Vh w;
    w[] = u;
    alpha = df(dx(w)*dx(w) + dy(w)*dy(w));
    varf au (uh, vh) = int2d(Th) {
        alpha*(dx(uh)*dx(vh) + dy(uh)*dy(vh))
        - b*vh
    } + on(1, 2, 3, 4, uh=0);
    u = au(0, Vh);
    return u; //warning: no return of local array
}

// Problem
alpha = df(dx(u)*dx(u) + dy(u)*dy(u));
varf alap (uh, vh) = int2d(Th) {
    alpha*(dx(uh)*dx(vh) + dy(uh)*dy(vh))
} + on(1, 2, 3, 4, uh=0);
varf amass(uh, vh) = int2d(Th) {
    uh*vh
} + on(1, 2, 3, 4, uh=0);
matrix Amass = amass(Vh, Vh, solver=CG);
matrix Alap = alap(Vh, Vh, solver=Cholesky, factorize=1);

// Preconditionner
func real[int] C(real[int] & u) {
    real[int] w = u;
    u = Alap^-1*w;
    return u; //warning: no return of local array variable
}

// Solve
int conv=0;
for(int i = 0; i < 20; i++) {
    conv = NLCG(dJ, u[], nbiter=10, precon=C, veps=eps, verbosity=5);
    if (conv) break;
    alpha = df(dx(u)*dx(u) + dy(u)*dy(u));
    Alap = alap(Vh, Vh, solver=Cholesky, factorize=1);
    cout << "Restart with new preconditionner " << conv << ", eps =" << eps << endl;
}

(continues on next page)
6.5.2 CMAES variational inequality

```plaintext
load "ff-cmaes"

// Parameters
int NN = 7;
func f1 = 1.;
func f2 = -1.;
func g1 = 0.;
func g2 = 0.1;
int iter = 0;
int nadapt = 1;
real starttol = 1e-10;
real bctol = 6.e-12;
real pena = 1000.;

// Mesh
mesh Th = square(NN, NN);

// Fespace
fespace Vh(Th, P1);
Vh ou1, ou2;

// Mesh adaptation loops
for (int al = 0; al < nadapt; ++al){
```

Fig. 6.21: Algorithms
// Problem
varf BVF (v, w)
   = int2d(Th)(
       0.5*dx(v)*dx(w)
     + 0.5*dy(v)*dy(w)
   )
;
varf LVF1 (v, w) = int2d(Th)(f1*w);
varf LVF2 (v, w) = int2d(Th)(f2*w);

matrix A = BVF(Vh, Vh);
real[int] b1 = LVF1(0, Vh);
real[int] b2 = LVF2(0, Vh);

varf Vbord (v, w) = on(1, 2, 3, 4, v=1);

Vh In, Bord;
Bord[] = Vbord(0, Vh, tgv=1);
In[] = Bord[] ? 0:1;
Vh gh1 = Bord*g1;
Vh gh2 = Bord*g2;

// Function which creates a vector of the search space type from
// two finite element functions
func int FEFToSSP (real[int] &fef1, real[int] &fef2, real[int] &ssp)
{
    int kX = 0;
    for (int i = 0; i < Vh.ndof; ++i){
        if (In[][i]){
            ssp[kX] = fef1[i];
            ssp[kX+In[]].sum = fef2[i];
            ++kX;
        }
    }
    return 1;
}

// Splits a vector from the search space and fills
// two finite element functions with it
func int SSPToFEF (real[int] &fef1, real[int] &fef2, real[int] &ssp)
{
    int kX = 0;
    for (int i = 0; i < Vh.ndof; ++i){
        if (In[][i]){
            fef1[i] = ssp[kX];
            fef2[i] = ssp[kX+In[]].sum;
            ++kX;
        } else{
            fef1[i] = gh1[][i];
            fef2[i] = gh2[][i];
        }
    }
    return 1;
}

func real IneqC (real[int] &X){
    real[int] constraints(In[]].sum);
    for (int i = 0; i < In[].sum; ++i){
      // Function which creates a vector of the search space type from
      // two finite element functions
      // Splits a vector from the search space and fills
      // two finite element functions with it
    }
81 constraints[i] = X[i] - X[i+In].sum;
82 constraints[i] = constraints[i] <= 0 ? 0. : constraints[i];
83 }
84 return constraints.l2;
85 }
86
87 func real J (real[int] &X)
88 {
89 Vh u1, u2;
90 SSPToFEF(u1[], u2[], X);
91 iter++;
92 real[int] Aul = A*u1[], Au2 = A*u2[];
93 Au1 -= b1;
94 Au2 -= b2;
95 real val = u1[]'*Aul + u2[]'*Au2;
96 val += pena * IneqC(X);
97 if (iter%200 == 199)
98 plot(u1, u2, nbiso=30, fill=1, dim=3, cmm="adapt level "+al+" - iteration "+iter+" - J = "+val, value=1);
99 return val;
100 }
101
102 // Solve
103 real[int] start(2*In[].sum);
104 if (al == 0)
105 start(0:In[].sum-1) = 0.;
106 start(In[].sum:2*In[].sum-1) = 0.1;
107 } else
108 FEFToSSP(ou1[], ou2[], start);
109
110 real mini = cmaes(J, start, stopMaxFunEval=10000*(al+1), stopTolX=1.e-3/(10*(al+1)), initialStdDev=(0.025/(pow(100.,al))));
111 Vh best1, best2;
112 SSPToFEF(best1[], best2[], start);
113
114 // Mesh adaptation
115 Th = adaptmesh(Th, best1, best2);
116 ou1 = best1;
117 ou2 = best2;
118 }
119
6.5.3 IPOPT minimal surface & volume
1 load "msh3";
2 load "medit";
3 load "ff-Ipopt";
4
5 // Parameters
6 int nadapt = 3;
7 real alpha = 0.9;
8 int np = 30;
9 real regtest;
10 int shapeswitch = 1;
11 real sigma = 2*pi/40.;
real treshold = 0.1;
real e = 0.1;
real r0 = 0.25;
real rr = 2-r0;
real E = 1./(e*e);
real RR = 1./(rr*rr);

// Mesh
mesh Th = square(2*np, np, [2*pi*x, pi*y]);

// Fespace
fespace Vh(Th, P1, periodic=[[2, y], [4, y]]);

// Initial shape definition
//outside of the mesh adaptation loop to initialize with the previous optimal shape found on further iterations
Vh startshape = 5;
Vh uz = 1., lz = 1.;

// Mesh adaptation loop
real[int] lm = [1];
for(int kkk = 0; kkk < nadapt; ++kkk){
  int iter=0;
  func sin2 = square(sin(y));

  // A function which transform Th in 3d mesh (r=rho)
  //a point (theta,phi) of Th becomes ( r(theta,phi)*cos(theta)*sin(phi) , r(theta,phi)*sin(theta)*sin(phi) , r(theta,phi)*cos(phi) )
  //then displays the resulting mesh with medit
  func int Plot3D (real[int] &rho, string cmm, bool ffplot){
    Vh rhoo;
    rhoo[] = rho;
    //mesh sth = square(np, np/2, [2*pi*x, pi*y]);
    //fespace svh(sth, P1);
//Vh rhoplot = rho0;

try{
  mesh3 Sphere = movemesh23(Th, transfo=[rho0(x,y)*cos(x)*sin(y), rho0(x,y)*sin(x)*sin(y), rho0(x,y)*cos(y)]);
  if(ffplot)
    plot(Sphere);
  else
    medit(cmm, Sphere);
  catch(...){
    cout << "PLOT ERROR" << endl;
  }
  return 1;
}

// Surface computation
//Maybe is it possible to use movemesh23 to have the surface function less complicated
//However, it would not simplify the gradient and the hessian
func real Area (real[int] &X){
  Vh rho;
  rho[] = X;
  Vh rho2 = square(rho);
  Vh rho4 = square(rho2);
  real res = int2d(Th)(sqrt(rho4*sin2 + rho2*square(dx(rho)) + rho2*sin2*square(dy(rho))));
  ++iter;
  if(1)
    plot(rho, value=true, fill=true, cmm="rho(theta,phi) on [0,2pi]x[0,pi] - S=", dim=3);
  else
    Plot3D(rho[], "shape_evolution", 1);
  return res;
}

func real[int] GradArea (real[int] &X){
  Vh rho, rho2;
  rho[] = X;
  rho2[] = square(X);
  Vh sqrtPsi, alpha;
  {
    Vh dxrho2 = dx(rho)*dx(rho), dyrho2 = dy(rho)*dy(rho);
    sqrtPsi = sqrt(rho2*rho2*sin2 + rho2*dxrho2 + rho2*dyrho2*sin2);
    alpha = 2.*rho2*rho*sin2 + rho*dxrho2 + rho*dyrho2*sin2;
  }
  varf dArea (u, v)
  = int2d(Th){
    1./sqrtPsi * (alpha*v + rho2*dx(rho)*dx(v) + rho2*dy(rho)*sin2*dy(v))
  };

  real[int] grad = dArea(0, Vh);
  return grad;
}

matrix hessianA;
func matrix HessianArea (real[int] &X){
  (continues on next page)
Vh rho, rho2;
rho[] = X;
rho2 = square(rho);
Vh sqrtPsi, sqrtPsi3, C00, C01, C02, C11, C12, C22, A;
{
    Vh C0, C1, C2;
    Vh dxrho2 = dx(rho) * dx(rho), dyrho2 = dy(rho) * dy(rho);
    sqrtPsi = sqrt(rho2 * rho2 + rho2 * dxrho2 + rho2 * dyrho2 + sin2);
    sqrtPsi3 = (rho2 * rho2 + rho2 * dxrho2 + rho2 * dyrho2 + sin2) * sqrtPsi;
    C0 = 2 * rho2 + rho * dxrho2 + rho * dyrho2 + sin2;
    C1 = rho2 * dx(rho);
    C2 = rho2 + sin2 * dy(rho);
    C00 = square(C0);
    C01 = C0 * C1;
    C02 = C0 * C2;
    C11 = square(C1);
    C12 = C1 * C2;
    C22 = square(C2);
    A = 6. * rho2 + sin2 + dxrho2 + dyrho2;
}

```
varf d2Area(w, v) = int2d(Th) {
    1./sqrtPsi * (A*w*v
    + 2*rho*dx(rho)*dx(w)*v
    + 2*rho*dx(rho)*w*dx(v)
    + 2*rho*dy(rho)*sin2*dx(v)*v
    + 2*rho*dy(rho)*sin2*w*dy(v)
    + rho2*dx(w)*dx(v)
    + rho2*sin2*dx(w)*dy(v)
    + rho2*sin2*dx(w)*dy(v)
    )
    + 1./sqrtPsi3 * (C00*w*v
    + C01*dx(w)*v
    + C01*w*dx(v)
    + C02*dy(w)*v
    + C02*w*dy(v)
    + C11*dx(w)*dx(v)
    + C12*dx(w)*dy(v)
    + C12*dy(w)*dx(v)
    + C22*dy(w)*dy(v)
    )
}
```

```
hessianA = d2Area(Vh, Vh);
return hessianA;
```

// Volume computation
```
func real Volume (real[int] &X) {
    Vh rho;
rho[] = X;
rho3 = rho*rho*rho;
real res = 1./3.*int2d(Th)(rho3*sin(y));
return res;
```

(continues on next page)
func real[|int|] GradVolume (real[|int|] &X)
    Vh rho;
    rho[] = X;
    varf dVolume(u, v) = int2d(Th)(rho*rho*sin(y)*v);
    real[|int|] grad = dVolume(0, Vh);
    return grad;
}

matrix hessianV;
func matrix HessianVolume(real[|int|] &X)
    Vh rho;
    rho[] = X;
    varf d2Volume(w, v) = int2d(Th)(2*rho*sin(y)*v*w);
    hessianV = d2Volume(Vh, Vh);
    return hessianV;
}

//if we want to use the volume as a constraint function
//we must wrap it in some freefem functions returning the appropriate type
//The lagrangian hessian also have to be wrapped since the Volume is not linear.

func real[|int|] ipVolume (real[|int|] &X)
    real[|int|] vol = Volume(X);
    return vol;
}

matrix mdV;
func matrix ipGradVolume (real[|int|] &X)
    real[|int|,|int|] dvol(1,Vh.ndof);
    dvol(0, -1) = GradVolume(X);
    mdV = dvol;
    return mdV;
}

matrix HLagrangian;
func matrix ipHessianLag (real[|int|] &X, real objfact, real[|int|] &lambda)
    HLagrangian = objfact*HessianArea(X) + lambda[0]*HessianVolume(X);
    return HLagrangian;
}

//building struct for GradVolume
int[|int|] gvi(Vh.ndof), gvj=0:Vh.ndof-1;
    gvi = 0;

Vh rc = startshape; //the starting value
Vh ub = 1.e19; //bounds definition
Vh lb = 0;

func real Gaussian (real X, real Y, real theta, real phi)
    real deltax2 = square((X-theta)*sin(Y)), deltay2 = square(Y-phi);
    return exp(-0.5 * (deltax2 + deltay2) / (sigma*sigma));
}

func disc1 = sqrt(1./(RR+(E-RR)*cos(y)*cos(y)))*(1+0.1*cos(7*x));
func disc2 = sqrt(1./(RR+(E-RR)*cos(x)*cos(x)*sin2));

if(1)
    lb = r0;
    for (int q = 0; q < 5; ++q)
        func f = rr*Gaussian(x, y, 2*q*pi/5., pi/3.);
        func g = rr*Gaussian(x, y, 2*q*pi/5.+pi/5., 2.*pi/3.);
        lb = max(max(lb, f), g);
    lb = max(lb, rr*Gaussian(x, y, 2*pi, pi/3));
    lb = max(lb, max(disc1, disc2));
real Vobj = Volume(lb[]);
real Vnvc = 4./3.*pi*pow(lb[].linfty,3);

if(1)
    Plot3D(lb[], "object_inside", 1);
real[int] clb = 0., cub = [(1-alpha)*Vobj + alpha*Vnvc];

// Call IPOPT
int res = IPOPT(Area, GradArea, ipHessianLag, ipVolume, ipGradVolume,
    rc[], ub=ub[], lb=lb[], clb=clb, cub=cub, checkindex=1, maxiter=kkk
    <nadapt-1 ? 40:150,
    warmstart=k, lm=lm, uz=uz[], lz=lz[], tol=0.00001, structjacc=[gvi,
    gvj]);
    cout << "IPOPT: res =" << res << endl;

// Plot
Plot3D(rc[], "Shape_at_"+kkk, 1);
Plot3D(GradArea(rc[]), "ShapeGradient", 1);

// Mesh adaptation
if (kkk < nadapt-1){
    Th = adaptmesh(Th, rc*cos(x)*sin(y), rc*sin(x)*sin(y), rc*cos(y),
        nbvx=50000, periodic=[[2, y], [4, y]]);
    plot(Th, wait=true);
    startshape = rc;
    uz = uz;
    lz = lz;
}
regtest = rc[]'*rc[];

Fig. 6.23: Mesh
6.5.4 CMAES MPI variational inequality

Command:

```bash
ff-mpirun -np 4 CMAESMPIVariationalInequality.edp -glut ffglut
```

```plaintext
load "mpi-cmaes"

// Parameters
int NN = 10;
func f1 = 1.;
func f2 = -1.;
func g1 = 0.;
func g2 = 0.1;
int iter = 0;
int nadapt = 1;
real starttol = 1e-10;
real bctol = 6.e-12;
real pena = 1000;

// Mesh
mesh Th = square(NN, NN);

// Fespace
fespace Vh(Th, P1);
Vh ou1, ou2;

// Mesh adaptation loop
for (int al = 0; al < nadapt; ++al){
    // Problem
    varf BVF (v, w) = int2d(Th)(
        0.5*dx(v)*dx(w) + 0.5*dy(v)*dy(w)
    );
    varf LVF1 (v, w) = int2d(Th)(f1*w);
    varf LVF2 (v, w) = int2d(Th)(f2*w);
    matrix A = BVF(Vh, Vh);
    real[int] b1 = LVF1(0, Vh);
    real[int] b2 = LVF2(0, Vh);
    varf Vbord (v, w) = on(1, 2, 3, 4, v=1);
    Vh In, Bord;
    Bord[] = Vbord(0, Vh, tgv=1);
    In[] = Bord[] ? 0:1;
    Vh gh1 = Bord*g1, gh2 = Bord*g2;

    // Function which create a vector of the search space type from
    // two finite element functions
    func int FEFToSSP (real[int] &fef1, real[int] &fef2, real[int] &ssp)
    {
        int kX = 0;
        for (int i = 0; i < Vh.ndof; ++i){
            if (In[i]){
                ssp[kX] = fef1[i];
                ssp[kX+In[i].sum] = fef2[i];
            }
        }
    }
```

(continues on next page)
++kX;
    }
  }
  return 1;
}

// Function splitting a vector from the search space and fills two finite element functions with it
func int SSPToFEF(real[int] &fef1, real[int] &fef2, real[int] &ssp){
  int kX = 0;
  for (int i = 0; i < Vh.ndof; ++i){
    if (In[i]){  
      fef1[i] = ssp[kX];
      fef2[i] = ssp[kX+In[].sum];
      ++kX;
    }
    else{
      fef1[i] = gh1[][i];
      fef2[i] = gh2[][i];
    }
  }
  return 1;
}

func real IneqC(real[int] &X){
  real[int] constraints(In[].sum);
  for (int i = 0; i < In[].sum; ++i){
    constraints[i] = X[i] - X[i+In[].sum];
    constraints[i] = constraints[i] <= 0 ? 0. : constraints[i];
  }
  return constraints.\text{L}_2;
}

func real J(real[int] &X){
  Vh u1, u2;
  SSPToFEF(u1[], u2[], X);
  iter++;
  real[int] Au1 = A*u1[], Au2 = A*u2[];
  Au1 -= b1;
  Au2 -= b2;
  real val = u1[]'*Au1 + u2[]'*Au2;
  val += pena * IneqC(X);
  plot(u1, u2, nbiso=30, fill=1, dim=3, cmm="adapt level "+al++" - iteration "+iter++" - J = "+val, value=1);
  return val ;
}

// Solve
real[int] start(2*In[].sum);

if (al==0){
  start(0:In[].sum-1) = 0.;
  start(In[].sum:2*In[].sum-1) = 0.1;
}
else
  FEFToSSP(ou1[], ou2[], start);

(continues on next page)
real mini = cmaesMPI(J, start, stopMaxFunEval=10000*(a1+1), stopTo1X=1.e-4/10*(a1+1)), initialStdDev=(0.025/pow(100.,(a1+1)));

Vh best1, best2;
SSPToFEF(best1[], best2[], start);

// Mesh adaptation
Th = adaptmesh(Th, best1, best2);
ou1 = best1;
ou2 = best2;

Fig. 6.24: Result

6.6 Parallelization

6.6.1 MPI-GMRES 2D

To launch this script, use for example:

ff-mpirun -np 12 MPIGMRES2D.edp -d 1 -k 1 -gmres 2 -n 50

//usage :
//arguments:
//-glut : to see graphically the process
//-n N: set the mesh cube split NxNxN
//-d D: set debug flag D must be one for mpiplot
//-k K: to refined by K all element
//-ns: remove script dump
//-gmres
//0: use iterative schwarz algo.

(continues on next page)
11 //1: Algo GMRES on residu of schwarz algo
12 //2: DDM GMRES
13 //3: DDM GMRES with coarse grid preconditionner (Good one)
14
15 load "MPICG"
16 load "medit"
17 load "metis"
18 include "getARGV.idp"
19 include "MPIplot.idp"
20 include "MPIGMRESPRmacro.idp"
21
22 searchMethod = 0; //more safe search algo (warning can be very expensive in case of ˓→lot of outside point)
23 assert(version >= 3.11); //need at least v3.11
24 real[int] ttt(10);
25 int ittt=0;
26 macro settt {ttt[ittt++] = mpiWtime();} //
27
28 // Arguments
29 verbosity = getARGV("-vv", 0);
30 int vdebug = getARGV("-d", 1);
31 int ksplit = getARGV("-k", 3);
32 int nloc = getARGV("-n", 10);
33 string sff = getARGV("-p", "");
34 int gmres = getARGV("-gmres", 2);
35 bool dplot = getARGV("-dp", 0);
36 int nC = getARGV("-N", max(nloc/10, 4));
37
38 if (mpirank==0 && verbosity){
39 cout << "ARGV: ";
40 for (int i = 0; i < ARGV.n; ++i)
41 cout << ARGV[i] << " ";
42 cout << endl;
43 }
44
45 if(mpirank==0 && verbosity)
46 cout << " vdebug: " << vdebug << ", ksplit " << ksplit << ", nloc " << nloc << " , sff " << ˓→< sff << "." << endl;
47
48 // Parameters
49 int withplot = 0;
50 bool withmetis = 1;
51 bool RAS = 1;
52 string sPk = "P2-2gd";
53 func Pk = P2;
54 int sizeoverlaps = 1; //size of overlap
55 int[int] l111 = [1, 1, 1, 1]; //mesh labels
56
57 // MPI function
58 func bool plotMPIall(mesh &Th, real[int] &u, string cm){
59 if(vdebug)
60 PLOTMPIALL(mesh, Pk, Th, u, {cm=cm, nbiso=20, fill=1, dim=3, value=1});
61 return 1;
62 }
63
64 // MPI
65 mpiComm comm(mpiCommWorld,0,0); //trick : make a no split mpiWorld

(continues on next page)
int npart = mpiSize(comm); // total number of partition
int ipart = mpiRank(comm); // current partition number

int nnpj = 0; // Number of partition with intersection (a jpart) with ipart without ipart
int [int] jpart(npart); // list of jpart

if (ipart==0)
    cout << " Final N = " << ksplit*nloc << ", nloc = " << nloc << " split = " << ksplit << endl;

// Mesh
mesh Thg = square(nloc, nloc, label=1111);
mesh ThC = square(nC, nC, label=1111); // Coarse mesh

mesh Thi, Thin; // with overlap, without overlap

// Fespace
fespace Phg(Thg, P0);
Phg part;

fespace Vhg(Thg, P1);
Vhg unssd; // boolean function: 1 in the subdomain, 0 elsewhere

fespace VhC(ThC, P1); // of the coarse problem

// Partitioning
{
    int [int] nupart(Thg.nt);
    nupart = 0;
    if (npart > 1 && ipart == 0)
        metisdual(nupart, Thg, npart);

    broadcast(processor(0, comm), nupart);
    for (int i = 0; i < nupart.n; ++i)
        part[][i] = nupart[i];
}

if (withplot > 1)
    plot(part, fill=1, cmm="dual", wait=1);

// Overlapping partition
Phg suppi = abs(part-ipart) < 0.1;

Thin = trunc(Thg, suppi>0, label=10); // non-overlapping mesh, interfaces have label 10

int nnn = sizeoverlaps*2; // to be sure
AddLayers(Thg, suppi[], nnn, unssd[]); // see above! suppi and unssd are modified
unssd[] *= nnn; // to put value nnn a 0
real nnn0 = nnn - sizeoverlaps*0.001;

Thi = trunc(Thg, unssd>nnn0, label=10); // overlapping mesh, interfaces have label 10

settt

// Fespace
fespace Vhi(Thi, P1);
int npij = npart;
Vhi[int] pij(npij); // local partition of unit + pii
Vhi pii;

real nnn1 = +0.001;
{
    /*
    construction of the partition of the unit,
    let phi_i P1 FE function 1 on Thin and zero outside of Thi and positive
    the partition is build with
    p_i = phi_i / \sum phi_i
    */
    // build a local mesh of thii such that all computation of the unit partition are
    // exact in thii
    mesh Thii = trunc(Thg, unssd>nnn1, label=10); // overlapping mesh, interfaces have
    // label 10

    {
        // find all j mes (supp(p_j) \cap supp(p_i)) >0
        // compute all phi_j on Thii
        // remark: supp p_i include in Thi

        // Fespace
        fespace Phi(Thii, P0);
        fespace Vhii(Thii, P1);
        Vhi sumphi = 0;
        Vhii phii = 0;

        jpart = 0;
        njpart = 0;
        int nlayer = RAS ? 1 : sizeoverlaps;
        if (ipart == 0)
            cout << "nlayer = " << nlayer << endl;
        pii = max(unssd-nnn1+nlayer, 0.)/nlayer;
        if(dplot)
            plot(pii, wait=1, cmm=" 0000");
        sumphi[] += pii[];
        if(dplot)
            plot(sumphi, wait=1, cmm=" summ 0000");

        real epsmes = 1e-10*Thii.area;
        for (int i = 0; i < npart; ++i)
            if (i != ipart) {
                Phii suppii = abs(i-part) < 0.2;
                if (suppii[].max > 0.5){
                    AddLayers(Thii, suppii[], nlayer, phii[]);
                    assert(phii[].min >= 0);
                    real interij = int2d(Thi)(phii);
                    if (interij > epsmes){
                        pij[njpart] = abs(phii);
                        if(vdebug > 2)
                            cout << " ***** " << int2d(Thi)(real(pij[njpart])<0) << " " <<
                            <pij[njpart][].min << " " << phii[].min << endl;
                    }
                    assert(int2d(Thi)(real(pij[njpart]) < 0) == 0);
                }
            }
    }

(continues on next page)
```cpp
if (dplot)
    plot(pij[njpart], wait=1, cmm=" j = " + i + " " + njpart);
sumphi[] += pij[njpart][];
if (dplot)
    plot(sumphi, wait=1, cmm=" sum j = " + i + " " + njpart);
jpart[njpart++] = i;
}
}
if (dplot)
    plot(sumphi, wait=1, dim=3, cmm="sum " fill=1);
pii[] = pii[] ./ sumphi[];
for (int j = 0; j < njpart; ++j)
    pij[j][] = pij[j][] ./ sumphi[];
jpart.resize(njpart);
for (int j = 0; j < njpart; ++j)
    assert(pij[j][].max <= 1);
{
    cout << ipart << " number of jpart " << njpart << " : ";
    for (int j = 0; j < njpart; ++j)
        cout << jpart[j] << " ";
    cout << endl;
}
sumphi[] = pii[];
for (int j = 0; j < njpart; ++j)
    sumphi[] += pij[j][];
if (vdebug > 2)
    cout << "sum min " << sumphi[].min << " " << sumphi[].max << "endl;"
    assert(sumphi[].min > 1.-1e-6 && sumphi[].max < 1.+1e-6);
}
//Thii is remove here
// end of the construction of the local partition of the unity ...
// on Thi
if (ipart == 0)
    cout << "End build partition" << endl;
// Computation of number of intersection
//here pii and the pij is the local partition of the unit on
//Thi (mesh with overlap)
if (dplot){
    plot(Thi, wait=1);
    for (int j = 0; j < njpart; ++j)
        plot(pij[j][], cmm=" j=" + j, wait=1);
}
//Partition of the unity on Thi
//computation of message
//data on intersection of the support of pij[0] and pij[j]
settt
if (vdebug)
    plotMPIall(Thi, pii[], "pi_i");
mesh[int] aThij(njpart);
matrix Pii;
```
matrix[int] sMj(njpart);  //M of send to j
matrix[int] rMj(njpart);  //M to recv from j
fespace Whi(Thi, Pk);
mesh Thij = Thi;
fespace Whij(Thij, Pk);

//construction of the mesh intersect i,j part
for(int jp = 0; jp < njpart; ++jp)
  aThij[jp] = trunc(Thi, pij[jp] > 1e-6, label=10);  //mesh of the supp of pij
for(int jp = 0; jp < njpart; ++jp)
  aThij[jp] = trunc(aThij[jp], 1, split=ksplit);
Thi = trunc(Thi, 1, split=ksplit);

if (ipart == 0)
  cout << "End build mesh intersection" << endl;

// Construction of transfert matrix
{
  Whi wpii = pii;
Pii = wpii[];
  for(int jp = 0; jp < njpart; ++jp){
    int j = jpart[jp];
    Thij = aThij[jp];
    matrix I = interpolate(Whij, Whi);  //Whji <- Whi
    sMj[jp] = I*Pii;  //Whi -> s Whij
    rMj[jp] = interpolate(Whij, Whi, t=1);  //Whji -> Whi
    if(vdebug > 10){
      {Whi uuu=1; Whij vvv=-1; vvv[]+=I*uuu[]; cout << jp << " %%% " << vvv[].linfty<< endl; assert(vvv[].linfty < 1e-6);}
      {Whi uuu=1; Whij vvv=-1; vvv[]+=rMj[jp]*uuu[]; cout << jp << " ### " << _vvv[].linfty << endl; assert(vvv[].linfty < 1e-6);}
    }
  }
}
if (ipart == 0)
  cout << "End build transfert matrix" << endl;

// Allocate array of send and recv data
InitU(njpart, Whij, Thij, aThij, Usend)  //initU(n, Vh, Th, aTh, U)
InitU(njpart, Whij, Thij, aThij, Vrecv)
if (ipart == 0)
  cout << "End init data for send/revc" << endl;

Whi ui, vi;
func bool Update(real[int] &ui, real[int] &vi){
  for(int j = 0; j < njpart; ++j)
    SendRecvUV(comm, jpart, Usend, Vrecv)
    vi = Pii*ui;
  for(int j = 0; j < njpart; ++j)
    vi += rMj[j]*Vrecv[j][];
  return true;
// Definition of the Problem
func G = x*0.1;
func F = 1.;
macro grad(u) [dx(u),dy(u)] //
varf vBC (U, V) = on(1, U=G);
varf vPb (U, V) = int2d(Thi)(grad(U)'*grad(V)) + int2d(Thi)(F*V) + on(10, U=0) + on(1, U=G);
varf vPbC (U, V) = int2d(ThC)(grad(U)'*grad(V)) + on(1, U=0);
varf vPbon (U, V) = on(10, U=1) + on(1, U=0);
varf vPbon10only (U, V) = on(10, U=1) + on(1, U=0);
//remark the order is important we want 0 part on 10 and 1

matrix Ai = vPb(Whi, Whi, solver=sparsesolver);
matrix AC, Rci, Pci;
if (mpiRank(comm) == 0)
    AC = vPbC(VhC, VhC, solver=sparsesolver);
Pci = interpolate(Whi, VhC);
Rci = Pci'*Pii;
real[int] onG10 = vPbon10only(0, Whi);
real[int] onG = vPbon(0, Whi);
real[int] Bi=vPb(0, Whi);

int kiter = -1;

func bool CoarseSolve(real[int] &V, real[int] &U, mpiComm &comm){
    //solving the coarse probleme
    real[int] Uc(Rci.n), Bc(Uc.n);
    Uc = Rci*U;
    mpiReduce(Uc, Bc, processor(0, comm), mpiSUM);
    if (mpiRank(comm) == 0)
        Uc = AC^-1*Bc;
    broadcast(processor(0, comm), Uc);
    V = Pci*Uc;
}

func real[int] DJ (real[int] &U){
    ++kiter;
    real[int] V(U.n);
    V = Ai*U;
    V = onG10 ? 0.: V; //remove internal boundary
    return V;
}

func real[int] PDJ (real[int] &U){
    real[int] V(U.n);
    real[int] b = onG10 ? 0. : U;
    V = Ai^-1*b;
    Update(V, U);
    return U;
}
```cpp
func real[int] PDJC (real[int] &U) {
    real[int] V(U.n);
    CoarseSolve(V, U, comm);
    V = -V; // -C2 * U0
    U += Ai*V; // U = (I - A C2) U0
    real[int] b = onG10 ? 0. : U;
    U = Ai^-1*b; // (C1 (I - A C2)) U0
    V = U - V;
    Update(V, U);
    return U;
}

func real[int] DJ0(real[int] &U) {
    ++kiter;
    real[int] V(U.n);
    real[int] b = onG .* U;
    b = onG ? b : Bi ;
    V = Ai^-1*b;
    Update(V, U);
    V -= U;
    return V;
}

while u = 0, v;
{
    // verification
    while u = 1, v;
    Update(u[], v[]);
    u[] -= v[];
    assert(U[].linfty < 1e-6);
}

settt
u[] = vBC(0, Whi, tgv=1); // set u with tgv BC value
real epss = 1e-6;
int rgmres = 0;
if (gmres == 1) {
    rgmres = MPIAffineGMRES(DJ0, u[], veps=epss, nbiter=300, comm=comm, dimKrylov=100,
    verbosity=ipart ? 0 : 50);
    real[int] b = onG .* U[];
    b = onG ? b : Bi;
    v[] = Ai^-1*b;
    Update(v[], u[]);
}
else if (gmres == 2) {
    rgmres = MPILinearGMRES(DJ, precon=PDJ, u[], Bi, veps=epss, nbiter=300, comm=comm,
    dimKrylov=100, verbosity=ipart ? 0 : 50);
} else if (gmres == 3) {
    rgmres = MPILinearGMRES(DJ, precon=PDJC, u[], Bi, veps=epss, nbiter=300, 
    comm=comm, dimKrylov=100, verbosity=ipart ? 0 : 50);
} else // algo Schwarz for demo
    for(int iter = 0; iter < 10; ++iter){
        real[int] b = onG .* U[];
        b = onG ? b : Bi;
        v[] = Ai^-1*b;
        Update(v[], u[]);
    }
(continues on next page)
```
if (vdebug)
   plotMPIall(Thi, u[], "u-"+iter);

v[] -= u[];

real err = v[].linfty;
real umax = u[].max;
real[int] aa = [err, umax], bb(2);
mpiAllReduce(aa, bb, comm, mpiMAX);
real errg = bb[0];
real umaxg = bb[1];

if (ipart == 0)
   cout << ipart << " err = " << errg << " u. max " << umaxg << endl;
   if (errg < 1e-5) break;

if (vdebug)
   plotMPIall(Thi, u[], "u-final");

settt

real errg = 1, umaxg;
{
   real umax = u[].max, umaxg;
   real[int] aa = [umax], bb(1);
   mpiAllReduce(aa, bb, comm, mpiMAX);
   errg = bb[0];
   if (ipart == 0)
      cout << "umax global = " << bb[0] << " Wtime = " << (ttt[ittt-1]-ttt[ittt-2])
          << " s " " " << kiter << endl;
}

if (sff != ""){
   ofstream ff(sff+.txt", append);
   cout << "++++ ";
   cout << mpirank << "/" << mpisize << " k=" << ksplit << " n= " << nloc << " " <<
       sizeoverlaps << " it= " << kiter;
   for (int i = 1; i < ittt; ++i)
      cout << " " << ttt[i]-ttt[i-1] << " ";
   cout << epss << " " << Ai.nbcоф << " " << Ai.n << endl;
/*
1 mpirank
2 mpisize
3 ksplit
4 nloc
5 sizeoverlaps
6 kiter
7 mesh & part build
8 build the partion
9 build mesh, transfer, and the fine mesh ..
10 build the matrix, the trans matrix, factorization
11 GMRES
*/
   ff << mpirank << " " << mpisize << " " << sPk << " ";
   ff << ksplit << " " << nloc << " " << sizeoverlaps << " " << kiter;

(continues on next page)
for (int i = 1; i < ittt; ++i)
    ff << " " << ttt[i]-ttt[i-1] << " ";
ff << epss << " " << Ai.nbcoef << " " << Ai.n << " " << gmres << endl;
}

Fig. 6.25: Results

6.6.2 MPI-GMRES 3D

Todo: todo

6.6.3 Direct solvers

load "MUMPS_FreeFem"
//default solver: real-> MUMPS, complex -> MUMPS
load "real_SuperLU_DIST_FreeFem"
default solver: real -> SuperLU_DIST, complex -> MUMPS
load "real_pastix_FreeFem"
//default solver: real-> pastix, complex -> MUMPS

// Solving with pastix
{
    matrix A =
        [[1, 2, 2, 1, 1],
         [ 2, 12, 0, 10 , 10],
         [ 2, 0, 1, 0, 2],
         [ 1, 10, 0, 22, 0 ],
         [ 1, 10, 2, 0., 22]];

    real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
    b = A*xx;
    cout << "b = " " << b << endl;
    cout << "xx = " " << xx << endl;
```
set(A, solver=sparssolver, datafilename="ffpastix_iparm_dparm.txt");
cout << "solve" << endl;
x = A^{-1}*b;
cout << "b = " << b << endl;
cout << "x = " << endl;
cout << x << endl;
di = xx - x;
if (mpirank == 0){
cout << "x-xx = " << endl;
cout << "Linf = " << di.linfy << ", L2 = " << di.l2 << endl;
}

// Solving with SuperLU_DIST
realdefaulttoSuperLudist();
//default solver: real-> SuperLU_DIST, complex -> MUMPS
{
   matrix A =
   [[1, 2, 2, 1, 1],
    [ 2, 12, 0, 10 , 10],
    [ 2, 0, 1, 0, 2],
    [ 1, 10, 0, 22, 0.],
    [ 1, 10, 2, 0., 22]];
   
   real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
   b = A*xx;
cout << "b = " << b << endl;
cout << "xx = " << xx << endl;
set(A, solver=sparssolver, datafilename="ffsuperlu_dist_fileparam.txt");
cout << "solve" << endl;
x = A^{-1}*b;
cout << "b = " << b << endl;
cout << "x = " << endl;
cout << x << endl;
di = xx - x;
if (mpirank == 0){
cout << "x-xx = " << endl;
cout << "Linf = " << di.linfy << ", L2 = " << di.l2 << endl;
}

// Solving with MUMPS
defaulttoMUMPS();
//default solver: real-> MUMPS, complex -> MUMPS
{
   matrix A =
   [[1, 2, 2, 1, 1],
    [ 2, 12, 0, 10 , 10],
    [ 2, 0, 1, 0, 2],
    [ 1, 10, 0, 22, 0.],
    [ 1, 10, 2, 0., 22]];
   
   real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
   b = A*xx;
cout << "b = " << b << endl;
cout << "xx = " << xx << endl;
```
set(A, solver=sparsetools, datafilename="ffmumps_fileparam.txt");
cout << "solving solution" << endl;
x = A^-1 * b;
cout << "b = " << b << endl;
cout << "x = " << endl;
cout << x << endl;
di = xx - x;
if (mpirank == 0) {
    cout << "x-xx = " << endl;
    cout << "Linf = " << di.linfy << ", L2 = " << di.l2 << endl;
}
}

6.6.4 Solver MUMPS

load "MUMPS_FreeFem"

// Parameters
int[int] ICNTL(40); //declaration of ICNTL parameter for MUMPS

//get value of ICNTL from file
if (mpirank == 0) {
    ifstream ff("ffmumps_fileparam.txt");
    string line;
    getline(ff, line);
    getline(ff, line);
    for (int iii = 0; iii < 40; iii++){
        ff >> ICNTL[iii];
        getline(ff, line);
    }
}
broadcast(processor(0), ICNTL);

// Given data of MUMPS solver in array iparams(SYM, PAR, ICNTL)
// There is no symmetric storage for a matrix associated with a sparse solver.
// Therefore, the matrix will be considered unsymmetric for parallel sparse solver.
{
    // Problem
    int SYM = 0;
    int PAR = 1;
    matrix A = [
        [40, 0, 45, 0, 0],
        [0, 12, 0, 0, 0],
        [0, 0, 40, 0, 0],
        [12, 0, 0, 22, 0],
        [0, 0, 20, 0, 22]
    ];

    // Construction of integer parameter for MUMPS
    int[int] MumpsLParams(42);
    MumpsLParams[0] = SYM;
}
MumpsLParams[1] = PAR;
for (int ii = 0; ii < 40; ii++)
    MumpsLParams[ii+2] = ICNTL[ii]; //ICNTL begin with index 0 here
real[int] xx = [1, 32, 45, 7, 2], x(5), b(5), di(5);
b = A*xx;
if (mpirank == 0)
    cout << "xx = " << xx << endl;

set(A, solver=sparse solver, lparams=MumpsLParams); //we take the default value  for CNTL MUMPS parameter

// Solve
if (mpirank == 0)
    cout << "Solve" << endl;
x = A^-1*b;
if (mpirank == 0)
    cout << "b = " << b << endl;
if (mpirank == 0)
    cout << "x = " << x << endl; cout << x << endl;
di = xx-x;
if (mpirank == 0){
    cout << "x-xx = " << endl;
    cout << "Linf = " << di.linfty <<", L2 = " << di.l2 << endl;
}

// Read parameter of MUMPS solver in file ffmumps_fileparam.txt
{
    // Problem
matrix A =
    [ [40, 0, 45, 0, 0],
      [0, 12, 0, 0, 0],
      [0, 0, 40, 0, 0],
      [12, 0, 0, 22, 0],
      [0, 0, 20, 0, 22] ];
real[int] xx = [1, 32, 45, 7000, 2], x(5), b(5), di(5);
b = A*xx;
if (mpirank == 0){
    cout << "b = " << b << endl;
    cout << "xx = " << xx << endl;
}

set(A, solver=sparse solver, datafilename="ffmumps_fileparam.txt");

// Solve
if (mpirank == 0)
    cout << "Solve" << endl;
x = A^-1*b;
if (mpirank == 0){
    cout << "b = " << b << endl;
    cout << "x = " << x << endl;
}
6.6.5 Solver superLU_DIST

_todo: write code (SuperLU_DIST seems to have a bug)

6.6.6 Solver PaStiX

_todo: write code (PaStiX seems to have a bug)

6.7 Developers

6.7.1 FFT

_load "dfft"

// Parameters
int nx = 32;
real ny = 16;
real N = nx*ny;
func f1 = cos(2*x*2*pi)*cos(3*y*2*pi);

// Mesh
//warning: the fourier space is not exactly the unit square due to periodic condition
mesh Th = square(nx-1, ny-1, [(nx-1)*x/nx, (ny-1)*y/ny]);
//warning: the numbering of the vertices (x,y) is
//given by i = x/nx + nx*y/ny

// Fespace
fespace Vh(Th, P1);
Vh<complex> u = f1, v;
Vh w = f1;
Vh ur, ui;

// FFT
//in fft the matrix n, m is in row-major order and array n, m is
//store j + m*i (the transpose of the square numbering)
v[] = dfft(u[], ny, -1);
u[] = dfft(v[], ny, +1);
    cout << "||u||\_linfty " << u[].linfty << endl;

(continues on next page)
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(continued from previous page)

28 \texttt{u[]} *= 1./\texttt{N};
29 \texttt{cout} << "||u||_{\infty} " << u[].\texttt{linfty} << endl;
30 \texttt{ur} = \texttt{real(u)};

33 \texttt{// Plot}
34 \texttt{plot(w, wait=1, value=1, cmm="w");}
35 \texttt{plot(ur, wait=1, value=1, cmm="u");}
36 \texttt{v} = w - u;
37 \texttt{cout} << "\texttt{diff} = " << v[].\texttt{max} << " " << v[].\texttt{min} << endl;
38 \texttt{assert( norm(v[].\texttt{max}) < 1e-10 && norm(v[].\texttt{min}) < 1e-10);}

39 \texttt{// Other example}
40 \texttt{//FFT Lapacian}
41 \texttt{//-\Delta u = f with biperiodic condition}
42 \texttt{func f = cos(3*2*\texttt{pi}*x)\times\cos(2*2*\texttt{pi}*y);}
43 \texttt{func ue = (1./(square(2*\texttt{pi})*13.))\times\cos(3*2*\texttt{pi}*x)\times\cos(2*2*\texttt{pi}*y); //the exact solution}
44 \texttt{Vh<complex> ff = f;}
45 \texttt{Vh<complex> fhat;}
46 \texttt{Vh<complex> wij;}

49 \texttt{// FFT}
50 \texttt{fhat[] = dfft(ff[], ny, -1);}
51 \texttt{//warning in fact we take mode between -nx/2, nx/2 and -ny/2, ny/2}
52 \texttt{//thanks to the operator ?}
53 \texttt{wij = square(2.*\texttt{pi})\times(square((x<0.5?x\times nx: (x-1)*nx)) + square((y<0.5?y\times ny: (y-1)*ny)));}
54 \texttt{wij[0] = 1e-5; //to remove div / 0}
55 \texttt{fhat[] = fhat[] ./ wij[];}
56 \texttt{u[] = dfft(fhat[], ny, 1);}
57 \texttt{u[] }=\texttt{ complex(N);}
58 \texttt{ur = real(u); //the solution}
59 \texttt{w = real(ue); //the exact solution}

62 \texttt{// Plot}
63 \texttt{plot(w, ur, value=1, cmm="ue", wait=1);}
66 \texttt{// Error}
67 \texttt{w[] }=\texttt{ ur[];}
68 \texttt{real err = abs(w[].\texttt{max}) + abs(w[].\texttt{min});}
69 \texttt{cout} << "err = " << err << endl;
70 \texttt{assert(err < 1e-6);}
71 \texttt{fftwplan p1 = plan(dfft(u[], v[], ny, -1));}
72 \texttt{fftwplan p2 = plan(dfft(u[], v[], ny, 1));}
73 \texttt{real ccc = square(2.*\texttt{pi});}
74 \texttt{cout} << "ny = " << ny << endl;
75 \texttt{map(wij[], ny, ccc*\texttt{x}\times\texttt{y}\times\texttt{y});}
76 \texttt{wij[0] = 1e-5;}
77 \texttt{plot(wij, cmm="wij");}

6.7.2 Complex

(continues on next page)
3
4
func string pc(complex z) {
    string r = "(" + real(z);
    if (imag(z) >= 0) r = r + " + ";
    return r + imag(z) + "i";
}

5
func string toPolar(complex z) {
    return "abs(z) + "*(cos(" + arg(z) + ")+i*sin(" + arg(z) + "))";
}

6
cout << "Standard output of the complex " << pc(z1) << " is the pair: " << z1 << endl;
cout << pc(z1) << " + " << pc(z2) << " = " << pc(z1+z2) << endl;
cout << pc(z1) << " - " << pc(z2) << " = " << pc(z1-z2) << endl;
cout << pc(z1) << " / " << pc(z2) << " = " << pc(z1/z2) << endl;
cout << "Real part of " << pc(z1) << " = " << real(z1) << endl;
cout << "Imaginary part of " << pc(z1) << " = " << imag(z1) << endl;
cout << "abs(" << pc(z1) << ") = " << abs(z1) << endl;
cout << "Polar coordinates of " << pc(z2) << " = " << toPolar(z2) << endl;
cout << "de Moivre formula: " << pc(z2) << "^3 = " << toPolar(z2^3) << endl;
cout << " and polar(2.82887, 0.523509) = " << pc(polar(abs(z2), -arg(z2))) << endl;
cout << "Conjugate of " << pc(z2) << " = " << pc(conj(z2)) << endl;
cout << pc(z1) << " ^ " << pc(z2) << " = " << pc(z1^z2) << endl;

Output of this script is:
Standard output of the complex (2.45+5.33i) is the pair: (2.45,5.33)
(2.45+5.33i) + (2.45+1.41421i) = (4.9+6.74421i)
(2.45+5.33i) - (2.45+1.41421i) = (0+3.91579i)
(2.45+5.33i) * (2.45+1.41421i) = (-1.53526+16.5233i)
(2.45+5.33i) / (2.45+1.41421i) = (1.692+1.19883i)
Real part of (2.45+5.33i) = 2.45
Imaginary part of (2.45+5.33i) = 5.33
abs((2.45+5.33i)) = 5.86612
Polar coordinates of (2.45+1.41421i) =
de Moivre formula: (2.45+1.41421i)^3 =
and polar(2.82887, 0.523509) = (2.45+1.41421i)
Conjugate of (2.45+1.41421i) = (2.45-1.41421i)
(2.45+5.33i) ^ (2.45+1.41421i) = (8.37072-12.7078i)

6.7.3 String

// Concatenation
string tt = "toto1" + 1 + " -- 77";

// Append
string t1 = "0123456789";
t1(4:3) = "abcdefghijk-";

// Sub string
string t55 = t1(4:14);
cout << "tt = " << tt << endl;

(continues on next page)
The output of this script is:

```
1  t = totoll -- 77
2  t1 = 0123abcdefhijk-456789
3  t1.find(abc) = 4
4  t1.rfind(abc) = 4
5  t1.find(abc, 10) = -1
6  t1.ffind(abc, 10) = 4
7  t1.length = 22
8  t55 = abcdefghijk
```

### 6.7.4 Elementary function

```
real b = 1.;
real a = b;
func real phix(real t){
    return (a+b)*cos(t) - b*cos(t*(a+b)/b);
}
func real phiy(real t){
    return (a+b)*sin(t) - b*sin(t*(a+b)/b);
}
```

```
border C(t=0, 2*pi){x=phix(t); y=phiy(t);}
mesh Th = buildmesh(C(50));
plot(Th);
```

**Fig. 6.26:** Mesh
6.7.5 Array

```cpp
real[int] tab(10), tab1(10); //2 array of 10 real
//real[int] tab2; //bug: array with no size

//set all the array to 1.03
real[int] tab2;
//real[int] tab2; //bug: array with no size

tab = 1.03; //set all the array to 1.03

cout << "tab: " << tab << endl;
cout << "min: " << tab.min << endl;
cout << "max: " << tab.max << endl;
cout << "sum: " << tab.sum << endl;

//change the size of array tab to 12 with preserving first value
//set values 10 & 11

real[int] a(5), b(5), c(5), d(5);

da = 1;
b = 2;
c = 3;
da[2] = 0;
d = (a ? b : c); //for i = 0, n-1 : d[i] = a[i] ? b[i] : c[i]

real[string] tt; //array with string index

int[int] ii(0:d.n-1); //set array ii to 0, 1, ..., d.n-1

sort(d, ii); //sort array d and ii in parallel
```

(continues on next page)
A1 = 1.:0.5:3.999;
cout << "1.:0.5:3.999 => " << A1 << endl;
}
{
complex[int] A1(2.+0i:10.+0i); //2, 3, 4, 5, 6, 7, 8, 9, 10
complex[int] A2(2.:3.:10.); //2, 5, 8
cout << " A1(2.+0i:10.+0i): " << A1 << endl;
A1.re real part array: " << A1.re << endl;
// he real part array of the complex array
cout << " A2(2.:3.:10.): " << A2 << endl;
A2.re real part array of the complex array
//the imaginary part array of the complex array
}

// Integer array operators
{
int N = 5;
real[int] a(N), b(N), c(N);
a = 1;
a(0:4:2) = 2;
a(3:4) = 4;
cout << " a: " << a << endl;
b = a + a;
cout<<"b = a + a: " << b << endl;
b += a;
cout<<"b += a: " << b << endl;
b += 2*a;
cout<<"b += 2*a: " << b << endl;
b /= 2;
cout<<"b /= 2: " << b << endl;
b .*= a; // same as b = b .* a
cout<<"b .*= a: " << b << endl;
b ./= a; //same as b = b ./ a
cout<<"b ./= a: " << b << endl;
c = a + b;
cout<<"c = a + b: " << c << endl;
c = 2*a + 4*b;
cout<<"c = 2*a + 4*b: " << c << endl;
c = a + 4*b;
cout<<"c = a + 4*b: " << c << endl;
c = -a + 4*b;
cout<<"c = -a + 4*b: " << c << endl;
c = -a - 4*b;
cout<<"c = -a - 4*b: " << c << endl;
c = a + b;
cout<<"c = a + b: " << c << endl;
c = 2*a + 4*b;
cout<<"c = 2*a + 4*b: " << c << endl;
c = a + 4*b;
cout<<"c = a + 4*b: " << c << endl;
c = -a + 4*b;
cout<<"c = -a + 4*b: " << c << endl;
c = -a - b;
cout<<"c = -a - b: " << c << endl;
c = a .* b;
cout<<"c = a .* b: " << c << endl;
c = a ./ b;
cout<<"c = a ./ b: " << c << endl;
c = 2 * b;
cout<<"c = 2 * b: " << c << endl;
c = b * 2;
cout<<"c = b * 2: " << c << endl;
//this operator do not exist
//c = b/2;
//cout << "c = b / 2: " << c << endl;

//Array methods
cout << "||a||_1 = " << a.l1 << endl;
cout << "||a||_2 = " << a.l2 << endl;
cout << "||a||_infty = " << a.linfy << endl;
cout << "max a_i = " << a.max << " a[ " << a.imax << " ] = " << a[a.imax] << endl;
cout << "min a_i = " << a.min << " a[ " << a.imin << " ] = " << a[a.imin] << endl;
cout << "a' * a = " << (a'*a) << endl;
cout << "a quantile 0.2 = " << a.quantile(0.2) << endl;

//Array mapping
int I = [2, 3, 4, -1, 3];
b = c = -3;
array a = c;
for (i = 0; i < b.n; i++) if (I[i] >= 0) b[i] = a[I[i]];
c = a;
array c = a;
for (i = 0; i < I.n; i++) if (I[i] >= 0) c[I[i]] = a[i];
cout << "b = a(I) : " << b << endl;
cout << "c(I) = a " << c << endl;
c(I) += a;
for (i = 0; i < I.n; i++) if (I[i] >= 0) C(I[i]) += a[i];
cout << "b = a(I) : " << b << endl;
cout << "c(I) = a " << c << endl;

// Array versus matrix
int N = 3, M = 4;
real A(N, M);
real b(N), c(M);
b = [1, 2, 3];
c = [4, 5, 6, 7];
complex C(N, M);
complex cb = [1, 2, 3], cc = [10i, 20i, 30i, 40i];
b = [1, 2, 3];
array I = [2, 0, 1];
array J = [2, 0, 1, 3];
A = 1; //set all the matrix
A(2, :) = 4; //the full line 2
A(:, 1) = 5; //the full column 1
A(0:N-1, 2) = 2; //set the column 2
A(1, 0:2) = 3; //set the line 1 from 0 to 2
cout << "A = " << A << endl;

//outer product
C = cb * cc';
C += 3 * cb * cc';
C -= 5i * cb * cc';
cout << "C = " << C << endl;
(continues on next page)
```cpp
// this transforms an array into a sparse matrix
matrix B;
B = A;
B = A(I, J);  // B(i, j) = A(I(i), J(j))
B = A(I^-1, J^-1);  // B(I(i), J(j)) = A(i, j)

// outer product
A = 2. * b * c';
cout << "A = " << A << endl;
B = b*c';  // outer product B(i, j) = b(i)*c(j)
B = b*c';  // outer product B(i, j) = b(i)*c(j)
B = (2*b*c')(I, J);  // outer product B(i, j) = b(I(i))*c(J(j))
B = (3.*b*c')(I^-1, J^-1);  // outer product B(I(i), J(j)) = b(i)*c(j)
cout << "B = (3.*b*c')(I^-1, J^-1) = " << B << endl;

// row and column of the maximal coefficient of A
int i, j, ii, jj;
ijmax(A, ii, jj);

i = A.imax;
j = A.jmax;
cout << "Max " << i << " " << j << ", = " << A.max << endl;

// row and column of the minimal coefficient of A
ijmin(A, i, j);

ii = A.imin;
jj = A.jmin;
cout << "Min " << ii << " " << jj << ", = " << A.min << endl;
```

The output of this script is:

```
tab: 10
  1.03  2.15  1.03  1.03  1.03
  1.03  1.03  1.03  1.03  1.03
min: 1.03
max: 2.15
sum: 11.42
resized tab: 12
  1.03  2.15  1.03  1.03  1.03
  1.03  1.03  1.03  1.03  1.03
  3.14  3.14
sorted tab: 12
  1.03  1.03  1.03  1.03  1.03
  1.03  1.03  1.03  1.03  2.15
  3.14  3.14
```

(continues on next page)
-5 -4 -3 -2 -1
ii: 5
  4  3  2  1  0
A1(2:10): 9
  2  3  4  5  6
  7  8  9 10
A2(2:3:10): 9
  2  3  4  5  6
  7  8  9 10
l:2:5 => 3
  1  3  5
A1(2:10): 9
  2  3  4  5  6
  7  8  9 10
A2(2:3:10): 9
  2  3  4  5  6
  7  8  9 10
l:0.5:3.999 => 6
  1  1.5  2  2.5  3
  3.5
A1(2.+0i:10.+0i): 9
  (2,0) (3,0) (4,0) (5,0) (6,0)
  (7,0) (8,0) (9,0) (10,0)
A2(2.:3.:10.)= 3
  (2,0) (5,0) (8,0)
A1.re real part array: 9
  2  3  4  5  6
  7  8  9 10
A1.im imag part array: 9
  0  0  0  0  0
  0  0  0  0
a: 5
  2  1  2  4  4
b = a + a: 5
  4  2  4  8  8
b += a: 5
  6  3  6 12 12
b += 2*a: 5
  10  5 10 20 20
b /= 2: 5
  5  2.5 5 10 10
b .*= a: 5
  10  2.5 10 40 40
b ./= a: 5
  5  2.5 5 10 10
c = a + b: 5
  7 3.5 7 14 14
(continues on next page)
```plaintext
\begin{verbatim}
c = 2*a + 4b: 5
     24 12 24 48 48

c = a + 4b: 5
     22 11 22 44 44

c = -a + 4b: 5
     18 9 18 36 36

c = -a - 4b: 5
     -22 -11 -22 -44 -44

c = a - b: 5
     -7 -3.5 -7 -14 -14

c = a .* b: 5
     10 2.5 10 40 40

c = a ./ b: 5
     0.4 0.4 0.4 0.4 0.4

c = 2 * b: 5
     10 5 10 20 20

c = b * 2: 5
     10 5 10 20 20

||a||_1 = 13
||a||_2 = 6.40312
||a||_\infty = 4
sum a_i = 13
max a_i = 4 a[ 3 ] = 4
min a_i = 1 a[ 1 ] = 1
a' * a = 41
a quantile 0.2 = 2
b = a(I) : 5
     2 4 4 -3 4
c(I) = a 5
     -3 -3 2 4 2
b = a(I) : 5
     2 4 4 -3 4
c(I) = a 5
     -3 -3 4 9 4
A = 3 4
     1 5 2 1
     3 3 3 1
     4 5 2 4
C = 3 4
     (-50,-40) (-100,-80) (-150,-120) (-200,-160)
     (-100,-80) (-200,-160) (-300,-240) (-400,-320)
     (-150,-120) (-300,-240) (-450,-360) (-600,-480)
\end{verbatim}
```
\[ A = \begin{bmatrix} 3 & 4 \\ 8 & 10 & 12 & 14 \\ 16 & 20 & 24 & 28 \\ 24 & 30 & 36 & 42 \end{bmatrix} \]

\[ B = (3.0 \times b \times c') (I^{-1}, J^{-1}) = \# \text{ Sparse Matrix (Morse)} \]

```plaintext
# first line: n m (is symmetric) nbcoef
# after for each nonzero coefficient: i j a_ij where (i,j) in \{1,...,n\}x\{1,...,m\}
3 4 0 12
  1  1 10
  1  2 12
  1  3  8
  1  4 14
  2  1 15
  2  2 18
  2  3 12
  2  4 21
  3  1  5
  3  2  6
  3  3  4
  3  4  7
```

### 6.7.6 Block matrix

```plaintext
// Parameters
real f1 = 1.;
real f2 = 1.5;

// Mesh
mesh Th1 = square(10, 10);
mesh Th2 = square(10, 10, [1+x, -1+y]);
plot(Th1, Th2);

// Fespace
fespace Uh1(Th1, P1);
Uh1 u1;

fespace Uh2(Th2, P2);
Uh2 u2;

// Macro
macro grad(u) [dx(u), dy(u)] //

// Problem
varf vPoisson1 (u, v)
  = int2d(Th1) (grad(u)' * grad(v))
  - int2d(Th1) (f1 * v) + on(1, 2, 3, 4, u=0);

varf vPoisson2 (u, v)
```

(continues on next page)
```c
int2d(Th2)(
    grad(u)' * grad(v)
) - int2d(Th2)(
    f1 * v
) + on(1, 2, 3, 4, u=0);

matrix<real> Poisson1 = vPoisson1(Uh1, Uh1);
real[int] Poisson1b = vPoisson1(0, Uh1);

matrix<real> Poisson2 = vPoisson2(Uh2, Uh2);
real[int] Poisson2b = vPoisson2(0, Uh2);

//block matrix
matrix<real> G = [[Poisson1, 0], [0, Poisson2]];
set(G, solver=sparsesolver);

//block right hand side
real[int] Gb = [Poisson1b, Poisson2b];

// Solve
real[int] sol = G^-1 * Gb;

// Dispatch
[u1[], u2[]] = sol;

// Plot
plot(u1, u2);
```

Fig. 6.27: Result
### Matrix operations

```cpp
// Mesh
mesh Th = square(2, 1);

// Fespace
fespace Vh(Th, P1);
Vh f, g;
f = x*y;
g = sin(pi*x);

Vh<complex> ff, gg; // a complex valued finite element function
ff = x*(y+1i);
gg = exp(pi*x*1i);

// Problem
varf mat (u, v) = int2d(Th) (1*dx(u)*dx(v) + 2*dx(u)*dy(v) + 3*dy(u)*dx(v) + 4*dy(u)*dy(v)) + on(1, 2, 3, 4, u=1);

varf mati (u, v) = int2d(Th) (1*dx(u)*dx(v) + 2i*dx(u)*dy(v) + 3*dy(u)*dx(v) + 4*dy(u)*dy(v)) + on(1, 2, 3, 4, u=1);

matrix A = mat(Vh, Vh);
matrix<complex> AA = mati(Vh, Vh); // a complex sparse matrix

// Operations
Vh m0; m0[] = A*f[];
Vh m01; m01[] = A'*f[];
Vh m1; m1[] = f[].*g[];
Vh m2; m2[] = f[]./g[];

// Display
cout << "f = " << f[] << endl;
cout << "g = " << g[] << endl;
cout << "A = " << A << endl;
cout << "m0 = " << m0[] << endl;
cout << "m01 = " << m01[] << endl;
cout << "m1 = " << m1[] << endl;
cout << "m2 = " << m2[] << endl;
cout << "dot Product = " << f[]'*g[] << endl;
cout << "hermitien Product = " << ff[]'*gg[] << endl;
cout << "outer Product = " << (A=f[]*g[]) << endl;
cout << "hermitien outer Product = " << (AA=ff[]*gg[]) << endl;
```

(continues on next page)
The output of this script is:

```
F = 6
0  0  0  0  0.5
1

G = 6
0  1  1.224646799e-16  0  1
1.224646799e-16
A = # Sparse Matrix (Morse)
# first line: n m (is symmetric) nbcoef
# after for each nonzero coefficient:  i j a_ij where (i,j) \in \{1,...,n\}x\{1,...,m\}
6 6 0 24
1 1 1.0000000000000000199e+30
1 2 0.49999999999999994449
1 4 0
1 5 -2.5
2 1 0
2 2 1.0000000000000000199e+30
2 3 0.49999999999999994449
2 5 0.499999999999999997796
2 6 -2.5
3 2 0
3 3 1.0000000000000000199e+30
3 6 0.499999999999999977796
4 1 0.499999999999999977796
4 4 1.0000000000000000199e+30
4 5 0
5 1 -2.5
5 5 0.499999999999999977796
5 6 0.49999999999999999449
6 2 -2.5
6 3 0
6 6 0.49999999999999999449
```
```
-1.25  -2.25  0.5  0 5e+29
le+30
m01 = 6
-1.25  -2.25  0 0.25 5e+29
le+30
m1 = 6
  0 0 0 0 0.5
  1.224646799e-16
m2 = 6
  -nan 0 0 -nan 0.5
  8.165619677e+15
dot Product = 0.5
hermitien Product = (1.11022e-16,2.5)
outer Product = # Sparse Matrix (Morse)
# first line: n m (is symmetric) nbcoef
# after for each nonzero coefficient: i j a_ij where (i,j) \in \{1,...,n\}\times\{1,...,m\}
6 6 0 8
  5 2 0.5
  5 3 6.123233957367660359e-17
  5 5 0.5
  6 2 1
  6 3 6.123233957367660359e-17
  6 5 1
  6 6 6.1246467991473532072e-16
hermitien outer Product = # Sparse Matrix (Morse)
# first line: n m (is symmetric) nbcoef
# after for each nonzero coefficient: i j a_ij where (i,j) \in \{1,...,n\}\times\{1,...,m\}
6 6 0 24
  2 1 (0,0.5)
  2 2 (0.5,3.061616997868330179e-17)
  2 3 (6.123233957367660359e-17,-0.5)
  2 4 (0,0.5)
  2 5 (0.5,3.061616997868330179e-17)
  2 6 (6.123233957367660359e-17,-0.5)
  3 1 (0,1)
  3 2 (1,6.123233957367660359e-17)
  3 3 (1.2246467991473532072e-16,-1)
  3 4 (0,1)
  3 5 (1,6.123233957367660359e-17)
  3 6 (1.2246467991473532072e-16,-1)
  5 1 (0.5,0.5)
  5 2 (0.5,-0.499999999999994449)
  5 3 (-0.4999999999999944449,-0.500000000000011102)
  5 4 (0.5,0.5)
  5 5 (0.5,-0.4999999999999944449)
  5 6 (-0.4999999999999944449,-0.500000000000011102)
  6 1 (1,1)
  6 2 (1,-0.999999999999888898)
  6 3 (-0.999999999999888898,-1.0000000000000222)
  6 4 (1,1)
  6 5 (1,-0.999999999999888898)
  6 6 (-0.999999999999888898,-1.0000000000000222)
I = 8
6 6 4 4 5
(continues on next page)
```

6.7. Developers
Warning: Due to Fortran indices starting at one, the output of a diagonal matrix $D$ is indexed from 1. but in FreeFEM, the indices start from 0.

6.7.8 Matrix inversion

```plaintext
load "lapack"
load "fflapack"

// Matrix
int n = 5;
real[int, int] A(n, n), A1(n, n), B(n, n);
for (int i = 0; i < n; ++i)
    for (int j = 0; j < n; ++j)
        A(i, j) = (i == j) ? n+1 : 1;
cout << A << endl;

// Inversion (lapack)
A1 = A^-1; //def in "lapack"
cout << A1 << endl;

B = 0;
for (int i = 0; i < n; ++i)
    for (int j = 0; j < n; ++j)
        for (int k = 0; k < n; ++k)
            B(i, j) += A(i, k) * A1(k, j);
cout << B << endl;

// Inversion (fflapack)
inv(A1); //def in "fflapack"
cout << A1 << endl;
```

The output of this script is:
### Tip:
To compile `lapack.cpp` and `fflapack.cpp`, you must have the `lapack` library on your system and compile the plugin with the command:

```bash
ff-c++ lapack.cpp -llapack  ff-c++ fflapack.cpp -llapack
```

#### 6.7.9 FE array

```cpp
// Mesh
mesh Th = square(20, 20, [2*x, 2*y]);

// Fespace
fespace Vh(Th, P1);
Vh u, v, f;

// Problem
problem Poisson (u, v)
  = int2d(Th) (dx(u)*dx(v) + dy(u)*dy(v))
  + int2d(Th) (-f*v)
  + on(1, 2, 3, 4, u=0);
```

(continues on next page)
Vh[int] uu(3); // an array of FE function
  // Solve problem 1
  f = 1;
Poisson;
uu[0] = u;
  // Solve problem 2
  f = sin(pi*x)*cos(pi*y);
Poisson;
uu[1] = u;
  // Solve problem 3
  f = abs(x-1)*abs(y-1);
Poisson;
uu[2] = u;

  // Plot
  for (int i = 0; i < 3; i++)
    plot(uu[i], wait=true);

Fig. 6.28: Finite element array
6.7.10 Loop

```cpp
for (int i = 0; i < 10; i=i+1)
    cout << i << endl;
real eps = 1.;
while (eps > 1e-5){
    eps = eps/2;
    if (i++ < 100)
        break;
    cout << eps << endl;
}
for (int j = 0; j < 20; j++){
    if (j < 10) continue;
    cout << "j = " << j << endl;
}
```

6.7.11 Implicit loop

```cpp
real [int, int] a(10, 10);
real [int] b(10);
for [i, bi : b]{
    bi = i+1;
    cout << i << " " << bi << endl;
}cout << "b = " << b << endl;
for [i, j, aij : a]{
    aij = 1./(2+i+j);
    if (abs(aij) < 0.2) aij = 0;
}cout << "a = " << a << endl;
matrix A = a;
string[string] ss; //a map
ss["1"] = 1;
ss["2"] = 2;
ss["3"] = 5;
for [i, bi : ss]
    bi = i + 6 + "-dddd";
cout << "ss = " << ss << endl;
int[string] si;
si[1] = 2;
si[50] = 1;
for [i, vi : si]{
    cout << "i = " << setw(3) << i << " " << setw(10) << vi << endl;
    vi = atoi(i)*2;
}cout << "si = " << si << endl;
for [i, j, aij : A]{
    cout << i << " " << j << " " << aij << endl;
}
```
aij = -aij;
}
cout << A << endl;

The output of this script is:

The output of this script is:

1 0 1
2 2
3 3
4 4
5 5
6 6
7 7
8 8
9 9
10 10

b = 10

1 1 2 3 4 5
6 7 8 9 10

a = 10 10

0.5 0.3333333333 0.25 0.2 0 0 0 0 0 0
0.3333333333 0.25 0.2 0 0 0 0 0 0 0
0.25 0.2 0 0 0 0 0 0 0 0
0.2 0 0 0 0 0 0 0 0 0
0 0 0 0 0 0 0 0 0 0
0 0 0 0 0 0 0 0 0 0
0 0 0 0 0 0 0 0 0 0
0 0 0 0 0 0 0 0 0 0
0 0 0 0 0 0 0 0 0 0
0 0 0 0 0 0 0 0 0 0

ss = 1 1
2 2
3 3

i 1 2
i 50 1
si = 1 2
50 100

0 0 0.5
0 1 0.333333
0 2 0.25
0 3 0.2
1 0 0.333333
1 1 0.25
1 2 0.2
2 0 0.25
2 1 0.2
3 0 0.2

# Sparse Matrix (Morse)
# first line: n m (is symmetric) nbcoef
# after for each nonzero coefficient: i j a_ij where (i,j) \in \{1,...,n\}\times\{1,...,m\}
10 10 0 10
1 1 -0.5
1 2 -0.3333333333333331483
6.7.12 I/O

```cpp
int i;
cout << "std-out" << endl;
cout << " enter i = ?";
cin >> i;
{
    ofstream f("toto.txt");
f << i << "hello world'\n";
} //close the file f because the variable f is delete

{
    ifstream f("toto.txt");
f >> i;
}
{
    ofstream f("toto.txt", append);
    //to append to the existing file "toto.txt"
f << i << "hello world'\n";
} //close the file f because the variable f is delete

cout << i << endl;
```

6.7.13 File stream

```cpp
int where;
real[int] f = [0, 1, 2, 3, 4, 5];
real[int] g(6);
{
    ofstream file("f.txt", binary);
    file.precision(16);
    file << f << endl;
    where = file.tellp();
    file << 0.1;
    cout << "Where in file " << where << endl;
    file << " # comment bla bla ... 0.3 \n";
    file << 0.2 << endl;
    file.flush; //to flush the buffer of file
}
```
17 // Function to skip comment starting with # in a file
18 func ifstream skipcomment(ifstream &ff) {
19     while(1) {
20         int where = ff.tellg(); // store file position
21         string comment;
22         ff >> comment;
23         if (!ff.good()) break;
24         if (comment(0:0) == "#") {
25             getline(ff, comment);
26             cout << " -- #" << comment << endl;
27         }
28         else {
29             ff.seekg(where); // restore file position
30             break;
31         }
32     } // while
33     return ff;
34 }
35
36 {
37     real xx;
38     ifstream file("f.txt", binary);
39     cout << "Where " << file.tellg << endl;
40     file >> xx;
41     cout << " xx = " << xx << " good ? " << file.good() << endl;
42     assert(xx == 0.1);
43     skipcomment(file) >> xx;
44     assert(xx == 0.2);
45     file.seekg(0); // rewind
46     cout << "Where " << file.tellg << " " << file.good() << endl;
47     file >> g;
48 }

6.7.14 Command line arguments

When using the command:

FreeFem++ script.edp arg1 arg2

The arguments can be used in the script with:

for (int i = 0; i < ARGV.n; i++)
    cout << ARGV[i] << endl;

When using the command:

FreeFem++ script.edp -n 10 -a 1. -d 42.

The arguments can be used in the script with:

include "getARGV.idp"

int n = getARGV("-n", 1);
real a = getARGV("-a", 1.);
real d = getARGV("-d", 1.);

### 6.7.15 Macro

```freefem
// Macro without parameters
macro xxx() {
    real i = 0;
    int j = 0;
    cout << i << " " << j << endl;
}

// Macro with parameters
macro toto(i) i //
toto({real i = 0; int j = 0; cout << i << " " << j << endl;})

// Macro as parameter of a macro
real[int,int] CC(7, 7), EE(5, 3), EEps(4, 4);
macro VIL6(v, i) [v(1,i), v(2,i), v(4,i), v(5,i), v(6,i)]
macro VIL3(v, i) [v(1,i), v(2,i)]
macro VV6(v, vv) [v(vv,1), v(vv,2), v(vv,4), v(vv,5), v(vv,6)]
macro VV3(v, vv) [v(vv,1), v(vv,2)]
func C5x5 = VV6(VIL6, CC);
func E5x2 = VV6(VIL3, EE);
func Eps = VV3(VIL3, EEps);

// Macro concatenation
mesh Th = square(2, 2);
fespace Vh(Th, P1);
Vh Ux=x, Uy=y;
macro div(V) (dx(V#x) + dy(V#y)) //
cout << int2d(Th)(div(U)) << endl;
// Verify the quoting
macro foo(i, j, k) i j k //
foo(, , )
foo({int[], {int} a[10], {}})

// IFMACRO - ENDIFMACRO
NewMacro grad(u) [dx(u), dy(u)] EndMacro
cout << int2d(Th)(grad(Ux)' * grad(Uy)) << endl;
// IFMACRO - ENDIFMACRO
macro AA CAS1 //
```

(continues on next page)
The output script generated with macros is:

```
// Macro without parameters
macro xxx {
  real i = 0;
  int j = 0;
  cout << i << " " << j << endl;
}

// Macro with parameters
macro toto(i) i

// Macro as parameter of a macro
real[int,int] CC(7, 7), EE(6, 3), EEps(4, 4);

// FILE - LINE
cout << "In " << FILE << ", line " << LINE << endl;
```
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(continued from previous page)

2 : [ CC(1,4), CC(2,4), CC(4,4), CC(5,4), CC(6,4)], [ CC(1, 
3 : -5), CC(2,5), CC(4,5), CC(5,5), CC(6,5)], 
27 : func E5x2 = 
1 : 2 : [ EE(1,1), EE(2,1)], [ EE(1,2), EE(2,2)], 
2 : [ EE(1,4), EE(2,4)], [ EE(1,5), EE(2,5)], 
3 : [ EE(1,6), EE(2,6)], 
28 : func Eps = [ [ EEps(1,1), EEps(2,1)], [ EEps(1,2), EEps(2,2)]]; 
29 : // Macro concatenation 
30 : mesh Th = square(2, 2); 
31 : fespace Vh(Th, P1); 
32 : Vh Ux=x, Uy=y; 
33 : 
34 : macro div(V) (dx(V#x) + dy(V#y)) // 
35 : cout << int2d(Th) ( (dx(Ux) + dy(Uy)) ) << endl; 
36 : // Verify the quoting 
37 : macro foo(i,j,k) i j k // 
38 : int[ int] a(10); 
39 : // NewMacro - EndMacro 
40 : macro grad(u) [dx(u), dy(u)]; 
41 : cout << int2d(Th) ( [dx(Ux), dy(Ux)]' * [dx(Uy), dy(Uy)] ) << endl; 
42 : // IFMACRO - ENDIFMACRO 
43 : macro AACS1 // 
44 : cout " " " AA = " " Stringification( CAS1 ) " " endl; 
45 : macro CASEfile1.edp 
46 : cout " " " CASE = " " Stringification(file1.edp) " " endl; 
47 : 
48 : include Stringification(file1.edp) cout " " "This is the file 1" " endl; 
49 : // FILE - LINE 
50 : cout " " " In " " " FILE " " " , line " " " LINE " " endl;

The output of this script is:

AA = CAS1
CASE = file1.edp
This is the file 1
In Macro.edp, line 59

6.7. Developers
6.7.16 Basic error handling

```cpp
real a;
try{
  a = 1./0.;
} catch (...){ //all exceptions can be caught
  cout << "Catch an ExecError" << endl;
  a = 0.;
}
```

The output of this script is:

```
l/0 : d d d
  current line = 3
Exec error : Div by 0
   -- number :1
Catch an ExecError
```

6.7.17 Error handling

```cpp
// Parameters
int nn = 5;
func f = 1; //right hand side function
func g = 0; //boundary condition function

// Mesh
mesh Th = square(nn, nn);

// Fespace
fespace Vh(Th, P1);
Vh uh, vh;

// Problem
real cpu = clock();
problem laplace (uh, vh, solver=Cholesky, tolpivot=1e-6)
  = int2d(Th)(
      dx(uh)*dx(vh)
      + dy(uh)*dy(vh)
  )
  + int2d(Th)(
      - f*vh
  )
try{
  cout << "Try Cholesky" << endl;
  // Solve
  laplace;
  // Plot
  plot(uh);
}
```

(continues on next page)
// Display
    cout << "laplacian Cholesky " << nn << ", x_" << nn << " : " << -cpu+clock() << "-
    s, max = " << uh[].max << endl;
}
catch(...) { //catch all error
    cout << " Catch cholesky PB " << endl;
}

The output of this script is:

try Cholesky
ERREUR choleskypivot (35)= -6.43929e-15 < 1e-06
    current line = 29
Exec error : FATAL ERREUR dans ./../femlib/MatriceCreuse_tpl.hpp
cholesky line:
    -- number :688
    catch an erreur in  solve  =>  set sol = 0 !!!!!!!!
Catch cholesky PB


