

# Setting up your workspace

FeenoX Tutorial #0

## Contents

<b>1</b>	<b>Foreword</b>	<b>2</b>
<b>2</b>	<b>FeenoX</b>	<b>3</b>
2.1	Compiling from source . . . . .	3
2.2	Executing an example . . . . .	4
2.3	Executing a test . . . . .	4
<b>3</b>	<b>Gmsh</b>	<b>4</b>
3.1	Checking that mesh refinement gives more accurate results . . . . .	5
<b>4</b>	<b>Text editor</b>	<b>6</b>
4.1	Vim . . . . .	6
4.2	Kate . . . . .	7
<b>5</b>	<b>Post-processors</b>	<b>7</b>
<b>6</b>	<b>Notes on hardware</b>	<b>8</b>
6.1	Memory . . . . .	8
6.2	CPU . . . . .	8

## 1 Foreword

FeenoX is a cloud-first engineering tool. Therefore, it runs natively on GNU/Linux platforms.

Theoretically, the tool could be compiled and run in other architectures such as Windows or MacOS in a non-cloud approach. However, this is *highly* discouraged because those two operating systems are

1. not cloud-friendly, let alone cloud-first; and
2. neither free-as-in-free-beer nor open source.

In order to take the tutorials that follow, it is then recommended to stick to **GNU/Linux** as explained below.

The best way to learn and to understand how FeenoX works is to use a native GNU/Linux distribution as the main operating system, either in a laptop or a desktop PC. This sentence from PETSc's Matt Kneppley from 2015 speaks for itself:

“It is really worth any amount of time and effort to get away from Windows if you are doing computational science.”

<https://lists.mcs.anl.gov/pipermail/petsc-users/2015-July/026388.html>

- If you already use GNU/Linux then you are almost set! Any distribution will do, although FeenoX is developed in Debian so `apt-get` (or `apt`) will be used as the package manager. Note that the names of the packages being installed as dependencies may vary from distribution to distribution.
- If you do not currently use GNU/Linux as your main operating system and still do not want to spend any time nor effort on doing things right, you can either
  - i. fire up a virtual GNU/Linux server in a cloud provider (e.g. AWS, Azure, DigitalOcean, Contabo, etc.) and connect through SSH, or
  - ii. use a containerized GNU/Linux (e.g. with `docker`), or
  - iii. use a GNU/Linux box through an virtual computer emulator (e.g. VirtualBox, VMWare, Vagrant, etc.)

Note that any of these three options is at least as difficult as using a native GNU/Linux box.

Up to this point, I assume you have access to a Unix-like shell (i.e. a GNU/Linux terminal, MacOS or even Cygwin or something of the like) as a regular user and that you have permissions to use `sudo`. If you do not know what this means, look it up in your favorite search engine, watch some videos online or—even better—ask for help. Spend some time (which will be really worth it) familiarizing with working with the terminal, issuing commands, etc.

In the following sections there will be terminal mimics. Lines starting with `$` show commands that ought to be typed into the command line. The `$` itself does not have to be typed as it is part of the prompt. Lines not starting with a dollar sign show the output of the invoked command. For example

```
$ feenox
FeenoX v0.2.144-g31d72de
a free no-fee no-X uniX-like finite-element(ish) computational engineering tool

usage: feenox [options] inputfile [replacement arguments] [petsc options]

-h, --help          display options and detailed explanations of command-line usage
-v, --version       display brief version information and exit
-V, --versions      display detailed version information
--pdes              list the types of PROBLEMs that FeenoX can solve, one per line
```

## Setting up your workspace

```
Run with --help for further explanations.  
$
```

When the output of a command is too long and not important for the tutorial, an ellipsis is denoted with [...]:

```
$ feenox  
[...]  
$
```

## 2 FeenoX

The most important thing to set up is FeenoX itself. Since there are still no Debian packages for FeenoX, the most straightforward way to go is to download the Linux binary tarball from <https://seamplex.com/feenox/dist/linux/> and copy the executable into `/usr/local/bin` so it is globally available. To download the tarball you need `wget` and to un-compress it `tar` and `gz` (which should be already installed anyway), so do

```
$ sudo apt-get install wget tar gzip  
$ wget https://seamplex.com/feenox/dist/linux/feenox-v0.2.142-g8f80cb9-linux-amd64.tar.gz  
[...]  
$ tar xvzf feenox-v0.2.142-g8f80cb9-linux-amd64.tar.gz  
[...]  
$ sudo cp feenox-v0.2.142-g8f80cb9-linux-amd64/bin/feenox /usr/local/bin/  
$
```

If you do not have root access, read the complete download or compilation instructions. Search for “root” and read along.

You should now be able to invoke FeenoX by executing `feenox` from any directory. See the “Invocation” section of the FeenoX Manual for details about how to invoke it. Check this is the case:

```
$ feenox  
FeenoX v0.2.144-g31d72de  
a free no-fee no-X uniX-like finite-element(ish) computational engineering tool  
  
usage: feenox [options] inputfile [replacement arguments] [petsc options]  
  
-h, --help          display options and detailed explanations of command-line usage  
-v, --version       display brief version information and exit  
-V, --versions      display detailed version information  
--pdes              list the types of PROBLEMS that FeenoX can solve, one per line  
  
Run with --help for further explanations.  
$
```

If you get stuck or get an error, please ask for help in the Github discussion page.

### 2.1 Compiling from source

If you want to tweak the compilation flags, use other libraries, modify the code or just learn how FeenoX works, follow the Compilation instructions. Again, do not hesitate to ask in the Github discussion page.

Setting up your workspace

## 2.2 Executing an example

The FeenoX examples are a set of annotated input files that can be found online at <https://www.seamless.com/feenox/examples/>. These examples range from a simple “Hello World” down to thermo-mechanical problems.

Let us run one of them to check FeenoX works. Find the `examples` directory and `cd` into it. If you have a binary version, it will be in `share/doc/examples`. If you have the source tarball or cloned the repository, it will be directly `examples`.

So far the only example that already comes with a mesh (so Gmsh is not needed to run it) is the “Parallelepiped...” one:

```
$ cd examples
$ feenox parallelepiped-thermal.fee
2.43384e-10
$ feenox parallelepiped-mechanical.fee
9.5239e-05
$
```

The smaller the numbers, the better the convergence with respect to the analytical solution. After installing Gmsh below, we will refine the meshes and check these numbers decrease.

## 2.3 Executing a test

By design, FeenoX has to have a set of tests such that if a regression is introduced in the code (say someone flips a sign by mistake), the bug can be quickly detected and eventually fixed. The tests are run with `make`  $\leftrightarrow$  check when compiling the source, as explained in the Compilation instructions.

Alternatively, go to the `tests` directory, pick any of the shell scripts ending in `.sh` and run them:

```
$ cd tests
$ ./nafems-le1.sh
nafems-le1.fee 1 ... ok
nafems-le1.fee 2 ... ok
nafems-le1.fee 3 ... ok
nafems-le1.fee 4 ... ok
nafems-le1.fee 5 ... ok
nafems-le1.fee 6 ... ok
nafems-le1.fee 7 ... ok
nafems-le1.fee 8 ... ok
$
```

If any of the lines does not say `ok` but something else, then the code contains at least one error.

## 3 Gmsh

To solve problems involving partial differential equations (i.e. elasticity, heat conduction, neutron transport, etc.) FeenoX needs a mesh in Gmsh’s MSH format. Any mesher whose output format can be converted to `.msh` should work, although of course the most natural way to create these meshes is to use Gmsh itself.

The easiest way to go is to install Gmsh from the `apt` repository:

```
$ sudo apt-get install gmsh
```

## Setting up your workspace

Check `gmsh` is globally available by calling it with `-info`:

```
$ gmsh -info
Version      : 4.10.5
License     : GNU General Public License
Build OS    : Linux64
Build date  : 20220701
Build host  : gmsh.info
Build options : 64Bit ALGLIB[contrib] ANN[contrib] Bamg Blas[petsc] Blossom Cgns DIntegration Dlopen ↔
              DomHex Eigen[contrib] Fltk Gmm[contrib] Hxt Jpeg Khipack Lapack[petsc] LinuxJoystick MathEx[contrib] ↔
              ] Med Mesh Metis[contrib] Mmg Mpeg Netgen ONELAB ONELABMetamodel OpenCASCADE OpenCASCADE-CAF OpenGL ↔
              OpenMP OptHom PETSc Parser Plugins Png Post QuadMeshingTools QuadTri Solver TetGen/BR Voro++[ ↔
              contrib] WinslowUntangler ZlibF
FLTK version : 1.4.0
PETSc version : 3.14.4 (real arithmetic)
OCC version  : 7.6.1
MED version  : 4.1.0
Packaged by  : geuzaine
Web site     : https://gmsh.info
Issue tracker : https://gitlab.onelab.info/gmsh/gmsh/issues
$
```

It should be noted that depending on the version of the base operating system, the Gmsh version in the apt repository might be old enough so as to fail with the examples provided in the tutorials, that are based on recent Gmsh versions. If this is the case, as with FeenoX above, you can always download newer Gmsh binaries:

```
$ wget http://gmsh.info/bin/Linux/gmsh-4.10.5-Linux64.tgz
[...]
$ tar xvzf gmsh-4.10.5-Linux64.tgz
[...]
$ sudo cp gmsh-4.10.5-Linux64/bin/gmsh /usr/local/bin
```

Also, Gmsh can be compiled from source by following the instructions in the documentation.

It is important to note that Gmsh creates its meshes by reading an input file with extension `.geo` without needing to use a graphical interface at all. Therefore, Gmsh can be used even through `ssh` or `docker` where there is no graphical device. However, it does provide a GUI that is very handy to create the `.geo` in the first place and to help identify the ids of the surfaces that will be subject to boundary conditions (or volumes for material properties). So you will be able to go through the tutorials in text-only mode but you will not be able to exploit its full potential nor even see what the mesh you are using look like.

### 3.1 Checking that mesh refinement gives more accurate results

Let us now go back to the `examples` directory and refine the mesh of the “Parallelepiped...” case.

```
$ gmsh -3 parallelepiped.geo -order 1 -clscale 1.2 -o parallelepiped-coarse.msh
[...]
$ gmsh -3 parallelepiped.geo -order 2 -clscale 0.6
[...]
$ feenox parallelepiped-thermal.fee
1.05973e-10
$ feenox parallelepiped-mechanical.fee
6.95446e-05
$
```

## Setting up your workspace

Don't worry if you do not understand the Gmsh command line. We will work out the details in the tutorials.

## 4 Text editor

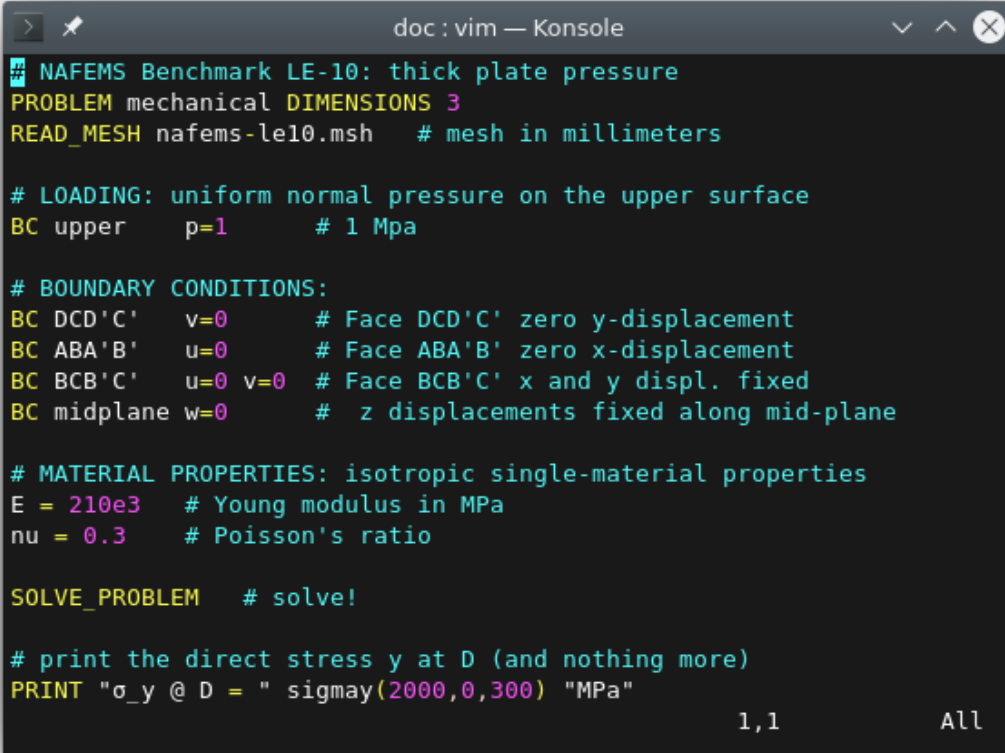
In order to create the both the FeenoX and the Gmsh input files, you will need a text editor. In principle any editor will do, but since FeenoX uses keywords to define the problem being solved, it is way more efficient (and aesthetically more pleasant) to use one that supports syntax highlighting.

So far, FeenoX supports highlighting for Vim and Kate, both discussed below. Other editors such as Emacs, Nano, Pico, Gedit, etc. can be used.

### 4.1 Vim

The `vim` text editor can be perfectly used to edit input files. It is text-only so it works through `ssh` and `docker`. It has a non-trivial learning curve but it is worth to learn its basics because at some point you will want to run FeenoX in an actual cloud server. Vim will be your friend there.

To enable syntax highlighting copy the file `fee.vim` into `~/.vim/syntax`.



```
doc : vim — Konsole
# NAFEMS Benchmark LE-10: thick plate pressure
PROBLEM mechanical DIMENSIONS 3
READ_MESH nafems-le10.msh # mesh in millimeters

# LOADING: uniform normal pressure on the upper surface
BC upper p=1 # 1 Mpa

# BOUNDARY CONDITIONS:
BC DCD'C' v=0 # Face DCD'C' zero y-displacement
BC ABA'B' u=0 # Face ABA'B' zero x-displacement
BC BCB'C' u=0 v=0 # Face BCB'C' x and y displ. fixed
BC midplane w=0 # z displacements fixed along mid-plane

# MATERIAL PROPERTIES: isotropic single-material properties
E = 210e3 # Young modulus in MPa
nu = 0.3 # Poisson's ratio

SOLVE_PROBLEM # solve!

# print the direct stress y at D (and nothing more)
PRINT "σ_y @ D = " sigmay(2000,0,300) "MPa"

1,1 All
```

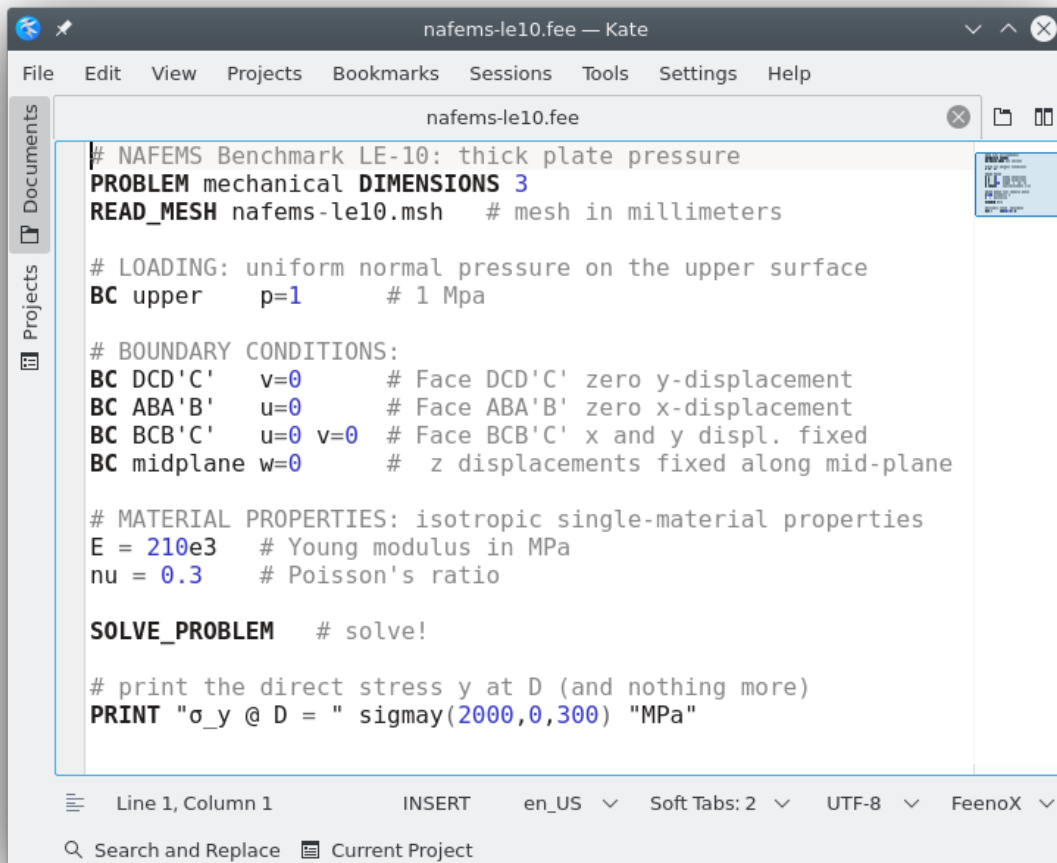
Figure 1: FeenoX input file edited in Vim on Konsole

Setting up your workspace

## 4.2 Kate

For those using a native GNU/Linux box with a graphical interface, the recommended editor is Kate.

To enable syntax highlighting copy the file `feenox.xml` into `~/.local/share/katepart5/syntax/`.



```
# NAFEMS Benchmark LE-10: thick plate pressure
PROBLEM mechanical DIMENSIONS 3
READ_MESH nafems-le10.msh # mesh in millimeters

# LOADING: uniform normal pressure on the upper surface
BC upper p=1 # 1 Mpa

# BOUNDARY CONDITIONS:
BC DCD'C' v=0 # Face DCD'C' zero y-displacement
BC ABA'B' u=0 # Face ABA'B' zero x-displacement
BC BCB'C' u=0 v=0 # Face BCB'C' x and y displ. fixed
BC midplane w=0 # z displacements fixed along mid-plane

# MATERIAL PROPERTIES: isotropic single-material properties
E = 210e3 # Young modulus in MPa
nu = 0.3 # Poisson's ratio

SOLVE_PROBLEM # solve!

# print the direct stress y at D (and nothing more)
PRINT "σ_y @ D = " sigmay(2000,0,300) "MPa"
```

Figure 2: FeenoX input file edited in Kate on Plasma

## 5 Post-processors

FeenoX can write mesh results either in `.msh` or `.vtk` format. The former can be read and postprocessed by Gmsh. The latter can be read and postprocessed by a few different tools, but Paraview is the flagship postprocessor. In general any version will do, so it can be installed with

```
$ sudo apt-get install paraview
```

Also binaries and source versions can be downloaded.

Setting up your workspace

Note that both Gmsh in post-processing mode and Paraview make sense only if you have access to a graphical device.

## 6 Notes on hardware

### 6.1 Memory

Solving PDEs with FeenoX might need a lot of memory (depending on the problem size). Mind the total RAM your system has. If you are using a laptop, FeenoX might thrash it if the problem is way too big.

### 6.2 CPU

Solving problems with FeenoX during a lot of time might increase the CPU temperature significantly. Make sure your hardware is properly cooled before executing FeenoX during a long period of time.