



*Version 0.5.2*

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*In nova fert animus mutatas dicere formas  
corpora; di, coeptis (nam vos mutastis et illas)  
adspirate meis primumque ab origine mundi  
ad mea perpetuum deducite tempora carmen!*

—Ovid, *Metamorphoses*

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# Chapter 1

## Overview

*Morpho* aims to solve the following class of problems. Consider a functional,

$$F = \int_C f(q, \nabla q, \nabla^2 q, \dots) d^n x + \int_{\partial C} g(q, \nabla q, \nabla^2 q, \dots) d^{n-1} x,$$

where  $q$  represents a set of fields defined on a manifold  $C$  that could include scalar, vector, tensor or other quantities and their derivatives  $\nabla^n q$ . The functional includes terms in the bulk and on the boundary  $\partial C$  and might also include geometric properties of the manifold such as local curvatures. This functional is to be minimized from an initial guess  $\{C_0, q_0\}$  with respect to the fields  $q$  and the shape of the manifold  $C$ . Global and local constraints may be imposed both on  $C$  and  $q$ .

*Morpho* is an object-oriented environment: all components of the problem, including the computational domain, fields, functionals etc. are all represented as objects that interact with one another. Much of the effort in writing a *morpho* program involves creating and manipulating these objects. The environment is flexible, modular, and users can easily create new kinds of object, or entirely change how *morpho* works.

## Chapter 2

# Installing *Morpho*

*Morpho* is hosted on a publicly available github repository <https://github.com/Morpho-lang/morpho>. *Morpho* is presently installed from source; streamlined installation will be provided in future releases. Instructions for different platforms are provided below.

### 2.1 Dependencies

*Morpho* leverages a few libraries to provide certain functionality:

**glfw** is used to provide gui functionality for an interactive visualization application, **morphoview**.

**blas/lapack** are used for dense linear algebra.

**suitesparse** is used for sparse linear algebra.

### 2.2 MacOS

1. Install the Homebrew package manager, following instructions on the homebrew site.
2. Install dependencies. Open the Terminal application and type:

```
brew update  
brew install glfw suite-sparse
```

3. Obtain the source by cloning the github public repository:

```
git clone https://github.com/Morpho-lang/morpho.git
```

4. Navigate to the **morpho5** folder within the downloaded repository and build the application<sup>1</sup>

```
cd morpho/morpho5  
make install
```

5. Navigate to the **morphoview** folder and build the viewer application

```
cd morpho/morpho5  
make install
```

---

<sup>1</sup>Some users may need to use `sudo make install`

6. Check that the application works by typing

```
morpho5
```

## 2.3 Linux

Note that the build script places `morpho5` and `morphoview` in the `/usr/local` file structure; this can easily be changed if a different location is preferred.

1. Install *morpho*'s dependencies using your distribution's package manager (or manually if you prefer). For example, on Ubuntu you would type:

```
sudo apt install libglfw3
sudo apt install libsuitesparse-dev
sudo apt install liblapacke
```

2. Obtain the source by cloning the github public repository:

```
git clone https://github.com/Morpho-lang/morpho.git
```

3. Navigate to the `morpho5` folder within the downloaded repository and build the application:

```
cd morpho/morpho5
sudo make -f Makefile.linux install
```

4. Navigate to the `morphoview` folder and build the viewer application:

```
cd ../morphoview
sudo make -f Makefile.linux install
```

5. Check that the application works by typing

```
morpho5
```

## 2.4 Windows via Windows Subsystem for Linux (WSL)

### 2.4 Install WSL

The instructions to install the Ubuntu App are [here](#). Once the Ubuntu terminal is working in Windows, you can install *morpho* through it in almost the same way as a Linux system, with the addition of an X windows manager to handle visualizations. Unless mentioned otherwise, all the commands below are run in the Ubuntu terminal.

### 2.4 Install Morpho

1. Install the dependencies. You can install the dependencies using the Advanced Package Tool or `apt`. First update the `apt` package list and then update existing packages.

```
sudo apt update
sudo apt upgrade
```

The dependencies for morpho can be then installed as follows:

```
sudo apt install libglfw3-dev
sudo apt install libsuitesparse-dev
sudo apt install liblapacke-dev
sudo apt install povray
```

To build the code you will also need to install build-essentials:

```
sudo apt install build-essential
```

2. Obtain the morpho source by cloning the Morpho repository:

```
git clone https://github.com/Morpho-lang/morpho.git
```

3. Navigate to the morpho5 folder within the downloaded repository and build the application:

```
cd morpho/morpho5
sudo make -f Makefile.linux install
```

4. Navigate to the morphoview folder and build the viewer application:

```
cd ../morphoview
sudo make -f Makefile.linux install
```

5. Check that the application works by typing:

```
morpho5
```

## 2.4 Get visualization working in WSL

1. A window manager must be installed so that the WSL can create windows. On Windows, install VcXsrv. It shows up as XLaunch in the Windows start menu.
2. Open Xlaunch. Then,
  - (a) choose 'Multiple windows', set display number to 0, and hit 'Next'
  - (b) choose 'start no client' and hit 'Next'
  - (c) **Unselect** 'native opengl' and hit 'Next'
  - (d) Hit 'Finish'
3. In Ubuntu download a package containing a full suite of desktop utilities that allows for the use of windows.

```
sudo apt install ubuntu-desktop mesa-utils
```

Tell ubuntu which display to use

```
export DISPLAY=localhost:0
```

To set the DISPLAY variable on launch add the line

```
DISPLAY=localhost:0
```



to `~/ .bashrc`.

4. Test that the window system is working by running `glxgears`.
5. Test the thomson example program! Navigate to the thomson example in the examples directory and run it. If you are in the `morphoview` directory.

```
cd ../examples/thomson  
morpho5 thomson.morpho
```

This example starts with randomly distributed charges on a sphere and minimizing electric potential. It should generate an interactive figure of points on a sphere.

## Chapter 3

# Using *Morpho*

*Morpho* is a command line application, like `python` or `lua`. It can be used to run scripts or programs, which are generally given the *.morpho* file extension, or run interactively responding to user commands.

### 3.1 Running a program

To run a program, simply run `morpho` with the name of the file,

```
morpho5 script.morpho
```

### 3.2 Interactive mode

To use *morpho* interactively, simply load the *Terminal* application (or equivalent on your system) and type

```
morpho5
```

As shown in Fig. 3.2.1, you'll be greeted by a brief welcome and a prompt `>` inviting you to enter *morpho* commands. For now, try a classic:

```
print "Hello World"
```

which will display `Hello World` as output. More information about the *morpho* language is provided in the Reference section, especially chapter 5; if you're familiar with C-like languages such as C, C++, Java, Javascript, etc. things should be quite familiar.

To assist the user, the contents of the reference manual are available to the user in interactive mode as online help. To get help, simply type:

```
help
```

or even more briefly,

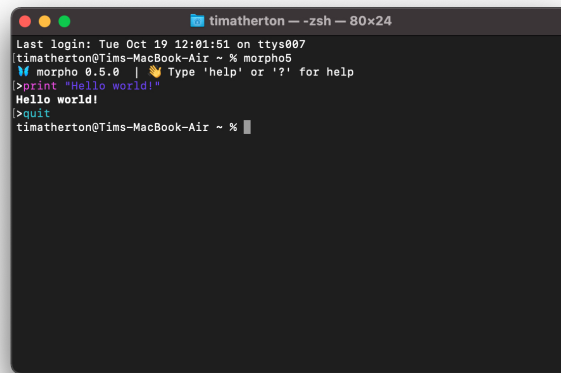
```
?
```

to see the list of main topics. To find help on a particular topic, for example `for` loops, simply type the topic name afterwards:

```
? for
```

Once you're done using *morpho*, simply type

```
quit
```



```
timatherton --zsh -- 80x24
Last login: Tue Oct 19 12:01:51 on ttys007
timatherton@Tims-MacBook-Air ~ % morpho5
morpho 0.5.0 | 🍌 Type 'help' or '?' for help
>print "Hello world!"
Hello world!
>quit
timatherton@Tims-MacBook-Air ~ %
```

Figure 3.2.1: Using *morpho* interactively from the command line.

to exit the program and return to the shell.

The interactive environment has a few other useful features to assist the user:

- **Autocomplete.** As you type, *morpho* will show you any suggested commands that it thinks you're trying to enter. For example, if you type `v` the command line will show the `var` keyword. To accept the suggestion, press the tab key. Multiple suggestions may be available; use the up and down arrow keys to rotate through them.
- **Command history.** Use the arrow keys to retrieve previously entered commands. You may then edit them before running them.
- **Line editing.** As you're typing a command, use the left and right arrows to move the cursor around; you can insert new characters at the cursor just by typing them or delete characters with the delete key. Hold down the shift key as you use the left and right arrow keys to select text; you can then use Ctrl-C to copy and Ctrl-V to paste. Ctrl-A moves to the start of the line and Ctrl-E the end.

## Chapter 4

# Tutorial

To illustrate how to use *morpho*, we will solve a problem involving nematic liquid crystals (NLCs), fluids composed of long, rigid molecules that possess a local average molecular orientation described by a unit vector field  $\hat{\mathbf{n}}$ . Droplets of NLC immersed in a host isotropic fluid such as water are called *tactoids* and, unlike droplets of, say, oil in water that form spheres, tactoids can adopt elongated shapes.

The functional to be minimized, the free energy of the system, is quite complex,

$$F = \underbrace{\frac{1}{2} \int_C K_{11} (\nabla \cdot \mathbf{n})^2 + K_{22} (\mathbf{n} \cdot \nabla \times \mathbf{n})^2 + K_{33} |\mathbf{n} \times \nabla \times \mathbf{n}|^2 dA}_{\text{Liquid crystal elastic energy}} + \underbrace{\sigma \int dl}_{s.t.} - \underbrace{\frac{W}{2} \int (\mathbf{n} \cdot \mathbf{t})^2 dl}_{\text{anchoring}} \quad (4.0.1)$$

where the three terms include **liquid crystal elasticity** that drives elongation of the droplet, **surface tension** (*s.t.*) that opposes lengthening of the boundary and an **anchoring term** that imposes a preferred orientation at the boundary. We need a local constraint,  $\mathbf{n} \cdot \mathbf{n} = 1$ , and will also impose a constraint on the volume of the droplet. For simplicity, we'll solve this problem in 2D. The complete code for this tutorial example is contained in the `examples/tactoid` folder in the repository.

### 4.1 Importing modules

*Morpho* is a modular system and hence we typically begin our program by telling *morpho* the modules we need so that they're available for us to use. To do so, we use the `import` keyword followed by the name of the module:

```
import meshtools
import optimize
import plot
```

We can also use the `import` keyword to import additional program files to assist in modularizing large programs. These are the modules we'll use for this example:

Module	Purpose
meshtools	Utility code to create and refine meshes
optimize	Perform optimization
plot	Visualize results

<pre> var a  declare a variable {  ← organize code in blocks   print foo print to the terminal } fn f(x) { define a function   return x^2 }      return a value  foo(f)  functions can be passed to other         functions!  parent class class Foo is Boo { define a class   init(p) { ...with methods like this initializer     self.prop = p   }      ← set and access object         properties }  a = Foo("hey") create an object a.foo() invoke a method </pre>	<pre> if (i&lt;1)  else  conditionally run code  for (a in b) {  loop over the contents                 of a collection  do {  } while (a&lt;b)  conventional loops while (a&lt;b) {  }    too  try {  } catch {  } deal with                     anticipated errors  a = [ 1, 2, 3 ]      lists b = { "a": 1, "b": 2 } ...and dictionaries  "Hello world \${i}"  strings (with interpolation)  m=Matrix(2,2)  dense and sparse matrices m[0,0]=1  set and access elements            of a collection by            indexing  import optimize  extensible,                   modular </pre>
--	---


Figure 4.2.1: Postcard-sized summary of the *morpho* language.

## 4.2 Morpho language

The *morpho* language is simple but expressive. If you're familiar with C-like languages (C, C++, Java, Javascript) you'll find it very natural. A much more detailed description is provided in Chapter 5, but a brief summary is provided in Fig. 4.2.1 and we provide an overview of key ideas to help you follow the tutorial:

- **Comments.** Any text after `//` or surrounded by `/*` and `*/` is a comment and not processed by morpho:

```
// This is a comment
/* This too! */
```

- **Variables.** To create a variable, use the `var` keyword; you can then assign and use the variable arbitrarily:

```
var a = 1
print a
```

- **Functions.** Functions may take parameters, and you call them like this:

```
print sin(x)
```

and declare them like this:

```
fn f(x,y) {
  return x^2+y^2
}
```

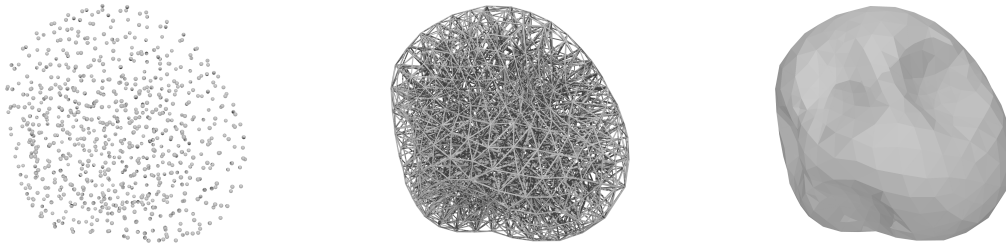


Figure 4.3.1: A *Mesh* object contains different kinds of element. In this example, the mesh contains points, lines and area elements referred to by their *grade*.

Some functions take optional arguments, which look like this:

```
var a = foo(quiet=true)
```

- **Objects.** *Morpho* is deeply object-oriented. Most things in *morpho* are represented as objects, which provide *methods* that you can use to control them. Objects are made by *constructor functions* that begin with a capital letter (and may take arguments):

```
var a = Object()
```

Method calls then look like this:

```
a.foo()
```

- **Collections.** *Morpho* provides a number of collection types—all of which are objects—including Lists,

```
var a = [1,2,3]
```

and Dictionaries:

```
var b = { "Massachusetts": "Boston", "California": "Sacramento" }
```

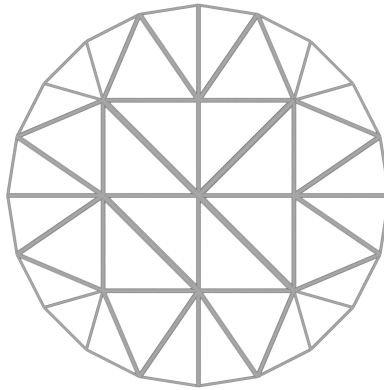
and Ranges (often used in loops):

```
var a = 0..10:2 # all even numbers 0-10 inclusive
```

There are many others, including Matrices, Sparse matrices, etc.

### 4.3 Creating the initial mesh

Mesheres are discretized regions of space. The very simplest region we can imagine is a *point* or *vertex* described by a set of coordinates  $(x_1, x_2, \dots, x_D)$  where the number of coordinates  $D$  defines the dimensionality of the space that the manifold is said to be *embedded* in. From more than one point, we can start constructing more complex regions. First, between two points we can imagine fixing an imaginary ruler and drawing a straight line or *edge* between them. Three points define a plane, and also a triangle; we can therefore identify the two dimensional area of the plane bounded by the triangle as a *face*, as in the face of a polyhedron. Using four points, we can define the volume bounded by a tetrahedron. Each of these **elements** has a different dimensionality—called a *grade*—and a complete **Mesh** may contain elements of many different grades as shown in Fig. 4.3.1.

Figure 4.3.2: The initial mesh, loaded from `disk.mesh`.

*Morpho* provides a number of ways of creating a mesh. One can load a mesh from a file, build one manually from a set of points, create one from a polyhedron, or from the level set (contours) of a function.

For this example, we'll use a predefined mesh file `disk.mesh`. To create a *Mesh* object from this file, we call the *Mesh* function with the file name:

```
var m = Mesh("disk.mesh")
```

Here, the **var** keyword tells *morpho* to create a new variable *m*, which now refers to the newly created *Mesh* object. The initial mesh is depicted in Fig. 4.3.2; we'll provide the code to perform the visualization in section 4.8.

If you open the file `disk.mesh`, which you can find in the same folder as `tactoid.morpho`, you'll find it has a simple human readable format:

```
vertices
```

```
1 -1. 0. 0
2 -0.951057 -0.309017 0
...
```

```
edges
```

```
1 8 2
2 2 4
...
```

```
faces
```

```
1 8 2 4
2 8 4 6
...
```

The file is broken into sections, each describing elements of a different grade. Each line begins either with a section delimiter such as *vertices*, *edges* or *faces*, or with an id. Vertices are then defined by a set of coordinates; edges and faces are defined by providing the respective vertex ids.

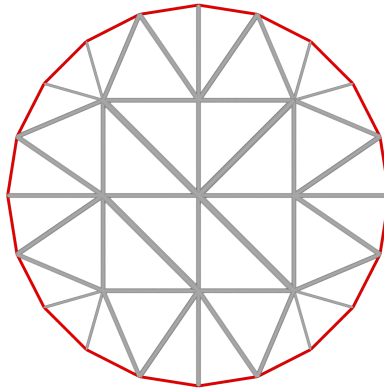


Figure 4.4.1: Selecting the boundary of the mesh.

## 4.4 Selections

Sometimes, we want to refer to specific parts of a `Mesh` object: elements that match some criterion, for example. `Selection` objects enable us to do this. Because selecting the boundary is a very common activity, the `Selection` constructor function takes an optional argument to do this:

```
var bnd=Selection(m, boundary=true)
```

By default, only the boundary elements are included in the `Selection`. For a mesh with at most grade 2 elements (facets), the boundaries are grade 1 elements (lines); for a mesh with grade 3 elements (volumes), the boundaries are grade 2 elements (facets). Quite often we want the vertices themselves as well, so we can call a method to achieve that:

```
bnd.addgrade(0)
```

Once a `Selection` has been created, it can be helpful to visualize it to ensure the correct elements are selected. We'll talk more about visualization in section 4.8, but for now the line

```
Show(plotselection(m, bnd, grade=1))
```

shows a visualization of the mesh with the selected grade 1 elements shaded red as displayed in Fig. 4.4.1.

## 4.5 Fields

Having created our initial computational domain, we will now create a `Field` object representing the director field  $\mathbf{n}$ :

```
var nn = Field(m, Matrix([1,0,0]))
```

As with the `Mesh` object earlier, we declare a variable,  $nn$ , to refer to the `Field` object. We have to provide two arguments to `Field`: the `Mesh` object on which the `Field` is defined, and something to initialize it. Here, we want the initial director to have a spatially uniform value, so we can just provide `Field` a constant `Matrix` object. By default, *morpho* stores a copy of this matrix on each vertex in the mesh; `Fields` can however store information on elements of any grade (and store both more than one quantity per grade and information on multiple grades at the same time).



It's possible to initialize a `Field` with spatially varying values by providing an *anonymous function* to `Field` like this:

```
var phi = Field(m, fn (x,y,z) x^2+y^2)
```

Here, *phi* is a scalar field that takes on the value  $x^2 + y^2$ . The **fn** keyword is used to define functions.

## 4.6 Defining the problem

We now turn to setting up the problem. Each term in the energy functional (4.0.1) is represented by a corresponding *functional* object, which acts on a `Mesh` (and possibly a `Field`) to calculate an integral quantity such as an energy; Functional objects are also responsible for calculating gradients of the energy with respect to vertex positions and components of `Fields`.

Let's take the terms in (4.0.1) one by one: To represent the nematic elasticity we create a `Nematic` object:

```
var lf=Nematic(nn)
```

The surface tension term involves the length of the boundary, so we need a `Length` object:

```
var lt=Length()
```

The anchoring term doesn't have a simple built in object type, but we can use a general `LineIntegral` object to achieve the correct result.

```
var la=LineIntegral(fn (x, n) n.inner(tangent())^2, nn)
```

Notice that we have to supply a function—the integrand—which will be called by `LineIntegral` when it evaluates the integral. Integrand functions are called with the local coordinates first (as a `Matrix` object representing a column vector) and then the local interpolated value of any number of `Fields`. We also make use of the special function `tangent()` that locally returns a local tangent to the line.

We also need to impose constraints. Any *functional* object can be used equally well as an energy or a constraint, and hence we create a `NormSq` (norm-squared) object that will be used to implement the local unit vector constraint on the director field:

```
var ln=NormSq(nn)
```

and an `Area` object for the global constraint. This is really a constraint fixing the volume of fluid in the droplet, but since we're in 2D that becomes a constraint on the area of the mesh:

```
var laa=Area()
```

Now we have a collection of functional objects that we can use to define the problem. So far, we haven't specified which functionals are energies and which are constraints; nor have we specified which parts of the mesh the functionals are to be evaluated over. All that information is collected in an `OptimizationProblem` object, which we will now create:

```
// Set up the optimization problem
```

```
var W = 1
```

```
var sigma = 1
```

```
var problem = OptimizationProblem(m)
```

```
problem.addenergy(lf)
```

```
problem.addenergy(la, selection=bnd, prefactor=-W/2)
```

```
problem.addenergy(lt, selection=bnd, prefactor=sigma)
```

```
problem.addconstraint(laa)
```

```
problem.addlocalconstraint(ln, field=nn, target=1)
```

Notice that some of these functionals only act on a selection such as the boundary and hence we use the optional `selection` parameter to specify this. We can also specify the prefactor of the functional.

## 4.7 Performing the optimization

We're now ready to perform the optimization, for which we need an `Optimizer` object. These come in two flavors: a `ShapeOptimizer` and a `FieldOptimizer` that respectively act on the shape and a field. We create them with the problem and quantity they're supposed to act on:

```
// Create shape and field optimizers
var sopt = ShapeOptimizer(problem, m)
var fopt = FieldOptimizer(problem, nn)
```

Having created these, we can perform the optimization by calling the `linesearch` method with a specified number of iterations for each:

```
// Optimization loop
for (i in 1..100) {
    fopt.linesearch(20)
    sopt.linesearch(20)
}
```

Each iteration of a `linesearch` evolves the field (or shape) down the gradient of the target functional, subject to constraints, and finds an optimal stepsize to reduce the value of the functional. Here, we alternate between optimizing the field and optimizing the shape, performing twenty iterations of each, and overall do this one hundred times. These numbers have been chosen rather arbitrarily, and if you look at the output you will notice that *morpho* doesn't always execute twenty iterations of each. Rather, at each iteration it checks to see if the change in energy satisfies,

$$|E| < \epsilon,$$

or,

$$\left| \frac{\Delta E}{E} \right| < \epsilon$$

where the value of  $\epsilon$ , the convergence tolerance can be changed by setting the `etol` property of the `Optimizer` object:

```
sopt.etol = 1e-7 // default value is 1e-8
```

Some other properties of an `Optimizer` that may be useful for the user to adjust are as follows:

Property	Default value	Purpose
<code>etol</code>	$1 \times 10^{-8}$	Energy tolerance (relative error)
<code>ctol</code>	$1 \times 10^{-10}$	Constraint tolerance (how well are constraints satisfied)
<code>stepsize</code>	0.1	Stepsize for <code>relax</code> (changed by <code>linesearch</code> )
<code>steplimit</code>	0.5	Largest stepsize a <code>linesearch</code> can take
<code>maxconstraintsteps</code>	20	Number of steps the optimizer may take to ensure constraints are satisfied
<code>quiet</code>	false	Whether to print output as the optimization happens

## 4.8 Visualizing results

*Morpho* provides a highly flexible graphics system, with an external viewer application *morphoview*, to enable rich visualizations of results. Visualizations typically involve one or more `Graphics` objects, which act as a container for graphical elements to be displayed. Various *graphics primitives*, such as spheres, cylinders, arrows, tubes, etc. can be added to a `Graphics` object to make a drawing.

We are now ready to visualize the results of the optimization. First, we'll draw the mesh. Because we're interested in seeing the mesh structure, we'll draw the edges (i.e. the grade 1 elements). The function to do this is provided as part of the `plot` module that we imported in section 4.1:

```
var g=plotmesh(m, grade=1)
```

Next, we'll create a separate `Graphics` object that contains the director. Since the director  $\mathbf{n}$  is a unit vector field, and the sign is not significant (the nematic elastic energy is actually invariant under  $\mathbf{n} \rightarrow -\mathbf{n}$ ), an appropriate way to display a single director is as a cylinder oriented along  $\mathbf{n}$ . We will therefore make a helper function that creates a `Graphics` object and draws such a cylinder at every mesh point:

```
// Function to visualize a director field
// m - the mesh
// nn - the director Field to visualize
// dl - scale the director
fn visualize(m, nn, dl) {
  var v = m.vertexmatrix()
  var nv = m.count() // Number of vertices
  var g = Graphics() // Create a graphics object
  for (i in 0...nv) {
    var x = v.column(i) // Get the ith vertex
    // Draw a cylinder aligned with nn at this vertex
    g.display(Cylinder(x-nn[i]*dl, x+nn[i]*dl, aspectratio=0.3))
  }
  return g
}
```

Once we've defined this function, we can use it:

```
var gnn=visualize(m, nn, 0.2)
```

The variables  $g$  and  $g_{nn}$  now refer to two separate `Graphics` objects. We can combine them using the  $+$  operator, and display them like so:

```
var gdisp = g+gnn
Show(gdisp)
```

The resulting visualization is shown in Fig. 4.8.1.

## 4.9 Refinement

We have now solved our first shape optimization problem, and the complete problem script is provided in the `examples/tutorial` folder inside the git repository as `tutorial.morpho`. The result we have obtained in Fig. 4.8.1 is, however, a very coarse, low resolution solution comprising only a relatively small number of elements. To gain an improved solution, we need to *refine* our mesh. Because modifying the mesh also

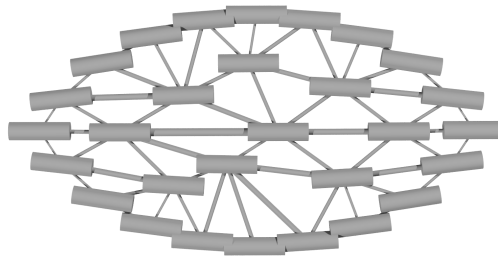


Figure 4.8.1: Optimized mesh and director field.

requires us to update other data structures like fields and selections, a special `MeshRefiner` object is used to perform the refinement.

To perform refinement we:

1. Create a `MeshRefiner` object, providing it a list of all the `Mesh`, `Field` and `Selection` objects (i.e. the mesh and objects that directly depend on it) that need to be updated:

```
var mr=MeshRefiner([m, nn, bnd]) // Set the refiner up
```

2. Call the `refine` method on the `MeshRefiner` object to actually perform the refinement. This method returns a `Dictionary` object that maps the old objects to potentially newly created ones.

```
var refmap=mr.refine() // Perform the refinement
```

3. Tell any other objects that refer to the mesh, fields or selections to update their references using `refmap`. For example, `OptimizationProblem` and `Optimizer` objects are typically updated at this step.

```
for (el in [problem, sopt, fopt]) el.update(refmap) // Update the problem
```

4. Update our own references

```
m=refmap[m]; nn=refmap[nn]; bnd=refmap[bnd] // Update variables
```

We insert this code after our optimization section, which causes *morpho* to successively optimize and refine<sup>1</sup>. The resulting optimized shapes are displayed in Fig. 4.9.1.

```
// Optimization loop
var refmax = 3
for (refiter in 1..refmax) {
  print "===Refinement level ${refiter}==="
  for (i in 1..100) {
```

<sup>1</sup>The complete code including refinement is in `examples/tutorial` folder inside the git repository as `tutorial2.morpho`

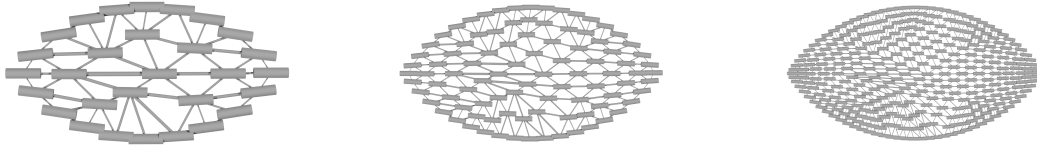


Figure 4.9.1: Optimized mesh and director field at three successive levels of refinement.

```

    fopt.linesearch(20)
    sopt.linesearch(20)
}

if (refiter==refmax) break

// Refinement
var mr=MeshRefiner([m, nn, bnd]) // Set the refiner up
var refmap=mr.refine() // Perform the refinement
for (el in [problem, sopt, fopt]) el.update(refmap) // Update the problem
m=refmap[m]; nn=refmap[nn]; bnd=refmap[bnd] // Update variables
}

```

## 4.10 Next steps

Having completed this tutorial, you may wish to explore the effect of changing some of the parameters in the file. What happens if you change  $\sigma$  and  $W$ , the coefficients in front of the terms in the energy? What happens if you take a different number of steps? Or change properties of the Optimizers like `stepsize` and `steplimit`?

You should look at other example files provided in the `examples` folder of the git repository. The remainder of the manual provides a detailed reference for *morpho* functionality, and a complete description of the scripting language.

## Reference

## Chapter 5

# Language

### 5.1 Syntax

Morpho provides a flexible object oriented language similar to other languages in the C family (like C++, Java and Javascript) with a simplified syntax.

Morpho programs are stored as plain text with the .morpho file extension. A program can be run from the command line by typing

```
morpho5 program.morpho
```

### 5.2 Comments

Two types of comment are available. The first type is called a ‘line comment’ whereby text after // on the same line is ignored by the interpreter.

```
a.dosomething() // A comment
```

Longer ‘block’ comments can be created by placing text between /\* and \*/. Newlines are ignored

```
/* This  
   is  
   a longer comment */
```

In contrast to C, these comments can be nested

```
/* A nested /* comment */ */
```

enabling the programmer to quickly comment out a section of code.

### 5.3 Symbols

Symbols are used to refer to named entities, including variables, classes, functions etc. Symbols must begin with a letter or underscore \_ as the first character and may include letters or numbers as the remainder. Symbols are case sensitive.

```
asymbol  
_alsoasymbol  
another_symbol  
EvenThis123
```

## YET\_ANOTHER\_SYMBOL

Classes are typically given names with an initial capital letter. Variable names are usually all lower case.

### 5.4 Newlines

Strictly, morpho ends statements with semicolons like C, but in practice these are usually optional and you can just start a new line instead. For example, instead of

```
var a = 1; // The ; is optional
```

you can simply use

```
var a = 1
```

If you want to put several statements on the same line, you can separate them with semicolons:

```
var a = 1; print a
```

There are a few edge cases to be aware of: The morpho parser works by accepting a newline anywhere it expects to find a semicolon. To split a statement over multiple lines, signal to morpho that you plan to continue by leaving the statement unfinished. Hence, do this:

```
print a +  
      1
```

rather than this:

```
print a    // < Morpho thinks this is a complete statement  
      + 1 // < and so this line will cause a syntax error
```

### 5.5 Booleans

Comparison operations like `==`, `<` and `>=` return `true` or `false` depending on the result of the comparison. For example,

```
print 1==2
```

prints `false`. The constants `true` or `false` are provided for you to use in your own code:

```
return true
```

### 5.6 Nil

The keyword `nil` is used to represent the absence of an object or value.

Note that in `if` statements, a value of `nil` is treated like `false`.

```
if (nil) {  
    // Never executed.  
}
```



## 5.7 Blocks

Code is divided into *blocks*, which are delimited by curly brackets like this:

```
{  
  var a = "Hello"  
  print a  
}
```

This syntax is used in function declarations, loops and conditional statements.

Any variables declared within a block become *local* to that block, and cannot be seen outside of it. For example,

```
var a = "Foo"  
{  
  var a = "Bar"  
  print a  
}  
print a
```

would print “Bar” then “Foo”; the version of *a* inside the code block is said to *shadow* the outer version.

## 5.8 Precedence

Precedence refers to the order in which morpho evaluates operations. For example,

```
print 1+2*3
```

prints 7 because  $2*3$  is evaluated before the addition; the operator  $*$  is said to have higher precedence than  $+$ .

You can always modify the order of evaluation by using parentheses:

```
print (1+2)*3 // prints 9
```

## 5.9 Print

The `print` keyword is used to print information to the console. It can be followed by any value, e.g.

```
print 1  
print true  
print a  
print "Hello"
```

## 5.10 Values

Values are the basic unit of information in morpho: All functions in morpho accept values as arguments and return values.

## 5.11 Int

Morpho provides integers, which work as you would expect in other languages, although you rarely need to worry about the distinction between floats and integers.

Convert a floating point number to an Integer:

```
print Int(1.3) // expect: 1
```

Convert a string to an integer:

```
print Int("10")+1 // expect: 11
```

## 5.12 Float

Morpho provides double precision floating point numbers.

Convert a string to an floating point number:

```
print Float("1.2e2")+1 // expect: 121
```

## 5.13 Ceil

Returns the smallest integer larger than or equal to its argument:

```
print ceil(1.3) // expect: 2
```

## 5.14 Floor

Returns the largest integer smaller than or equal to its argument:

```
print floor(1.3) // expect: 1
```

## 5.15 Variables

Variables are defined using the `var` keyword followed by the variable name:

```
var a
```

Optionally, an initial assignment may be given:

```
var a = 1
```

Variables defined in a block of code are visible only within that block, so

```
var greeting = "Hello"
{
  var greeting = "Goodbye"
  print greeting
}
print greeting
```

will print

*Goodbye Hello*

Multiple variables can be defined at once by separating them with commas

```
var a, b=2, c[2]=[1,2]
```

where each can have its own initializer (or not).

## 5.16 Indexing

Morpho provides a number of collection objects, such as `List`, `Range`, `Array`, `Dictionary`, `Matrix` and `Sparse`, that can contain more than one value. Index notation (sometimes called subscript notation) is used to access elements of these objects.

To retrieve an item from a collection, you use the `[` and `]` brackets like this:

```
var a = List("Apple", "Bag", "Cat")
print a[0]
```

which prints *Apple*. Note that the first element is accessed with `0` not `1`.

Similarly, to set an entry in a collection, use:

```
a[0]="Adder"
```

which would replace the first element in `a` with `"Adder"`.

Some collection objects need more than one index,

```
var a = Matrix([[1,0],[0,1]])
print a[0,0]
```

and others such as `Dictionary` use non-numerical indices,

```
var b = Dictionary()
b["Massachusetts"]="Boston"
b["California"]="Sacramento"
```

as in this dictionary of state capitals.

## 5.17 Control Flow

Control flow statements are used to determine whether and how many times a selected piece of code is executed. These include:

- `if` - Selectively execute a piece of code if a condition is met.
- `else` - Execute a different block of code if the test in an `if` statement fails.
- `for` - Repeatedly execute a section of code with a counter
- `while` - Repeatedly execute a section of code while a condition is true.

## 5.18 If

If allows you to selectively execute a section of code depending on whether a condition is met. The simplest version looks like this:

```
if (x<1) print x
```

where the body of the loop, `print x`, is only executed if `x` is less than 1. The body can be a code block to accommodate longer sections of code:

```
if (x<1) {
    ... // do something
}
```

If you want to choose between two alternatives, use `else`:

```
if (a==b) {
    // do something
} else {
    // this code is executed only if the condition is false
}
```

You can even chain multiple tests together like this:

```
if (a==b) {
    // option 1
} else if (a==c) {
    // option 2
} else {
    // something else
}
```

## 5.19 While

While loops repeat a section of code while a condition is true. For example,

```
var k=1
while (k <= 4) { print k; k+=1 }
    ^cond    ^body
```

prints the numbers 1 to 4. The loop has two sections: `cond` is the condition to be executed and `body` is the section of code to be repeated.

Simple loops like the above example, especially those that involve counting out a sequence of numbers, are more conveniently written using a `for` loop,

```
for (k in 1..4) print k
```

Where while loops can be very useful is where the state of an object is being changed in the loop, e.g.

```
var a = List(1,2,3,4)
while (a.count()>0) print a.pop()
```

which prints 4,3,2,1.

## 5.20 Do

A `do...while` loop repeats code while a condition is true—similar to a `while` loop—but the test happens at the end:

```
var k=1
do {
  print k;
  k+=1
} while (k<5)
```

which prints 1,2,3,4

Hence this type of loop executes at least one iteration

## 5.21 For

For loops allow you to repeatedly execute a section of code. They come in two versions: the simpler version looks like this,

```
for (var i in 1..5) print i
```

which prints the numbers 1 to 5 in turn. The variable `i` is the *loop variable*, which takes on a different value each iteration. `1..5` is a range, which denotes a sequence of numbers. The *body* of the loop, `print i`, is the code to be repeatedly executed.

Morpho will implicitly insert a `var` before the loop variable if it's missing, so this works too:

```
for (i in 1..5) print i
```

If you want your loop variable to count in increments other than 1, you can specify a stepsize in the range:

```
for (i in 1..5:2) print i
      ^step
```

Ranges need not be integer:

```
for (i in 0.1..0.5:0.1) print i
```

You can also replace the range with other kinds of collection object to loop over their contents:

```
var a = Matrix([1,2,3,4])
for (x in a) print x
```

Morpho iterates over the collection object using an integer *counter variable* that's normally hidden. If you want to know the current value of the counter (e.g. to get the index of an element as well as its value), you can use the following:

```
var a = [1, 2, 3]
for (x, i in a) print "${i}: ${x}"
```

Morpho also provides a second form of `for` loop similar to that in C:

```
for (var i=0; i<5; i+=1) { print i }
      ^start  ^test ^inc.  ^body
```

which is executed as follows: *start*: the variable `i` is declared and initially set to zero. *test*: before each iteration, the test is evaluated. If the test is `false`, the loop terminates. *body*: the body of the loop is executed. *inc*: the variable `i` is increased by 1.

You can include any code that you like in each of the sections.

## 5.22 Break

Break is used inside loops to finish the loop early. For example

```
for (i in 1..5) {
  if (i>3) break // --.
  print i        // | (Once i>3)
}                // |
...             // <-'
```

would only print 1, 2 and 3. Once the condition `i>3` is true, the `break` statement causes execution to continue after the loop body.

Both `for` and `while` loops support `break`.

## 5.23 Continue

Continue is used inside loops to skip over the rest of an iteration. For example

```
for (i in 1..5) { // <-.
  print "Hello"   |
  if (i>3) continue // --'
  print i
}
```

prints “Hello” five times but only prints 1, 2 and 3. Once the condition `i>3` is true, the `continue` statement causes execution to transfer to the start of the loop body.

Traditional `for` loops also support `continue`:

```
                // v increment
for (var i=0; i<5; i+=1) {
  if (i==2) continue
  print i
}
```

Since `continue` causes control to be transferred *to the increment section* in this kind of loop, here the program prints 0..4 but the number 2 is skipped.

Use of `continue` with `while` loops is possible but isn’t recommended as it can easily produce an infinite loop!

```
var i=0
while (i<5) {
  if (i==2) continue
  print i
  i+=1
}
```

In this example, when the condition `i==2` is true, execution skips back to the start, but `i` *isn’t* incremented. The loop gets stuck in the iteration `i==2`.

## 5.24 Try

A `try` and `catch` statement allow you handle errors. For example

```
try {
  // Do something
} catch {
  "Tag" : // Handle the error
}
```

Code within the block after the `try` keyword is executed. If an error is generated then Morpho looks to see if the tag associated with the error matches any of the labels in the `catch` block. If it does, the code after the matching label is executed. If no error occurs, the `catch` block is skipped entirely.

## 5.25 Functions

A function in morpho is defined with the `fn` keyword, followed by the function's name, a list of parameters enclosed in parentheses, and the body of the function in curly braces. This example computes the square of a number:

```
fn sqr(x) {
  return x*x
}
```

Once a function has been defined you can evaluate it like any other morpho function.

```
print sqr(2)
```

## 5.26 Return

The `return` keyword is used to exit from a function, optionally passing a given value back to the caller. `return` can be used anywhere within a function. The below example calculates the  $n$ th Fibonacci number,

```
fn fib(n) {
  if (n<2) return n
  return fib(n-1) + fib(n-2)
}
```

by returning early if  $n < 2$ , otherwise returning the result by recursively calling itself.

## 5.27 Closures

Functions in morpho can form *closures*, i.e. they can enclose information from their local context. In this example,

```
fn foo(a) {
  fn g() { return a }
  return g
}
```

the function `foo` returns a function that captures the value of `a`. If we now try calling `foo` and then calling the returned functions,

```
var p=foo(1), q=foo(2)
print p() // expect: 1
print q() // expect: 2
```

we can see that `p` and `q` seem to contain different copies of `g` that encapsulate the value that `foo` was called with.

Morpho hints that a returned function is actually a closure by displaying it with double brackets:

```
print foo(1) // expect: <<fn g>>
```

## 5.28 Classes

Classes are defined using the `class` keyword followed by the name of the class. The definition includes methods that the class responds to. The special `init` method is called whenever an object is created.

```
class Cake {
  init(type) {
    self.type = type
  }

  eat() {
    print "A delicious "+self.type+" cake"
  }
}
```

Objects are created by calling the class as if it was a function:

```
var c = Cake("carrot")
```

Methods are called using the `.` operator:

```
c.eat()
```

## 5.29 Self

The `self` keyword is used to access an object's properties and methods from within its definition.

```
class Vehicle {
  init (type) { self.type = type }

  drive () { print "Driving my ${self.type}." }
}
```

## 5.30 Super

The keyword `super` allows you to access methods provided by an object's superclass rather than its own. This is particularly useful when the programmer wants a class to extend the functionality of a parent class, but needs to make sure the old behavior is still maintained.

For example, consider the following pair of classes:

```
class Lunch {
  init(type) { self.type=type }
}
```



```
class Soup < Lunch {
    init(type) {
        print "Delicious soup!"
        super.init(type)
    }
}
```

The subclass Soup uses `super` to call the original initializer.

### 5.31 Modules

Morpho is extensible and provides a convenient module system that works like standard libraries in other languages. Modules may define useful variables, functions and classes, and can be made available using the `import` keyword. For example,

```
import color
```

loads the `color` module that provides functionality related to color.

You can create your own modules; they're just regular morpho files that are stored in a standard place. On UNIX platforms, this is `/usr/local/share/morpho/modules`.

### 5.32 Import

Import provides access to the module system and including code from multiple source files.

To import code from another file, use `import` with the filename:

```
import "file.morpho"
```

which immediately includes all the contents of `"file.morpho"`. Any classes, functions or variables defined in that file can now be used, which allows you to divide your program into multiple source files.

Morpho provides a number of built in modules—and you can write your own—which can be loaded like this:

```
import color
```

which imports the `color` module.

You can selectively import symbols from a modules by using the `for` keyword:

```
import color for HueMap, Red
```

which imports only the `HueMap` class and the `Red` variable.

### 5.33 Help

Morpho provides an online help system. To get help about a topic called `topicname`, type

```
help topicname
```

A list of available topics is provided below and includes language keywords like `class`, `fn` and `for`, built in classes like `Matrix` and `File` or information about functions like `exp` and `random`.

Some topics have additional subtopics: to access these type

```
help topic subtopic
```

For example, to get help on a method for a particular class, you could type

```
help Classname.methodname
```

Note that `help` ignores all punctuation.

You can also use `?` as a shorthand synonym for `help`

```
? topic
```

A useful feature is that, if an error occurs, simply type `help` to get more information about the error.

## 5.34 Builtin functions

Morpho provides a number of built-in functions.

### 5.35 arctan

Returns the arctangent of an input value that lies from `-Inf` to `Inf`. You can use one argument:

```
print arctan(0) // expect: 0
```

or use two arguments to return the angle in the correct quadrant:

```
print arctan(x, y)
```

Note the order `x, y` differs from some other languages.

### 5.36 isnil

Returns `true` if a value is `nil` or `false` otherwise.

### 5.37 isint

Returns `true` if a value is an integer or `false` otherwise.

### 5.38 isfloat

Returns `true` if a value is a floating point number or `false` otherwise.

### 5.39 isbool

Returns `true` if a value is a boolean or `false` otherwise.

### 5.40 isobject

Returns `true` if a value is an object or `false` otherwise.

### 5.41 isstring

Returns `true` if a value is a string or `false` otherwise.

### 5.42 isclass

Returns `true` if a value is a class or `false` otherwise.

### 5.43 isrange

Returns `true` if a value is a range or `false` otherwise.

### 5.44 isdictionary

Returns `true` if a value is a dictionary or `false` otherwise.

### 5.45 islist

Returns `true` if a value is a list or `false` otherwise.

### 5.46 isarray

Returns `true` if a value is an array or `false` otherwise.

### 5.47 ismatrix

Returns `true` if a value is a matrix or `false` otherwise.

### 5.48 issparse

Returns `true` if a value is a sparse matrix or `false` otherwise.

### 5.49 isinf

Returns `true` if a value is infinite or `false` otherwise.

### 5.50 isnan

Returns `true` if a value is a Not a Number or `false` otherwise.

### 5.51 iscallable

Returns `true` if a value is callable or `false` otherwise.

## 5.52 Apply

Apply calls a function with the arguments provided as a list:

```
apply(f, [0.5, 0.5]) // calls f(0.5, 0.5)
```

It's often useful where a function or method and/or the number of parameters isn't known ahead of time. The first parameter to apply can be any callable object, including a method invocation or a closure.

You may also instead omit the list and use apply with multiple arguments:

```
apply(f, 0.5, 0.5) // calls f(0.5, 0.5)
```

There is one edge case that occurs when you want to call a function that accepts a single list as a parameter. In this case, enclose the list in another list:

```
apply(f, [[1,2]]) // equivalent to f([1,2])
```

## Chapter 6

# Data Types

### 6.1 Array

Arrays are collection objects that can have any number of indices. Their size is set when they are created:

```
var a[5]
var b[2,2]
var c[nv,nv,nv]
```

Values can be retrieved with appropriate indices:

```
print a[0,0]
```

Array can be indexed with slices:

```
print a[[0,2,4],2]
print a[1,0..2]
```

Any morpho value can be stored in an array element

```
a[0,0] = [1,2,3]
```

### 6.2 Dimensions

Get the dimensions of an Array object:

```
var a[2,2]
print a.dimensions() // expect: [ 2, 2 ]
```

### 6.3 Dictionary

Dictionaries are collection objects that associate a unique *key* with a particular *value*. Keys can be any kind of morpho value, including numbers, strings and objects.

An example dictionary mapping states to capitals:

```
var dict = { "Massachusetts" : "Boston",
             "New York" : "Albany",
             "Vermont" : "Montpelier" }
```

Look up values by a given key with index notation:

```
print dict["Vermont"]
```

You can change the value associated with a key, or add new elements to the dictionary like this:

```
dict["Maine"]="Augusta"
```

Create an empty dictionary using the Dictionary constructor function:

```
var d = Dictionary()
```

Loop over keys in a dictionary:

```
for (k in dict) print k
```

## 6.4 List

Lists are collection objects that contain a sequence of values each associated with an integer index.

Create a list like this:

```
var list = [1, 2, 3]
```

Look up values using index notation:

```
list[0]
```

Indexing can also be done with slices: `list[0..2]` `list[[0,1,3]]`

You can change list entries like this:

```
list[0] = "Hello"
```

Create an empty list:

```
var list = []
```

Loop over elements of a list:

```
for (i in list) print i
```

## 6.5 Append

Adds an element to the end of a list:

```
var list = []  
list.append("Foo")
```

## 6.6 Insert

Inserts an element into a list at a specified index:

```
var list = [1,2,3]  
list.insert(1, "Foo")  
print list // prints [ 1, Foo, 2, 3 ]
```

## 6.7 Pop

Remove the last element from a list, returning the element removed:

```
print list.pop()
```

If an integer argument is supplied, returns and removes that element:

```
var a = [1,2,3]
print a.pop(1) // prints '2'
print a        // prints [ 1, 3 ]
```

## 6.8 Sort

Sorts the contents of a list into ascending order:

```
list.sort()
```

Note that this sorts the list “in place” (i.e. it modifies the order of the list on which it is invoked) and hence returns `nil`.

You can provide your own function to use to compare values in the list

```
list.sort(fn (a, b) a-b)
```

This function should return a negative value if  $a < b$ , a positive value if  $a > b$  and `0` if  $a$  and  $b$  are equal.

## 6.9 Order

Returns a list of indices that would, if used in order, would sort a list. For example

```
var list = [2,3,1]
print list.order() // expect: [2,0,1]

    would produce [2,0,1]
```

## 6.10 Remove

Remove any occurrences of a value from a list:

```
var list = [1,2,3]
list.remove(1)
```

## 6.11 ismember

Tests if a value is a member of a list:

```
var list = [1,2,3]
print list.ismember(1) // expect: true
```

## 6.12 Add

Join two lists together:

```
var l1 = [1,2,3], l2 = [4, 5, 6]
print l1+l2 // expect: [1,2,3,4,5,6]
```

## 6.13 Tuples

Generate all possible n-tuples from a list:

```
var t = [ 1, 2, 3].tuples(2)
    produces [ [ 1, 1 ], [ 1, 2 ], [ 1, 3 ] ... ].
```

## 6.14 Sets

Generate all possible sets of order n from a list.

```
var t = [ 1, 2, 3 ].tuples(2)
    produces [ [ 1, 2 ], [ 1, 3 ], [ 2, 3 ] ].
```

Note that sets include only distinct elements from the list (no element is repeated) and ordering is unimportant, hence only one of [ 1, 2 ] and [ 2, 1 ] is returned.

## 6.15 Matrix

The Matrix class provides support for matrices. A matrix can be initialized with a given size,

```
var a = Matrix(nrows,ncols)
```

where all elements are initially set to zero. Alternatively, a matrix can be created from an array,

```
var a = Matrix([[1,2], [3,4]])
```

You can create a column vector like this,

```
var v = Matrix([1,2])
```

Once a matrix is created, you can use all the regular arithmetic operators with matrix operands, e.g.

```
a+b
a*b
```

The division operator is used to solve a linear system, e.g.

```
var a = Matrix([[1,2],[3,4]])
var b = Matrix([1,2])

print b/a
```

yields the solution to the system  $a \cdot x = b$ .



## 6.16 Range

Ranges represent a sequence of numerical values. There are two ways to create them depending on whether the upper value is included or not:

```
var a = 1..5 // inclusive version, i.e. [1,2,3,4,5]
var b = 1...5 // exclusive version, i.e. [1,2,3,4]
```

By default, the increment between values is 1, but you can use a different value like this:

```
var a = 1..5:0.5 // 1 - 5 with an increment of 0.5.
```

You can also create Range objects using the appropriate constructor function:

```
var a = Range(1,5,0.5)
```

Ranges are particularly useful in writing loops:

```
for (i in 1..5) print i
```

They can easily be converted to a list of values:

```
var c = List(1..5)
```

To find the number of elements in a Range, use the `count` method

```
print (1..5).count()
```

## 6.17 Sparse

The `Sparse` class provides support for sparse matrices. An empty sparse matrix can be initialized with a given size,

```
var a = Sparse(nrows,ncols)
```

Alternatively, a matrix can be created from an array of triplets,

```
var a = Sparse([[row, col, value] ...])
```

For example,

```
var a = Sparse([[0,0,2], [1,1,-2]])
```

creates the matrix

```
[ 2 0 ]
[ 0 -2 ]
```

Once a sparse matrix is created, you can use all the regular arithmetic operators with matrix operands, e.g.

```
a+b
a*b
```

## 6.18 String

Strings represent textual information. They are written in Morpho like this:

```
var a = "hello world"
```

Unicode characters including emoji are supported.

You can also create strings using the constructor function `String`, which takes any number of parameters:

```
var a = String("Hello", "World")
```

A very useful feature, called *string interpolation*, enables the results of any morpho expression can be interpolated into a string. Here, the values of `i` and `func(i)` will be inserted into the string as it is created:

```
print "${i}: ${func(i)}"
```

To get an individual character, use index notation

```
print "morpho"[0]
```

You can loop over each character like this:

```
for (c in "morpho") print c
```

Note that strings are immutable, and hence

```
var a = "morpho"  
a[0] = 4
```

raises an error.

## 6.19 split

The `split` method splits a `String` into a list of substrings. It takes one argument, which is a string of characters to use to split the string:

```
print "1,2,3".split(",")
```

gives

```
[ 1, 2, 3 ]
```

## Chapter 7

# Computational Geometry

### 7.1 Field

Fields are used to store information, including numbers or matrices, associated with the elements of a `Mesh` object.

You can create a `Field` by applying a function to each of the vertices,

```
var f = Field(mesh, fn (x, y, z) x+y+z)
```

or by supplying a single constant value,

```
var f = Field(mesh, Matrix([1,0,0]))
```

Fields can then be added and subtracted using the `+` and `-` operators.

To access elements of a `Field`, use index notation:

```
print f[grade, element, index]
```

where `* grade` is the grade to select `* element` is the element id `* index` is the element index

As a shorthand, it's possible to omit the grade and index; these are then both assumed to be `0`:

```
print f[2]
```

### 7.2 Grade

To create fields that include grades other than just vertices, use the `grade` option to `Field`. This can be just a grade index,

```
var f = Field(mesh, 0, grade=2)
```

which creates an empty field with `0` for each of the facets of the mesh `mesh`.

You can store more than one item per element by supplying a list to the `grade` option indicating how many items you want to store on each grade. For example,

```
var f = Field(mesh, 1.0, grade=[0,2,1])
```

stores two numbers on the line (grade 1) elements and one number on the facets (grade 2) elements. Each number in the field is initialized to the value `1.0`.

### 7.3 Shape

The `shape` method returns a list indicating the number of items stored on each element of a particular grade. This has the same format as the list you supply to the `grade` option of the `Field` constructor. For example,

```
[1,0,2]
```

would indicate one item stored on each vertex and two items stored on each facet.

### 7.4 Op

The `op` method applies a function to every item stored in a `Field`, returning the result as elements of a new `Field` object. For example,

```
f.op(fn (x) x.norm())
```

calls the `norm` method on each element stored in `f`.

Additional `Field` objects may be supplied as extra arguments to `op`. These must have the same shape (the same number of items stored on each grade). The function supplied to `op` will now be called with the corresponding element from each field as arguments. For example,

```
f.op(fn (x,y) x.inner(y), g)
```

calculates an elementwise inner product between the elements of Fields `f` and `g`.

### 7.5 Functionals

A number of functionals are available in `Morpho`. Each of these represents an integral over some `Mesh` and `Field` objects (on a particular `Selection`) and are used to define energies and constraints in an `OptimizationProblem` provided by the `optimize` module.

Many functionals are built in. Additional functionals are available by importing the `functionals` module:

```
import functionals
```

Functionals provide a number of standard methods:

- `total(mesh)` - returns the value of the integral with a provided mesh, selection and fields
- `integrand(mesh)` - returns the contribution to the integral from each element
- `gradient(mesh)` - returns the gradient of the functional with respect to vertex motions.
- `fieldgradient(mesh, field)` - returns the gradient of the functional with respect to components of the field

Each of these may be called with a mesh, a field and a selection.

### 7.6 Length

A `Length` functional calculates the length of a line element in a mesh.

Evaluate the length of a circular loop:

```
import constants
import meshtools
var m = LineMesh(fn (t) [cos(t), sin(t), 0], 0...2*Pi:Pi/20, closed=true)
var le = Length()
print le.total(m)
```

## 7.7 AreaEnclosed

An AreaEnclosed functional calculates the area enclosed by a loop of line elements.

```
var la = AreaEnclosed()
```

## 7.8 Area

An Area functional calculates the area of the area elements in a mesh:

```
var la = Area()
print la.total(mesh)
```

## 7.9 VolumeEnclosed

A VolumeEnclosed functional is used to calculate the volume enclosed by a surface. Note that this estimate may become inaccurate for highly deformed surfaces.

```
var lv = VolumeEnclosed()
```

## 7.10 Volume

A Volume functional calculates the volume of volume elements.

```
var lv = Volume()
```

## 7.11 ScalarPotential

The ScalarPotential functional is applied to point elements.

```
var ls = ScalarPotential(potential, gradient)
```

You must supply two functions (which may be anonymous) that return the potential and gradient respectively.

This functional is often used to constrain the mesh to the level set of a function. For example, to confine a set of points to a sphere:

```
import optimize
fn sphere(x,y,z) { return x^2+y^2+z^2-1 }
fn grad(x,y,z) { return Matrix([2*x, 2*y, 2*z]) }
var lsph = ScalarPotential(sphere, grad)
problem.addlocalconstraint(lsph)
```

See the thomson example for use of this technique.

## 7.12 LinearElasticity

The `LinearElasticity` functional measures the linear elastic energy away from a reference state.

You must initialize with a reference mesh:

```
var le = LinearElasticity(mref)
```

Manually set the poisson's ratio and grade to operate on:

```
le.poissonratio = 0.2
le.grade = 2
```

## 7.13 EquiElement

The `EquiElement` functional measures the discrepancy between the size of elements adjacent to each vertex. It can be used to equalize elements for regularization purposes.

## 7.14 LineCurvatureSq

The `LineCurvatureSq` functional measures the integrated curvature squared of a sequence of line elements.

## 7.15 LineTorsionSq

The `LineTorsionSq` functional measures the integrated torsion squared of a sequence of line elements.

## 7.16 MeanCurvatureSq

The `MeanCurvatureSq` functional computes the integrated mean curvature over a surface.

## 7.17 GaussCurvature

The `GaussCurvature` computes the integrated gaussian curvature over a surface.

## 7.18 GradSq

The `GradSq` functional measures the integral of the gradient squared of a field. The field can be a scalar, vector or matrix function.

Initialize with the required field:

```
var le=GradSq(phi)
```

## 7.19 Nematic

The `Nematic` functional measures the elastic energy of a nematic liquid crystal.

```
var lf=Nematic(nn)
```

There are a number of optional parameters that can be used to set the splay, twist and bend constants:

```
var lf=Nematic(nn, ksplay=1, ktwist=0.5, kbend=1.5, pitch=0.1)
```

These are stored as properties of the object and can be retrieved as follows:

```
print lf.ksplay
```

## 7.20 NematicElectric

The `NematicElectric` functional measures the integral of a nematic and electric coupling term  $\int (n \cdot E)^2$  where the electric field  $E$  may be computed from a scalar potential or supplied as a vector.

Initialize with a director field `nn` and a scalar potential `phi`: `var lne = NematicElectric(nn, phi)`

## 7.21 NormSq

The `NormSq` functional measures the elementwise L2 norm squared of a field.

## 7.22 LineIntegral

The `LineIntegral` functional computes the line integral of a function. You supply an integrand function that takes a position matrix as an argument.

To compute  $\int (x^2 + y^2)$  over a line element:

```
var la=LineIntegral(fn (x) x[0]^2+x[1]^2)
```

The function `tangent()` returns a unit vector tangent to the current element:

```
var la=LineIntegral(fn (x) x.inner(tangent()))
```

You can also integrate functions that involve fields:

```
var la=LineIntegral(fn (x, n) n.inner(tangent()), n)
```

where `n` is a vector field. The local interpolated value of this field is passed to your integrand function. More than one field can be used; they are passed as arguments to the integrand function in the order you supply them to `LineIntegral`.

## 7.23 AreaIntegral

The `AreaIntegral` functional computes the area integral of a function. You supply an integrand function that takes a position matrix as an argument.

To compute  $\int (x \cdot y)$  over an area element:

```
var la=AreaIntegral(fn (x) x[0]*x[1])
```

You can also integrate functions that involve fields:

```
var la=AreaIntegral(fn (x, phi) phi^2, phi)
```

More than one field can be used; they are passed as arguments to the integrand function in the order you supply them to `AreaIntegral`.

## 7.24 FloryHuggins

The FloryHuggins functional computes the Flory-Huggins mixing energy over an element:

```
a*phi*log(phi) + b*(1-phi)+log(1-phi) + c*phi*(1-phi)
```

where a, b and c are parameters you can supply. The value of phi is calculated from a reference mesh that you provide on initializing the Functional:

```
var lfh = FloryHuggins(mref)
```

Manually set the coefficients and grade to operate on:

```
lfh.a = 1; lfh.b = 1; lfh.c = 1;
lfh.grade = 2
```

## 7.25 Mesh

The Mesh class provides support for meshes. Meshes may consist of different kinds of element, including vertices, line elements, facets or area elements, tetrahedra or volume elements.

To create a mesh, you can import it from a file:

```
var m = Mesh("sphere.mesh")
```

or use one of the functions available in `meshtools` or `implicitmesh` packages.

Each type of element is referred to as belonging to a different grade. Point-like elements (vertices) are *grade 0*; line-like elements (edges) are *grade 1*; area-like elements (facets; triangles) are *grade 2* etc.

The `plot` package includes functions to visualize meshes.

## 7.26 Save

Saves a mesh as a `.mesh` file.

```
m.save("new.mesh")
```

## 7.27 Vertexposition

Retrieves the position of a vertex given an id:

```
print m.vertexposition(id)
```

## 7.28 Setvertexposition

Sets the position of a vertex given an id and a position vector:

```
print m.setvertexposition(1, Matrix([0,0,0]))
```



## 7.29 Addgrade

Adds a new grade to a mesh. This is commonly used when, for example, a mesh file includes facets but not edges. To add the missing edges:

```
m.addgrade(1)
```

## 7.30 Addsymmetry

Adds a symmetry to a mesh. Experimental in version 0.5.

## 7.31 Maxgrade

Returns the highest grade element present:

```
print m.maxgrade()
```

## 7.32 Count

Counts the number of elements. If no argument is provided, returns the number of vertices. Otherwise, returns the number of elements present of a given grade:

```
print m.count(2) // Returns the number of area-like elements.
```

## 7.33 Selection

The Selection class enables you to select components of a mesh for later use. You can supply a function that is applied to the coordinates of every vertex in the mesh, or select components like boundaries.

Create an empty selection:

```
var s = Selection(mesh)
```

Select vertices above the  $z=0$  plane using an anonymous function:

```
var s = Selection(mesh, fn (x,y,z) z>0)
```

Select the boundary of a mesh:

```
var s = Selection(mesh, boundary=true)
```

Selection objects can be composed using set operations:

```
var s = s1.union(s2)
```

or `var s = s1.intersection(s2)`

To add additional grades, use the addgrade method. For example, to add areas: `s.addgrade(2)`

### 7.34 addgrade

Adds elements of the specified grade to a Selection. For example, to add edges to an existing selection, use

```
s.addgrade(1)
```

By default, this only adds an element if *all* vertices in the element are currently selected. Sometimes, it's useful to be able to add elements for which only some vertices are selected. The optional argument `partials` allows you to do this:

```
s.addgrade(1, partials=true)
```

Note that this method modifies the existing selection, and does not generate a new Selection object.

### 7.35 removegrade

Removes elements of the specified grade from a Selection. For example, to remove edges from an existing selection, use

```
s.removegrade(1)
```

Note that this method modifies the existing selection, and does not generate a new Selection object.

### 7.36 idlistforgrade

Returns a list of element ids included in the selection.

To find out which edges are selected:

```
var edges = s.idlistforgrade(1)
```

### 7.37 isselected

Checks if an element id is selected, returning `true` or `false` accordingly.

To check if edge number 5 is selected:

```
var f = s.isselected(1, 5))
```

## Chapter 8

# I/O

### 8.1 File

The `File` class provides the capability to read from and write to files, or to obtain the contents of a file in convenient formats.

To open a file, create a `File` object with the filename as the argument

```
var f = File("myfile.txt")
```

which opens "myfile.txt" for *reading*. To open a file for writing or appending, you need to provide a mode selector

```
var g = File("myfile.txt", "write")
```

or

```
var g = File("myfile.txt", "append")
```

Once the file is open, you can then read or write by calling appropriate methods:

```
f.lines()           // reads the contents of the file into an array of lines.
f.readline()        // reads a single line
f.readchar()         // reads a single character.
f.write(string)      // writes the arguments to the file.
```

After you're done with the file, close it with

```
f.close()
```

### 8.2 lines

Returns the contents of a file as an array of strings; each element corresponds to a single line.

Read in the contents of a file and print line by line:

```
var f = File("input.txt")
var s = f.lines()
for (i in s) print i
f.close()
```

### 8.3 readline

Reads a single line from a file; returns the result as a string.

Read in the contents of a file and print each line:

```
var f = File("input.txt")
while (!f.eof()) {
    print f.readline()
}
f.close()
```

### 8.4 readchar

Reads a single character from a file; returns the result as a string.

### 8.5 write

Writes to a file.

Write the contents of a list to a file:

```
var f = File("output.txt", "w")
for (k, i in list) f.write("${i}: ${k}")
f.close()
```

### 8.6 close

Closes an open file.

### 8.7 eof

Returns true if at the end of the file; false otherwise

## Chapter 9

# Modules

### 9.1 Color

The `color` module provides support for working with color. Colors are represented in morpho by `Color` objects. The module predefines some colors including `Red`, `Green`, `Blue`, `Black`, `White`.

To use the module, use `import` as usual:

```
import color
```

Create a `Color` object from an RGB pair:

```
var col = Color(0.5,0.5,0.5) // A 50% gray
```

### 9.2 Colormap

The `color` module provides `ColorMaps` which are subclasses of `Color` that map a single parameter in the range 0..1 onto a continuum of colors. These include `GradientMap`, `GrayMap` and `HueMap`. Colors and `Colormaps` have the same interface.

Get the red, green or blue components of a color or colormap:

```
var col = HueMap()
print col.red(0.5) // argument can be in range 0..1
```

Get all three components as a list:

```
col.rgb(0)
```

Create a grayscale:

```
var c = Gray(0.2) // 20% gray
```

### 9.3 RGB

Gets the `rgb` components of a `Color` or `ColorMap` object as a list. Takes a single argument in the range 0..1, although the result will only depend on this argument if the object is a `ColorMap`.

```
var col = Color(0.1,0.5,0.7)
print col.rgb(0)
```

## 9.4 Constants

The `constants` module contains a number of useful mathematical and physical constants. Import it like any other module:

```
import constants
```

Available constants:

- `E` the base of natural logarithms.
- `Pi` ratio of the perimeter of a circle to its diameter.

## 9.5 Delaunay

The `delaunay` module creates Delaunay triangulations from point clouds. It is dimensionally independent, so generates tetrahedra in 3D and higher order simplices beyond.

To use the module, first import it:

```
import delaunay
```

To create a Delaunay triangulation from a list of points:

```
var pts = []
for (i in 0...100) pts.append(Matrix([random(), random()])))
var del=Delaunay(pts)
print del.triangulate()
```

The module also provides `DelaunayMesh` to directly create meshes from Delaunay triangulations.

## 9.6 Triangulate

The `triangulate` method performs the delaunay triangulation. To use it, first construct a `Delaunay` object with the point cloud of interest:

```
var del=Delaunay(pts)
```

Then call `triangulate`:

```
var tri = del.triangulate()
```

This returns a list of triangles `[ [i, j, k], ... ]`.

## 9.7 DelaunayMesh

The `DelaunayMesh` constructor function creates a `Mesh` object directly from a point cloud using the Delaunay triangulator.

```
var pts = []
for (i in 0...100) pts.append(Matrix([random(), random()])))
var m=DelaunayMesh(pts)
Show(plotmesh(m))
```

You can control the output dimension of the mesh (e.g. to create a 2D mesh embedded in 3D space) using the optional `outputdim` property.

```
var m = DelaunayMesh(pts, outputdim=3)
```

## 9.8 Circumsphere

The `Circumsphere` class calculates the circumsphere of a set of points, i.e. a sphere such that all the points are on the surface of the sphere. It is used internally by the `delaunay` module.

Create a `Circumsphere` from a list of points and a triangle specified by indices into that list:

```
var sph = Circumsphere(pts, [i,j,k])
```

Test if an arbitrary point is inside the `Circumsphere` or not:

```
print sph.pointinsphere(pt)
```

## 9.9 Graphics

The `graphics` module provides a number of classes to provide simple visualization capabilities. To use it, you first need to import the module:

```
import graphics
```

The `Graphics` class acts as an abstract container for graphical information; to actually launch the display see the `Show` class. You can create an empty scene like this,

```
var g = Graphics()
```

Additional elements can be added using the `display` method.

```
g.display(element)
```

Morpho provides the following fundamental Graphical element classes:

```
TriangleComplex
```

You can also use functions like `Arrow`, `Tube` and `Cylinder` to create these elements conveniently.

To combine graphics objects, use the add operator:

```
var g1 = Graphics(), g2 = Graphics()
// ...
Show(g1+g2)
```

## 9.10 Show

`Show` is used to launch an interactive graphical display using the external `morphoview` application. `Show` takes a `Graphics` object as an argument:

```
var g = Graphics()
Show(g)
```

## 9.11 TriangleComplex

A `TriangleComplex` is a graphical element that can be used as part of a graphical display. It consists of a list of vertices and a connectivity matrix that selects which vertices are used in each triangle.

To create one, call the constructor with the following arguments:

```
TriangleComplex(position, normals, colors, connectivity)
```

- `position` is a `Matrix` containing vertex positions as *columns*.
- `normals` is a `Matrix` with a normal for each vertex.
- `colors` is the color of the object.
- `connectivity` is a `Sparse` matrix where each column represents a triangle and rows correspond to vertices.

You can also provide optional arguments:

- `transmit` sets the transparency of the object. This parameter is only used by the `povray` module as of now. Default is 0.
- `filter` sets the transparency of the object using a filter effect. This parameter is only used by the `povray` module as of now. Default is 0. For the difference between `transmit` and `filter`, checkout the `POV-Ray` documentation.

Add to a `Graphics` object using the `display` method.

## 9.12 Arrow

The `Arrow` function creates an arrow. It takes two arguments:

```
arrow(start, end)
```

- `start` and `end` are the two vertices. The arrow points `start -> end`.

You can also provide optional arguments:

- `aspectratio` controls the width of the arrow relative to its length
- `n` is an integer that controls the quality of the display. Higher `n` leads to a rounder arrow.
- `color` is the color of the arrow. This can be a list of `RGB` values or a `Color` object
- `transmit` sets the transparency of the arrow. This parameter is only used by the `povray` module as of now. Default is 0.
- `filter` sets the transparency of the arrow using a filter effect. This parameter is only used by the `povray` module as of now. Default is 0. For the difference between `transmit` and `filter`, checkout the `POV-Ray` documentation.

Display an arrow:

```
var g = Graphics([])
g.display(Arrow([-1/2,-1/2,-1/2], [1/2,1/2,1/2], aspectratio=0.05, n=10))
Show(g)
```



### 9.13 Cylinder

The `Cylinder` function creates a cylinder. It takes two required arguments:

```
cylinder(start, end)
```

- `start` and `end` are the two vertices.

You can also provide optional arguments:

- `aspectratio` controls the width of the cylinder relative to its length.
- `n` is an integer that controls the quality of the display. Higher `n` leads to a rounder cylinder.
- `color` is the color of the cylinder. This can be a list of RGB values or a `Color` object.
- `transmit` sets the transparency of the cylinder. This parameter is only used by the povray module as of now. Default is 0.
- `filter` sets the transparency of the cylinder using a filter effect. This parameter is only used by the povray module as of now. Default is 0. For the difference between `transmit` and `filter`, checkout the POV-Ray documentation.

Display an cylinder:

```
var g = Graphics()  
g.display(Cylinder([-1/2,-1/2,-1/2], [1/2,1/2,1/2], aspectratio=0.1, n=10))  
Show(g)
```

### 9.14 Tube

The `Tube` function connects a sequence of points to form a tube.

```
Tube(points, radius)
```

- `points` is a list of points; this can be a list of lists or a `Matrix` with the positions as columns.
- `radius` is the radius of the tube.

You can also provide optional arguments:

- `n` is an integer that controls the quality of the display. Higher `n` leads to a rounder tube.
- `color` is the color of the tube. This can be a list of RGB values or a `Color` object.
- `closed` is a `bool` that indicates whether the tube should be closed to form a loop.
- `transmit` sets the transparency of the tube. This parameter is only used by the povray module as of now. Default is 0.
- `filter` sets the transparency of the tube using a filter effect. This parameter is only used by the povray module as of now. Default is 0. For the difference between `transmit` and `filter`, checkout the POV-Ray documentation.

Draw a square:

```
var a = Tube([[ -1/2, -1/2, 0], [1/2, -1/2, 0], [1/2, 1/2, 0], [-1/2, 1/2, 0]], 0.1, closed=true)
var g = Graphics()
g.display(a)
```

## 9.15 Sphere

The `Sphere` function creates a sphere.

```
Sphere(center, radius)
```

- `center` is the position of the center of the sphere; this can be a list or column `Matrix`.
- `radius` is the radius of the sphere

You can also provide an optional argument:

- `color` is the color of the sphere. This can be a list of RGB values or a `Color` object.
- `transmit` sets the transparency of the sphere. This parameter is only used by the `povray` module as of now. Default is 0.
- `filter` sets the transparency of the sphere using a filter effect. This parameter is only used by the `povray` module as of now. Default is 0. For the difference between `transmit` and `filter`, checkout the `POV-Ray` documentation.

Draw some randomly sized spheres:

```
var g = Graphics()
for (i in 0...10) {
  g.display(Sphere([random()-1/2, random()-1/2, random()-1/2], 0.1*(1+random()),
    color=Gray(random())))
}
Show(g)
```

## 9.16 ImplicitMesh

The `implicitmesh` module allows you to build meshes from implicit functions. For example, the unit sphere could be specified using the function  $x^2+y^2+z^2-1 == 0$ .

To use the module, first import it:

```
import implicitmesh
```

To create a sphere, first create an `ImplicitMeshBuilder` object with the implicit function you'd like to use:

```
var impl = ImplicitMeshBuilder(fn (x,y,z) x^2+y^2+z^2-1)
```

You can use an existing function (or method) as well as an anonymous function as above.

Then build the mesh,

```
var mesh = impl.build(stepsize=0.25)
```

The build method takes a number of optional arguments:

- **start** - the starting point. If not provided, the value `Matrix([1,1,1])` is used.
- **stepsize** - approximate lengthscale to use.
- **maxiterations** - maximum number of iterations to use. If this limit is exceeded, a partially built mesh will be returned.

## 9.17 KDTree

The `kdtree` module implements a k-dimensional tree, a space partitioning data structure that can be used to accelerate computational geometry calculations.

To use the module, first import it:

```
import kdtree
```

To create a tree from a list of points:

```
var pts = []
for (i in 0...100) pts.append(Matrix([random(), random(), random()]))
var tree=KDTree(pts)
```

Add further points:

```
tree.insert(Matrix([0,0,0]))
```

Test whether a given point is present in the tree:

```
tree.ismember(Matrix([1,0,0]))
```

Find all points within a given bounding box:

```
var pts = tree.search([-1,1], [-1,1], [-1,1])
for (x in pts) print x.location
```

Find the nearest point to a given point:

```
var pt = tree.nearest(Matrix([0.1, 0.1, 0.5]))
print pt.location
```

## 9.18 Insert

Inserts a new point into a k-d tree. Returns a `KDTreeNode` object.

```
var node = tree.insert(Matrix([0,0,0]))
```

Note that, for performance reasons, if the set of points is known ahead of time, it is generally better to build the tree using the constructor function `KDTree` rather than one-by-one with `insert`.

## 9.19 Ismember

Checks if a point is a member of a k-d tree. Returns `true` or `false`.

```
print tree.ismember(Matrix([0,0,0]))
```

## 9.20 Nearest

Finds the point in a k-d tree nearest to a point of interest. Returns a `KDTreeNode` object.

```
var pt = tree.nearest(Matrix([0.1, 0.1, 0.5]))
```

To get the location of this nearest point, access the `location` property:

```
print pt.location
```

## 9.21 Search

Finds all points in a k-d tree that lie within a cuboidal bounding box. Returns a list of `KDTreeNode` objects.

Find and display all points that lie in a cuboid  $0 \leq x \leq 1$ ,  $0 \leq y \leq 2$ ,  $1 \leq z \leq 2$ :

```
var result = tree.search([[0,1], [0,2], [1,2]])
for (x in result) print x.location
```

## 9.22 KDTreeNode

An object corresponding to a single node in a k-d tree. To get the location of the node, access the `location` property:

```
print node.location
```

## 9.23 Meshgen

The `meshgen` module is used to create `Mesh` objects corresponding to a specified domain. It provides the `MeshGen` class to perform the meshing, which are created with the following arguments:

```
MeshGen(domain, boundingbox)
```

Domains are specified by a scalar function that is positive in the region to be meshed and locally smooth. For example, to mesh the unit disk:

```
var dom = fn (x) -(x[0]^2+x[1]^2-1)
```

A `MeshGen` object is then created and then used to build the `Mesh` like this:

```
var mg = MeshGen(dom, [-1..1:0.2, -1..1:0.2])
var m = mg.build()
```

A bounding box for the mesh must be specified as a `List` of `Range` objects, one for each dimension. The increment on each `Range` gives an approximate scale for the size of elements generated.

To facilitate convenient creation of domains, a `Domain` class is provided that provides set operations `union`, `intersection` and `difference`.

`MeshGen` accepts a number of optional arguments:

- `weight` A scalar weight function that controls mesh density.
- `quiet` Set to `true` to suppress `MeshGen` output.
- `method` a list of options that controls the method used.

Some method choices that are available include:

- "FixedStepSize" Use a fixed step size in optimization.
- "StartGrid" Start from a regular grid of points (the default).
- "StartRandom" Start from a randomly generated collection of points.

There are also a number of properties of a `MeshGen` object that can be set prior to calling `build` to control the operation of the mesh generation:

- `stepsize`, `steplimit` Stepsize used internally by the Optimizer
- `fscale` an internal "pressure"
- `ttol` how far the vertices are allowed to move before retriangulation
- `etol` energy tolerance for optimization problem

`MeshGen` picks default values that cover a reasonable range of uses.

## 9.24 Domain

The `Domain` class is used to conveniently build a domain by composing simpler elements.

Create a `Domain` from a scalar function that is positive in the region of interest:

```
var dom = Domain(fn (x) -(x[0]^2+x[1]^2-1))
```

You can pass it to `MeshGen` to specify the region to mesh:

```
var mg = MeshGen(dom, [-1..1:0.2, -1..1:0.2])
```

You can combine `Domain` objects using set operations `union`, `intersection` and `difference`:

```
var a = CircularDomain(Matrix([-0.5,0]), 1)
var b = CircularDomain(Matrix([0.5,0]), 1)
var c = CircularDomain(Matrix([0,0]), 0.3)
var dom = a.union(b).difference(c)
```

## 9.25 CircularDomain

Conveniently constructs a `Domain` object corresponding to a disk. Requires the position of the center and a radius as arguments.

Create a domain corresponding to the unit disk:

```
var c = CircularDomain([0,0], 1)
```

## 9.26 HalfSpaceDomain

Conveniently constructs a Domain object corresponding to a half space defined by a plane at  $x_0$  and a normal  $n$ :

```
var hs = HalfSpaceDomain(x0, n)
```

Note  $n$  is an “outward” normal, so points into the *excluded* region.

Half space corresponding to the allowed region  $x < 0$ :

```
var hs = HalfSpaceDomain(Matrix([0,0,0]), Matrix([1,0,0]))
```

Note that HalfSpaceDomains cannot be meshed directly as they correspond to an infinite region. They are useful, however, for combining with other domains.

Create half a disk by cutting a HalfSpaceDomain from a CircularDomain:

```
var c = CircularDomain([0,0], 1)
var hs = HalfSpaceDomain(Matrix([0,0]), Matrix([-1,0]))
var dom = c.difference(hs)
var mg = MeshGen(dom, [-1..1:0.2, -1..1:0.2], quiet=false)
var m = mg.build()
```

## 9.27 MshGnDim

The MeshGen module currently supports 2 and 3 dimensional meshes. Higher dimensional meshing will be available in a future release; please contact the developer if you are interested in this functionality.

## 9.28 Meshtools

The Meshtools package contains a number of functions and classes to assist with creating and manipulating meshes.

## 9.29 AreaMesh

This function creates a mesh composed of triangles from a parametric function. To use it:

```
var m = AreaMesh(function, range1, range2, closed=boolean)
```

where

- `function` is a parametric function that has one parameter. It should return a list of coordinates or a column matrix corresponding to this parameter.
- `range1` is the Range to use for the first parameter of the parametric function.
- `range2` is the Range to use for the second parameter of the parametric function.
- `closed` is an optional parameter indicating whether to create a closed loop or not. You can supply a list where each element indicates whether the relevant parameter is closed or not.

To use AreaMesh, import the meshtools module:

```
import meshtools
```

Create a square:

```
var m = AreaMesh(fn (u,v) [u, v, 0], 0..1:0.1, 0..1:0.1)
```

Create a tube:

```
var m = AreaMesh(fn (u, v) [v, cos(u), sin(u)], -Pi...Pi:Pi/4,
                        -1..1:0.1, closed=[true, false])
```

Create a torus:

```
var c=0.5, a=0.2
var m = AreaMesh(fn (u, v) [(c + a*cos(v))*cos(u),
                           (c + a*cos(v))*sin(u),
                           a*sin(v)], 0...2*Pi:Pi/16, 0...2*Pi:Pi/8, closed=true)
```

### 9.30 LineMesh

This function creates a mesh composed of line elements from a parametric function. To use it:

```
var m = LineMesh(function, range, closed=boolean)
```

where

- **function** is a parametric function that has one parameter. It should return a list of coordinates or a column matrix corresponding to this parameter.
- **range** is the Range to use for the parametric function.
- **closed** is an optional parameter indicating whether to create a closed loop or not.

To use **LineMesh**, import the **meshtools** module:

```
import meshtools
```

Create a circle:

```
import constants
var m = LineMesh(fn (t) [sin(t), cos(t), 0], 0...2*Pi:2*Pi/50, closed=true)
```

### 9.31 PolyhedronMesh

This function creates a mesh from a polyhedron specification.

### 9.32 Equiangular

Attempts to equiangularate a mesh.

### 9.33 MeshBuilder

The `MeshBuilder` class simplifies user creation of meshes. To use this class, begin by creating a `MeshBuilder` object:

```
var build = MeshBuilder()
```

You can then add vertices, edges, etc. one by one using `addvertex`, `addedge` and `addtriangle`. Each of these returns an element id:

```
var id1=build.addvertex(Matrix([0,0,0]))
var id2=build.addvertex(Matrix([1,1,1]))
build.addedge([id1, id2])
```

Once the mesh is ready, call the `build` method to construct the `Mesh`:

```
var m = build.build()
```

### 9.34 MeshRefiner

The `MeshRefiner` class is used to refine meshes, and to correct associated data structures that depend on the mesh.

### 9.35 Optimize

The `optimize` package contains a number of functions and classes to perform shape optimization.

### 9.36 OptimizationProblem

An `OptimizationProblem` object defines an optimization problem, which may include functionals to optimize as well as global and local constraints.

Create an `OptimizationProblem` with a mesh:

```
var problem = OptimizationProblem(mesh)
```

Add an energy:

```
var la = Area()
problem.addenergy(la)
```

Add an energy that operates on a selected region, and with an optional prefactor:

```
problem.addenergy(la, selection=sel, prefactor=2)
```

Add a constraint:

```
problem.addconstraint(la)
```

Add a local constraint (here a onesided level set constraint):

```
var ls = ScalarPotential(fn (x,y,z) z, fn (x,y,z) Matrix([0,0,1]))
problem.addlocalconstraint(ls, onesided=true)
```



### 9.37 Optimizer

Optimizer objects are used to optimize Meshes and Fields. You should use the appropriate subclass: ShapeOptimizer or FieldOptimizer respectively.

### 9.38 ShapeOptimizer

A ShapeOptimizer object performs shape optimization: it moves the vertex positions to reduce an overall energy.

Create a ShapeOptimizer object with an OptimizationProblem and a Mesh:

```
var sopt = ShapeOptimizer(problem, m)
```

Take a step down the gradient with fixed stepsize:

```
sopt.relax(5) // Takes five steps
```

Linesearch down the gradient:

```
sopt.linesearch(5) // Performs five linesearches
```

Perform conjugate gradient (usually gives faster convergence):

```
sopt.conjugategradient(5) // Performs five conjugate gradient steps.
```

Control a number of properties of the optimizer:

```
sopt.stepsize=0.1 // The stepsize to take  
sopt.steplimit=0.5 // Maximum stepsize for optimizing methods  
sopt.etol = 1e-8 // Energy convergence tolerance  
sopt.ctol = 1e-9 // Tolerance to which constraints are satisfied  
sopt.maxconstraintsteps = 20 // Maximum number of constraint steps to use
```

### 9.39 FieldOptimizer

A FieldOptimizer object performs field optimization: it changes elements of a Field to reduce an overall energy.

Create a FieldOptimizer object with an OptimizationProblem and a Field:

```
var sopt = FieldOptimizer(problem, fld)
```

Field optimizers provide the same options and methods as Shape optimizers: see the ShapeOptimizer documentation for details.

### 9.40 Plot

The plot module provides visualization capabilities for Meshes, Selections and Fields. These functions produce Graphics objects that can be displayed with Show.

To use the module, first import it:

```
import plot
```

## 9.41 Plotmesh

Visualizes a Mesh object:

```
var g = plotmesh(mesh)
```

Plotmesh accepts a number of optional arguments to control what is displayed:

- `selection` - Only elements in a provided `Selection` are drawn.
- `grade` - Only draw the specified grade. This can also be a list of multiple grades to draw.
- `color` - Draw the mesh in a provided `Color`.
- `filter` and `transmit` - Used by the `povray` module to indicate transparency.

## 9.42 Plotselection

Visualizes a `Selection` object:

```
var g = plotselection(mesh, sel)
```

Plotselection accepts a number of optional arguments to control what is displayed:

- `grade` - Only draw the specified grade. This can also be a list of multiple grades to draw.
- `filter` and `transmit` - Used by the `povray` module to indicate transparency.

## 9.43 Plotfield

Visualizes a scalar `Field` object:

```
var g = plotfield(field)
```

Plotfield accepts a number of optional arguments to control what is displayed:

- `grade` - Draw the specified grade.
- `colormap` - A `Colormap` object to use. The field is automatically scaled.
- `style` - Plot style. See below.
- `filter` and `transmit` - Used by the `povray` module to indicate transparency.

Supported plot styles:

- `default` - Color Mesh elements by the corresponding value of the `Field`.
- `interpolated` - Interpolate `Field` quantities onto higher elements.

## 9.44 POVRay

The `povray` module provides integration with POVRay, a popular open source ray-tracing package for high quality graphical rendering. To use the module, first import it:

```
import povray
```

To raytrace a graphic, begin by creating a `POVRaytracer` object:

```
var pov = POVRaytracer(graphic)
```

Create a `.pov` file that can be run with POVRay:

```
pov.write("out.pov")
```

Create, render and display a scene using POVRay:

```
pov.render("out.pov")
```

This also creates the `.png` file for the scene.

The `POVRaytracer` constructor supports a number of optional arguments:

- `antialias` - whether to antialias the output or not
- `width` - image width
- `height` - image height
- `viewangle` - camera angle (higher means wider view)
- `viewpoint` - position of camera

The `render` method supports two optional boolean arguments:

- `quiet` - whether to suppress the parser and render statistics from `povray` or not (`false` by default)
- `display` - whether to turn on the graphic display while rendering or not (`true` by default)

## 9.45 VTK

The `vtk` module contains classes to allow I/O of meshes and fields using the VTK Legacy Format.

## 9.46 VTKExporter

This class can be used to export the field(s) and/or at a given state to a single `.vtk` file. To use it, import the `vtk` module:

```
import vtk
```

Initialize the `VTKExporter`

```
var vtkE = VTKExporter(obj)
```

where `obj` can either be

- A `Mesh` object: This prepares the `Mesh` for exporting.

- A `Field` object: This prepares both the `Field` and the `Mesh` associated with it for exporting.

Use the `export` method to export to a VTK file.

```
vtkE.export("output.vtk")
```

Optionally, use the `addfield` method to add one or more fields before exporting:

```
vtkE.addfield(f, fieldname="f")
```

where,

- `f` is the field object to be exported
- `fieldname` is an optional argument that assigns a name to the field in the VTK file. This name is required to be a character string without embedded whitespace. If not provided, the name would be either “scalars” or “vectors” depending on the field type\*\*.

\*\* Note that this currently only supports scalar or vector (column matrix) fields that live on the vertices (shape `[1,0,0]`). Support for tensorial fields and fields on cells coming soon.

Minimal example:

```
import vtk
import meshtools

var m1 = LineMesh(fn (t) [t,0,0], -1..1:2)

var vtkE = VTKExporter(m1) // Export just the mesh

vtkE.export("mesh.vtk")

var f1 = Field(m1, fn(x,y,z) x)

var g1 = Field(m1, fn(x,y,z) Matrix([x,2*x,3*x]))

vtkE = VTKExporter(f1, fieldname="f") // Export fields

vtkE.addfield(g1, fieldname="g")

vtkE.export("data.vtk")
```

## 9.47 VTKImporter

This class can be used to import the field(s) and/or the mesh at a given state from a single `.vtk` file. To use it, import the `vtk` module:

```
import vtk
```

Initialize the `VTKImporter` with the filename

```
var vtkI = VTKImporter("output.vtk")
```

Use the `mesh` method to get the mesh:

```
var mesh = vtkI.mesh()
```

Use the `field` method to get the field:

```
var f = vtkI.field(fieldname)
```

Use the `fieldlist` method to get the list of the names of the fields contained in the file:

```
print vtkI.fieldlist()
```

Use the `containsfield` method to check whether the file contains a field by a given `fieldname`:

```
if (vtkI.containsfield(fieldname)) {  
    ...  
}
```

where `fieldname` is the name assigned to the field in the `.vtk` file

Minimal example:

```
import vtk  
import meshtools  
  
var vtkI = VTKImporter("data.vtk")  
  
var m = vtkI.getmesh()  
  
var f = vtkI.getfield("f")  
  
var g = vtkI.getfield("g")
```