PRECEPT 09

CEE 361-513: Introduction to Finite Element Methods

Monday Nov. 27

Poisson's Equation with FEniCS

1. We are going to solve the simplest 2-D problem, the Poisson Equation. The problem reads: find u: $[0,1] \times [0,1] \to \mathbb{R}$ such that

$$-\Delta u = f \quad \text{in } \Omega$$
$$u = 0 \quad \text{on } \Gamma_D$$

We solve the problem on the domain $\Omega = [0,1] \times [0,1]$. Γ_D represents the Dirichlet boundary of the domain. For this problem we employ the method of manufactured solution to study the convergence. The manufactured solution is:

$$u = exp(-((x - 0.5)^2 + (y - 0.5)^2)/0.125^2))$$

First and foremost we use sympy to create the analytical solution ue and obtain the expressions for g and f.

```
# Create manufactured solution
x = sp.symbols('x[0]')
y = sp.symbols('x[1]')
solution = sp.exp(-((x-0.5)**2+(y-0.5)**2)/0.125**2)
ue_code = sp.ccode(solution).replace('M_PI','pi')
f_code = sp.ccode( -sp.diff(solution,x,2) -sp.diff(solution,y,2))\
.replace('M_PI','pi')
7
```

Next we create the mesh/subdivision of the domain:

```
V = FunctionSpace(mesh, 'Lagrange', poly_order)
1
2
```

The mesh looks like:

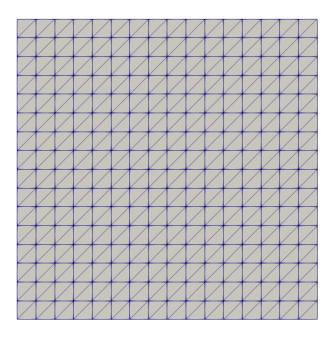


Figure 1: Mesh

We now create the subspace of piecewise linear polynomials as we saw in class.

```
# Create the function space
poly_order = 1
V = FunctionSpace(mesh, 'Lagrange', poly_order)
1
2
3
```

Next we identify the parts of the boundary of our domain that are Dirichlet Γ_D . We do not have a Neumann Boundary in this problem

```
# Define the dirichlet boundary
class dirichlet_boundary(SubDomain):
   def inside(self, x, on_boundary):
    return on_boundary

4
```

We define the exact solution as FEniCS expression

```
# Define the analaytical solution
ue = Expression(ue_code, pi=np.pi, degree=5)
1
```

and we define the Dirichlet boundary conditions

```
# Define the dirichlet boundary conditions
bc = DirichletBC(V, ue, dirichlet_boundary())
```

Now we are ready to define the trial function and the test function, the expression for the source term

```
# Define the trial and test function
u = TrialFunction(V)
v = TestFunction(V)

1
2
```

```
# Define the forcing function
f = Expression(f_code,pi=np.pi, degree=5)
1
2
```

We now define the bilinear form and the forcing functional

```
# Define the bilinear form
a = dot(grad(u),grad(v))*dx

# Define the forcing
F = f*v*dx
1
2
3
4
5
```

and we finally solve the problem

```
# Compute solution
uh = Function(V, name='displacement')
solve(a == F, uh, bc)
1
2
3
```

Since we know the exact the solution we can compute the L2-norm, H1-norm of error , and the maximum error.

```
# Compute maximum error at vertices
vertex_values_ue = ue.compute_vertex_values(mesh)
vertex_values_uh = uh.compute_vertex_values(mesh)

error_L2 = errornorm(ue, uh, 'L2')
error_H1 = errornorm(ue, uh, 'H1')

error_max = np.max(np.abs(vertex_values_ue - vertex_values_uh))

8
```

The maximum error was 0.0204

We can solve the problem on successively refined meshes and obtain the rate of convergence.

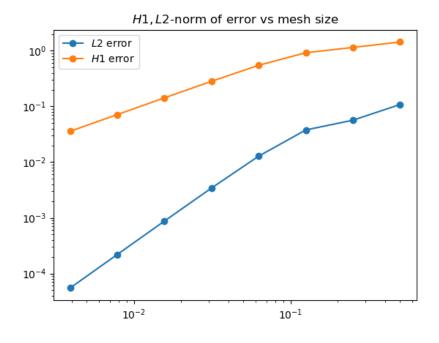


Figure 2: Convergence plot

The rate of convergence using L2-norm was 2 The rate of convergence using H1-norm was 1

2. Saving the solution in XDMF format

```
# Save as xdmf format
# The file in which the solution is stored
file_xdmf = XDMFFile('poisson.xdmf')
file_xdmf.write(uh,0)
4
```

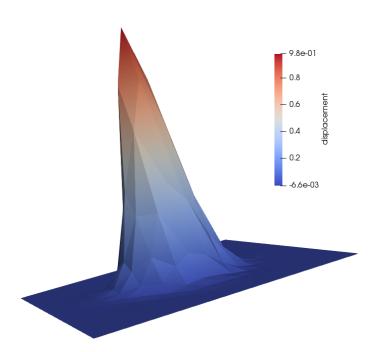


Figure 3: Solution with initial mesh

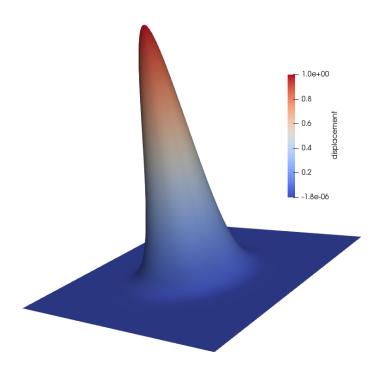


Figure 4: Converged Solution

Poisson's Equation with Quad Elements

Here we modify the poisson's problem file provided along with the homework code to add the function for computing L1-norm and H1-norm of the errors and determine the rate of convergence.

1. Compute L2 - norm of error

```
def compute_12_norm_error(local_to_global_map, elements, uh, ue):
                                                                           1
                                                                           2
 12_{error} = 0.
                                                                           3
                                                                           4
                                                                           5
 for element in elements:
    # Get the element local degress of freedom
                                                                           6
    num_dofs = element.get_num_dofs( )
                                                                           7
                                                                           8
                                                                           9
    # Get the quadrature rule
    gauss_points, gauss_weights = element.get_quadrature()
                                                                           10
                                                                           11
    # Fill in the stiffness matrix
                                                                           12
    for q in range(len(gauss_points)):
                                                                           13
                                                                           14
      # Get the jacobian at the quadrature point
                                                                           15
      jacobian = element.get_dmap( gauss_points[q], jacobian=True )
                                                                           16
                                                                           17
      # Get the image of the gauss point on the physical domain
                                                                           18
                                                                           19
      x_g = element.get_map( gauss_points[q] )
                                                                           20
      uh_g = 0.
                                                                           21
                                                                           22
                                                                           23
      # Loop over all degree of freedoms
      for i in range(num_dofs):
                                                                           24
        # Get the global index of the local i dof
                                                                           25
        i_global = local_to_global_map.get_global_dof( \
                                                                           26
                                                                           27
          element.element_index , i )
                                                                           28
        # If we specified a source term
                                                                           29
        uh_g += element.get_base_function_val(i,\
                                                                           30
          gauss_points[q])*uh[i_global]
                                                                           31
                                                                           32
      # Add contribution of intergral at gauss point
                                                                           33
      12_error += pow( ue( x_g ) - uh_g , 2 )*jacobian*gauss_weights[q]
                                                                           34
                                                                           35
 return np.sqrt(12_error)
                                                                           36
```

2. Compute H1 - norm of error

```
num_dofs = element.get_num_dofs( )
                                                                        8
                                                                        g
  # Get the quadrature rule
                                                                        10
  gauss_points, gauss_weights = element.get_quadrature()
                                                                        11
                                                                        12
                                                                        13
  # Fill in the stiffness matrix
  for q in range(len(gauss_points)):
                                                                        14
                                                                        15
    # Get the jacobian at the quadrature point
                                                                        16
    jacobian = element.get_dmap( gauss_points[q], jacobian=True )
                                                                        17
                                                                        18
    # Get the image of the gauss point on the physical domain
                                                                        19
    x_g = element.get_map( gauss_points[q] )
                                                                        20
                                                                        21
                                                                        22
    # Get the value of gradient of all basis at quadrature point
                                                                        23
    uh_g = 0.
                                                                        24
    duh_g = 0.
                                                                        25
                                                                        26
    # Loop over all degree of freedoms
                                                                        27
    for i in range(num_dofs):
                                                                        28
      # Get the global index of the local i dof
                                                                        29
      i_global = local_to_global_map.get_global_dof( \
                                                                        30
        element.element_index , i )
                                                                        31
                                                                        32
      # If we specified a source term
                                                                        33
                                                                        34
      uh_g += element.get_base_function_val(i,gauss_points[q])*\
                                                                        35
          uh[i_global]
      duh_g +=element.get_base_function_grad(i,gauss_points[q])*\
                                                                        36
                                                                        37
          uh[i_global]
                                                                        38
    # Add contribution of intergral at gauss point
                                                                        39
    h1_{error} += pow(ue(x_g) - uh_g, 2)*jacobian*gauss_weights[q] \ 40
          + np.dot((due(x_g) - duh_g), (due(x_g) - duh_g))
                                                                        41
          *jacobian*gauss_weights[q]
                                                                        42
                                                                        43
                                                                        44
return np.sqrt(h1_error)
```

3. Compute rate of convergence.

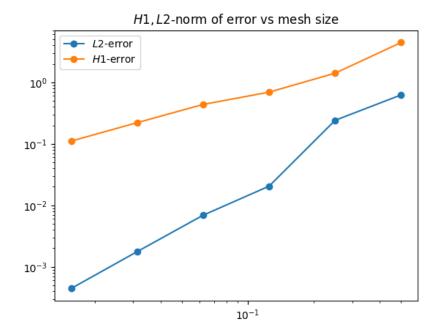


Figure 5: Convergence plot for quad elements

The rate of convergence for L2-norm of error is 2 and the rate of convergence of H1-norm of error is 1