Finite Element Modelling for Geosciences: Code Verification and the Method of Manufactured **Solutions**

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How do I know the code is working?

- \triangleright As you've no doubt noticed, there are many, many ways to make mistakes in your implementation
- \blacktriangleright How do you show that the code "works"?
- \triangleright One way we've seen is to confirm that our finite element solution looks "close" to an exact solution that we know beforehand
- \blacktriangleright But ...
	- \blacktriangleright What does "close" mean?
	- \triangleright What if I don't have an exact solution to compare with?

\blacktriangleright Verification

- \blacktriangleright Is your code solving the equations correctly?
- \triangleright Does the code produce approximate solutions that converge at the correct rate?
- \blacktriangleright The subject of Chapter 9 in the course notes
- \blacktriangleright Validation
	- \triangleright Does the code produce physically-meaningful results?
	- \blacktriangleright Not covered in this course
- If we can construct an exact solution to the continuous problem, we can use it to verify our code
- \triangleright The FEM is well-known to converge, if implemented correctly!
- \triangleright Convergence is expressed in terms of the grid spacing and an error norm

$$
||u-u^h||
$$

If an analytical expression is available, one can compute the solution at different spacings and examine how quickly it converges

$$
||u-u^h||\leq Ch^k
$$

- \blacktriangleright The rate of convergence can be computed by fitting a straight line to a log-log plot of the grid spacing versus the error
- ▶ Obvious issue: what if we don't know μ ?

Computing errors

The L^2 error

$$
\left(\int_{\Omega}\left(u^h(x,y)-u(x,y)\right)^2 d\vec{x}\right)^{\frac{1}{2}}
$$

 \triangleright An integral over the entire domain; compute it with quadrature!

In Sometimes one sees "superconvergence" if only comparing nodal values.

err = sqrt $(sum ((T-T_exact) .^2))$ % Not an L 2 norm!

The Method of Manufactured Solutions (MMS)

- \triangleright General approach works for most PDE.
- Intuition: choose the solution you want and adjust the forcing and boundary terms to match.
- \triangleright Choose ("manufacture") your solution
- \blacktriangleright Plug it into your equations (physics $+$ material properties)
- \triangleright Extract the required boundary and forcing terms (CAS tools like SAGE/SymPy/MAPLE/Mathematica/etc. are helpful)
- \triangleright Now you can use your FEM code to compute an approximate solution, and compare.
- \blacktriangleright How to choose your manufactured solution
	- \blacktriangleright Make it complex enough
	- \blacktriangleright In particular, avoid low-order polynomials which might be exactly captured by your finite element basis functions
	- \triangleright Good choices are combinations of sinusoidal or exponential functions

 \triangleright Assume we're interested in approximate solutions to the Poisson equation (steady-state heat equation) on the unit square

$$
\nabla \cdot (\kappa(x, y) \nabla T) = f(x, y), \quad \kappa(x, y) = 1 + x
$$

- \blacktriangleright Note that we specified the coefficient structure, but not the boundary conditions or forcing
- \triangleright Choose our own solution (depending on both x and y, not just a polynomial)

$$
T_{MMS} \doteq \sin(x)\cos(2y)
$$

Insert T_{MMS} into our equation¹ to obtain the forcing

$$
f(x,y) = \frac{\partial}{\partial x} \left((1+x) \frac{\partial T_{MMS}}{\partial x} \right) + \frac{\partial}{\partial y} \left((1+x) \frac{\partial T_{MMS}}{\partial y} \right)
$$

$$
= \cos(2y) \left(\cos(x) - 5(x+1) \sin(x) \right)
$$

 \triangleright Use T_{MMS} to define boundary conditions, for example Dirichlet conditions

> $T(x, 0) = T_{MMS}(x, 0), \quad T(x, 1) = T_{MMS}(x, 1), \quad \forall x \in [0, 1]$ $T(0, y) = T_{MMS}(0, y), \quad T(1, y) = T_{MMS}(1, y), \quad \forall y \in [0, 1]$

 $^{\text{\tiny 1}}$ I lazily just plugged this into Wolfram Alpha: $\text{\rm Div}{\text{\small [}}(1\text{+x)}\text{*Grad}{\text{\small [}}\text{\small Sin}{\text{\small [}} x\text{\small]}\text{*Cos}{\text{\small [}} 2\text{*y}{\text{\small]}}\text{\small]}$

L2 error computation MATLAB code snippet

```
L2_error_2 = 0;for i<sub>-el</sub> = 1: el<sub>-tot</sub>
     nodes= el = elementnodemap(:, iel);
     coord_e1 = good(:,nodes_e1)T_e1 = T(nodes_e1);for i<sub>-pt</sub> = 1: pts-per_el
          \texttt{jac} = dNdxi(:, :, i\_pt) * coord_el;
          det_iac = det(iac)coord_p t = N(i_p t, :) * coord_e 1;T_{\text{mm}}s_pt = T_{\text{MMS}} (coord_pt (1), coord_pt (2));
          T_f = pt = N(i_p t, :) * T_e1;
          L2_error_2 = L2_error_2 + \(T_{\texttt{mms\_pt}} - T_{\texttt{fe\_pt}})^2 * det<sub>riac</sub> * w(i_pt);
     end
L2<sub>-</sub>error = sqrt(L2<sub>-</sub>error<sub>-2</sub>)
```
 \blacktriangleright The expected convergence rate, from the course notes, is

$$
\left[\int_{\Omega} (T-T^h)^2 d\vec{x}\right]^{\frac{1}{2}} \leq Ch^{p+1}
$$

 \triangleright Our modified code shows slopes of approximately 2 and 3 on a log-log plot, so we are now much more confident that the implementation is correct.

MMS: Wrinkle - non-physical forcing

- If using MMS for something like the Stokes equations (Friday), you will perhaps run into the following issue
- \triangleright As an example, let's say that I have written a FE code to solve this (very) simple PDE:

$$
\frac{d}{dx}\left(k(x)\frac{du}{dx}\right) = 0, \quad u \in [0,1], \quad u(0) = 1, u(1) = 2
$$

- \triangleright What issue might I run into if I apply the MMS?
- I may have written my code assuming zero forcing!
- \triangleright So I may have to **modify my weak form** to take the additional forcing into account. Don't forget to consider terms like

$$
\mathbf{F} = \int_{\Omega^e} \mathbf{N}^T f \ d\vec{x}
$$

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