Computer Modeling and Finite Difference Methods

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1 Philosophy of Computer Modeling

Before starting to write a program to solve a problem, one must ask oneself important questions such as,

- Why do I want to model this system?
 - To test relationships between variables and gain some new physical insight that would be difficult to attain with reasoning or analytical solutions alone?
 - To accurately represent a real physical system, including efforts to try to match data and their associated uncertainties?
 - To make a prediction of the future and/or past?
 - Other?
- What is the best way to model it?
 - As a set of continuum equations?
 - As a number of individual agents that take actions in their virtual environment?
 - With full-realism physics or simplified rules?
 - Other?
- Should I wish to, how do I compare the "real world" and the data? Or perhaps improve the fit between data and model?

The first algorithm

The first computer program was written in 1843 (or maybe 1842) by Lady Ada Lovelace, with a numerical method to calculate the Bernoulli Numbers. She was 28!

2 Mathematics Review

2.1 Vectors

A vector is a one-dimensional matrix. It is thus a special case of the more general class of matrices. It may be a row,

$$\vec{a} = \begin{bmatrix} a_1 & a_2 & \cdots & a_n \end{bmatrix}, \tag{1}$$

or a column,

$$\vec{a} = \begin{bmatrix} a_1 \\ a_2 \\ \vdots \\ a_m \end{bmatrix} . \tag{2}$$

Vectors can define a direction. For example, in three dimensions:

$$\vec{r} = \begin{bmatrix} x \\ y \\ z \end{bmatrix} . \tag{3}$$

In addition to the top arrow notation, a boldface (but italicized) notation is also common for vectors:

$$\vec{a} = a \tag{4}$$

2.2 Two-dimensional matrices

A two-dimensional matrix of m rows and n columns can be represented in a way that is an extension of our vector notation; these are represented by boldface Roman (non-italic) characters. They are often, but not always, represented by capital letters:

$$\mathbf{A} = \begin{bmatrix} a_{1,1} & a_{1,2} & \cdots & a_{1,n} \\ a_{2,1} & a_{2,2} & \cdots & a_{2,n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{m,1} & a_{m,2} & \cdots & a_{m,n} \end{bmatrix}$$
 (5)

A 3×3 matrix is often used to represent a *tensor* that can define the orientation and dimensions in a three-dimensional shape (ellipsoid). Such a matrix can also describe stress-strain relationships, mineral symmetry, or any other material anisotropy (or isotropy, if the 3×3 matrix is the identity matrix).

One special matrix is the *Identity* matrix, which is an expansion of "1" for an array:

$$\mathbf{I} = \begin{bmatrix} 1 & 0 & \cdots & 0 \\ 0 & 1 & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & 1 \end{bmatrix} \tag{6}$$

2.3 Matrix operations

2.3.1 Matrix multiplication

To multiply two matrices, $\mathbf{A} \times \mathbf{B}$, follow these steps:

- 1. Multiply each number in the first column of the matrix on the left (A) by each number in the first row of the matrix on the right (**B**), in order: $\mathbf{A}_{1,1} \times \mathbf{B}_{1,1}$, $\mathbf{A}_{2,1} \times \mathbf{B}_{1,2}, ..., \mathbf{A}_{1,n} \times \mathbf{B}_{m,1}$. Note that in order for this to work, the number of rows in matrix **A** must equal the number of columns in matrix **B**.
- 2. Sum all of these products that you have just calculated.
- 3. The result is the values for the resultant matrix in location (1,1).
- 4. Repeat this for each row in Matrix A and column in Matrix B. If the sizes of these arrays are $(m_{\mathbf{A}} \times n_{\mathbf{A}})$ and $(m_{\mathbf{B}} \times n_{\mathbf{B}})$, then your resultant array will have dimensions $(m_{\mathbf{A}} \times n_{\mathbf{B}})$.

An example of matrix multiplication follows:

$$\mathbf{AB} = \begin{bmatrix} a_{1,1} & a_{1,2} & a_{1,3} \\ a_{2,1} & a_{2,2} & a_{2,3} \end{bmatrix} \begin{bmatrix} b_{1,1} & b_{1,2} \\ b_{2,1} & b_{2,2} \\ b_{3,1} & b_{3,2} \end{bmatrix} = \begin{bmatrix} a_{1,1}b_{1,1} + a_{1,2}b_{2,1} + a_{1,3}b_{3,1} & a_{1,1}b_{1,2} + a_{1,2}b_{2,2} + a_{1,3}b_{3,2} \\ a_{2,1}b_{1,1} + a_{2,2}b_{2,1} + a_{2,3}b_{3,1} & 4d_{2}, 1b_{1,2} + a_{2,2}b_{2,2} + a_{2,3}b_{3,2} \end{bmatrix}$$
(8)

$$\begin{bmatrix} a_{1,1}b_{1,1} + a_{1,2}b_{2,1} + a_{1,3}b_{3,1} & a_{1,1}b_{1,2} + a_{1,2}b_{2,2} + a_{1,3}b_{3,2} \\ a_{2,1}b_{1,1} + a_{2,2}b_{2,1} + a_{2,3}b_{3,1} & 4d2, 1b_{1,2} + a_{2,2}b_{2,2} + a_{2,3}b_{3,2} \end{bmatrix}$$
(8)

I usually think of matrix multiplication as "over... down". But some better memonics include "karate... chop!" (the "chop" being obviously down) and the visual way of multiplying A and B, as follows:

$$\mathbf{AB} = \tag{9}$$

$$\begin{bmatrix} b_{1,1} & b_{1,2} \\ b_{2,1} & b_{2,2} \\ b_{3,1} & b_{3,2} \end{bmatrix}$$
 (10)

$$\begin{bmatrix} b_{1,1} & b_{1,2} \\ b_{2,1} & b_{2,2} \\ b_{3,1} & b_{3,2} \end{bmatrix}$$

$$\begin{bmatrix} a_{1,1} & a_{1,2} & a_{1,3} \\ a_{2,1} & a_{2,2} & a_{2,3} \end{bmatrix} \begin{bmatrix} & & & & & & & & \\ & & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & \\ & & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & & \\ & &$$

Here, the empty matrix can be filled with the values by multiplying the rows and columns that project (on a straight line) into that cell. (Thanks to Jack for both of these.)

2.3.2 Determinants

The determinant of a matrix provides information that can help us with:

• Systems of linear equations

• Matrix inversions

The determinant of a matrix, \mathbf{A} , is denoted:

$$|\mathbf{A}|. \tag{12}$$

For a 2×2 matrix, the determinant is:

$$\mathbf{A} = \begin{vmatrix} a_{1,1} & a_{1,2} \\ a_{2,1} & a_{2,2} \end{vmatrix} = a_{1,1}a_{2,2} - a_{1,2}a_{2,1} \tag{13}$$

While I prefer to include subscripts with a single letter, this is more commonly displayed in math textbooks as:

$$\begin{vmatrix} a & b \\ c & d \end{vmatrix} = ad - bc \tag{14}$$

For a 3×3 matrix, one takes an additional step. I am showing the method called Laplace Expansion, because it has an easy-to-remember pattern.

$$\mathbf{A} = \begin{vmatrix} a_{1,1} & a_{1,2} & a_{1,3} \\ a_{2,1} & a_{2,2} & a_{2,3} \\ a_{3,1} & a_{3,2} & a_{3,3} \end{vmatrix} = a_{1,1} \begin{vmatrix} a_{1,2} & a_{1,3} \\ a_{2,2} & a_{2,3} \end{vmatrix} - a_{1,2} \begin{vmatrix} a_{1,1} & a_{1,3} \\ a_{2,1} & a_{2,3} \end{vmatrix} + a_{1,3} \begin{vmatrix} a_{1,1} & a_{1,2} \\ a_{2,1} & a_{2,2} \end{vmatrix}$$
 (15)

Note the pattern – eliminate the rows and columns in straight lines from each term on one row (you may pick any row... or column, I think, though I haven't tried this), and then solve.

Following this same nested-determinant and +, -, +, -, ... pattern, you can solve any determinant to an arbitrarily large matrix. But perhaps you see one reason why we have computers.

2.3.3 Dot products

The dot product between two arrays can be visualized as the magnitude the product of two n-dimensional vectors. It is therefore also called the "scalar product", because it gives a scalar result.

$$\vec{a} \cdot \vec{b} = \begin{bmatrix} a_1 & a_2 & \cdots & a_n \end{bmatrix} \begin{bmatrix} b_1 \\ b_2 \\ \vdots \\ b_m \end{bmatrix} = \sum_{i=1}^{m(=n)} a_i b_i$$
 (16)

If you know the angle, $\theta_{\vec{a},\vec{b}}$, between the two vectors, a dot product can also be represented as:

$$|\vec{a}| |\vec{b}| \cos \left(\theta_{\vec{a}, \vec{b}}\right) \tag{17}$$

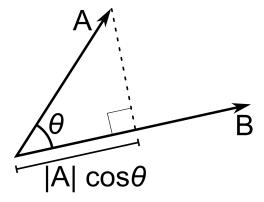


Figure 1: The dot product can be thought of as the scalar product of the magnitude of one vector, times the magnitude of another vector that has been projected onto it. https://commons.wikimedia.org/wiki/File:Dot_Product.svg.

2.3.4 Cross products

A cross product, or "vector product", gives the magnitude and direction of a product of two vectors, \vec{a} and \vec{b} , where the magnitude is the area of the parallelogram created by the two vectors and the direction is given by the right-hand rule.

Cross-products are useful for:

• Vector

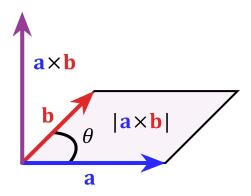


Figure 2: Graphical definition of the cross-product of two vectors. https://commons.wikimedia.org/wiki/File:Cross_product_parallelogram.svg.

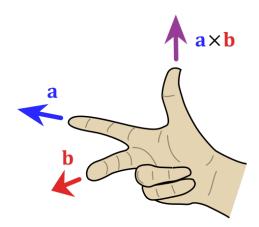


Figure 3: Direction of a cross-product of vectors $\vec{a} \times \vec{b}$ as given by the right-hand rule. https://commons.wikimedia.org/wiki/File:Right_hand_rule_cross_product.svg.

The cross product can be solved in a 3-space (so for two orthogonal 2-D vectors) by solving the determinant of the following equation, here given in a 3-space (\mathbb{R}^3):

$$\vec{u} \times \vec{v} = \begin{vmatrix} \hat{i} & \hat{j} & \hat{k} \\ u_1 & u_2 & u_3 \\ v_1 & v_2 & v_3 \end{vmatrix}$$
 (18)

The scalar value (i.e., magnitude) of the cross product can also be obtained as follows:

$$\left\| \vec{a} \times \vec{b} \right\| = \left\| \vec{a} \right\| \left\| \vec{b} \right\| \sin \theta \tag{19}$$

2.3.5 Dot and cross products

The relationship between the dot and cross products is that:

$$\|\vec{a} \times \vec{b}\|^2 = \|\vec{a}\|^2 \|\vec{b}\|^2 - (\vec{a} \cdot \vec{b})^2$$
 (20)

2.3.6 Trace

The trace operator is given by the sum of the diagonal elements:

$$Tr(\mathbf{A}) = \sum \mathbf{AI} \tag{21}$$

2.3.7 Einstein notation

@TODO:

2.4 Derivatives

A derivative is the change in a function (f(x)) with respect to the change in the independent variable (x) as the interval (Δx) approaches 0:

2.4.1 Ordinary

$$\frac{d}{dx}f(x) = \lim_{\Delta x \to 0} \frac{f(x + \Delta x) - f(x)}{\Delta x}$$
 (22)

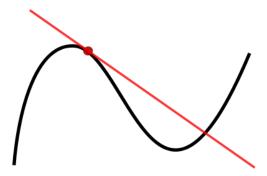


Figure 4: The graph of a function, drawn in black, and a tangent line to that function, drawn in red. The slope of the tangent line is equal to the derivative of the function at the marked point. (Text from https://en.wikipedia.org/wiki/Derivative on 2015.05.07; borrowing it because I couldn't find a better way to say it!)

@TODO: Create a better figure for derivative definition and finite difference, with shown Δx

2.4.2 Partial

@TODO:

2.5 Integrals

Antiderivatives, or integrals, provide the sum of the area under a given curve. These are very useful for solving problems involving processes that occur over space, time, or other (more mathematical, less tangible) dimensions. An example of a non-tangible dimensions would be one that relates to a degree of freedom of some equation and/or error analysis.

Indefinite integrals can be defined as:

Definite integrals can be defined as:

@TODO: Check

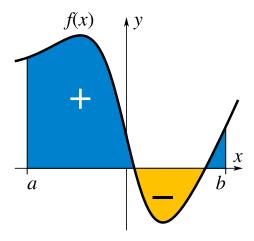


Figure 5: An integral is the sum of all area under a curve. (Contributed to Wikimedia Commons by User:KSmrq)

$$\int_{a}^{b} f(x) dx = F(b) - F(a)$$

$$\int_{a}^{b} f(x) dx$$
(23)

$$\int_{a}^{b} f(x) \, dx \tag{24}$$



Figure 6: This "S" in the Berlin-Wannsee station sign is written in the old style. It is used for integrals to stand for "sum", as the area under a curve can be imagined to be a sum (\sum) of every infinites simally thin column of space under a curve.

2.6 Taylor series

The Taylor series of any complex function f(x) (i.e., $f(x) \in \mathbb{C}$) that is infinitely differentiable at a point x_0 approximates that function as a power series:

$$f(x_0) + \frac{f'(x_0)}{1!}(x - x_0) + \frac{f''(x_0)}{2!}(x - x_0)^2 + \frac{f^{(3)}(x_0)}{3!}(x - x_0)^3 + \cdots$$
 (25)

Or, in the more-compact summation notation, this is:

$$\sum_{n=0}^{\infty} \frac{f^{(n)}(x_0)}{n!} (x - x_0)^n \tag{26}$$

where n is the number of the derivative of f.

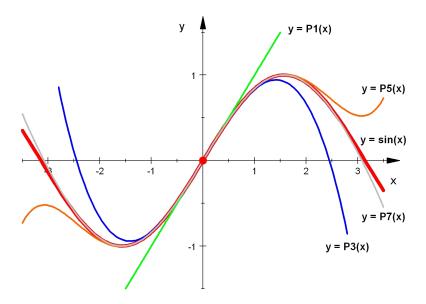


Figure 7: Increasing orders of the Taylor series expansion (and thus its derivative), denoted Pn, where $n=1,2,3,\cdots$, show the increasing approximation of a sum of polynomials to the sine function.

3 Finite difference

One of the most useful functions of computers is their ability to approximate the values of derivatives at points. This allow us to directly compute approximations of the values of derivatives in differential equations and therefore solve them, typically through space and/or time.

You may have wondered why I covered derivatives and Taylor series right after one another. This is because these provide two ways to generate finite difference solutions to differential equations. The definition of a derivative provides a way to think of a derivative as the linear tangent to a curve at a given point – and hence, also, to be well-approximated by a straight line over a finite distance so long as that distance is small compared to the nonlinearity of the curve that it is approximating. This is the same as the second term in the Taylor series expansion.

The Taylor series provides a way to consider higher-order equations that could fit a curve near a certain point. A straight line often is not a poor fit, but if we want to replace an arbitrary function with an infinitely differentiable polynomial, generating a Taylor series approximation is an easy way to do so – and thus, to numerically solve differential equations.

The result of this section will be a formalilization of the **finite difference method**, by which a set of differential equations is estimated on a grid with constant (but not necessarily uniform) spacings between grid cells (hence finite difference) and then solved. By this, I mean that the distances between grid points stays constant across the variable over which you are integrating, but may vary with respect to any other variables.

In this class, we will most likely truncate all expansions after the linear first terms: the higher-order terms are more work to include, for both you and the computer, and smaller time-steps or more clever solution nethods can often make up for not having them.

3.1 Where this fits in the spectrum of modeling

To provide a bit more background, finite difference is just one of several major methods for modeling with continuum equations. Other methods are finite volume (similar to finite difference except considering full cell volume, finite element (which allows a more irregular grid) and spectral (using how something varies in terms of periodic functions in order to write solutions to equations – this generally requires Fourier and/or other spectral transformations of the equations). Agent-based models that follow the actions of individual "agets" in a computational space also exist. These agents are often given simple rules and then allowed to interact with one another.

3.2 Thermal diffusion

In the example given here, we will solve the diffusion equation. First, I will present a derivation.

Heat is conducted at a rate that is proportional to the:

• Temperature difference (directly proportional): hotter things more quickly make other things hotter and colder things more quickly make other things colder. The contents of a bottle placed in the freezer will become cold more quickly than will those of one placed in the refrigerator

• Distance along which heat must be conducted (inversely proportional): if there is a thicker wall to a thermos, cooler, or refrigerator, it will better insulate the heat

The constant of proportionality, k, is called the "thermal conductivity" and tells one how easily heat is conducted through a medium. It is an analog to electrical conductivity.

Putting all of these together, we can write (in one dimension):

$$q_x = -k\frac{dT}{dx} \tag{27}$$

Here, q_x is the heat flux (units of Watts per meter squared [W m⁻²]). x indicates that we are solving the equation in the x-direction; we will work with a one-dimensional solution here, though you can imagine higher-dimensional solutions to require 2-D instead of 1-D arrays of values for our numerical implementation. The negative sign indicates that heat is flowing downgradient, from hotter items to colder ones.

This expression for thermal conduction (Fourier's Law of Thermal Conduction @TODO: check) can be paired with an expression that describes how the transfer of heat from one object to another results in changes in temperature. This change in temperature produces a *feedback*, by which an object that is conducting heat to another object cools (while the other object warms. This decreases the gradient in temperature between the two objects, and therefore decreases the heat flux.

As one might expect, temperature increases with heat flux into an object and decreases with heat flux out of an object. The negative spatial derivative of heat flux provides the negative divergence (i.e. convergence) of heat at the point of interest. Assuming no phase changes (e.g., solid to liquid), this occurrs at a rate that is inversely proportional to the specific heat capacity of the material and its density.

$$\frac{\partial T}{\partial t} = -\frac{1}{c_p \rho} \frac{\partial q_x}{\partial x} \tag{28}$$

We can plug Equation 27 into Equation 28 to produce a single equation for thermal conduction

$$\frac{\partial T}{\partial t} = \frac{1}{c_p \rho} \frac{\partial}{\partial x} \left(k \frac{\partial q_x}{\partial x} \right) \tag{29}$$

If we assume that the thermal conductivity, k, is spatially uniform, then we can pull it outside the derivative. This yields the most common form of the *diffusion equation*, one of the most widely-used equations in science.

$$\frac{\partial T}{\partial t} = \frac{k}{c_p \rho} \frac{\partial^2 q_x}{\partial x^2} \tag{30}$$

Redefining the coefficients on the right-hand side of the equation as κ , which is known as "thermal diffusivity" and for many Earth materials $\approx 10^{-6} \text{ m}^2 \text{ s}^{-1}$ yields:

$$\frac{\partial T}{\partial t} = \kappa \frac{\partial^2 q_x}{\partial x^2} \tag{31}$$

In addition to heat transfer, the diffusion equation cab be used to model any process in which a rate is linearly dependent on the difference in values of some state variable. This includes chemical fluxes, temperature fluxes, groundwater flow, hillslope geomorphic change, ...

3.3 Discretization

The simplest way to "discretize" an equation, or turn it into something that can represent changes over finite (i.e., $\Delta x = \text{some number} > 0$) distances, is to replace the dx or ∂x terms with a finite Δx . In doing so, we need to take into account a couple of considerations:

- 1. What do our grids of variables look like?
- 2. Do we want to generate differences forwards or backwards in time? (I am focusing on changes in time here, but you can really model a function in this case, integrating it numerically over any variable.)

3.3.1 Grids

For our thermal diffusion example, there are two (or three) important grids, each of which I will represent as a column vector (not to waste space, but to prepare you for matrix methods to be introduced in a later section). The first is our temperatures:

$$T = \begin{bmatrix} T_1 \\ T_2 \\ \vdots \\ T_m \end{bmatrix}$$
 (32)

And the second is position:

$$T = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_m \end{bmatrix} \tag{33}$$

3.3.2 Differences

In order to do finite difference calculations, we must make, well, finite differences! But how should we define our differences? There are a few ways. Using the example of T(x):

• Left-handed or "upwind" difference scheme:

$$1 (34)$$

$$\frac{\Delta T}{\Delta t} = \kappa \frac{\partial^2 q_x}{\partial x^2} \tag{35}$$

Forward difference and "implicit" on both structured and unstructured meshes Stencils

Linearizing equations (is this the right word? turning them into a set of linear equations)

Differential equations and linear algebra review Include examples

4 Example: hillslope diffusion

This example is from something I wrote for CSDMS, http://csdms.colorado.edu/wiki/Introduction_to_Python

4.1 Theory

Numerical models are fundamentally a computer-readable way to interpret rules that we have provided about how the natural world should work. Here, we are creating a simple model of a hillslope profile with rivers incising on either end. We don't worry about regolith development or any specific hillslope processes. In our case, these rules take the form of partial differential equations. It is extremely important to have a full grasp of and—insofar as it is possible—analytical solution for these equations before beginning.

Here, we make the simple assumption that regolith discharge is a linear function of slope:

$$Q = -kS \tag{36}$$

Here, Q is regolith discharge [square meters per year], S is slope [unitless], and k is a constant of proportionality that scales slope to discharge. The negative sign is present because we want positive discharge where there are negative (i.e. downhill) slopes (or, in other words, elevation decreasing as x-distance increases).

We want all of our equations in terms of elevation z, and distance across the hillslope profile, x. Substituting these into S gives us:

$$Q = -k\frac{dz}{dx} \tag{37}$$

Hillslope evolution occurs when we combine these discharges to conserve mass. We use a continuity equation for conservation of volume to state that:

$$\frac{\partial z}{\partial t} = -\frac{\partial Q}{\partial x} \tag{38}$$

Combining these two terms together gives us the diffusion equation:

$$\frac{\partial z}{\partial t} = \frac{\partial}{\partial x} \left(k \frac{\partial z}{\partial x} \right) \tag{39}$$

If we assume that the hillslope diffusivity, k, is uniform across the landscape, we obtain the diffusion equation in its more familiar form:

$$\frac{\partial z}{\partial t} = k \frac{\partial^2 z}{\partial x^2} \tag{40}$$

4.2 Source code

First, I will write the whole source code. Then I will break it down into sections where I explain each component. Then I will present a new model that uses an implicit matrix-style solution instead of a foward-time-stepping solution, and write about numerical (finite difference) methods.

```
/usr/bin/python
  # Import modules
  import numpy as np
  from matplotlib import pyplot as plt
  Copyright 2012 Andrew D. Wickert
      This program is free software: you can redistribute it and/or modify it under the terms of the GNU General Public License as published by
      the Free Software Foundation, either version 3 of the License, or
      (at your option) any later version.
      This program is distributed in the hope that it will be useful,
      but WITHOUT ANY WARRANTY; without even the implied warranty of
      MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE.
      GNU General Public License for more details.
      You should have received a copy of the GNU General Public License along with this program. If not, see <http://www.gnu.org/licenses/>.
21
23
      hill.py
25
    HILLSLOPE PROFILE
27
    Center of hill is at 0, two incising channels at steady rate at edges
    no channel-hillslope feedback and assuming entire hill is made of mobile
    regolith
    This is a very simple example code in which variables are defined inline,
    and the equations are solved by forward difference methods without
    consideration for stability
    Written by Andrew D. Wickert; August, 2013
```

```
# We will have the simple boundary condition of rivers on either end of the
    # hillslope that are incising at the same, steady rate.
    zdot channel = 2E-4 # [m], 0.5 mm/yr boundary incision
43
    # time
    \mathrm{dt} = 100.~\# [years], the time step (from guessing, not from stability analysis
45
       final = 4E5 # [years], the final time for hillslope evolution
    timesteps = np.arange(0, t_final + dt/10., dt) # [years]
    # Set up domain
    dx = 10 \# [m]

xmax = 500 \# [m]
51
    x = np.arange(-xmax, xmax + dx/10., dx) \# [m], +dx/10. to make sure that edges
            are included
    z = t\_final * zdot\_channel * np.ones(x.shape) \# [m], elevation in meters - set such that the edges are 0 at t\_final
    \# We're using a very simplified assumption that the rate of hillslope \# material transport is linearly proportional to local slope, and that this \# constant of proportionality is constant across the hill
55
    \dot{k}=5E-3~\#~Slope~-
                                   - soil flux scaling
59
    # Loop through time and evolve elevation
    for t in timesteps:
    # The channels set the boundaries
61
       # The channels set the boundaries z[0] = z \det_{channel} * \det_{class} z[0] = z \det_{channel} * \det_{class} z[-1] = z \det_{class} z \det_{class} z[-1] = a[n+1] - a[n]
# We are calculating the slopes between each of the cells, then using
63
       # these to solve for discharge of material between interior cells. 
 S = np. diff(z) \# [-] slope   Q = -k*S \# [m**2/yr] discharge, goes downslope, hence the negative sign # POSITIVE - WHY? CANCELLED OUT AGAIN!
67
69
       \# The change in internal elevation, due to conservation of mass, is equal
       # once again to the x-defivative
      # once again to the x-defivative dzdt_interior = np. diff(Q) z[1:-1] -= dzdt_interior * dt # (It would have been quicker to take both derivatives at once, but would be # somewhat less intuitive for descriptive purposes)
    # Plot
    # Plot plt.plot(x,z,'k-',linewidth=3) plt.title('Final hillslope profile', fontsize=20, weight='bold') plt.xlabel('Distance across hillside [m]', fontsize=16) plt.ylabel('Elevation [m]', fontsize=16) plt.xlim((x[0],x[-1]))
    plt.show()
```

4.2.1 The "Shebang"

This script starts with:

```
#! /usr/bin/python
This is a "shebang" that tells the program where to look for the Python
interpreter on Unix-like systems (the most popular being Mac and Linux,
the former more popular for desktop applications and the latter more
popular for supercomputers). You can find where Python is located by
typing:

3 <syntaxhighlight lang=bash>
which python
```

at the Unix command prompt (i.e. terminal).

4.2.2 Module Import (and Namespaces)

```
# Import modules
import numpy as np
from matplotlib import pyplot as plt
```

Python is a general-purpose language, which means that for specific functionalities, you have to import "modules". These modules are sets of functions that share a namespace ("np" and "plt", respectively). "'Numpy" does general mathematical and "n"-dimensional array operations, and "'Pyplot" does basic plotting. "Sharing a namespace" means that to call functions within these modules, we must use the "np" and "plt" prefixes, e.g., to plot a parabola:

```
a = np.array([[0,1,2,3,4],[0,1,4,9,16]]) # Create an array with 2 rows and 3 columns - nested sets of brackets add dimensions to the array plt.plot(a[0,:], a[1,:], 'ko-') # Plot the first row of "a" as ''x''-values and the second row as ''y''-values with a dashed line between points plt.show() # Show the plot in a display screen
```

Namespaces are important, because it means that you cannot easily accidentally overwrite the name of a function by giving a variable that name. For example, typing "np.mean = 1.5" seems a less-likely accident than writing "mean = 1.5".

4.2.3 Copyright Notice

The aforementioned code is trivial, but lately (2012) it seems that big, ugly patent wars are being fought over trivial chunks of code. So better to be safe than sorry. Below is text included for version 3 of the [http://www.gnu.org GNU] General Public License (GPL). This is an open-source "copyleft" license, meaning that if this code is included in a project, the project must be open-source with a "copyleft" license as well. This is a very good license for science, which is a community-driven collaborative effort. To give your code the GNU GPL v3, include the following text (though replacing my name with yours in the copyright, unless you are feeling generous).

In order for your code to be included in CSDMS, you need to include some type of open-source license, even if it is not as strong a license as the GPL v3.

```
"""
Copyright 2012 Andrew D. Wickert

This program is free software: you can redistribute it and/or modify it under the terms of the GNU General Public License as published by the Free Software Foundation, either version 3 of the License, or (at your option) any later version.
```

```
This program is distributed in the hope that it will be useful, but WITHOUT ANY WARRANTY; without even the implied warranty of MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE. See the GNU General Public License for more details.

You should have received a copy of the GNU General Public License along with this program. If not, see <a href="https://www.gnu.org/licenses/">http://www.gnu.org/licenses/</a>.
```

On a technical note, the triple quotes at the beginning and end of the block are used in Python to create blocks of text as strings. These can be used later inside functions, blocks of code designed for a single purpose, to provide accessible documentation. For now, since we are not assigning this string anywhere, it simply functions as a comment.

4.2.4 Introductory Comments and Authorship

This section is huge overkill for this short of a model, but I put it in here to be complete.

4.2.5 Initialization

```
# We will have the simple boundary condition of rivers on either end of the # hillslope that are incising at the same, steady rate. zdot_channel = 2E-4 # [m], 0.5 mm/yr boundary incision  

# time dt = 100. # [years], the time step (from guessing, not from stability analysis )

t_final = 4E5 # [years], the final time for hillslope evolution timesteps = np.arange(0,t_final + dt/10.,dt) # [years]

# Set up domain dx = 10 # [m]  
xmax = 500 # [m]
```

```
13 x = np.arange(-xmax, xmax + dx/10., dx) \# [m], +dx/10. to make sure that edges are included z = t\_final * zdot\_channel * np.ones(x.shape) \# [m], elevation in meters - set such that the edges are 0 at <math>t\_final 15 \# We're using a very simplified assumption that the rate of hillslope 17 \# material transport is linearly proportional to local slope, and that this \# constant of proportionality is constant across the hill \# 5E-3 \# Slope — soil flux scaling
```

4.2.6 Main Loop

```
# Loop through time and evolve elevation
for t in timesteps:
# The channels set the boundaries

z [0] -= zdot_channel * dt

z [-1] -= zdot_channel * dt

# for np.diff, out [n] = a [n+1] - a [n]

# We are calculating the slopes between each of the cells, then using
# these to solve for discharge of material between interior cells.

S = np.diff(z) # [-] slope
Q = -k*S # [m**2/yr] discharge, goes downslope, hence the negative sign #
POSITIVE - WHY? CANCELLED OUT AGAIN!

# The change in internal elevation, due to conservation of mass, is equal
# once again to the x-defivative
dzdt_interior = np.diff(Q)
z [1:-1] -= dzdt_interior * dt
# (It would have been quicker to take both derivatives at once, but would be
# somewhat less intuitive for descriptive purposes)
```

4.2.7 Plotting

```
# Plot
plt.plot(x,z,'k-',linewidth=3)
plt.title('Final hillslope profile', fontsize=20, weight='bold')
plt.xlabel('Distance across hillside [m]', fontsize=16)
plt.ylabel('Elevation [m]', fontsize=16)
plt.xlim((x[0],x[-1]))
plt.show()
```

4.3 Numerical stability

Run hillslope_diffusion_no_matrix.py or hillslope_diffusion_no_matrix_plot_through_time.py, found in the code/NumericalMethods folder. You get a nice, rounded hillslope!

Now, run the same code, but change dx to 101 years. What happens? It goes unstable. You can play around with this.

What is the problem? Well, we are projecting out into the future along a straight line. sometimes this is a good approximation. Sometimes this isn't.

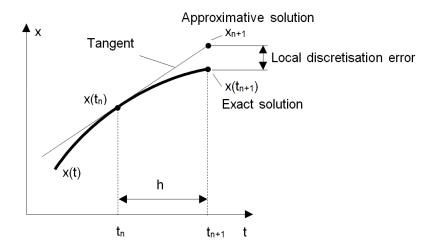


Figure 8: Euler forward numerical integration discretization error; figure from course notes at http://moodle.autolab.uni-pannon.hu/Mecha_tananyag/mechatronikai_modellezes_angol/ch07.html

When is it a good approximation? And how can we improve the approximation? Well, because our solution (hillslope diffusion) is directly analogous to the classic thermal diffusion problem, we can follow a scheme so classic that it is even on Wikipedia (https://en.wikipedia.org/wiki/Von_Neumann_stability_analysis).

How do we improve stability?

- Reduce the time-step
- Increase the spatial discretization
- Write a local approximation that is closer to the curve (Taylor series approximation)
- Find a clever mathematical way around our "shooting forward in time [or other variable]" method of numerical integration.