
- Numerical simulation of a stratigraphic model - single lithology case

Single lithology stratigraphic model

Our objective is to simulate the infill of sedimentary basins at large space and time scales.

- $\Omega = (0, L_x)$: horizontal extension of the basin
- $(0, T)$ is the time interval of the simulation with $T > 0$
- $h(x, t)$ is the sediment thickness for $(x, t) \in \Omega \times (0, T)$
- $h_{\text{sea}}(t)$ is the given sea level for $t \in (0, T)$
- $b(x, t) = h_{\text{sea}}(t) - h(x, t)$ is the bathymetry

The model accounts for the following sediment thickness conservation and boundary and initial conditions

$$\left\{ \begin{array}{l} \partial_t h(x, t) + \text{div}(\nabla \psi(b(x, t))) = 0 \text{ on } \Omega \times (0, T) \\ h(x, 0) = h_{\text{init}}(x) \text{ on } \Omega \\ \nabla \psi(b(x, t)) \cdot \mathbf{n} = g_0 \text{ at } x = 0 \\ \nabla \psi(b(x, t)) \cdot \mathbf{n} = g_1 \text{ at } x = L_x \end{array} \right. \quad (1)$$

where $\psi(b) = \int_0^b k(u) du$ with $k(b) > 0$ the diffusion coefficient of the sediments measuring their ability to be transported by gravity. This coefficient is modeled by a nonlinear function of the bathymetry as follows

$$k(b) = \begin{cases} k^m & \text{if } b \geq 0 \\ k^c & \text{otherwise} \end{cases}$$

Finite Volume discretization

The model is discretized using a Two Points Flux Approximation (TPFA) on an orthogonal mesh. We obtain at each time step $n > 0$ and for each cell K

$$\left\{ \begin{array}{l} |K| \frac{h_K^n - h_K^{n-1}}{\Delta t^n} + \sum_{\sigma=K|L \in \Sigma_K \cap \Sigma_{\text{int}}} T_\sigma (\psi(b_L^n) - \psi(b_K^n)) + \sum_{\sigma \in \Sigma_K \cap \Sigma_b} g_\sigma = 0 \\ \text{with } T_{\sigma=K|L} = \frac{|\sigma|}{\text{dist}(x_K, x_L)} \\ b_K^n = h_{\text{sea}}(t^n) - h_K^n. \end{array} \right. \quad (2)$$

The initial condition is computed by

$$h_K^0 = h_{\text{init}}(x_K).$$

Scilab implementation

The data structure needed for the implementation of the scheme for the given uniform 1D mesh with N cells of the domain $(0, L_x)$ is already given.

You need to implement the scheme (2) by completing the given Scilab file.

At each time step (inside the time loop), solve using a Newton algorithm the nonlinear system (2) denoted by

$$R(h^n) = 0$$

where $R : \mathbb{R}^N \rightarrow \mathbb{R}^N$ is the so-called “residual” function representing the finite volume scheme conservation equations in all cells. To do this, the Newton algorithm is applied: set $\epsilon = 10^{-6}$, $h^{0,n} = h^{n-1}$, and for $l \geq 0$ until $\|R(h^{l,n})\| \leq \epsilon \|R(h^{0,n})\|$ compute

$$\frac{\partial R}{\partial h}(h^{l,n}) dh = -R(h^{l,n}),$$

$$h^{l+1,n} = h^{l,n} + dh.$$

We underline that the Newton algorithm has a quadratic convergence if the initial solution $h^{0,n}$ is closed enough to the solution h^n , ie there exist $\alpha > 0$ and $\beta > 0$ such that if $\|h^{0,n} - h^n\| \leq \alpha$ then

$$\|R(h^{l+1,n})\| \leq \beta \|R(h^{l,n})\|^2. \quad (3)$$

Note also that if the Newton algorithm is not converged in less than *Newtmax* iterations, then the time step is restarted using a reduced time step by a factor 2. If the time step is converged in less than *Newtmax* iterations we can increase the time step by a factor 1.2 until the maximum time step is reached.

- (1) Write the Scilab function computing the **residual** $R(h^n)$ given h^{n-1} and $h_{sea}(t^n)$. This computation is achieved using one loop on the cells, one loop on the inner faces, and one loop on the boundary faces.
- (2) Write the Scilab function computing the **jacobian** $\frac{\partial R}{\partial h}(h^n)$ given $h_{sea}(t^n)$. This computation is achieved using one loop over the cells, one loop over the inner faces, and one loop over the boundary faces.
- (3) Complete the Newton and the time loops. By using a logarithm scale, check graphically the quadratic convergence of the Newton algorithm.
- (4) Plot h and b function of x for the computed discrete times (one plot for h and one plot for b).
- (5) Write the algorithm to compute

$$h_s(x, t) = \min_{\{t \leq q \leq T\}} h(x, q).$$

Comment on the geological meaning of h_s ?

Then, plot $h_s(x, t)$ function of x for all the discrete times t^n , $n = 0, \dots$, nb time steps, of the simulation (in a single plot).

- (6) (More difficult) We want to compute the paleobathymetry $b_p(x, z, t)$ in the basin defined as the bathymetry of the sediment at the time it has been deposited.

It is the solution of the equation:

$$\begin{cases} \partial_t b_p(x, z, t) = 0 \text{ on } \mathcal{B} \\ b_p(x, h(x, t), t) = b(x, t) \text{ if } \partial_t h(x, t) > 0 \text{ (sedimentation)} \\ b_p(x, z, 0) = b_{p,\text{init}}(x) \text{ for } z \in (-\infty, h_{\text{init}}(x)). \end{cases} \quad (4)$$

with

$$\mathcal{B} = \{(x, z, t) \mid (x, t) \in \Omega \times (0, T), z \in (-\infty, h(x, t))\}.$$

The initial paleobathymetry $b_{p,\text{init}}(x)$ will be assumed constant in the following to fix ideas.

The above equation is discretized using the following algorithm: let at time $t = 0$ $b_{p,K}^0(z) = b_{p,\text{init}}$ (constant) for all $z < h_K^0$. Then we compute at each time step:

If $h_K^n \geq h_K^{n-1}$ one has

$$\begin{cases} b_{p,K}^n(z) = b_{p,K}^{n-1}(z) \text{ for all } z < h_K^{n-1}, \\ b_{p,K}^n(z) = b_K^n = h_{\text{sea}}(t^n) - h_K^n \text{ for all } h_K^{n-1} < z < h_K^n, \end{cases}$$

Else if $h_K^n < h_K^{n-1}$ one has

$$b_{p,K}^n(z) = b_{p,K}^{n-1}(z) \text{ for all } z < h_K^n.$$

Code the algorithm in scilab (initialization before the time loop and at each time step in the time loop after the computation of h^n) using the following arrays:

The paleobathymetry $b_{p,K}^n(z)$ is piecewise constant on the intervals

$$]-\infty, h_{\text{col}}(K, 1)],]h_{\text{col}}(K, 1), h_{\text{col}}(K, 2)], \dots]h_{\text{col}}(K, n_{\text{col}}(K) - 1), h_{\text{col}}(K, n_{\text{col}}(K))],$$

where $n_{\text{col}}(K)$ is the number of intervals. The value of the concentration $b_{p,K}^n(z)$ in each interval is stored in $b_{\text{col}}(K, l)$ for $l = 1, \dots, n_{\text{col}}(K)$.

The arrays n_{col} , h_{col} and b_{col} have to be updated according to the above algorithm at each time step and in each cell K .

Question: Write the corresponding code in scilab and plot the paleobathymetry at final time T in 3 wells located at one third, one half and two third of the basin as functions of z .

- (8) Comment the physical results that you obtain. What do you think of this model? Why is it used in oil exploration?
- (9) Look at the plots of the solution obtained for different mesh sizes $N = 10, 20, 50, 100, 200, 400$ and $\Delta t = 0.02$ and comment the results.

- (10) Look at the plot of the solutions obtained for different time steps and $N = 100$ and comment the results.
- (11) Comment the behavior of the Newton algorithm for the different mesh sizes and time steps values