

## Exercise 6. Iterative Solution of Matrix Problems.

1. Start with your code that solved the diffusion equation explicitly. Adjust it to always take timesteps at the stability limit  $\Delta t = \Delta x^2/2D$ , so that:

$$\psi_j^{(n+1)} - \psi_j^{(n)} = \left(\frac{1}{2}\psi_{j+1}^{(n)} - \psi_j^{(n)} + \frac{1}{2}\psi_{j-1}^{(n)} + \frac{s_j^{(n)}}{2D}\Delta x^2\right).$$

Now it is a Jacobi iterator for solving the steady-state elliptical matrix equation. Implement a convergence test that finds the maximum absolute *change in  $\psi$*  and divides it by the maximum absolute  $\psi$ , giving the normalized  $\psi$ -change. Consider the iteration to be converged when the normalized  $\psi$ -change is less than (say)  $10^{-5}$ . Use it to solve

$$\frac{d^2\psi}{dx^2} = -1$$

on the domain  $x = [0, 1]$  with boundary conditions  $\psi = 0$  at  $x = 0$ ,  $\frac{\partial\psi}{\partial x} = 0$  at  $x = 1$ , with a total of  $N_x$  equally-spaced nodes. Find how many iterations it takes to converge, starting from an initial state  $\psi = 0$ , when

- (a)  $N_x = 10$
- (b)  $N_x = 30$
- (c)  $N_x = 100$

Compare the number of iterations you require with the analytic estimate in the notes. How good is the estimate?

Now we want to check how accurate the solution really is.

- (d) Solve the equation analytically, and find the value of  $\psi$  at  $x = 1$ ,  $\psi(1)$ .
- (e) For the three  $N_x$  values, find the relative error<sup>1</sup> in  $\psi(1)$ .
- (f) Is the actual relative error the same as the convergence test value  $10^{-5}$ ? Why?

**Enrichment only**, optional and not for credit. Turn your iterator into a SOR solver by splitting the iteration matrices up into red and black (odd and even) advancing parts. Each part-iterator then uses the latest values of  $\psi$ , that has just been updated by the other part-iterator. Also provide yourself an over-relaxation parameter  $\omega$ . Explore how fast the iterations converge as a function of  $N_x$  and  $\omega$ .

Note. Although in Octave/MATLAB® it is convenient to implement the matrix multiplications of the advancing using a literal multiplication by a big sparse matrix, one does not do that in practice. There are far more efficient ways of doing the multiplication, that avoid all the irrelevant multiplications by zero.

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<sup>1</sup>the difference between the “converged” iterative  $\psi(1)$  and the analytic  $\psi(1)$  normalized to the analytic value.

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