# Lecture 5: Systems of nonlinear equations

COMP5930M Scientific Computation

# Today

Notation

Newton's method

The Jacobian matrix

Efficiency

Next

### **Notation**

► Single equation in a single unknown

$$F(x) = 0$$

n equations in n unknowns

$$F(x) = 0$$

 $\underline{\mathbf{x}}$  is a vector  $\{x_1, x_2, ..., x_n\}$  of n unknown values  $\underline{\mathbf{F}}$  is a set  $\{F_1(\mathbf{x}), F_2(\mathbf{x}), ... F_n(\mathbf{x})\}$  of n nonlinear equations

# The nonlinear problem

Find the *n*-dimensional point  $\{x_j^*\}$ , j = 1, 2, ...n such that the set of functions

$$F_i(x_j^*) = 0$$
,  $i = 1, 2, ...n$  simultaneously

## Example: n = 2 system

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$$\frac{\partial F_1}{\partial x_1} = 6 x_1 + 4 x_2, \qquad \frac{\partial F_1}{\partial x_2} = 4 x_1$$

$$\frac{\partial F_2}{\partial x_1} = 2x_1 + 2 x_2^2, \qquad \frac{\partial F_2}{\partial x_2} = 4 x_1 x_2$$

## Newton's Method

- Assume we have access to the set of functions F(x)
- Assume we know an initial state x<sub>0</sub>

These first two are the minimum information necessary

Assume we have access to all the partial derivatives of **F** with respect to **x** 

$$\frac{\partial F_i}{\partial x_i}$$
,  $i = 1, 2, ..., n$ ,  $j = 1, 2, ..., n$ 

## Derivation of Newton's method in the vectorial case

Given current iterate  $\mathbf{x}_k$ , find an increment  $\delta \in \mathbb{R}^n$  s.t.

$$\mathbf{F}(\mathbf{x}_k + \delta) = 0. \tag{1}$$

If the function **F** is differentiable at  $\mathbf{x}_k$ , we can linearise it:

$$\mathbf{F}(\mathbf{x}_k + \delta) = \mathbf{F}(\mathbf{x}_k) + \frac{\partial \mathbf{F}}{\partial \mathbf{x}}(\mathbf{x}_k)\delta + o(|\delta|^2).$$

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Dropping the second order terms, we can solve for  $\delta$  from (1):

$$\mathbf{x}_{k+1} - \mathbf{x}_k = \delta = -\left[\frac{\partial \mathbf{F}}{\partial \mathbf{x}}(\mathbf{x}_k)\right]^{-1} \mathbf{F}(\mathbf{x}_k).$$

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# The algorithm

Generate a sequence of approximations  $\boldsymbol{x}_1, \ \boldsymbol{x}_2, \ ...$  using

$$\mathbf{x}_{k+1} = \mathbf{x}_k - \left(\frac{\partial \mathbf{F}}{\partial \mathbf{x}}(\mathbf{x}_k)\right)^{-1} \mathbf{F}(\mathbf{x}_k)$$

starting from an initial guess  $\mathbf{x}_0$ .

### The Jacobian Matrix

The  $n \times n$  matrix of partial derivatives is called the Jacobian, J, where  $J_{ij} = \frac{\partial F_i}{\partial x_i}$ 

Newton's Method is more compactly written as

$$\mathbf{x}_{k+1} = \mathbf{x}_k - \mathbf{J}(\mathbf{x}_k)^{-1}\mathbf{F}(\mathbf{x}_k)$$

This implies the solution of an  $n \times n$  linear system at every step

This is a costly part of the algorithm

## Rewrite as 2-step algorithm

$$\begin{array}{ccc}
\mathbf{J}(\mathbf{x}_k)\delta & = & -\mathbf{F}(\mathbf{x}_k) \\
\mathbf{x}_{k+1} & = & \mathbf{x}_k + \delta
\end{array}$$

#### Notes:

- Step 2 is trivial in this basic form
- Step 1 is a linear algebra problem: system of  $n \times n$  linear equations with J a known  $n \times n$  matrix at each iteration .

# Computational cost

	Function calls	Algorithmic
Evaluate <b>J</b>	?	n <sup>2</sup>
Solve $\mathbf{J}\delta = -\mathbf{F}$	1	$n^3$
Update $\mathbf{x}_{k+1}$	0	n

# Evaluating the Jacobian in practice

- Analytical Jacobian may be expensive to evaluate
- Numerical approximation of Jacobian requires  $2n^2$  evaluations
- Quasi-Newton methods rely on approximation of Jacobian that is updated at each step

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# Evaluating the Jacobian numerically

- We require  $\frac{n^2}{n^2}$  values to fill the Jacobian matrix
- Numerical approximation term-by-term

$$\frac{\partial F_i}{\partial x_j} \approx \frac{F_i(x_1, ..., x_j + \delta_j, ..., x_n) - F_i(x_1, ..., x_n)}{\delta_j}$$

- Efficient implementation
  - $\triangleright$  Can simultaneously perturb all the  $F_i$  with respect to  $x_i$
  - Equivalent to n Matlab function calls
     where a function call evaluates all the F<sub>i</sub>

## Jacobian for the n=2 system

```
% Vectorised version of the function F(x1,x2)
F=0(x1,x2)([3*x1.^2 + 4.*x1.*x2;
          x1.^2 + 2*x1.*x2.^2);
\% Optimal choice of perturbation parameter h
h = 10 * sqrt(eps);
% Call F(x1,x2) once and store
Fx = F(x1, x2):
% Numerical Jacobian based on difference approximation
dFnum = [ (F(x1+h,x2) - Fx) / h ...
          (F(x1.x2+h) - Fx) / h:
```

# Computational cost



	Function calls	Algorithmic
Evaluate <b>J</b>	n	n <sup>2</sup>
Solve $\mathbf{J}\delta = -\mathbf{F}$	1	n <sup>3</sup>
Update $\mathbf{x}_{k+1}$	0	n

## Problems?

- More difficult to make the algorithm robust overall
  - There is no bisection method in higher dimensions



- For this reason damped Newton-like methods are preferred
  - Take steps in the direction  $\frac{\delta}{||\delta||}$  but control the step size to avoid divergence when the Jacobian  $\partial \mathbf{F}/\partial \mathbf{x} \approx \mathbf{0}$ .
  - In some circumstances this allows global convergence of Newton-type algorithms

#### Next time...

#### **Tutorial**

- ► Solving nonlinear systems with MATLAB
- Problems with convergence of non-damped Newton's method

#### Lecture

- Line-search algorithms for Newton's method
- Computational algorithms for systems