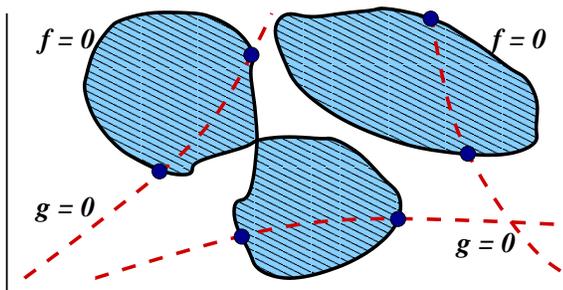


# Class 9. Root Finding in Multi-D, and Numerical Differentiation

## Nonlinear Systems of Equations

- Consider the system  $f(x, y) = 0, g(x, y) = 0$ . Plot zero contours of  $f$  and  $g$ :



- No information about  $f$  in  $g$ , and *vice versa*.
  - In general, no good method for finding roots.

- If you are near root, best bet is NR.

E.g., For  $\mathbf{F}(\mathbf{x}) = \mathbf{0}$ , choose  $\mathbf{x}_{i+1} = \mathbf{x}_i + \boldsymbol{\delta}$ , where  $\mathbf{F}'(\mathbf{x})\boldsymbol{\delta} = -\mathbf{F}(\mathbf{x})$ .

- This is a matrix equation:  $\mathbf{F}'(\mathbf{x})$  is a matrix with elements  $\partial F_i / \partial x_j$ . The matrix is called the Jacobian.

- Written out (2-D example):

$$\begin{aligned}\frac{\partial f}{\partial x} \delta_x + \frac{\partial f}{\partial y} \delta_y &= -f(x, y), \\ \frac{\partial g}{\partial x} \delta_x + \frac{\partial g}{\partial y} \delta_y &= -g(x, y).\end{aligned}$$

- Given initial guess, must evaluate matrix elements and RHS, solve system for  $\boldsymbol{\delta}$ , and compute next iteration  $\mathbf{x}_{i+1}$ . Then repeat (must solve  $2 \times 2$  linear system each time).
- Essentially the non-linear system has been linearized to make it easier to work with.
- *NRiC* §9.7 discusses a global convergence strategy that combines multi-D NR with “backtracking” to improve chances of finding solutions.

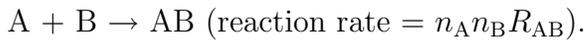
## Example: Interstellar Chemistry

- ISM is multiphase plasma consisting of electrons, ions, atoms, and molecules.
- Originally, the ISM was thought to be too hostile for molecules.

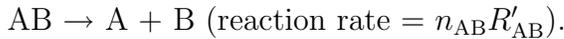
- But in 1968-69, radio observations discovered absorption/emission lines of  $\text{NH}_3$ ,  $\text{H}_2\text{CO}$ ,  $\text{H}_2\text{O}$ , ...
- Lots of organic molecules, e.g.,  $\text{CH}_3\text{CH}_2\text{OH}$  (ethanol), etc.
- In some places, all atoms have been incorporated into molecules.
- E.g., molecular clouds: dense, cold clouds of gas composed primarily of molecules.

$$(T \sim 30 \text{ K}, n \sim 10^6 \text{ cm}^{-3}, M \sim 10^{5-6} M_\odot, R \sim 10-100 \text{ pc.})$$

- How do we predict what the abundances of different molecules should be, given  $n$  and  $T$ ?
- Need to solve a chemical reaction network.
- Consider reaction between two species A and B:



- Reverse also possible:

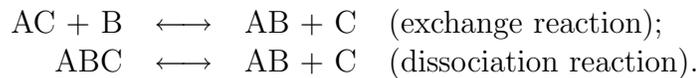


- In equilibrium:

$$\begin{aligned} n_{\text{A}}n_{\text{B}}R_{\text{AB}} &= n_{\text{AB}}R'_{\text{AB}}; \\ n_{\text{A}} + n_{\text{AB}} &= n_{\text{A}}^0; \\ n_{\text{B}} + n_{\text{AB}} &= n_{\text{B}}^0. \end{aligned}$$

where  $n_{\text{A}}^0$  and  $n_{\text{B}}^0$  are normalizations so that A and B are conserved.

- Substitute normalization equations into reaction equation to get quadratic in  $n_{\text{AB}}$ , easily solved.
- However, many more possible reactions:



- Wind up with large nonlinear system describing all forward/reverse reactions, involving known reaction rates  $R$ , plus normalizations. Must solve given fixed  $n^0$  and  $T$ .

## Numerical Derivatives

- For NR and function minimization, often need derivatives of functions. It's always better to use an analytical derivative if it's available.

- If you're stuck, could try:

$$f'(x) \simeq \frac{f(x+h) - f(x)}{h},$$

where  $|h|$  is small.

- However, this is very susceptible to RE. Better:

$$f'(x) \simeq \frac{f(x+h) - f(x-h)}{2h}.$$

(This version cancels the second-derivative term in the Taylor series expansion of  $f(x+h) - f(x-h)$ , leaving just the third- and higher-order terms.)

- Read *NRiC* §5.7 before trying this!