

## Numerical Solution of Laplace's Eq.

$$\nabla^2 \phi = -\frac{P}{\epsilon} \quad (\text{Poisson's Eq.})$$

$$\nabla^2 \phi = 0 \quad (\text{Laplace's Eq.})$$

notice that Laplace's equation is a homogeneous P.D.E.  $\therefore \phi = 0$  is automatically a solution. Where do the non-trivial solutions come from?

{ Ans: imposed boundary conditions determine the form of the solution.

example (1-D) :

$$(O.D.E.) \quad \frac{d^2 \phi}{dx^2} = 0 \quad 0 \leq x \leq 1$$

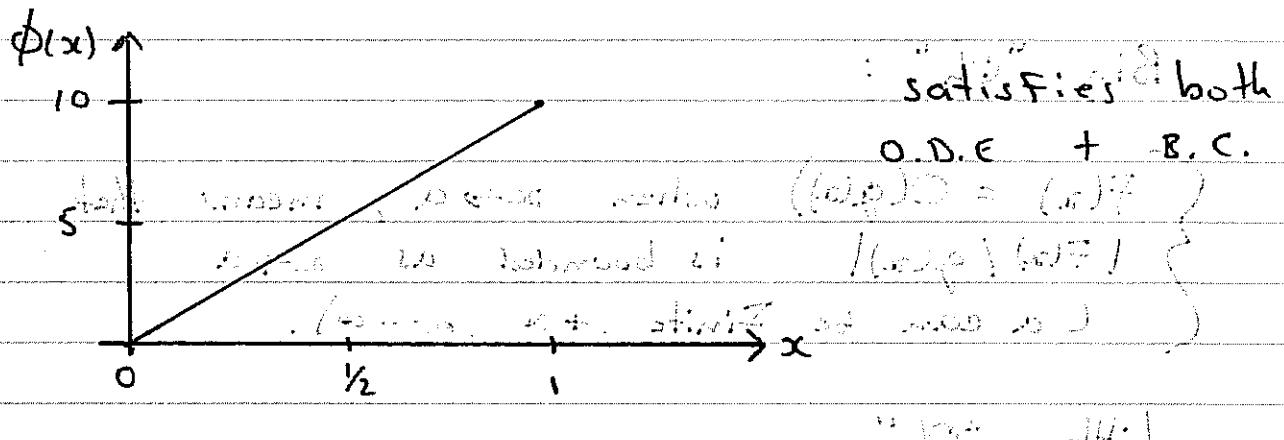
the general solution is

$$\phi(x) = k_1 x + k_2$$

$k_1, k_2$  - arbitrary constants.

$$(B.C.) \quad \phi(0) = 0, \quad \phi(1) = 10 \quad (\text{Dirichlet}).$$

$$\phi(0) = k_2 = 0 \quad \phi(1) = 10 = k_1 \\ \therefore \boxed{\phi(x) = 10x}$$



How would we apply some numerical approximations to the boundary value problem (B.V.P)

$$\frac{d^2\phi}{dx^2} = 0 \quad \text{B.C. } \phi(0) = 0, \phi(1) = 10 ?$$

Main idea in finite difference techniques is to approximate the fcn we're looking for  $\phi(x)$  by a grid fcn

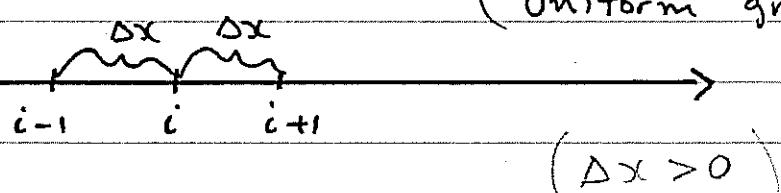
$$\phi_i \approx \phi(i\Delta x)$$

and then discretize the differential operators. (usually using Taylor's expansion).

Given the grid fcn  $f_i \approx f(i\Delta x)$

$$(i \in \mathbb{Z})$$

(uniform grid)



## Taylor's expansion

$$f_{i+1} = f_i + \Delta x f'_i + \frac{\Delta x^2}{2} f''_i + \frac{\Delta x^3}{3!} f'''_i + \dots \quad (1)$$

Solving for  $f'_i \approx f'(i \Delta x)$

$$f'_i = \frac{f_{i+1} - f_i}{\Delta x} - f''_i \frac{\Delta x}{2} - f'''_i \frac{\Delta x^2}{3!} - \dots \quad (1)$$

Forward difference approximation: (F. d.)

$$f'_i = \frac{f_{i+1} - f_i}{\Delta x} + O(\Delta x)$$

$O(\Delta x) \rightarrow$  "of order  $\Delta x$ "  
 $\rightarrow$  truncation error due  
 to replacing a continuous  
 process by a discrete one.

continuous derivative

$\Rightarrow$

Discrete finite  
difference.

$(\Delta x > 0)$

$$f_{i+1} = f_i - \Delta x f'_i + \frac{\Delta x^2}{2} f''_i - f'''_i \frac{\Delta x^3}{6} + \dots \quad (2)$$

$$f'_i = \frac{f_i - f_{i-1}}{\Delta x} + \frac{\Delta x}{2} f''_i - f'''_i \frac{\Delta x^2}{6} \quad (2')$$

backward difference (b.d.)

$$f'_i = \frac{f_i - f_{i-1}}{\Delta x} + O(\Delta x)$$

now take eq. (1) - eq (2) :

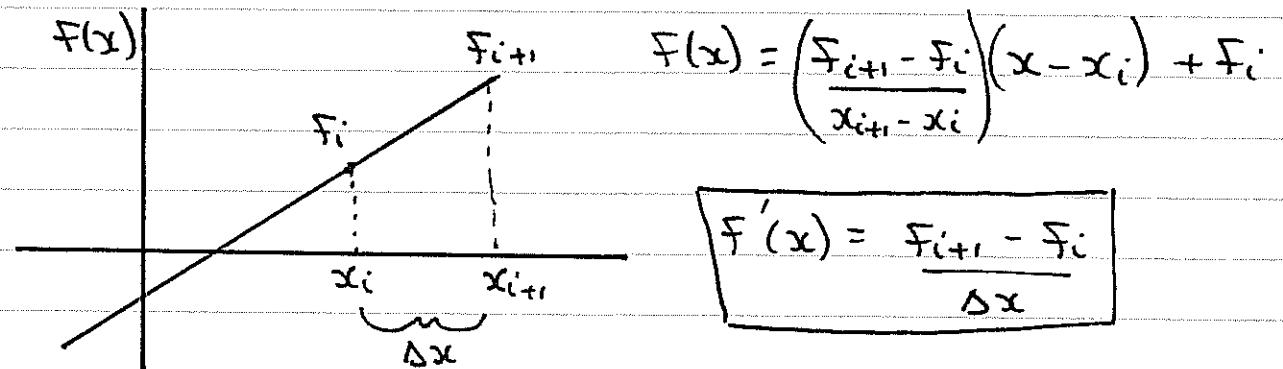
$$f_{i+1} - f_{i-1} = f'_i 2\Delta x + f'''_i \frac{\Delta x^3}{3} + \dots$$

$$f'_i = \frac{f_{i+1} - f_{i-1}}{2\Delta x} - f'''_i \frac{\Delta x^2}{6} + \dots$$

central difference

$$f'_i = \frac{f_{i+1} - f_{i-1}}{2\Delta x} + O(\Delta x^2)$$

1) Let us show that the F.d. and b.d. operators are exact for linear fns.



(obvious - slope)

(For b.d. just reliable  $i+1 \rightarrow i$   $i \rightarrow i-1$ ).

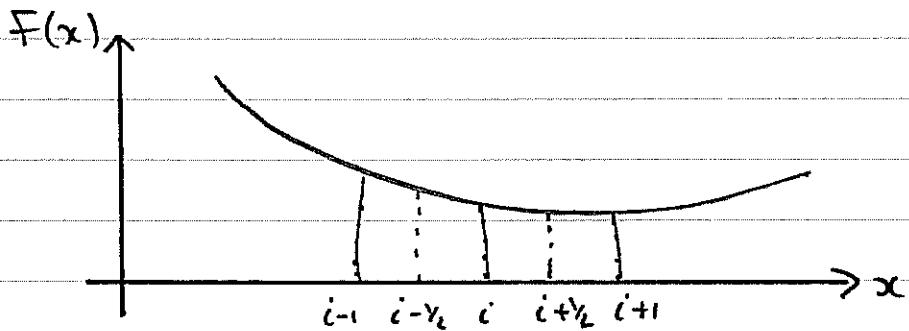
2) Show that the central difference operator is exact for quadratic fns.

$$f'_i = \frac{f_{i+1} - f_{i-1}}{2\Delta x} - f''' \frac{\Delta x^2}{6} + \dots$$

if  $f(x) = a + bx + cx^2$  (arbitrary quadratic).  
 $\Rightarrow f'''(x) = 0$

$\therefore$  c.d. exact.

approximations for higher order derivatives:



using central differences at  $i-\gamma_2$ ,  $i+\gamma_2$ :

$$f'_i + \gamma_2 = \frac{f_{i+1} - f_i}{\Delta x} + O(\Delta x^2)$$

$$f'_{i-\gamma_2} = \frac{f_i - f_{i-1}}{\Delta x} + O(\Delta x^2)$$

$$\therefore f''_i = \frac{f_{i+\gamma_2} - f'_{i-\gamma_2}}{\Delta x} + O(\Delta x^2)$$

$$= \frac{f_{i+1} - f_i}{\Delta x} - \frac{f_i - f_{i-1}}{\Delta x} + O(\Delta x^2)$$

$$f''_i = \frac{f_{i+1} - 2f_i + f_{i-1}}{\Delta x^2} + O(\Delta x^2)$$

Note it turns out that the two  $O(\Delta x^2)$  errors cancel to see this use

$$(1) - (2) ! \\ \text{solve for } f'' - 6$$

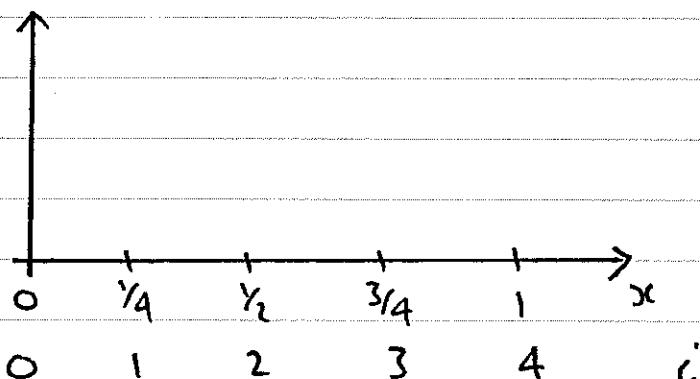
Now we can apply this to Laplace's eq in 1-D: ( $h = \Delta x$ )

$$\frac{\phi_{i+1} - 2\phi_i + \phi_{i-1}}{h^2} = 0$$

$$\phi_{i+1} - 2\phi_i + \phi_{i-1} = 0$$

let  $h = \frac{1}{4}$

$$\phi(x)$$



B.C.'s.

$$\phi_0 \approx \phi(0 \Delta x) = 0$$

$$\phi_4 \approx \phi(4 \Delta x) = \phi(1) = 10$$

$$\left. \begin{array}{l} \phi_0 - 2\phi_1 + \phi_2 = 0 \\ \phi_1 - 2\phi_2 + \phi_3 = 0 \\ \phi_2 - 2\phi_3 + \phi_4 = 0 \end{array} \right\} \begin{array}{l} \text{three equations} \\ \text{from applying C.D.} \\ \text{at 3 internal points.} \end{array}$$

3 unknowns  $\phi_1, \phi_2, \phi_3$ .

$$\begin{bmatrix} -2 & 1 & 0 \\ 1 & -2 & 1 \\ 0 & 1 & -2 \end{bmatrix} \begin{bmatrix} \phi_1 \\ \phi_2 \\ \phi_3 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ -10 \end{bmatrix} \leftarrow \text{From B.C.}$$

by Gaussian elimination:

$$\phi_1 = \frac{5}{2} \quad \phi_2 = 5 \quad \phi_3 = \frac{15}{2}$$

recall exact solution:  $\phi(x) = 10x$

$$\phi\left(\frac{1}{4}\right) = 10\left(\frac{1}{4}\right) = \frac{5}{2} = \phi_1$$

$$\phi\left(\frac{1}{2}\right) = 10\left(\frac{1}{2}\right) = 5 = \phi_2$$

$$\phi\left(\frac{3}{4}\right) = 10\left(\frac{3}{4}\right) = \frac{15}{2} = \phi_3$$

∴ the second order approximation gives the exact solution on the grid points. (as it should since exact sol<sup>m</sup> is linear).

Try for more grid points!

Now what have we solved for? " $\phi$ "

This is like a nodal voltage.

$$\bar{E} = -\frac{d\phi}{dx} \hat{a}_x = -10 \hat{a}_x$$

$$\text{if } \phi(0) = 10 \quad \phi(1) = 20 \Rightarrow \phi(x) = 10x + 10$$

$$\bar{E} = -10 \hat{a}_x \text{ still!}$$

{  $\bar{E}$  Field or voltage differences is what we can measure physically not  $\phi$ !

## Other types of Boundary Conditions

Say we were given  $E$  at one boundary.

$$-\frac{d\phi}{dx} = E = 10 \quad @ \quad x = 1$$

$$\cdot | \cdot - \frac{d\phi}{dx} \approx \frac{\phi_{N-1} - \phi_{N+1}}{2\Delta x} = 10$$

$\uparrow$  boundary.

$\therefore$  at the point  $N$ :

$$\frac{d^2\phi}{dx^2} \approx \frac{\phi_{N-1} + \phi_{N+1} - 2\phi_N}{\Delta x^2} = 0$$

we don't want to introduce new node to solve for " $\phi_{N+1}$ "

$$\therefore \text{use } \phi_{N+1} = \phi_{N-1} - 20\Delta x$$

$$\therefore \phi_{N-1} + \phi_{N-1} - 20\Delta x - 2\phi_N = 0$$

$$2\phi_{N-1} - 20\Delta x - 2\phi_N = 0$$



$$\phi_0 = 0$$

$$\Delta x = \frac{1}{4}$$

- @ i=1  $\phi_0 - 2\phi_1 + \phi_2 = 0$
- @ i=2  $\phi_1 - 2\phi_2 + \phi_3 = 0$
- @ i=3  $\phi_2 - 2\phi_3 + \phi_4 = 0$
- @ i=4  $2\phi_3 - 5 - 2\phi_4 = 0$

$$\begin{bmatrix} -2 & 1 & 0 & 0 \\ 1 & -2 & 1 & 0 \\ 0 & 1 & -2 & 1 \\ 0 & 0 & 2 & -2 \end{bmatrix} \begin{bmatrix} \phi_1 \\ \phi_2 \\ \phi_3 \\ \phi_4 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ 0 \\ 5 \end{bmatrix}$$

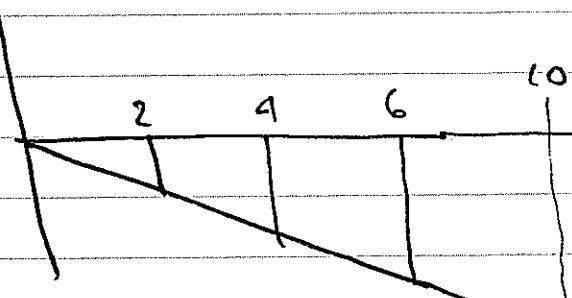
$$\begin{bmatrix} -2 & 1 & 0 & 0 & 0 \\ 0 & -\frac{3}{2} & 1 & 0 & 0 \\ 0 & 0 & -\frac{5}{3} & 1 & 0 \\ 0 & 0 & 2 & -2 & 5 \end{bmatrix} \rightarrow \begin{bmatrix} -2 & 1 & 0 & 0 & 0 \\ 0 & -\frac{3}{2} & 1 & 0 & 0 \\ 0 & 0 & -1 & \frac{3}{5} & 0 \\ 0 & 0 & 0 & -\frac{4}{5} & 5 \end{bmatrix}$$

$$\phi_4 = -10$$

$$\phi_3 = -6$$

$$\phi_2 = -4$$

$$\phi_1 = -2$$



check that this is exact.

$$\phi(x) = k_1(x) + k_2 \quad \phi_0 = 0 = k_2 \quad \left. \frac{d\phi}{dx} \right|_{x=1} = -10 = k_1 \quad \boxed{\phi(x) = -10x - 2}$$

# Application of Iterative Methods to Solve Matrix Equations

take same example as before:

$$\begin{bmatrix} -2 & 1 & 0 \\ 1 & -2 & 1 \\ 0 & 1 & -2 \end{bmatrix} \begin{bmatrix} \phi_1 \\ \phi_2 \\ \phi_3 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ -10 \end{bmatrix}$$

solve for diagonal components:

$$\begin{aligned} -2\phi_1 &= -\phi_2 \Rightarrow \phi_1^{(k)} = \frac{1}{2}\phi_2^{(k-1)} \\ -2\phi_2 &= -\phi_1 - \phi_3 \Rightarrow \phi_2^{(k)} = \frac{1}{2}(\phi_1^{(k-1)} + \phi_3^{(k-1)}) \\ -2\phi_3 &= -\phi_2 - 10 \Rightarrow \phi_3^{(k)} = \frac{1}{2}\phi_2^{(k-1)} + 5 \end{aligned} \quad \left. \right\}$$

K	$\phi_1^{(k)}$	$\phi_2^{(k)}$	$\phi_3^{(k)}$	$\  \underline{\delta} \  = \sum  \delta_i $
0	1.00	1.00	1.00	0
1	0.5	1.0	5.5	5
2	0.5	3.0	5.5	2
3	1.5	3.0	6.5	2
4	1.5	4.0	6.5	1
5	2.0	4.0	7.0	1
:	:	:	:	:
14	2.46875	4.96875	7.46875	0.03125

exact: 2.5 5.0 7.5

A). This is called Jacobi's method is usually slow to converge.

in matrix form :  $Ax = b$

then choose  $D$  to be the diagonal of  $A$

$$Dx^{(k+1)} = (D - A)x^{(k)} + b$$

in algebraic form:

$$[a_{ij}]_{n \times n} (x_i)_n = (b_i)_n$$

$$(x_i)^{(k+1)} = \frac{b_i}{a_{ii}} - \sum_{j=1}^{i-1} \frac{a_{ij}}{a_{ii}} x_j^{(k)} - \sum_{j=i+1}^n \frac{a_{ij}}{a_{ii}} x_j^{(k)}$$

### B) Gauss-Seidel Method (G-S)

succinctly: use the new  $(k+1)$  values of  $x$  as you find them!

Same example:

$$\begin{aligned} -2\phi_1 &= -\phi_2 \Rightarrow \phi_1^{(k+1)} = \frac{1}{2}\phi_2^{(k)} \\ -2\phi_2 &= -\phi_1 - \phi_3 \Rightarrow \phi_2^{(k+1)} = \frac{1}{2}(\phi_1^{(k+1)} + \phi_3^{(k)}) \\ -2\phi_3 &= -\phi_2 - 10 \Rightarrow \phi_3^{(k+1)} = \frac{1}{2}\phi_2^{(k+1)} + 5 \end{aligned}$$

K	$\phi_1^{(k)}$	$\phi_2^{(k)}$	$\phi_3^{(k)}$	$\ \bar{\delta}\  = \sum  \delta_i $
0	1.00	1.00	1.00	0
1	0.5	0.75	5.375	5.125
2	0.375	2.875	6.4375	3.3125
⋮	⋮	⋮	⋮	⋮
10	2.4916993	4.9916992	7.4958496	~0
Exact:	2.5	5.0	7.5	

in matrix form :  $A \mathbf{x} = \mathbf{b}$

choose  $E$  to be the lower triangular part of  $A$  (including diagonal)

$$E \mathbf{x}^{(k+1)} = (E - A) \mathbf{x}^{(k)} + \mathbf{b}$$

in algebraic form :  $[a_{ij}]_{n \times n} (x_i)_n = (b_i)_n$

$$(x_i)^{(k+1)} = \frac{b_i}{a_{ii}} - \sum_{j=1}^{i-1} \frac{a_{ij}}{a_{ii}} x_j^{(k+1)} - \sum_{j=i+1}^n \frac{a_{ij}}{a_{ii}} x_j^{(k)}$$

This method is still relatively slow to converge.

### c) Successive Overrelaxation (SOR)

Succinctly : take the difference between the G-S calculated point and the original value of the point and multiply by the over-relaxation factor  $1 < \omega < 2$  then add to the original value.

same example :

$$-2\phi_1 = -\phi_2 \Rightarrow \left\{ \begin{array}{l} \phi_1^{(k+1)} = \omega \left[ \frac{1}{2} \phi_2^{(k)} - \phi_1^{(k)} \right] + \phi_1^{(k)} \end{array} \right.$$

$$-2\phi_2 = -\phi_1 - \phi_3 \Rightarrow \left\{ \begin{array}{l} \phi_2^{(k+1)} = \omega \left[ \frac{1}{2} (\phi_1^{(k+1)} + \phi_3^{(k)}) - \phi_2^{(k)} \right] + \phi_2^{(k)} \end{array} \right.$$

$$-2\phi_3 = -\phi_2 - 10 \Rightarrow \left\{ \begin{array}{l} \phi_3^{(k+1)} = \omega \left[ \frac{1}{2} \phi_2^{(k+1)} + 5 - \phi_3^{(k)} \right] + \phi_3^{(k)} \end{array} \right.$$

in matrix form:  $Ax = b$

choose  $M$  to be the lower triangular part of  $A$  as in G-S but divide the diagonal entries by  $\omega$  (overrelaxation factor)

$$M \overset{(k+1)}{x} = (M-A) \overset{(k)}{x} + b$$

in algebraic form:  $[a_{ij}]_{n \times n} (x_i)_n = (b_i)_n$

$$\frac{a_{ii}}{\omega} (x_i)^{(k+1)} = b_i - \sum_{j=1}^{i-1} a_{ij} (x_j)^{(k+1)} - \sum_{j=i+1}^n a_{ij} (x_j)^{(k)} + (\frac{1}{\omega} - 1) a_{ii} (x_i)^{(k)}$$

$$(x_i)^{(k+1)} = \omega \left[ \frac{b_i}{a_{ii}} - \sum_{j=1}^{i-1} \frac{a_{ij}}{a_{ii}} (x_j)^{(k+1)} - \sum_{j=i+1}^n \frac{a_{ij}}{a_{ii}} (x_j)^{(k)} \right] + (1-\omega) (x_i)^{(k)}$$

$$(x_i)^{(k+1)} = \omega \left[ (x_i)_{G-S}^{(k+1)} - (x_i)^{(k)} \right] + (x_i)^{(k)}$$

for optimum " $\omega$ " this method is usually much faster to converge.

we find this optimum " $\omega$ " experimentally or check theory for specific P.D.E's.

## Stability of Iterative Methods

consider the linear equation

$$Ax = b \quad (1)$$

and an arbitrary simpler matrix  $M$  which is inserted into (1) as:

$$\begin{aligned} Mx + Ax &= Mx + b \\ Mx &= (M - A)x + b \end{aligned} \quad (2)$$

now use (2) to develop an iterative scheme starting from an initial guess  $x^0$ :

$$Mx^{k+1} = (M - A)x^k + b \quad (3)$$

now if the series  $x^0, x^1, x^2, \dots$  converges to say  $x^\infty$  then

$$Mx^\infty = (M - A)x^\infty + b$$

$$\Rightarrow Ax^\infty = b$$

$\therefore x = x^\infty$  is the solution  
of (1)

the convergence of the series to the "fixed point"  $x^\infty$  depends on the properties of  $A$  and the chosen  $M$ . (note:  $M$  has been arbitrary until now).

If we subtract ③ from ② we can define the error at step  $k$  as

$$Mx - Mx^{k+1} = (M-A)x - (M-A)x^k + b - b \\ M(x - x^{k+1}) = (M-A)(x - x^k)$$

$$e^k = x - x^k \quad \text{error at } k^{\text{th}} \text{ step}$$

$$\therefore \text{error equation : } M e^{k+1} = (M-A) e^k \quad ④$$

this is a one-step difference equation and the R.H.S of the original system (b) is not involved,

$$e^{k+1} = M^{-1}(M-A) e^k = B e^k \quad ⑤$$

$\therefore$  at each new iteration the current error vector is multiplied by the matrix  $B$

If  $e^0$  is the error in the initial guess then we can write:

$$e^k = B^k e_0 \quad ⑥$$

thus we see From ⑥ that

if  $B^k \rightarrow 0$  as  $k \rightarrow \infty$  then  
 $e^k \rightarrow 0$  as  $k \rightarrow \infty$   
and :  $x^k \rightarrow x$

∴ To test that our iterative scheme will converge we must test that

$B^k \rightarrow 0 \Rightarrow$  that our difference equation is stable.

A useful theorem for our purposes is the following:

defn: spectral radius,  $\rho(A)$ , of a matrix A is its largest eigenvalue, in magnitude:  
$$\rho(A) = \max_{\lambda \in \Lambda(A)} |\lambda|$$

where  $\Lambda(A)$  is the set of eigenvalues of the matrix A.

Theorem:

The powers  $B^k$  approach zero iff

$$\rho(B) < 1$$

The rate of convergence is governed by the size of  $\rho(B)$

$|\lambda_i| < 1 \Rightarrow$  that the eigenvalues must lie inside the unit circle in the complex plane.

Examples: check convergence of the following:

a)  $B = \begin{bmatrix} \gamma_2 & 10 \\ 0 & \gamma_2 \end{bmatrix}$   $(\lambda - \gamma_2)(\lambda - \gamma_2) = 0$   
 $\lambda_1 = \frac{1}{2}, \lambda_2 = \gamma_2 \Rightarrow$  convergence

b)  $B = \begin{bmatrix} \gamma_2 & \gamma_2 \\ \gamma_2 & \gamma_2 \end{bmatrix}$   $(\lambda - \gamma_2)(\lambda - \gamma_2) - \gamma_4 = 0$   
 $\lambda^2 - \lambda + \gamma_4 - \gamma_4 = 0$   
 $\lambda(\lambda - 1) = 0$   
 $\lambda = 0, \lambda = 1$   
 $\Rightarrow$  divergent

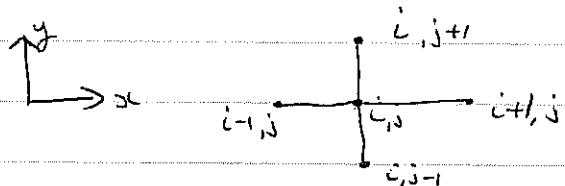
in fact

$$B^k = \begin{bmatrix} \gamma_2 & \gamma_2 \\ \gamma_2 & \gamma_2 \end{bmatrix}$$

## 2-D Laplace Equation

$$\nabla^2 \phi = \frac{\partial^2 \phi}{\partial x^2} + \frac{\partial^2 \phi}{\partial y^2} = 0$$

computational molecule :

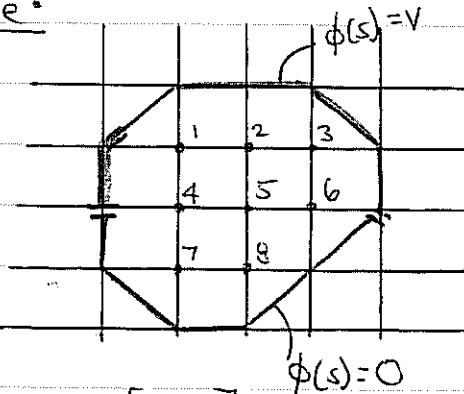


$$\frac{\partial^2 \phi}{\partial x^2} + \frac{\partial^2 \phi}{\partial y^2} \approx \frac{\phi_{i+1,j} - 2\phi_{i,j} + \phi_{i-1,j}}{(\Delta x)^2} + \frac{\phi_{i,j+1} - 2\phi_{i,j} + \phi_{i,j-1}}{(\Delta y)^2}$$

If  $\Delta x = \Delta y = h$  (i.e square grid)

$$\nabla^2 \phi \approx \frac{\phi_{i+1,j} + \phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1} - 4\phi_{i,j}}{h^2} = 0$$

example:



$$A = \begin{bmatrix} -4 & 1 & 0 & 1 & 0 & 0 & 0 & 0 & 0 \\ 1 & -4 & 1 & 0 & 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & -4 & 0 & 0 & 1 & 0 & 0 & 0 \\ 1 & 0 & 0 & -4 & 1 & 0 & 1 & 0 & 0 \\ 0 & 1 & 0 & 1 & -4 & 1 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 & 1 & -4 & 0 & 0 & 1 \\ 0 & 0 & 0 & 1 & 0 & 0 & -4 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 & 1 & -4 & 1 \end{bmatrix}$$

$$\underline{\phi} = \begin{bmatrix} \phi_1 \\ \phi_2 \\ \phi_3 \\ \phi_4 \\ \phi_5 \\ \phi_6 \\ \phi_7 \\ \phi_8 \end{bmatrix}$$

$$\underline{b} = \begin{bmatrix} -2V \\ -V \\ -2V \\ -V \\ 0 \\ -V \\ 0 \\ 0 \end{bmatrix}$$

$$A \underline{\phi} = \underline{b}$$

"A" will be a sparse matrix and so iteration techniques are appropriate.

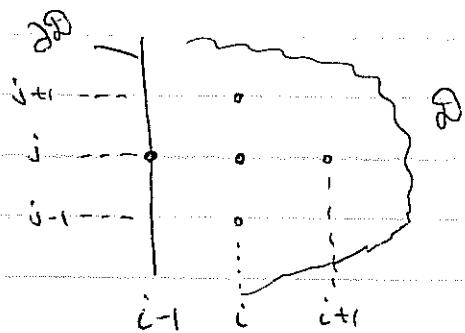
Notice that A will be diagonally dominant which is good for convergence.

In two space dimensions it is more convenient to label the nodes with their appropriate double indices  $\phi_{ij}$

and as long as all boundary points lie on the "integer" locations then we can use

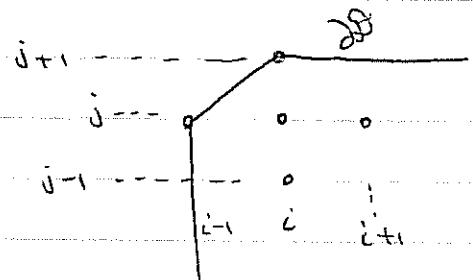
$$\phi_{i+1,j} + \phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1} - 4\phi_{ij} = 0$$

at all the interior points of the domain as well as the ones which have neighbouring boundaries in which the value of  $\phi$  is specified



↗ Dirichlet boundary conditions.

if on  $\partial\Omega$   $\phi = V$ , say, then

$$\phi_{i+1,j} + V + \phi_{i,j+1} + \phi_{i,j-1} - 4\phi_{ij} = 0$$


$$\phi_{i+1,j} + V + V + \phi_{i,j-1} - 4\phi_{ij} = 0$$

now if the normal derivative is specified at a boundary point:

$\partial\phi/\partial n$



$$\text{if on } \partial\Omega \quad \frac{\partial\phi}{\partial n} = 0 \quad \left( \frac{\partial\phi}{\partial x} = 0 \right)$$

(Neumann boundary condition)

then we can consider a fictitious grid point to the left of the boundary which does not belong in  $\Omega$ .

Laplace's equation at (i,j) gives.

$$\phi_{i+1,j} + \phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1} - 4\phi_{i,j} = 0 \quad (1)$$

Neumann's B.C. at (i,j) expressed as a central difference approximation gives:

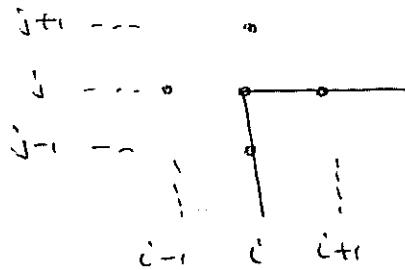
$$\frac{\phi_{i+1,j} - \phi_{i-1,j}}{2h} = 0 \quad O(h^2)$$

$$\text{or } \phi_{i+1,j} = \phi_{i-1,j} \quad (2)$$

using (2) in (1) gives:

$$2\phi_{i+1,j} + \phi_{i,j+1} + \phi_{i,j-1} - 4\phi_{i,j} = 0 \quad (3)$$

at a corner with two Neumann B.C's we get:



$$\phi_{i+1,j} = \phi_{i,j+1}, \quad \phi_{i-1,j} = \phi_{i,j-1}$$

$$2\phi_{i+1,j} + 2\phi_{i,j+1} - 4\phi_{i,j} = 0 \quad (4)$$

Now that we have discussed the special cases of the boundary conditions, we can formulate the SOR scheme:

at a regular point:

$$\phi_{ij} = \frac{\omega}{4} [\phi_{i+1,j} + \phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1}] + (1-\omega)\phi_{ij}$$

at an interior point bounded by a Dirichlet point:

$$\phi_{ij} = \frac{\omega}{4} [\phi_{i+1,j} + V + \phi_{i,j+1} + \phi_{i,j-1}] + (1-\omega)\phi_{ij}$$

etc.

Notice we are using the most upto date  $\phi_{ij}$ 's and over-writing  $\phi_{ij}$ .

in implementation only one array  $\phi(n,n)$  is required

## \* Odd-even updating.

For the Laplace equation notice that the updating of a point  $(i,j)$  involves only the points  $(i+1,j)$   $(i-1,j)$   $(i,i+1)$   $(i,i-1)$

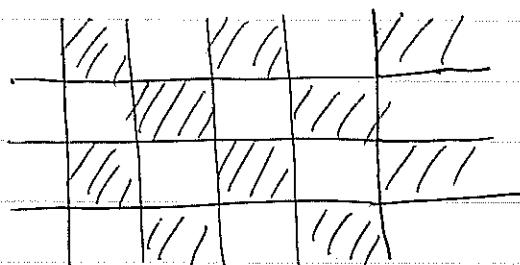
Thus we can update all points with

$$(i,j) \bmod 2 = 0$$

First and then scan through the points with

$$(i,j) \bmod 2 \neq 0$$

(checker board effect)



Stopping condition:

we define a displacement norm as,

$$\delta = \sum_{i=1}^N |\phi_i^{(m+1)} - \phi_i^{(m)}| = \|\Delta\phi\|$$

and a relative displacement norm as

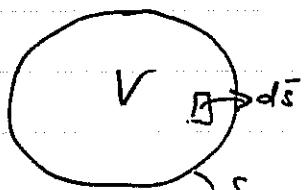
$$\epsilon = \frac{\delta}{\|\phi\|} = \frac{\delta}{\sum_{i=1}^n |\phi_i^{(m+1)}|}$$

we stop when the calculated error is below a specified  $\epsilon$  (say  $10^{-8}$ )

## Integration of Electric Flux Density

From Gauss' Law we know that

$$\Phi_0 = \iint_S \bar{D} \cdot d\bar{s} = \iiint_V q \, dv = Q_T$$

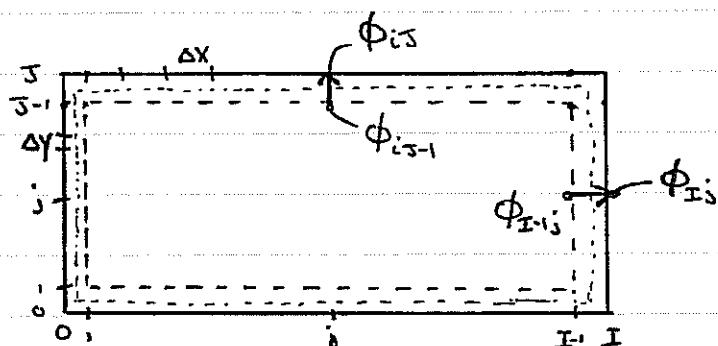


For the assigned problem:  $Q_T = 0$

$$\therefore \iint_S \bar{D} \cdot d\bar{s} = 0$$

and we have  $\epsilon = \epsilon_0$  inside the volume

$$\therefore \iint_S \bar{E} \cdot d\bar{s} = 0$$

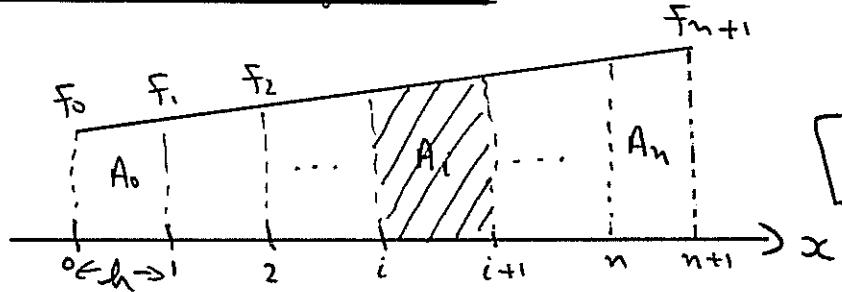


$$E_{y_{i,j-\frac{1}{2}}} \approx \frac{\phi_{ij} - \phi_{i,j-1}}{\Delta y} + O(\Delta y)$$

$$E_{x_{i-\frac{1}{2},j}} \approx \frac{\phi_{ij} - \phi_{i-1,j}}{\Delta x} + O(\Delta x^2)$$

$$\iint_S \bar{E} \cdot d\bar{s} = - \int_{j=0}^{j=j} E_{x_{0,j}} dy + \int_{i=0}^{i=i} E_{y_{i,j}} dx + \int_{j=0}^{j=j} E_{x_{i,j}} dy - \int_{i=0}^{i=i} E_{y_{i,0}} dx$$

## Numerical integration:



Trapezoidal Rule

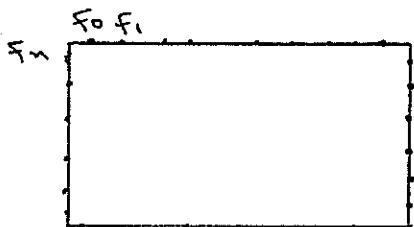
$$A_i \approx \frac{1}{2}(f_i + f_{i+1})h$$

$$I = \int f(x) dx = \sum_{i=0}^n \frac{1}{2}(f_i + f_{i+1})h + O(h)$$

$$I \approx h \left( \frac{1}{2}f_0 + f_1 + \dots + f_n + \frac{1}{2}f_{n+1} \right)$$

Around a closed loop:

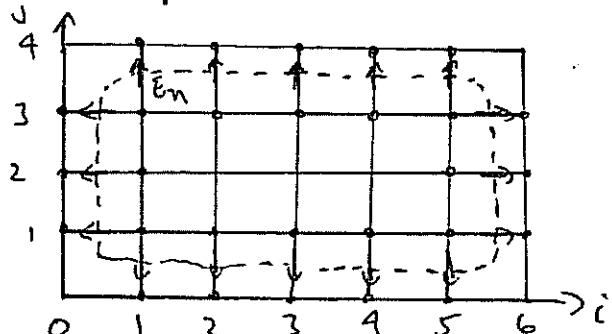
$$f_0 = f_{n+1}$$



$$\therefore I \approx (f_0 + f_1 + \dots + f_n)h$$

$$I \approx \left( \sum_{i=0}^n f_i \right)h$$

For our case the  $f$  are the normal component of electric field. ( $h = \Delta x = \Delta y$ )



$$\begin{aligned} \Phi_0 &\approx \sum_{i=1}^5 \left( \frac{\phi_{i,4} - \phi_{i,3}}{h} \right) h \\ &+ \sum_{i=1}^5 \left( \frac{\phi_{i,0} - \phi_{i,1}}{h} \right) h \\ &+ \sum_{j=1}^3 \left( \frac{\phi_{0,j} - \phi_{1,j}}{h} \right) h \\ &+ \sum_{j=1}^3 \left( \frac{\phi_{6,j} - \phi_{5,j}}{h} \right) h \end{aligned}$$

$$\bar{\Phi}_D \approx (\text{Sum outside potentials}) - (\text{Sum inside potentials} + \text{inside corners})$$

$$\approx \frac{\text{zero}}{=}$$

This should be checked in the program.

say  $i = 0(1)6$   $j = 0(1)6$  i.e. square.

$$\text{flux} = 0$$

For  $j = 1(1)5$

(outer boundary)

$$\text{flux} = \text{flux} + \phi(0, j)$$

Left

$$\text{flux} = \text{flux} + \phi(j, 0)$$

Bottom

$$\text{flux} = \text{flux} + \phi(6, j)$$

Right

$$\text{flux} = \text{flux} + \phi(j, 6)$$

Top

(inner boundary)

$$\text{flux} = \text{flux} - \phi(1, j)$$

Left

$$\text{flux} = \text{flux} - \phi(j, 1)$$

Bottom

$$\text{flux} = \text{flux} - \phi(5, j)$$

Right

$$\text{flux} = \text{flux} - \phi(j, 5)$$

Top.

end for