In a previous chapter  $\phi$  was approximted by a continuous function defined over the whole domain  $\Omega$  $\Omega \in \mathbb{R}^n$  with  $\Gamma \in \mathbb{R}^{n-1}$  its boundary  $\phi \simeq \hat{\phi} = \psi(x) + \sum a_m N_m(x)$  (completeness requirement)  $N_m$ ;  $m = 1,2,3,\cdots$  such that  $N_m|_{\Gamma} = 0$  $|\psi|_{\Gamma} = \phi|_{\Gamma}$  $N_{\scriptscriptstyle m}$  and  $\psi$  defined globally over  $\Omega$ Main drawback: treatment of arbitrary geometries boundary conditions to be imposed

Now 
$$\Omega = \bigcup_{e} \Omega^{e}$$
 ;  $\bigcap_{e} \Omega^{e} = \{ \}$  empty set

 $\phi$  is approximated piecewise over each subdomain (locally)

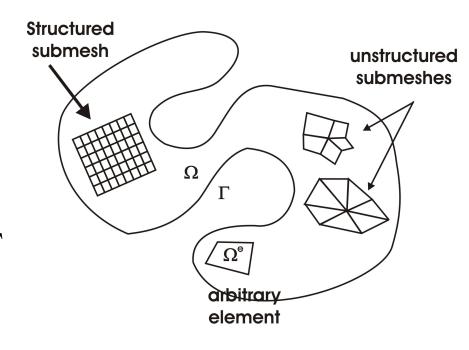
$$\int_{\Omega} W_l R_{\Omega} dx = \sum_{e=1}^{E} \int_{\Omega^e} W_l R_{\Omega} d\Omega$$

$$\int_{\Gamma} \overline{W_l} R_{\Gamma} dx = \sum_{e=1}^{E} \int_{\Gamma^e} \overline{W_l} R_{\Gamma} d\Gamma$$

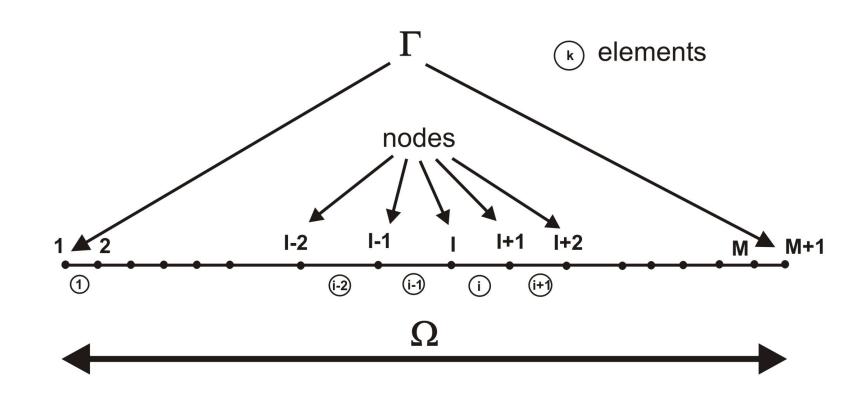
$$\sum_{e=1}^{E} \Omega^{e} = \Omega \qquad ; \qquad \sum_{e=1}^{E} \Gamma^{e} = \Gamma$$

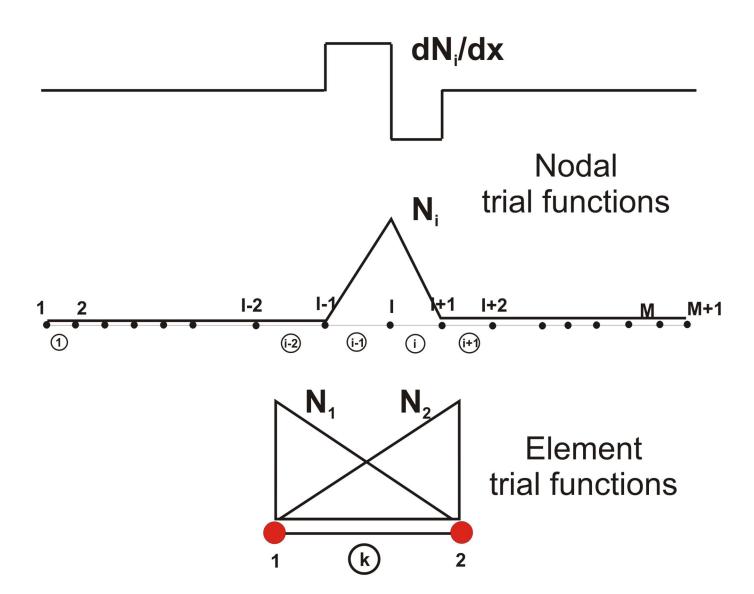
 $\Gamma^e$  boundary of each  $\Omega^e$ 

 $\Omega^e$  arbitrary, but may be simple enough



However piecewise approximation of trial function produces a loss of regularity for the globally assembled functions. In 1D the elements are intervals





## Some typical locally defined narrow-base shape functions

Approximate a continuous function  $\phi$  defined over  $\Omega = [0, L_x]$  discretized by a set of points  $[x_l; l = 1, 2, ..., M_n]$  with

$$x_1 = 0; x_{M_n} = L_x$$

$$\Omega = \bigcup_{e} \Omega^{e} = \bigcup_{e=1}^{M_n-1} [x_{e-1}, x_e]$$

$$\phi \cong \hat{\phi} = \psi + \sum_{m=1}^{M_n-1} \phi_m N_m(x)$$
 with  $\hat{\phi}$  piecewise constant

### Approximating a given function in 1D

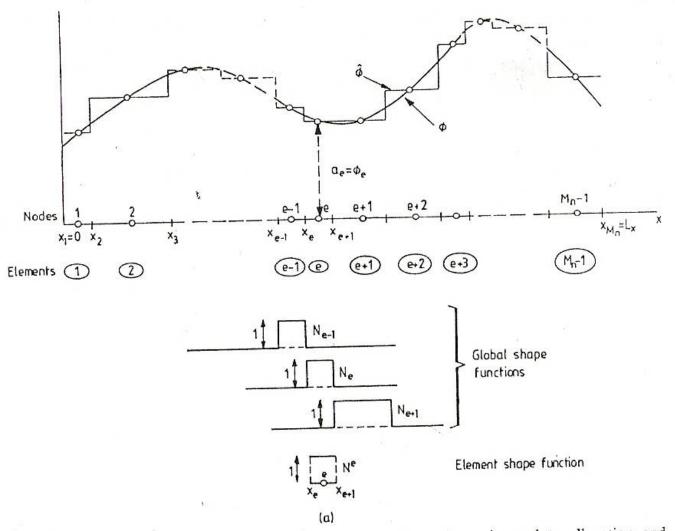
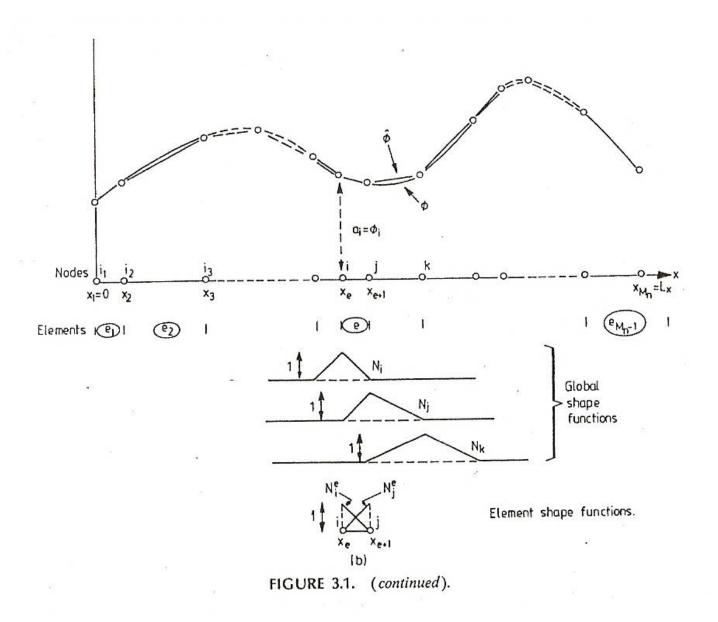


FIGURE 3.1. Approximating a given function in one dimension using point collocation and (a) piecewise constant elements and (b) piecewise linear elements.

### Approximating a given function in 1D



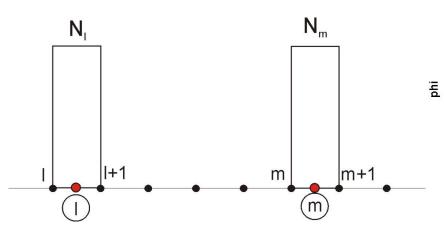
## Point collocation & piecewise constant approximation of 1D functions (see Ej\_3\_1 routine)

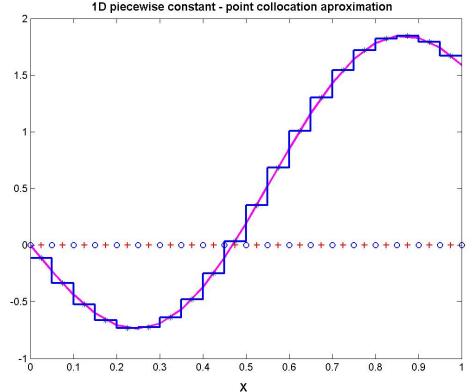
Using Weighted Residual Method (WRM) with  $\psi = 0$ 

$$\int_{\Omega} W_l R_{\Omega} d\Omega = \int_{\Omega} \delta(x - x_l) \left( \phi - \hat{\phi} \right) d\Omega = \int_{\Omega} \delta(x - x_l) \left( \phi - \sum_{m=1}^{M_n-1} \phi_m N_m(x) \right) d\Omega = 0$$

 $K_{lm} = \int_{\Omega} \delta(x - x_l) N_m(x) d\Omega = N_m(x_l) = \delta_{lm}$  (by orthonormality of trial functions)

$$f_l = \int \delta(x - x_l) \, \phi(x) \, d\Omega = \phi(x_l)$$





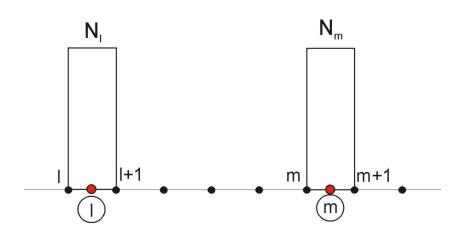
### Galerkin & piecewise constant approximation of 1D functions (see Ej\_3\_1\_2 routine)

Using Weighted Residual Method (WRM)

$$\int_{\Omega} W_l R_{\Omega} d\Omega = \int_{\Omega} W_l \left( \phi - \hat{\phi} \right) d\Omega = \int_{\Omega} N_l \left( \phi - \sum_{m=1}^{M_n-1} \phi_m N_m(x) \right) d\Omega = 0$$

$$K_{lm} = \int_{\Omega} N_l N_m d\Omega = \delta_{lm} |\Omega^m|$$
 (by orthogonality of trial functions)

$$f_l = \int_{\Omega} N_l \, \phi(x) \, d\Omega = \int_{\Omega^l} \phi(x) \, d\Omega$$



#### Point collocation

a = [-0.1159 -0.3365 -0.5244 -0.6608 -0.7308 -0.7249 -0.6396 -0.4776 -0.2480 0.0351 0.3531 0.6847 1.0077 1.3002 1.5431 1.7211 1.8239 1.8474 1.7936 1.6709]

#### **Galerkin solution**

a = [-0.1154 -0.3351 -0.5223 -0.6580 -0.7276 -0.7215 -0.6363 -0.4748 -0.2458 0.0365 0.3536 0.6844 1.0064 1.2982 1.5404 1.7179 1.8205 1.8441 1.7907 1.6686 l

### Point collocation & piecewise linear approximation of 1D functions (see Ej\_3\_1\_3 routine)

Using Weighted Residual Method (WRM) with  $\psi=0$ 

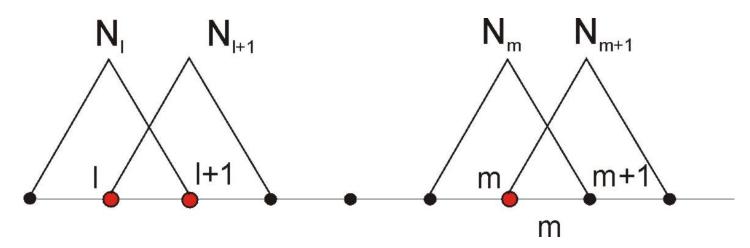
$$\int_{\Omega} W_l R_{\Omega} d\Omega = \int_{\Omega} \delta(x - x_l) \left( \phi - \hat{\phi} \right) d\Omega = \int_{\Omega} \delta(x - x_l) \left( \phi - \sum_{m=1}^{M_n} \phi_m N_m(x) \right) d\Omega = 0$$

the unknowns are now nodal values instead of element - wise values

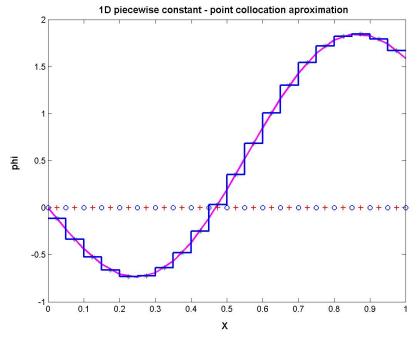
$$K_{lm} = \int_{\Omega} \delta(x - x_l) N_m(x) d\Omega = N_m(x_l) = \delta_{lm}$$
 (by orthonormality of trial functions)

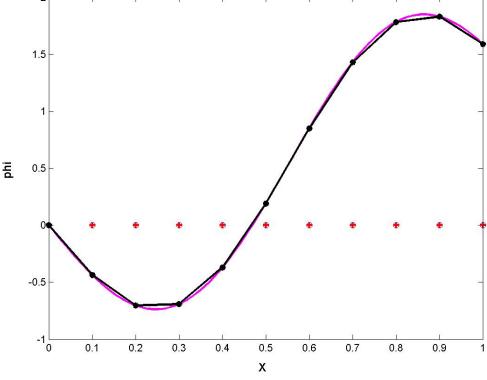
$$f_l = \int_{\Omega} \delta(x - x_l) \, \phi(x) \, d\Omega = \phi(x_l)$$

Ka = f is the same as using point collocation with piecewise constant!



### piecewise constant vs linear





1D piecewise linear - point collocation aproximation

## Galerkin & piecewise linear approximation of 1D functions

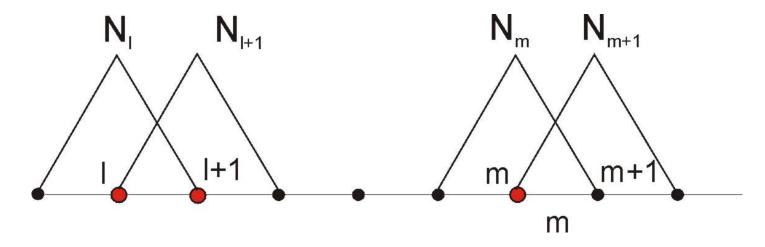
Using Weighted Residual Method (WRM)

$$\int_{\Omega} W_l R_{\Omega} d\Omega = \int_{\Omega} W_l \left( \phi - \hat{\phi} \right) d\Omega = \int_{\Omega} N_l \left( \phi - \sum_{m=1}^{M_n} \phi_m N_m(x) \right) d\Omega = 0$$

$$K_{lm} = \int_{\Omega} N_l N_m d\Omega \quad \neq 0 \quad \text{si} \quad |l-m| \leq 1 \text{ (for linear (*))}$$

$$f_l = \int_{\Omega} N_l \phi(x) d\Omega$$

(\*) assuming a natural numbering of nodes in 1D



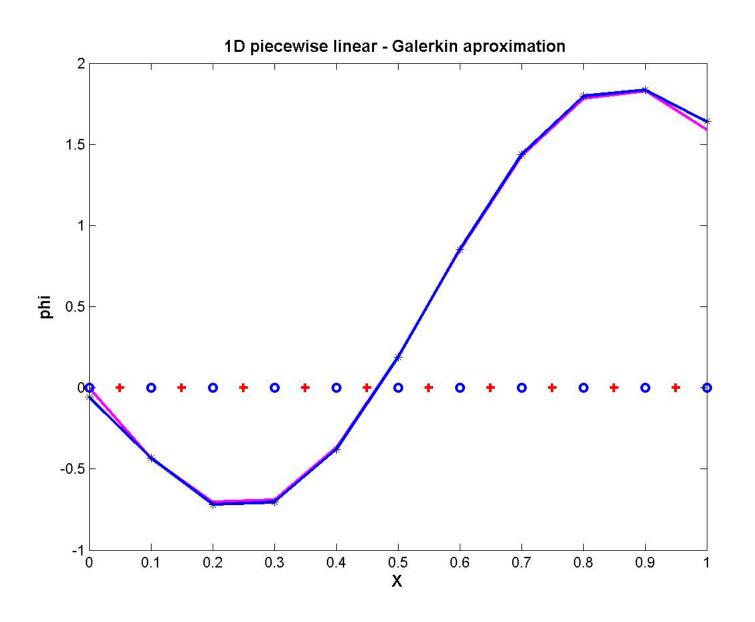
## Galerkin & piecewise linear integral kernel for Galerkin

```
K = zeros(Nx+1,Nx+1);
f = zeros(Nx+1,1);
                                                              piecewise linear Galerkin integrands
for I=2:Nx
  psi = (0:2)'/2;
  xx = psi*(x(l)-x(l-1))+x(l-1);
  Na = shape_Ej_3_1_4(xx,x(I-1),x(I));
  Nb = shape_{Ej_3_1_4(xx,x(l),x(l-1))};
  K(I,I-1) = K(I,I-1) + trapz(xx,Nb.*Na);
  K(I,I) = K(I,I) + trapz(xx,Nb.*Nb);
  xx = psi*(x(l+1)-x(l))+x(l);
  Na = shape_Ej_3_1_4(xx,x(l+1),x(l));
                                                 0.2
  Nb = shape_Ej_3_1_4(xx,x(l),x(l+1));
  K(I,I) = K(I,I) + trapz(xx,Na.*Na);
  K(I,I+1) = K(I,I+1) + trapz(xx,Na.*Nb);
                                                  0.25
                                                                0.35
                                                                              0.45
                                                                                            0.55
                                                                        Х
  psi = (0:20)'/20;
  xx = psi*(x(l)-x(l-1))+x(l-1);
  f(l,1) = f(l,1) + trapz(xx,shape_Ej_3_1_4(xx,x(l-1),x(l)).*ffun_Ej_3_1_4(xx));
  xx = psi*(x(l+1)-x(l))+x(l);
  f(I,1) = f(I,1) + trapz(xx,shape_Ej_3_1_4(xx,x(I+1),x(I)).*ffun_Ej_3_1_4(xx));
end
```

## Galerkin & piecewise linear boundary terms

```
I=Nx+1;
psi = (0:2)'/2;
xx = psi*(x(l)-x(l-1))+x(l-1);
Na = shape_Ej_3_1_4(xx,x(l-1),x(l));
Nb = shape_Ej_3_1_4(xx,x(l),x(l-1));
K(I,I-1) = K(I,I-1) + trapz(xx,Nb.*Na);
K(I,I) = K(I,I) + trapz(xx,Nb.*Nb);
I=1;
xx = psi*(x(l+1)-x(l))+x(l);
Na = shape_Ej_3_1_4(xx,x(l+1),x(l));
Nb = shape_Ej_3_1_4(xx,x(l),x(l+1));
K(I,I) = K(I,I) + trapz(xx,Na.*Na);
K(I,I+1) = K(I,I+1) + trapz(xx,Na.*Nb);
I=Nx+1;
psi = (0:20)'/20;
xx = psi*(x(l)-x(l-1))+x(l-1);
f(I,1) = f(I,1) + trapz(xx,shape_Ej_3_1_4(xx,x(I-1),x(I)).*ffun_Ej_3_1_4(xx));
I=1;
xx = psi*(x(l+1)-x(l))+x(l);
f(I,1) = f(I,1) + trapz(xx,shape_Ej_3_1_4(xx,x(I+1),x(I)).*ffun_Ej_3_1_4(xx));
```

### piecewise linear Galerkin results



```
Galerkin & piecewise linear
                                                                                  icone = [ ...
                                                                                      1
                       Element-wise assembling
                                                                                      3
                                                                                          4
for k=1:numel
                                                                                          5
  psi = (0:10)'/10;
                                                                                          6
  node1 = icone(k,1);
  node2 = icone(k,2);
                                                                                          7
  xx = psi*(xnod(node2,1)-xnod(node1,1))+xnod(node1,1);
                                                                                          8
  Na = shape_Ej_3_1_4(xx,xnod(node2,1),xnod(node1,1));
                                                                                          9
  Nb = shape Ei 3 1 4(xx,xnod(node1,1),xnod(node2,1));
                                                                                      9
                                                                                          10
  Ke(k,1,1) = Ke(k,1,1) + trapz(xx,Na.*Na);
                                                                                     10
                                                                                          11
  Ke(k,1,2) = Ke(k,1,2) + trapz(xx,Na.*Nb);
                                                                                     11
                                                                                          12
  Ke(k,2,1) = Ke(k,2,1) + trapz(xx,Nb.*Na);
                                                                                     12
                                                                                          13
  Ke(k,2,2) = Ke(k,2,2) + trapz(xx,Nb.*Nb);
  fe(k,1) = fe(k,1) + trapz(xx,Na.*ffun_Ej_3_1_4(xx));
                                                                                     13
                                                                                          14
  fe(k,2) = fe(k,2) + trapz(xx,Nb.*ffun Ei 3 1 4(xx));
                                                                                     14
                                                                                          15
end
                                                                                     15
                                                                                          16
                                                                       node numbering
% gather Ke and fe in Kg and fg
                                                                                          17
                                                                                     16
Kg = zeros(Nx+1,Nx+1); fg = zeros(Nx+1,1);
                                                     0.3
                                                                                          18
                                                                                     17
for k=1:numel
                                                                                          19
                                                                                     18
  node1 = icone(k,1):
                                                     0.2
                                                                                          20
                                                                                     19
  node2 = icone(k,2);
  Kg(node1,node1)=Kg(node1,node1)+Ke(k,1,1);
                                                     0.1
                                                                                          21]
                                                                                     20
  Kg(node1,node2)=Kg(node1,node2)+Ke(k,1,2);
  Kg(node2,node1)=Kg(node2,node1)+Ke(k,2,1);
                                                                 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21
  Kg(node2,node2)=Kg(node2,node2)+Ke(k,2,2);
                                                     -0.1
  fg(node1,1) = fg(node1,1) + fe(k,1);
  fg(node2,1) = fg(node2,1) + fe(k,2);
                                                     -0.2
end
                                                     -0.3
% solver
a = Kg \setminus fg;
                                                          0.1
                                                              0.2
                                                                  0.3
                                                                       0.4
                                                                           0.5
                                                                               0.6
                                                                                   0.7
                                                                                       0.8
                                                                                           0.9
                                                                           Х
```

### Approximating a given function in 2D

Using piecewise constant on triangles (Fig 3.2 (a))

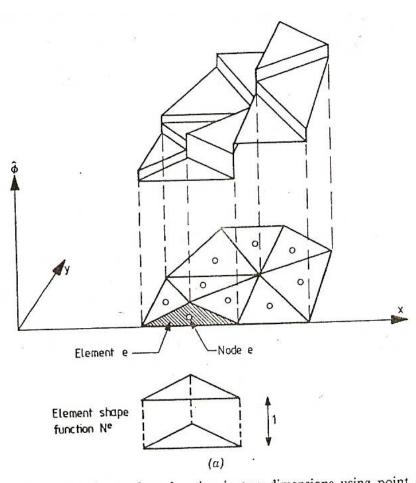


FIGURE 3.2. Approximating a given function in two dimensions using point collocation and (a) piecewise constant triangular elements and (b) piecewise linear triangular elements.

### Approximating a given function in 2D

Using piecewise linear on triangles (Fig 3.2 (b))

