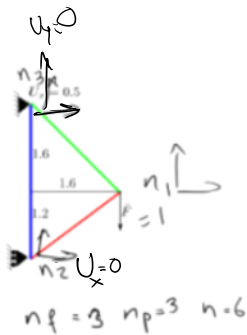


From last time



node	P	pos	v	F
1	0		0?	0
	0		0?	0
2	1		0.0	0?
	0		0?	-
3	1		0.5	0 p
	1		0.0	0?

$$U_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

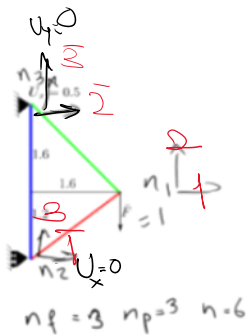
$$F_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

$$K_{ref} F_n = \begin{bmatrix} \\ \\ \end{bmatrix}$$

e	LEM	dof Map	dofs	fe

Steps 6 & 7

Step 6: dof positions; Step 7: Set F(F_f)



node	P	pos	v	F
1	0	1	0?	0
	0	2	0?	0
2	1	1	0.0	0?
	0	3	0?	-
3	1	1	0.5	0 p
	1	3	0.0	0?

$$U_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

$$F_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

$$K_{ref} F_n = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix}$$

e	LEM	dof Map	dofs	fe

free center

prescribed center



in C++, Python
dof positions start from 0
in Fortran, Matlab
= 1

posf = 0, posp = 0

for n = 1:nNodes

for dofi = 1: node(n).ndof num dof for node (n)

if node(n).ndof(dofi).p == true prescribed dof

posp = posp - 1;

node(n).ndof(dofi).pos = posp;

else free dof

```

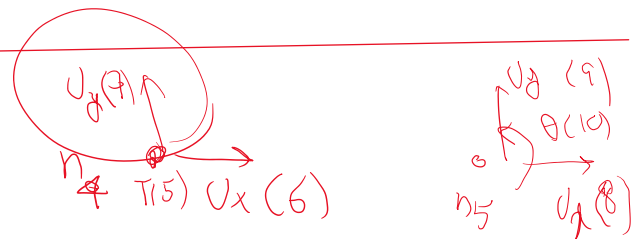
posf = 0, posp = 0
for n = 1:nNodes
  for dofi = 1: node(n).nndof num dof for node (n)
    if node(n).ndof(dofi).p == true prescribed dof
      posp = posp - 1;
      node(n).ndof(dofi).pos = posp;
    else free dof
      posf = posf + 1;
      node(n).ndof(dofi).pos = posf;
      F(posf) = node(n).ndof(dofi).f
    end
  end
end
end
end

```

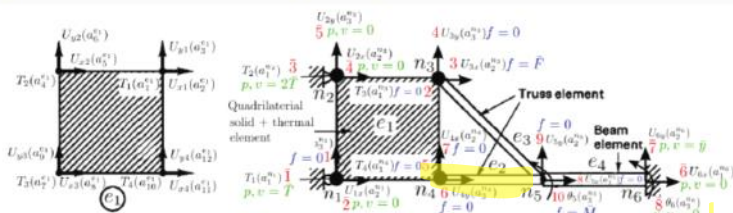
in Fortran, Matlab

~ 7

Step 8:
Setting dofMap of elements



Step 8: Element dof maps M_t^e



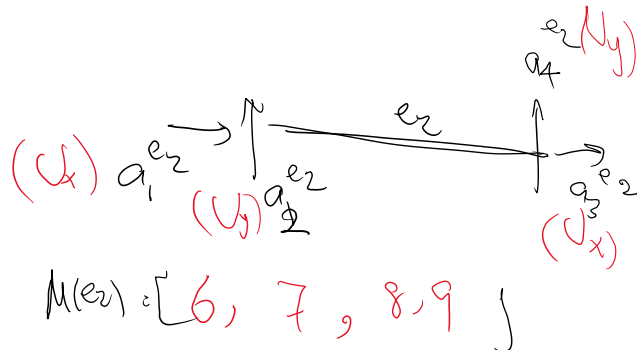
- As mentioned, M_t^e is a vector of size n_{dof}^e that maps element dofs to global positions.
- For element 1, dofs are ordered as (loop over nodes, then loop over dofs for the node):

$$a_1^e = [a_1^{e1} \quad a_2^{e1} \quad \dots \quad a_{12}^{e1}]$$

$$= [T_1 \quad U_{x1} \quad U_{y1} \mid T_2 \quad U_{x2} \quad U_{y2} \mid T_3 \quad U_{x3} \quad U_{y3} \mid T_4 \quad U_{x4} \quad U_{y4}]$$

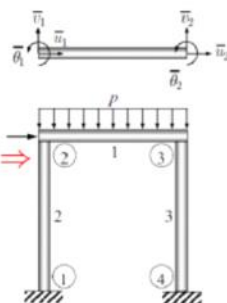
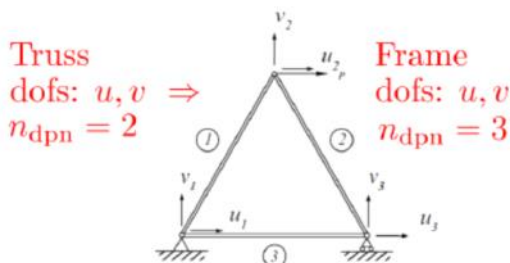
- We need to map these dofs to global dofs and have their position in M_t^e vector. For example, 1st dof of node 1 ($a_1^{e1} = T_1$) is mapped to first dof of n_3 which has position 2.
- 2nd dof of node 3 ($a_8^{e1} = U_{y2}$) is mapped to 2nd dof of n_1 which has position 2(-2).
- The map for element e_1 is:

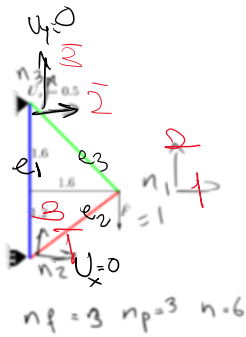
$$M^{e1}_t = [2 \quad 3 \quad 4 \quad \bar{3} \quad \bar{4} \quad \bar{5} \quad \bar{1} \quad \bar{2} \quad 1 \quad 5 \quad 6 \quad 7]$$



Not needed for the team project

Step 8: Element dof maps M_t^e : Simplified limited case





node	P	pos	v	f
1	0	1	0?	0
	0	2	0?	0
2	1	1	0.0	0?
	0	3	0?	-
3	1	2	0.5	0?
	1	3	0.0	0?

$$U_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix} \quad K_{ref} \quad F_n = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix}$$

$$F_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

e	LEM	dof Map	dofs	fe
1	[2,3]	[1 3 2 3]		
2	[2,1]	[1 3 1 2]		
3	[3,1]	[2 3 1 2]		

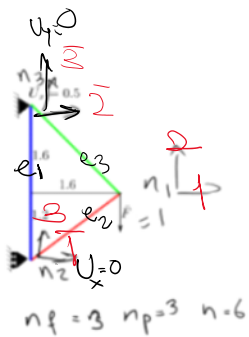
Be careful with indices stored in dofMap, in C++ I subtract 1 from prescribed dofs. For example, for e2 without the subtraction, I would have had

0 2 0 1, this becomes -1, 2 0 1

Similarly element 1 will have dofMap -1, 2, -2, -3

Step 9: Set element dofs a^e

needed for $f_D^e = K^e a^e$



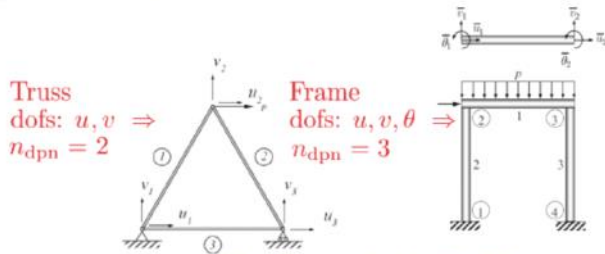
node	P	pos	v	f
1	0	1	0?	0
	0	2	0?	0
2	1	1	0.0	0?
	0	3	0?	-
3	1	2	0.5	0?
	1	3	0.0	0?

$$U_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix} \quad K_{ref} \quad F_n = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix}$$

$$F_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

e	LEM	dof Map	dofs	fe
1	[2,3]	[1 3 2 3]	[0.0 0 .5 0]	
2	[2,1]	[1 3 1 2]	[0.0 0 0 0]	
3	[3,1]	[2 3 1 2]	[.5 0 0 0]	

Step 9: Set element dofs a^e: Simplified limited case



- Similar to steps 1, 2, and 8, step 9 can be greatly simplified if we assume all nodes share exactly the same set of dofs.
- Noting n_{dofpn} (ndofpn) = Number of dof. per node, simplified merged steps 8 & 9 are:
dofs = zeros(ndof) element dofs (edof) resized to number of element dofs and zeroed
ecdf = 1 dof counter for element
for en = 1: neNodes number of element nodes
 gn = LEM(en) global node number for element node en
 for endof = 1: ndofpn This number is fixed now, e.g., 2 for 2D trusses
 if (node(gn).dof(endof).p == true) gndof = endof, we bypass some steps here
 dofs(ecdf) = node(gn).dof(endof).value; e dof val = corresponding global val
 end
 dofMap(ecdf) = node(gn).dof(endof).pos
 ecdf = ecdf + 1 increment counter
 end
end

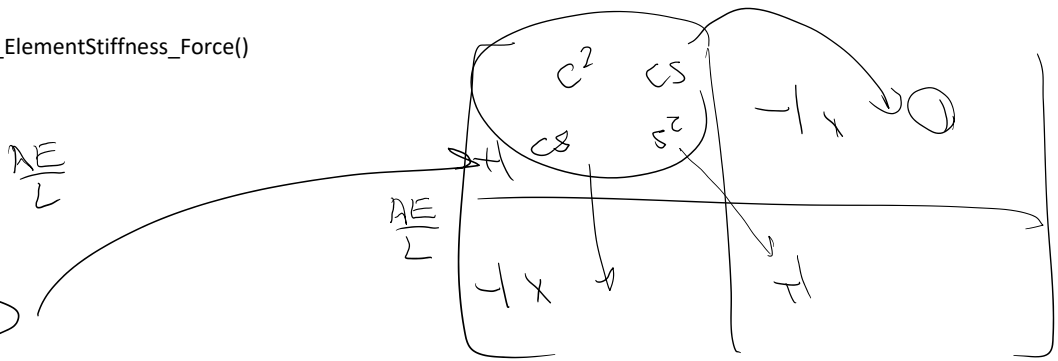
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Step 10: Compute element stiffness/force

Example: Truss element

```
class PhyElementTruss : public PhyElement
{
public:
    virtual void setInternalMaterialProperties(PhyMaterial*
    pMat);
    virtual void Calculate_ElementStiffness_Force();
};
```

```
void PhyElementTruss::Calculate_ElementStiffness_Force()
{
    // compute stiffness matrix:
    ke.resize(4, 4);
    double factor = A * E / L;
    for (int I = 0; I < 2; ++I)
        for (int J = 0; J < 2; ++J)
        {
            double f2 = factor;
            if ((I + J) % 2 != 0)
                f2 = -factor;
            ke(2 * I, 2 * J) = c * c * f2;
            ke(2 * I + 1, 2 * J) = ke(2 * I, 2 * J + 1) = c * s * f2;
            ke(2 * I + 1, 2 * J + 1) = s * s * f2;
        }
    cout << "ke\n" << ke << endl;
}
```



we need to have a few things for this element

@ A, E → Material properties

ⓑ L, c, s

ⓐ A, E :

no matID para
1 2 10 10000
2 3 0.01 100 0.3

- In the sample data input, material refers to truss element material parameters
A = 10, E = 10000 while material 2 is for 2D solid with
t(thickness) = 0.01, E = 100, ν(Poisson ratio) = 0.3.
- At element level a matID refers to global {mats}; e.g.,

For e3, matID = 1, ⇒ A = {mats}(1)(1) = 10, E = {mats}(1)(2) = 10000 (450)

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eType = 1 : bar element (451)
eType = 2 : beam element (452)
eType = 3 : truss element (453)
eType = 4 : frame element (454)
eType = 5 : 2D thermal element (455)
eType = 6 : 2D solid element (456)

eType = 1	Bar	mat{en} = [E A]	(454a)
eType = 2	Beam	mat{en} = [E I]	(454b)
eType = 3	Truss	mat{en} = [E A]	(454c)
eType = 4	Frame	mat{en} = [E A I]	(454d)
eType = 5	2D thermal	mat{en} = [κ t]	(454e)
eType = 6	2D solid	mat{en} = [E ν t]	(454f)

matID ≠ 1

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Materials
nMat 1
id numPara Paras
1 2 1 0 1 0
E, A
2 properties

id elementType matID neNodes eNodes
1 3 1 2 2 3
2 3 1 2 2 1
3 3 1 2 3 1

ⓑ L, C, S

How do we calculate these?

The diagram shows a truss element with nodes 1 and 2. Node 1 is at the origin (0,0) and node 2 is at (1.6, 1.2). The element length is L = 2.0. The cross-sectional area is A = 10 and the modulus of elasticity is E = 10000. The forces at node 1 are F1x = 1.6 and F1y = 1.2. The forces at node 2 are F2x = 1.6 and F2y = 1.2. The table below summarizes the node properties:

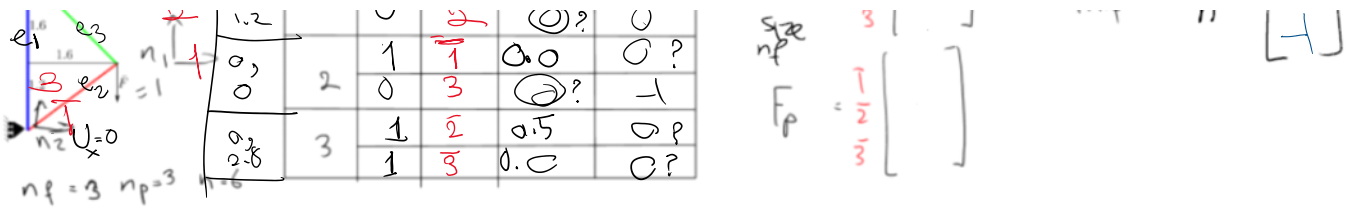
node	x	y	u	v	Fx	Fy
1	0	0	0	0	1.6	1.2
2	1.6	1.2	0	0	1.6	1.2

The stiffness matrix K is calculated as:

$$K = \frac{EA}{L} \begin{bmatrix} \cos^2 \theta & \cos \theta \sin \theta & -\cos \theta & 0 \\ \cos \theta \sin \theta & \sin^2 \theta & \sin \theta & 0 \\ -\cos \theta & \sin \theta & \cos \theta & -\sin \theta \\ 0 & 0 & -\sin \theta & -\cos \theta \end{bmatrix}$$

The force vector F is:

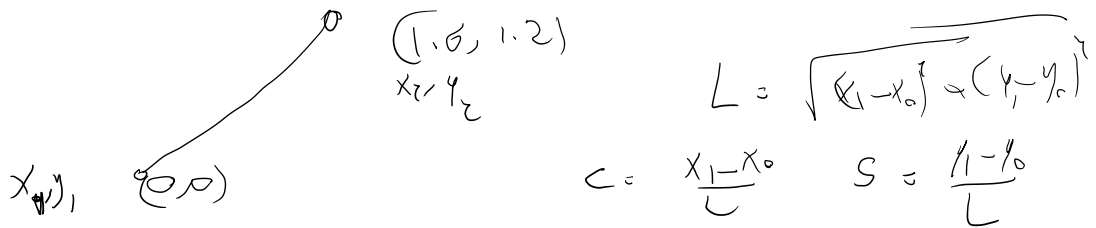
$$F = \begin{bmatrix} 0 \\ 0 \\ 1.6 \\ 1.2 \end{bmatrix}$$



e	LEM	def Map	def s	fe
1	[2,3]	[1 3 2 3]	[0.0 0 .5 0]	
2	[2,1]	[1 3 1 2]	[0.0 0 0 0]	
3	[3,1]	[2 3 1 2]	[.5 0 0 0]	

LEM

e2 node(1)_{e2} = 2 crd 0, 0
 node(2)_{e2} = 1 crd 1.6, 1.2



First, we add the extra data for each particular element type:

class PhyElementTruss : public PhyElement

```

{
public:
  virtual void setInternalMaterialProperties(PhyMaterial*
  pMat);
  virtual void Calculate_ElementStiffness_Force();
  virtual void SpecificOutput(ostream& out) const;

```

```

double L;
double A;
double E;
double c;
double s;

```

```

void PhyElementTruss::setInternalMaterialProperties(PhyMaterial* pMat)

```

```

{
  A = pMat->paras(mpb_A);
  E = pMat->paras(mpb_E);
}

```

```

void PhyElementTruss::setGeometry()

```

```

{
  VECTOR *crd0, *crd1;
  crd1 = &eNodePtrs[1]->coordinate;
  crd0 = &eNodePtrs[0]->coordinate;

  int sz = crd1->size();
}

```

```

if (sz != 2)
    THROW("implementation only for 2D truss");
double delX, delY;
delX = (*crd1)(0) - (*crd0)(0);
delY = (*crd1)(1) - (*crd0)(1);
L = sqrt(delX * delX + delY * delY);
c = delX / L;
s = delY / L;
}

```

Non-object oriented calculation of stiffness matrix

if (type == 1) v. bar

$$k = \frac{AE}{L} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix}$$

elseif (type == 2) v. brace

$$k = \begin{bmatrix} & & & \\ & & & \\ & & & \\ & & & \end{bmatrix}_{4 \times 4}$$

or
cases

Step 11: Assembly from local to global system

Truss example: Assembly of global system

e_1 	e_2 	e_3 	
$k^{e_1} = \frac{(1)(1)}{1.6} \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & -1 \\ 0 & 0 & 0 & 0 \\ 0 & -1 & 0 & 1 \end{bmatrix}$	$k^{e_2} = \frac{(1)(1)}{2} \begin{bmatrix} 0.64 & 0.48 & -0.64 & -0.48 \\ 0.48 & 0.36 & -0.48 & -0.36 \\ -0.64 & -0.48 & 0.64 & 0.48 \\ -0.48 & -0.36 & 0.48 & 0.36 \end{bmatrix}$	$k^{e_3} = \frac{(1)(1)}{1.6\sqrt{2}} \begin{bmatrix} 0.5 & -0.5 & -0.5 & 0.5 \\ -0.5 & 0.5 & 0.5 & -0.5 \\ -0.5 & 0.5 & 0.5 & -0.5 \\ 0.5 & -0.5 & -0.5 & 0.5 \end{bmatrix}$	
$f_D^{e_1} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.3571 & 0 & -0.3571 \\ 0 & 0 & 0 & 0 \\ 0 & -0.3571 & 0 & 0.3571 \end{bmatrix}$	$f_D^{e_2} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.24 & 0.18 & -0.24 & -0.18 \\ 0 & -0.32 & -0.24 & 0.32 & 0.24 \\ 0 & -0.24 & -0.18 & 0.24 & 0.18 \end{bmatrix}$	$f_D^{e_3} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.221 & -0.221 & -0.221 & 0.221 \\ 0 & -0.221 & 0.221 & 0.221 & -0.221 \\ 0 & -0.221 & 0.221 & 0.221 & -0.221 \\ 0 & 0.221 & -0.221 & -0.221 & 0.221 \end{bmatrix}$	
$f_e^{e_1} = f_r^{e_1} + f_N^{e_1} - f_D^{e_1} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.3571 & 0 & -0.3571 \\ 0 & 0 & 0 & 0 \\ 0 & -0.3571 & 0 & 0.3571 \end{bmatrix}$	$f_e^{e_2} = f_r^{e_2} + f_N^{e_2} - f_D^{e_2} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.24 & 0.18 & -0.24 & -0.18 \\ 0 & -0.32 & -0.24 & 0.32 & 0.24 \\ 0 & -0.24 & -0.18 & 0.24 & 0.18 \end{bmatrix}$	$f_e^{e_3} = f_r^{e_3} + f_N^{e_3} - f_D^{e_3} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.1105 & -0.1105 & -0.1105 & 0.1105 \\ 0 & -0.1105 & 0.1105 & 0.1105 & -0.1105 \\ 0 & -0.1105 & 0.1105 & 0.1105 & -0.1105 \\ 0 & 0.1105 & -0.1105 & -0.1105 & 0.1105 \end{bmatrix}$	
$K = \frac{1}{2} \begin{bmatrix} 0.32+0.221 & 0.24 & -0.221 & -0.24 \\ 0.24 & 0.18+0.221 & 0.18 & -0.18 \\ -0.24 & 0.18 & 0.3571+0.18 & 0 \end{bmatrix} = \begin{bmatrix} 0.5410 & 0.019 & -0.24 \\ 0.019 & 0.401 & -0.18 \\ -0.24 & -0.18 & 0.5371 \end{bmatrix}$			
$F = F_N + F_e = \begin{bmatrix} 0 \\ -1 \\ 0 \end{bmatrix} \Rightarrow U = K^{-1}F = \begin{bmatrix} U_1 \\ U_2 \\ U_3 \end{bmatrix} = \begin{bmatrix} -0.2123 \\ -3.2980 \\ -1.300 \end{bmatrix}$			

row column
 only k^e go to K (global)
 which one contribute to f_D^e

only P matters

ff
 fp
 pf
 pp

ff
 fp
 pf
 pp

Numbers encircled in the computation of essential BC force are displacements corresponding to free dofs. As mentioned before, in reality we do not consider them in computation of this force, but in hand calculation we just put zero for those values.

Function
 (rows) for $i = 1 : \text{ne dof}$ # dof
 $I = \text{dofMap}(i)$
 if $I < 0$ % row is prescribed
 continue;
 end

(columns) for $j = 1 : \text{ne dof}$ free rows
 $J = \text{dofMap}(j)$
 if $J > 0$ free columns
 $K(I, J) = K(I, J) + k_e(i, j) \cdot A_e(i)$ ff
 else ($J < 0$)
 $f_D(i) = f_D(i) + k_e(i, j) \cdot A_e(j)$ fp case
 end

$$\begin{aligned}
 & \text{end} \\
 & f_{ee}(i) = f_{ee}(i) - f_{eD}(i) \\
 & F(I) = F(I) + f_{ee}(i) \\
 & \text{end}
 \end{aligned}$$

Step 11: Assembly from local to global system

- K and F (global stiffness and force) are already sized and set to zero.
 - Element level (local) stiffness and force is calculated (Step 10).
 - Element local to global dof M_t^e is already set (Step 8).
 - Using dof map, we assemble local values to global values.
 - Clearly, only free dofs are added to stiffness matrix and force vector.
 - Element dof values (dofs: a^e is also set (Step 9).
 - $f_D^e = k^e a^e$ may not need to be formed and can be directly added to f_e^e .
- ```

for e = 1:ne loop over elements
 fee = feo element total force = element all forces except essential force
 for i = 1:nedof loop over rows of ke; nedof = element # dof
 I = dofMap(i) local to global dof map M_t^e
 if (I > 0) I corresponds to a free dof, we skip prescribed dofs
 for j = 1:nedof loop over columns of ke
 J = dofMap(j) global dof corresponding to j
 if (J > 0) now both I and J are free and can add ke(i,j) to global K
 K(I, J) = K(I, J) + ke(i, j)
 else J < 0, prescribed dof j; add contributions of $f_D^e = k^e a^e$ to f_e^e
 fee(i) = fee(i) - ke(i, j) * edofs(j) edofs: element dofs = a^e
 end
 end
 end
 F(I) = F(I) + fee(i) element's total force fee component i'th is computed → added to F(I)
 end
end
end
end

```

## Step 11: Assembly from local to global system

- **Rows:** In loop over edof (element dofs) we skip prescribed dofs  $i$ .
- **Columns:** free columns  $j$ : Stiffness is assembled  

$$K(I, J) = K(I, J) + ke(i, j).$$
- **Columns:** prescribed columns  $j$ :  $f_D^e$  is added to total element force  
 $f_e^e$ :  $fee(i) = fee(i) - ke(i, j) * edofs(j).$
- **Force:** Once  $fee$  is updated with  $fde$  (essential BD force), it is assembled to global  $F$ :  $F(I) = F(I) + fee(i).$
- **Both  $K$  and  $k$  are symmetric:**
  - 1  $K$  can be stored into a symmetric matrix  $\rightarrow$  almost half the storage.
  - 2  $ke$  symmetric: can be stored in symmetric matrix  $\rightarrow$  almost half the storage.
  - 3  $ke$  symmetric: half the loop computation: instead of looping ( $i: 1 \rightarrow nedof; j: 1 \rightarrow nedof$ ) we can loop ( $i: 1 \rightarrow nedof; j: i \rightarrow nedof$ ).  
How?

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## Step 12: Solve global (free) dof $a$ from $Ka = F$

- Two major computational costs during FEM solve are:
  - 1 **Assembly:** Refers to: all node, element, and dof set up; computation of local  $ke$  and  $fee$ ; assembly of those to global system. This step scales linearly versus  $n_e$  ( $ne$ )
  - 2 **Linear algebra solution:**  $Ka = F$ : We solve for unknown  $a$ . Although conceptually simple, this step is a major source of computational cost. It scales higher than linear versus  $n_e \Rightarrow$  As the problem size increases this term becomes more dominant.
- Solution of  $Ka = F$ :
  - WE DO NOT OBTAIN  $a$  from  $a = K^{-1}F$ : We do not invert  $K$ .
  - We only solve the problem for the specific RHS of  $F$ .
  - In Comparison  $K^{-1}$  corresponds to the solution of  $Ka = F$  for  $n_f$  RHS of  $F = e_i, i = 1, \dots, n_f$  where  $n_f$  is the number of rows (and columns) of  $K$ .
  - We employ methods such as LU factorization that computationally only solve the problem for the given RHS  $F$ .
  - We take advantage of the structure of stiffness matrix: symmetry, bandedness, sparsity in choosing the right solution technique.
  - order of free dofs affects band of the matrix  $\rightarrow$  various algorithms reorder free dofs such that the matrix band get smaller and the solution cost is optimized.
  - In your term projects you can simply employ simply compute

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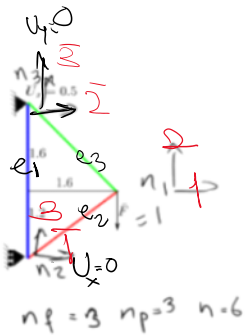
$$a = K^{-1}F \quad \swarrow \quad \searrow \quad K \setminus F$$

4.0  
2.0

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 67 68 69 70 71 72 73 74 75 76 77 78 79 80 81 82 83 84 85 86 87 88 89 90 91 92 93 94 95 96 97 98 99 100

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 67 68 69 70 71 72 73 74 75 76 77 78 79 80 81 82 83 84 85 86 87 88 89 90 91 92 93 94 95 96 97 98 99 100

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 67 68 69 70 71 72 73 74 75 76 77 78 79 80 81 82 83 84 85 86 87 88 89 90 91 92 93 94 95 96 97 98 99 100



| node | p | pos | v   | f  |
|------|---|-----|-----|----|
| 1    | 0 | 1   | 0?  | 0  |
|      | 0 | 2   | 0?  | 0  |
| 2    | 1 | 1   | 0.0 | 0? |
|      | 0 | 3   | 0?  | -1 |
| 3    | 1 | 2   | 0.5 | 0? |
|      | 1 | 3   | 0.0 | 0? |

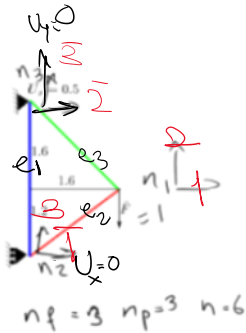
$$U_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix} \begin{bmatrix} -21.33 \\ -3.29 \\ -1.2 \end{bmatrix}$$

$$F_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

$$K_{ref} \quad F_n = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix}$$

| e | LEM   | dof Map   | dofs         | fe |
|---|-------|-----------|--------------|----|
| 1 | [2,3] | [1 3 2 3] | [0.0 0 .5 0] |    |
| 2 | [2,1] | [1 3 1 2] | [0.0 0 0 0]  |    |
| 3 | [3,1] | [2 3 1 2] | [.5 0 0 0]   |    |

### Step 13: Assign a to nodes and elements



| node | p | pos | v      | f  |
|------|---|-----|--------|----|
| 1    | 0 | 1   | -21.33 | 0  |
|      | 0 | 2   | -3.29  | 0  |
| 2    | 1 | 1   | 0.0    | 0? |
|      | 0 | 3   | -1.2   | -1 |
| 3    | 1 | 2   | 0.5    | 0? |
|      | 1 | 3   | 0.0    | 0? |

$$U_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix} \begin{bmatrix} -21.33 \\ -3.29 \\ -1.2 \end{bmatrix}$$

$$F_p = \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

$$K_{ref} \quad F_n = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix}$$

| e | LEM   | dof Map   | dofs                    | fe |
|---|-------|-----------|-------------------------|----|
| 1 | [2,3] | [1 3 2 3] | [0.0 -1.2 .5 0]         |    |
| 2 | [2,1] | [1 3 1 2] | [0.0 -1.2 -21.33 -3.29] |    |
| 3 | [3,1] | [2 3 1 2] | [.5 0 -21.33 -3.29]     |    |

for n = 1:nNodes

for dofi = 1: node(n).nndof num dof for node (n)

if node(n).ndof(dofi).p == false free dof

posn = node(n).ndof(dofi).pos position of dof in global free F

node(n).ndof(dofi).v = dofs(posn) set free dof val to corresponding val in global dofs (a)

node dofs

end

end

end

```
for e = 1:ne loop over elements
 for i = element(e).nedof loop over element dofs; nedof = # dof (n_{dof}^e)
 posn = element(e).dofMap(i) corresponding global position using dofMat (M_i^e)
 if (posn > 0) free dof
 element(e).edofs(i) = dofs(posn)
 set free element dof a^e to corresponding val in global dofs (a)
 end
 end
end
```

element  $e$

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