

much better than displacement & stress approaches

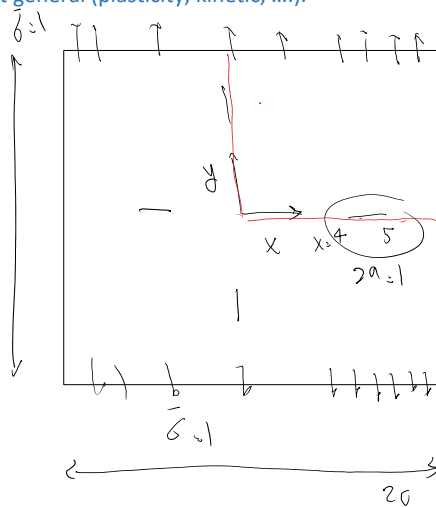
So EDI method is by far the most accurate, most robust, and most general (plasticity, kinetic, ...).

Analytical  $K_I$   
assuming infinite domain approx.

$$K_I = \delta \sqrt{\pi a} = 1 \sqrt{\pi \times 5} = 1.2533$$

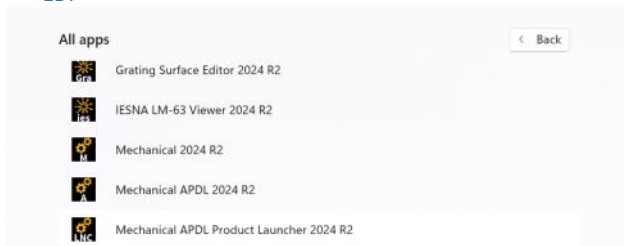
Eol  $v = 0$

$$G = J = \frac{K_I^2}{E} = 1.5708$$

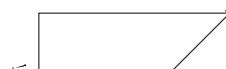
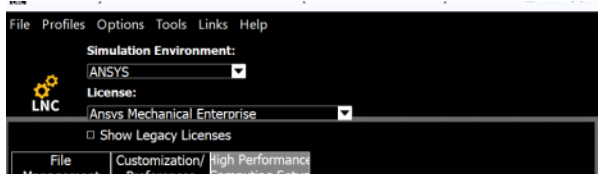


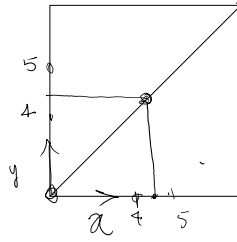
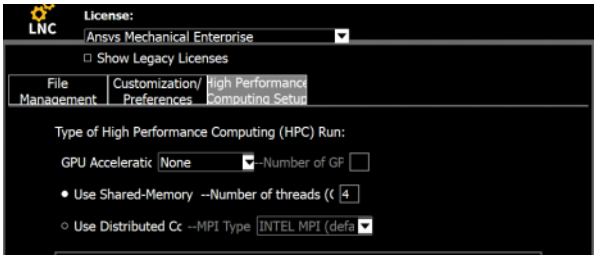
Computing K and J with Ansys:

- Displacement approach
- EDI

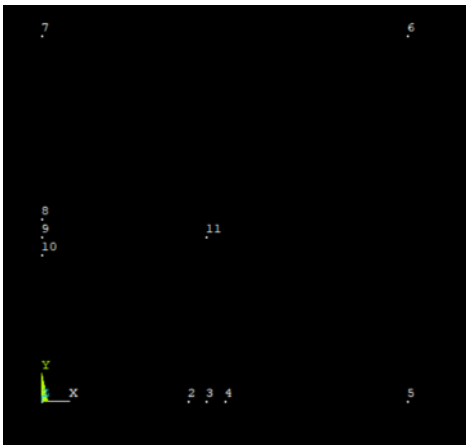
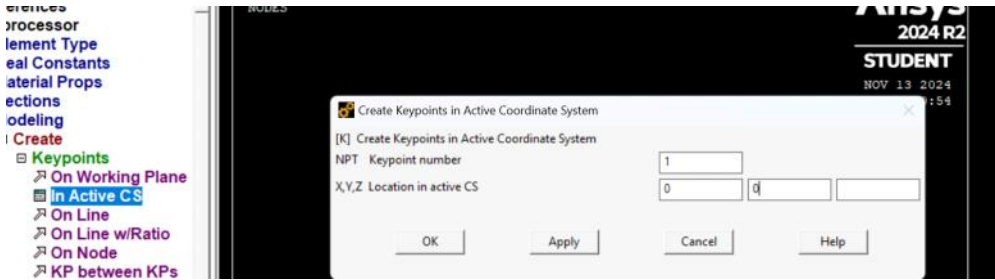


Change this to shared memory:

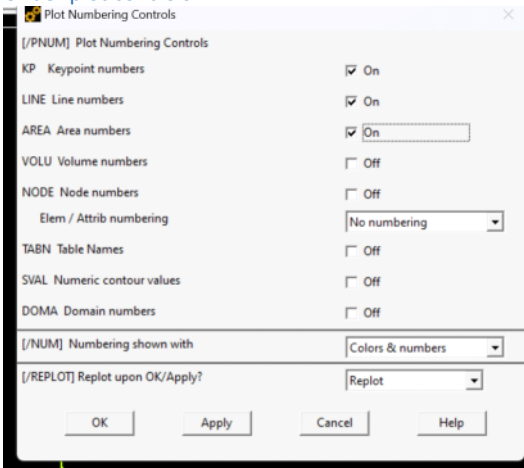




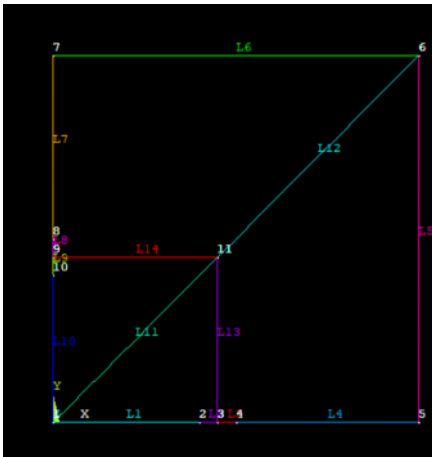
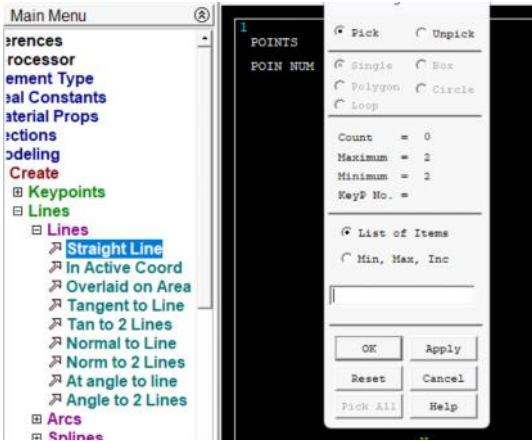
Creating the geometry



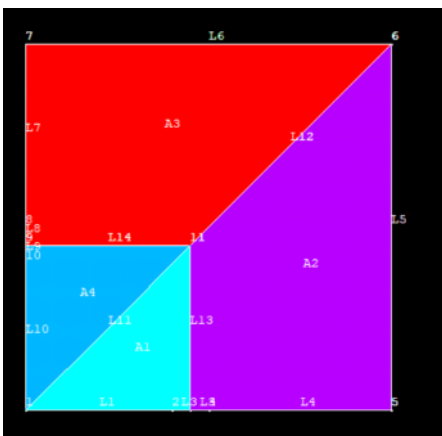
Under plot controls



Next we create lines



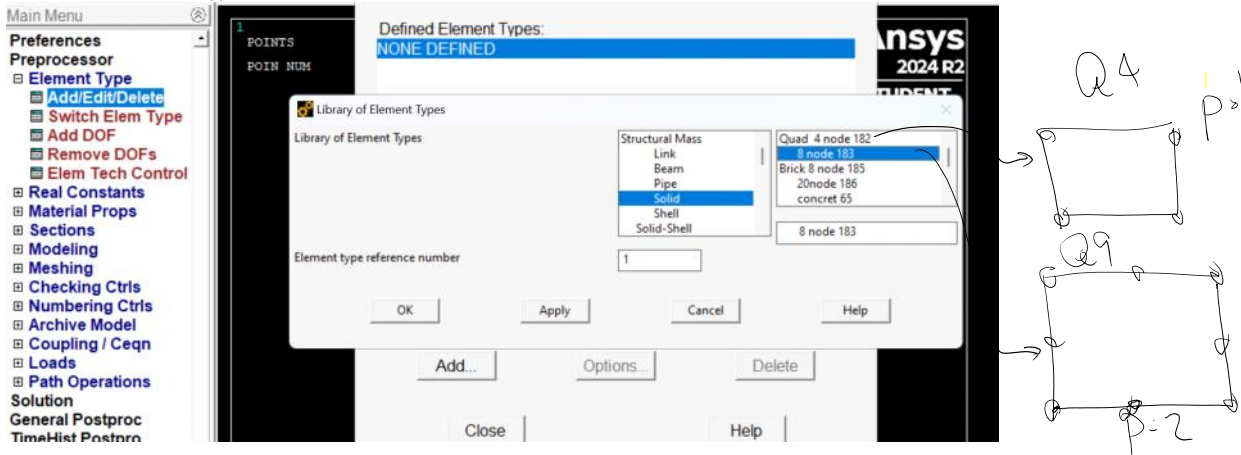
Create areas:



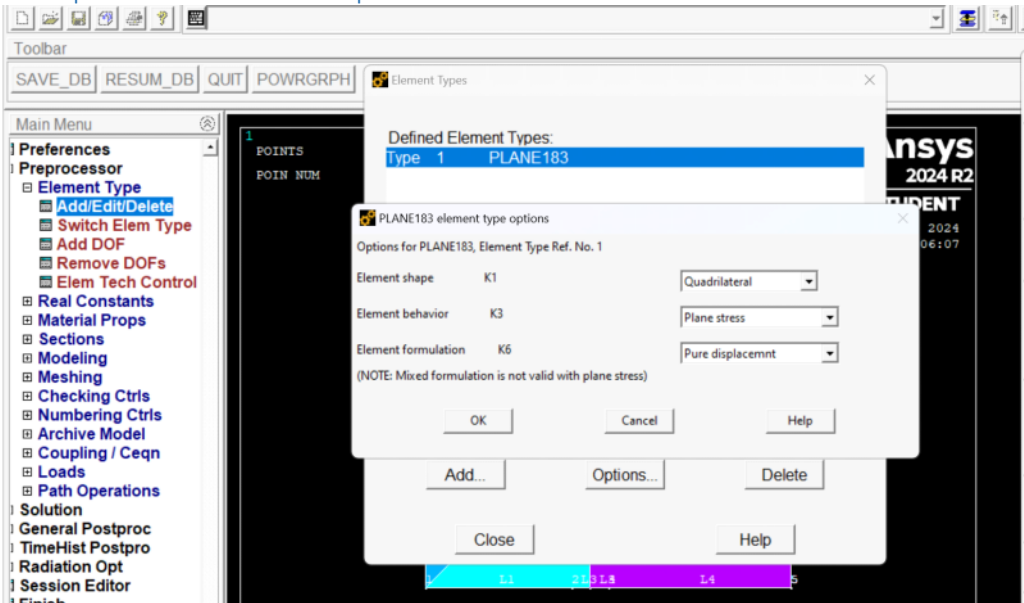
Create the element type



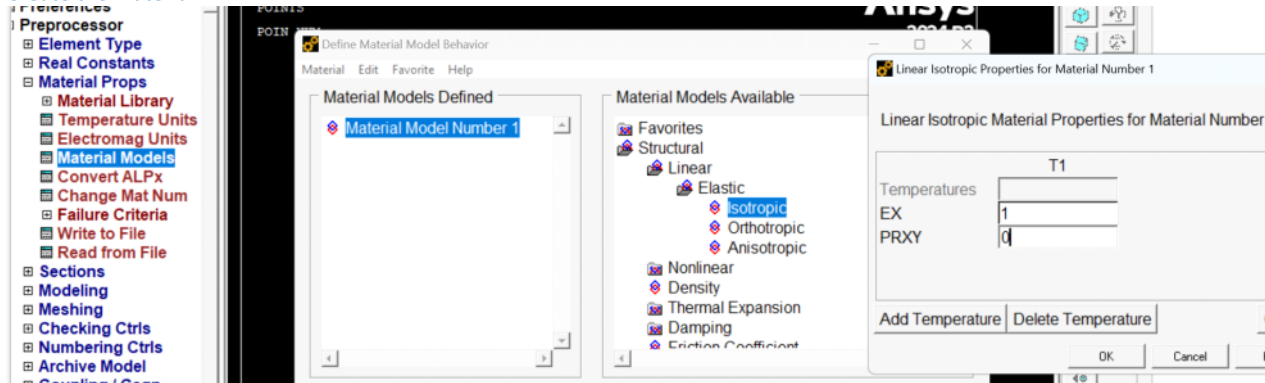
Create the element type



Choose plane stress or strain under options

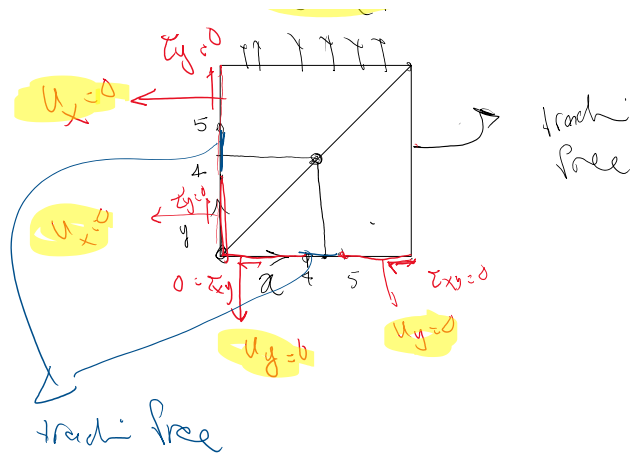


Create the material

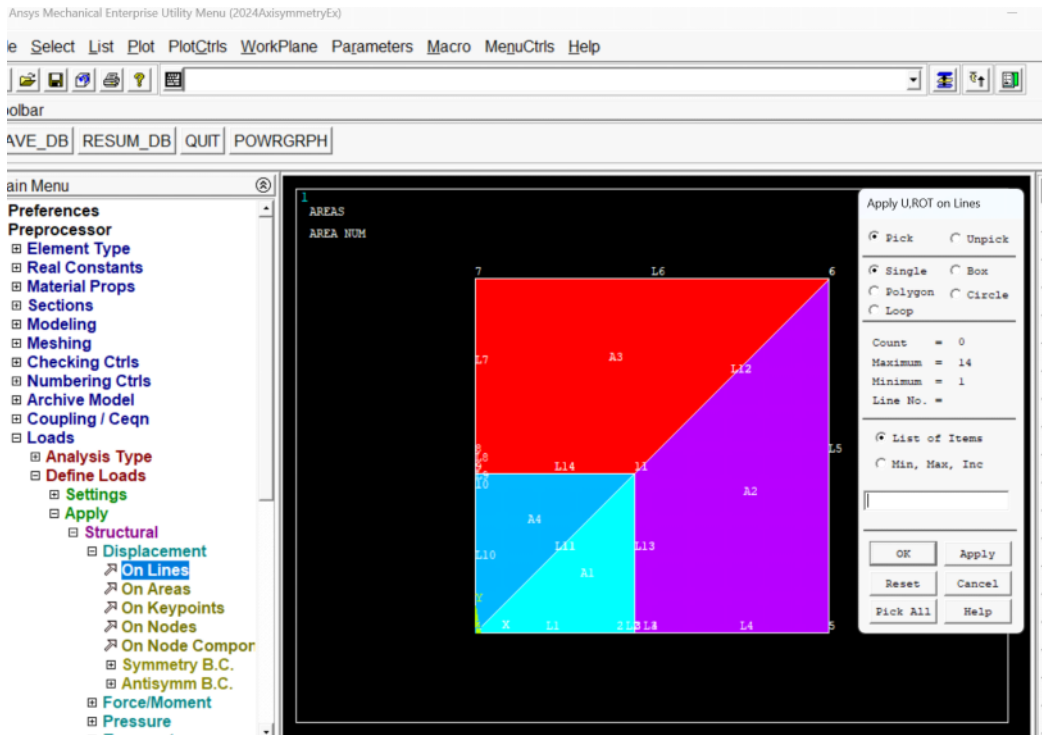


Apply the BCs on the geometry

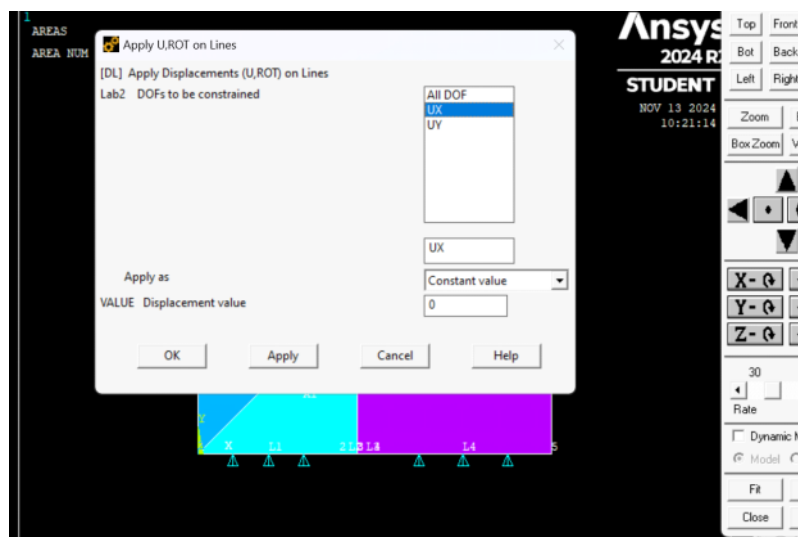
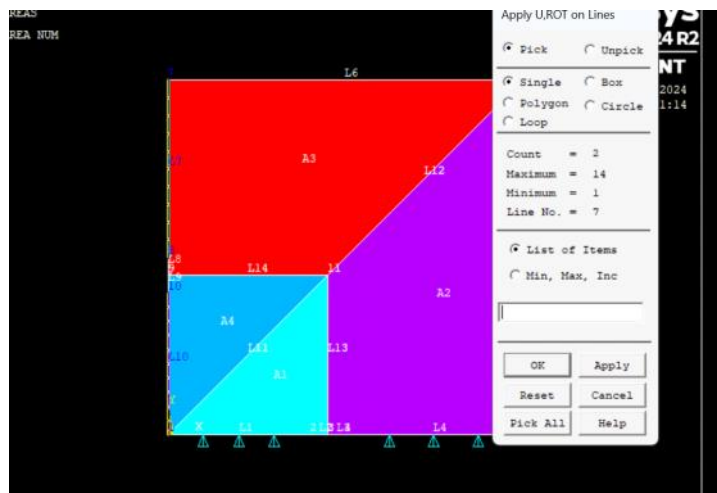
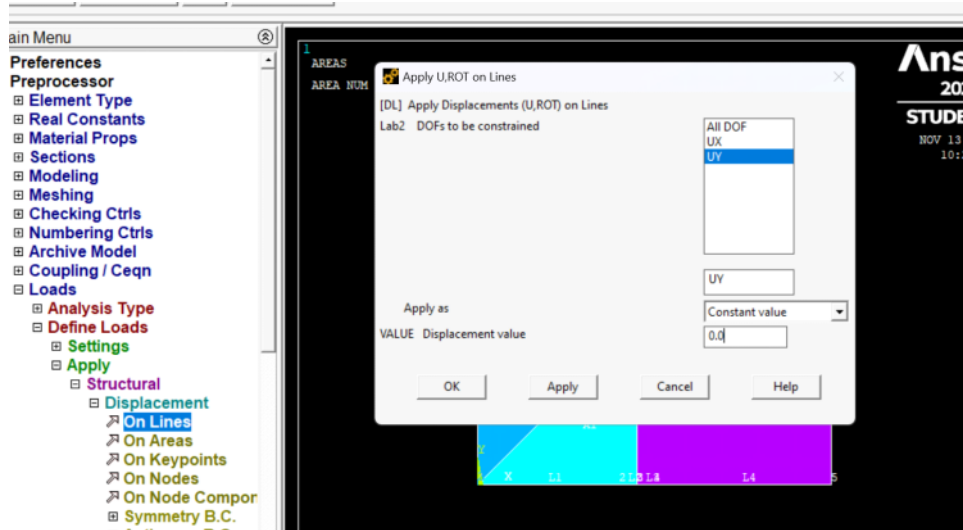




In FEM, the default is traction free. So, if for a direction traction is zero, we don't do anything

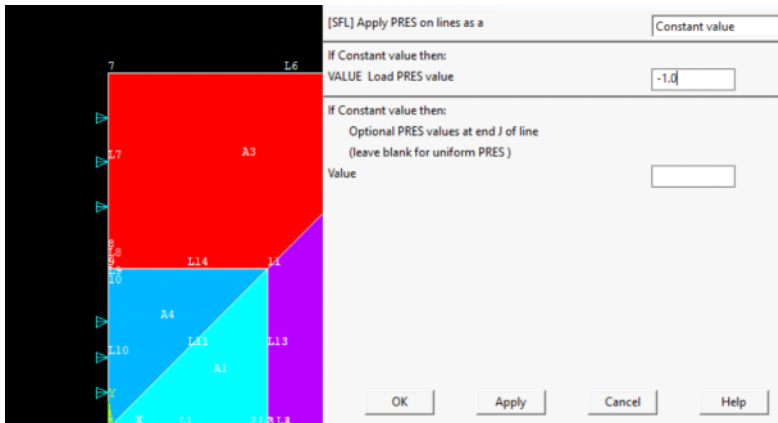


Apply  $u_y = 0$  on the left and right segments (bottom edge)

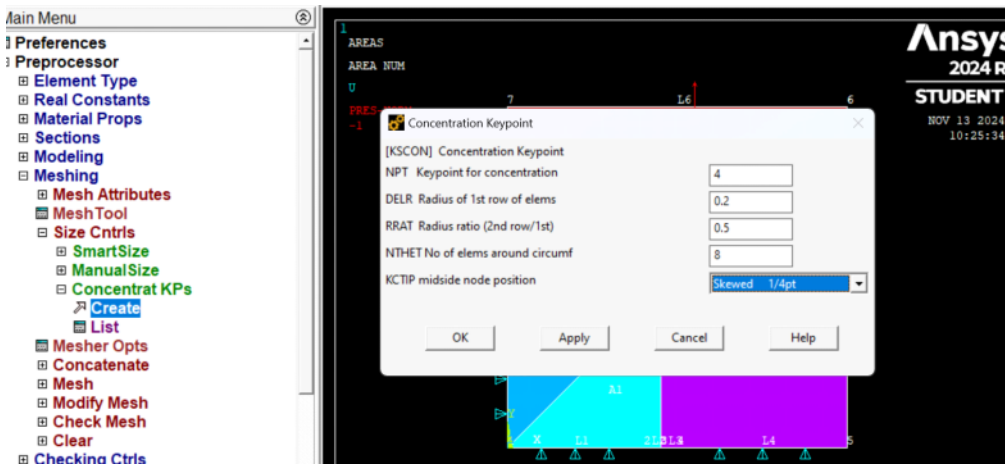
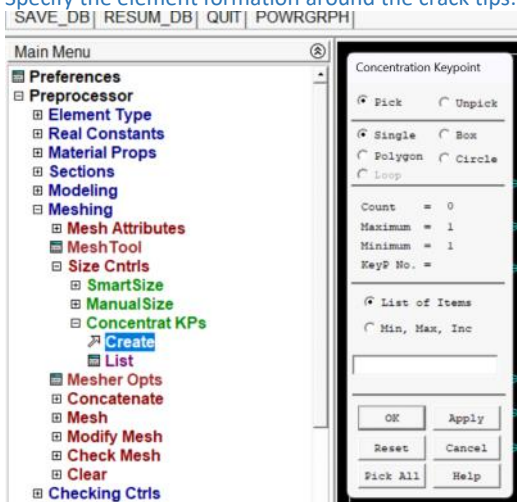


Pressure on the top line:

- Analysis Type
- Define Loads
- Settings
- Apply
  - Structural
  - Displacement
  - Force/Moment
  - Pressure
    - On Lines
    - On Areas

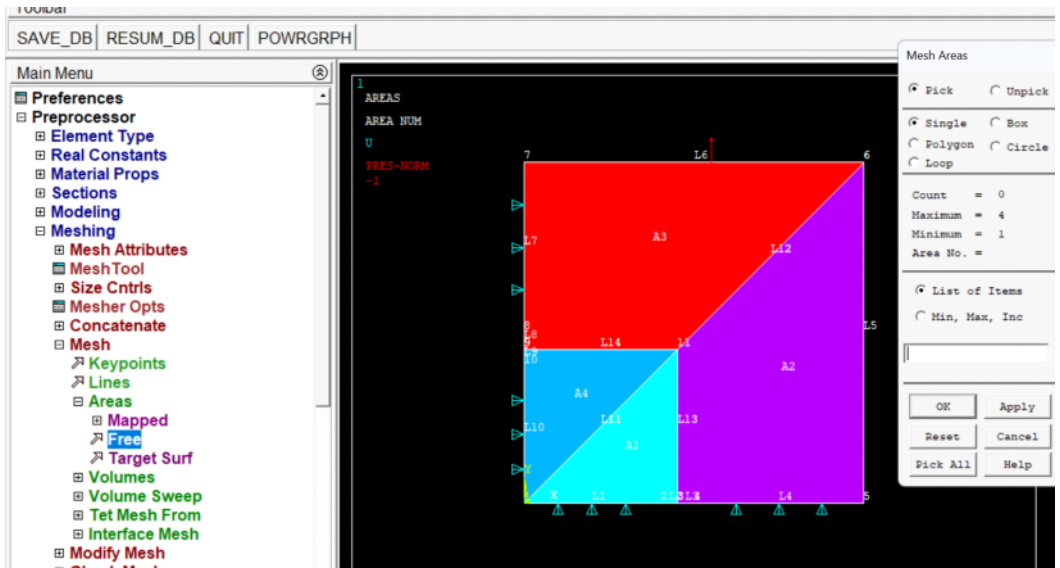


Specify the element formation around the crack tips:

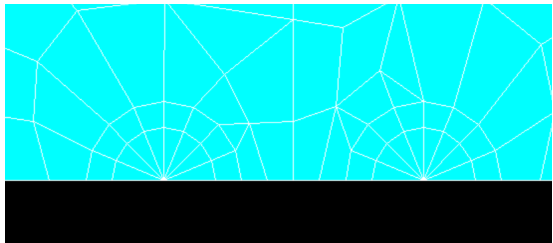
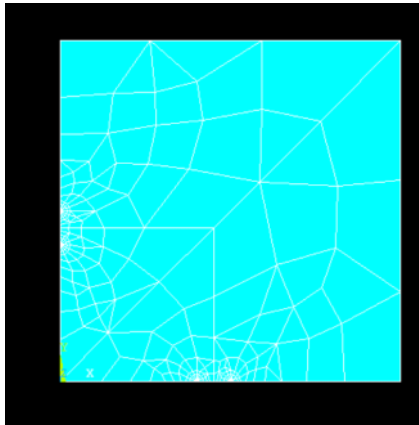


Repeat this for the other 3 CTs

Next we mesh the domain

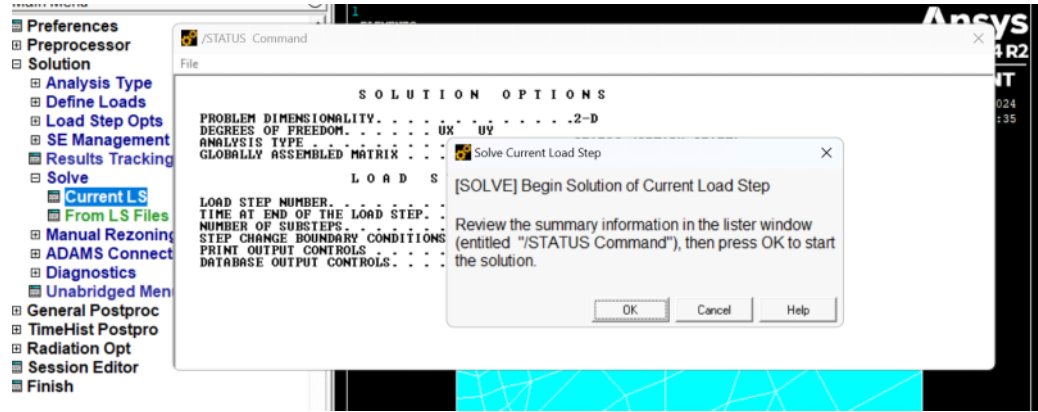


Pick all to mesh the whole domain

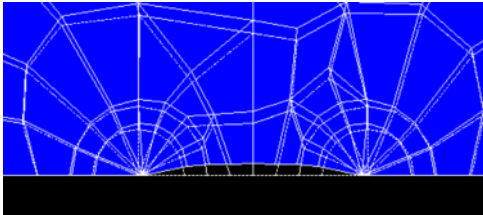
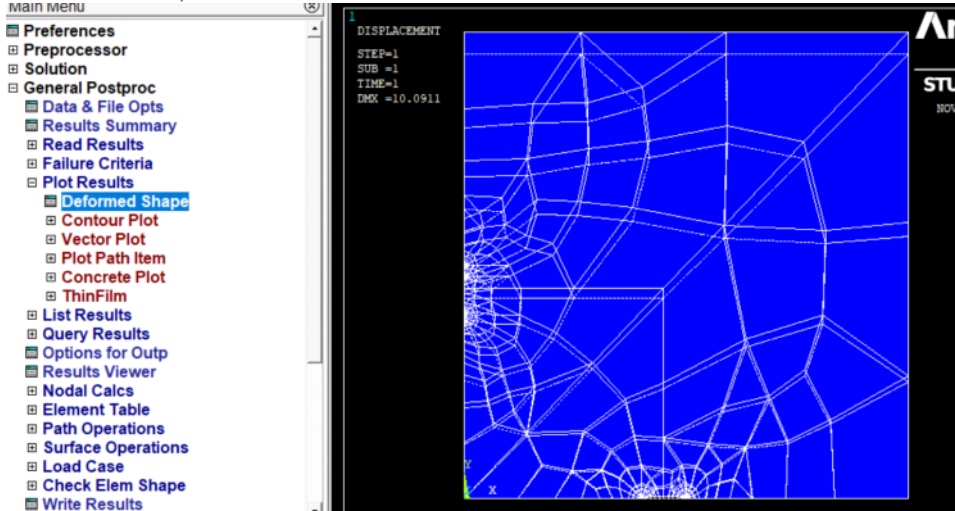


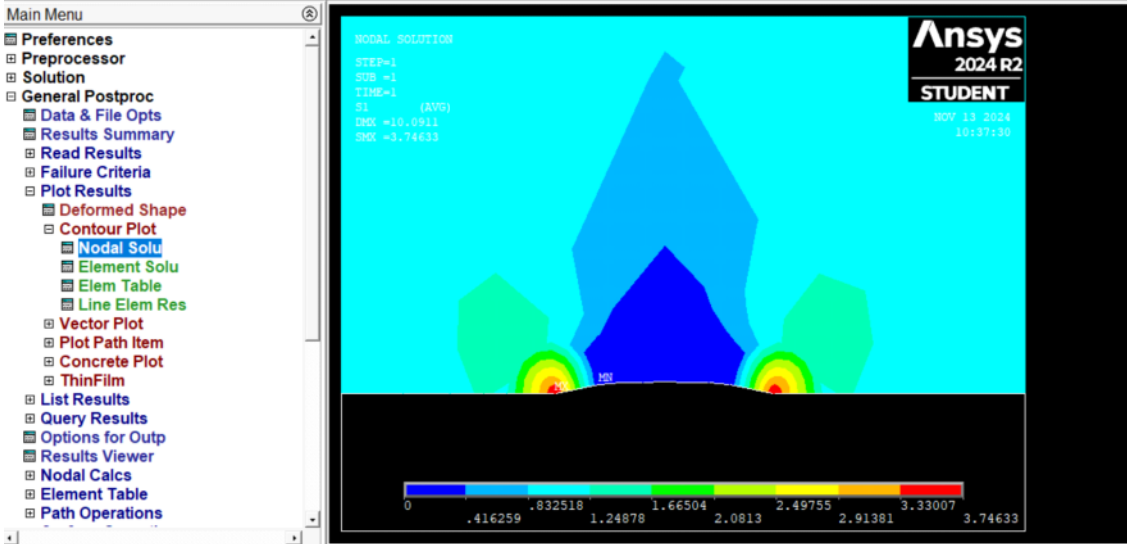


solution



Post processing  
- Deformed shape





Calculating SIF:

- Method 1: Displacement

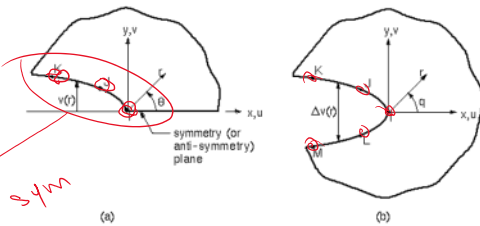


Figure 4) Nodes used for the approximate crack-tip displacements in (a) Half model and (b) Full model

or alternatively from the first quarter point element:

$$v = K_I \frac{\kappa + 1}{2G} \sqrt{\frac{r}{2\pi}}$$

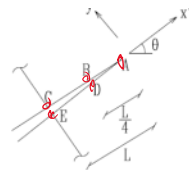
$$\left. \begin{aligned} u' &= \bar{u}'_A + (-3\bar{u}'_A + 4\bar{u}'_B - \bar{u}'_C) \sqrt{\frac{r}{L}} + (2\bar{u}'_A + 2\bar{u}'_C - 4\bar{u}'_B) \frac{r}{L} \\ v' &= \bar{v}'_A + (-3\bar{v}'_A + 4\bar{v}'_B - \bar{v}'_C) \sqrt{\frac{r}{L}} + (2\bar{v}'_A + 2\bar{v}'_C - 4\bar{v}'_B) \frac{r}{L} \end{aligned} \right\} \rightarrow K_I = \frac{2G}{\kappa + 1} \sqrt{\frac{2\pi}{L}} (-3\bar{v}'_A + 4\bar{v}'_B - \bar{v}'_C)$$

Recall for 1D

$$u = u_1 + \frac{\sqrt{x}}{\sqrt{L}} (-3u_1 - u_2 + 4u_3) + \frac{2x}{L} (u_1 + u_2 - 2u_3)$$

$$\left\{ \begin{matrix} K_I \\ K_{II} \end{matrix} \right\} = \frac{1}{2} \frac{2G}{\kappa + 1} \sqrt{\frac{2\pi}{L}} \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix} \begin{bmatrix} -3\bar{u}'_A + 4(\bar{u}'_B - \bar{u}'_D) - (\bar{u}'_C - \bar{u}'_E) \\ -3\bar{v}'_A + 4(\bar{v}'_B - \bar{v}'_D) - (\bar{v}'_C - \bar{v}'_E) \end{bmatrix}$$

278 Mixed mode generalization:



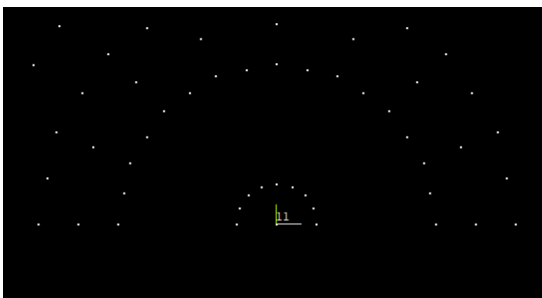
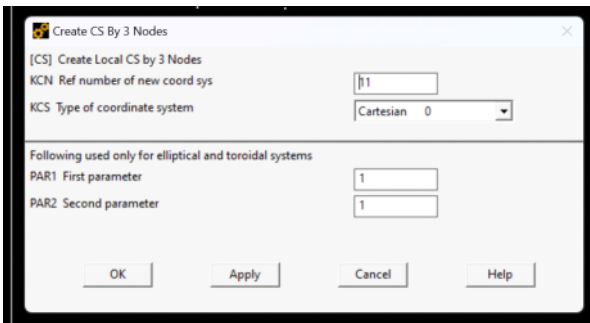
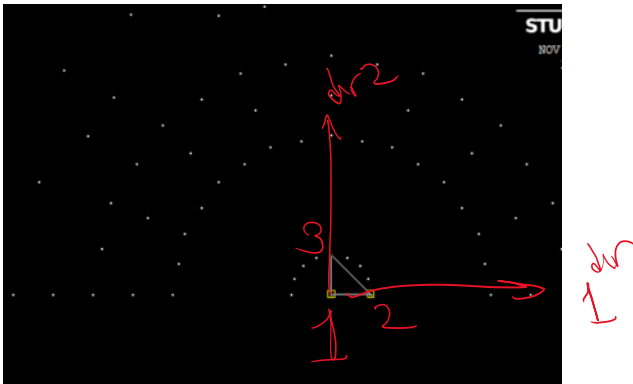
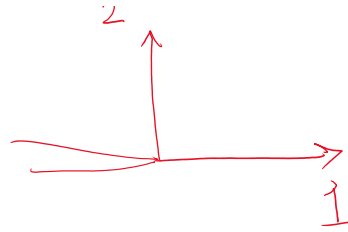
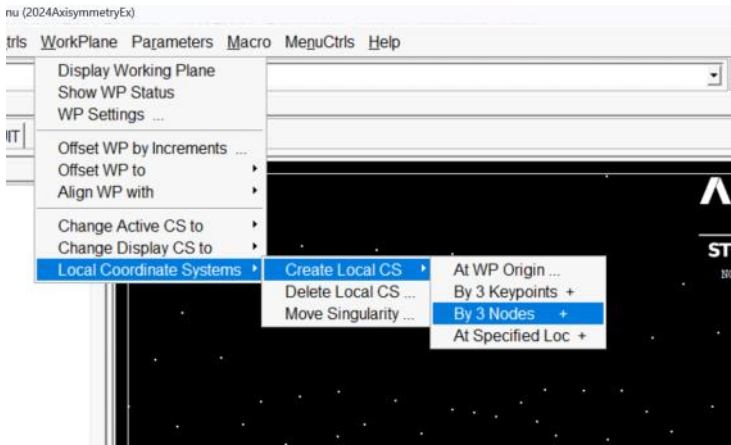
To get K from this approach, we need to do the following:

- a. Define a local coordinate system
- b. Define points 1 to 3 (or 1 to 5) in the figure above.

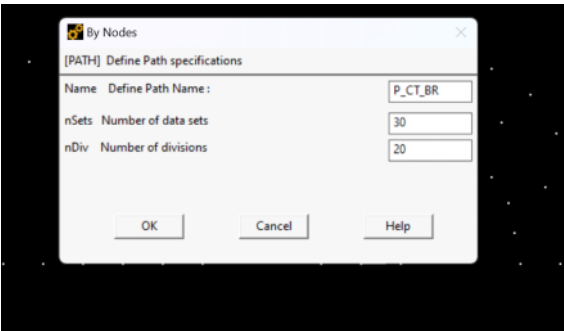
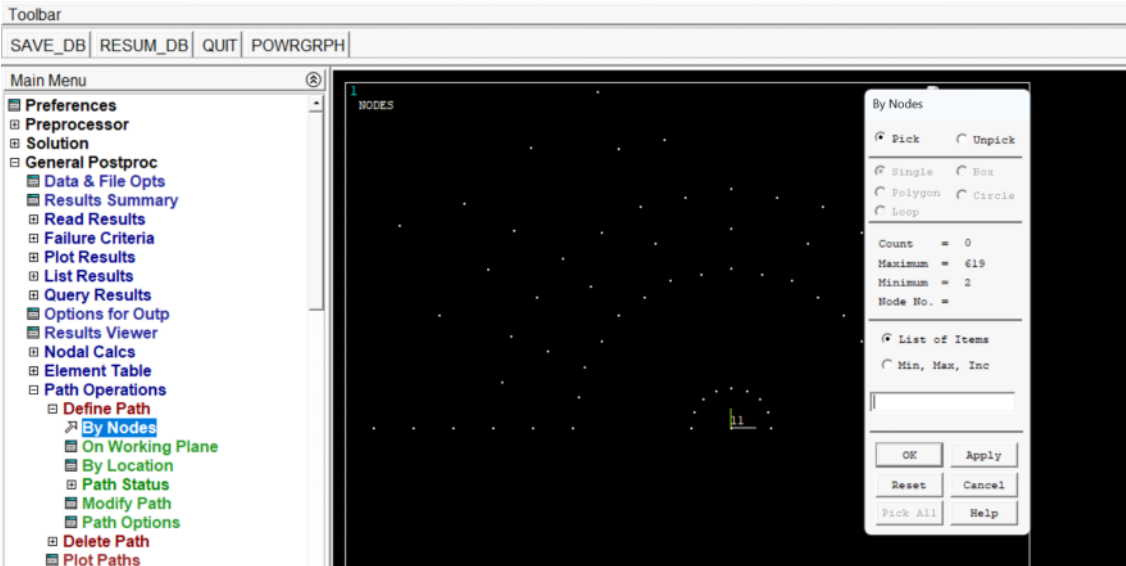
Plot -> nodes  
Create a local coordinate system



### Create a local coordinate system



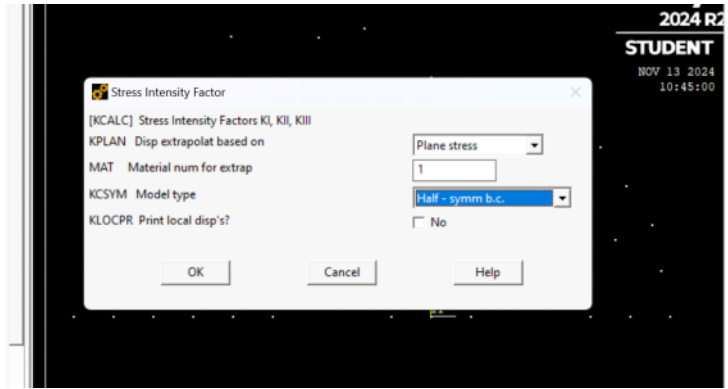
-----  
Define the three nodes used to calculate K:  
This is done using the "path"



\*\*\*\*\* PATH STATUS \*\*\*\*\*

Path	nPts	nSets	nDiv
P_CT_BR	3	30	20

- Preprocessor
- Solution
- General Postproc
  - Data & File Opts
  - Results Summary
  - Read Results
  - Failure Criteria
  - Plot Results
  - List Results
  - Query Results
  - Options for Outp
  - Results Viewer
  - Nodal Calcs
    - Total Force Sum
    - Sum @ Each Node
    - Summation Pt
    - Stress Int Factr
  - Element Table
  - Path Operations
  - Surface Operations



\*\*\*\* CALCULATE MIXED-MODE STRESS INTENSITY FACTORS \*\*\*\*

ASSUME PLANE STRESS CONDITIONS

ASSUME A HALF-CRACK MODEL WITH SYMMETRY BOUNDARY CONDITIONS (USE 3 NODES)

EXTRAPOLATION PATH IS DEFINED BY NODES: 147 154 153  
WITH NODE 147 AS THE CRACK-TIP NODE

USE MATERIAL PROPERTIES FOR MATERIAL NUMBER 1  
EX = 1.0000 NUXY = 0.0000 AT TEMP = 0.0000

\*\*\*\* KI = 1.2627 , KII = 0.0000 , KIII = 0.0000 \*\*\*\*

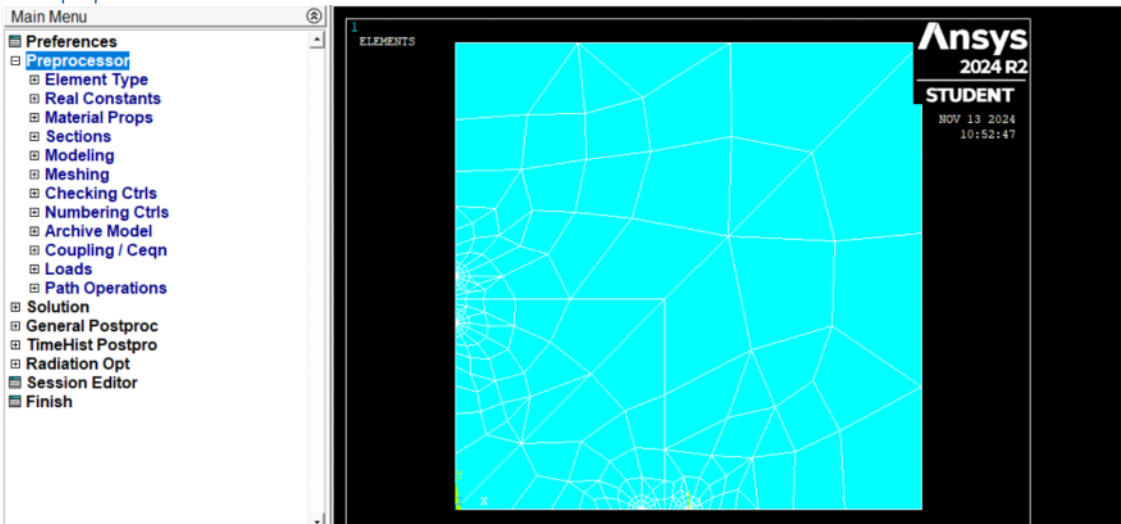
Analytical K  
assuming infinite domain approx.

$$K_I = \sqrt{\pi a} = \sqrt{\pi \times 5} = 1.2533$$

Method 2:

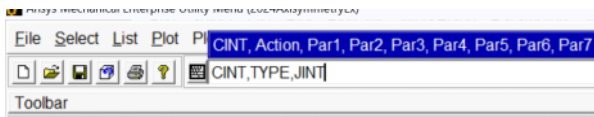
Calculating K through J (Equivalent Domain Integral: EDI)

Go to preprocessor

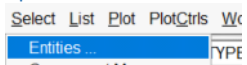


1. Define an integral  
CINT,NEW,1

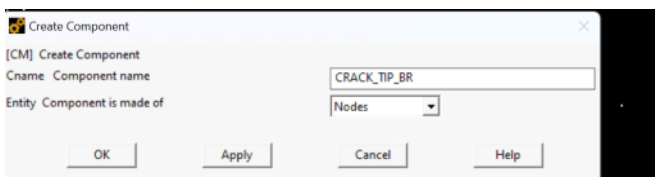
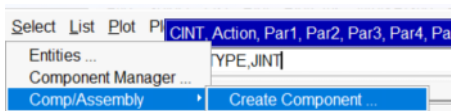
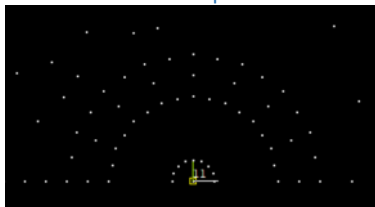
2. Specify the type of contour integral to be done (here it is the J-integral)



3. We need to create an assembly containing only the crack tip:



Select the crack tip



CINT,CTNC,CRACK\_TIP

-----