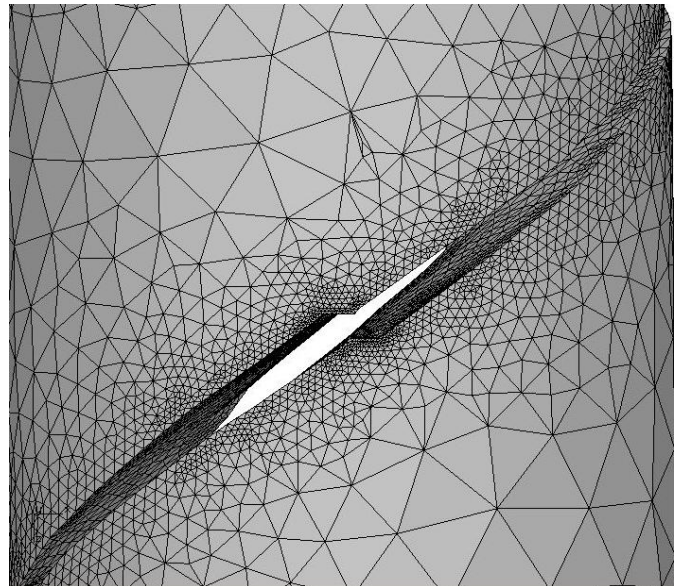


FRANC3D

ANSYSTM Tutorial

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Version 9.2



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<https://www.fracanalysis.com/>

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1.0 Introduction

This tutorial introduces the fracture simulation capabilities of FRANC3D Version 9 and ANSYS 2025 R2; older versions of ANSYS should work also. FRANC3D is introduced by analyzing a single surface crack in an ANSYS finite element (FE) model of a cube.

Subsequent tutorials (see the FRANC3D Tutorials #2-15 document) build on this first example and describe additional capabilities and features of the software. It is intended that the user perform operations as they are presented, but you can experiment and consult other reference documentation whenever necessary.

The FRANC3D menu and dialog button selections are indicated by bold text, such as **File** menu. Window regions along with dialog options, fields and labels are underlined. Model names and file names are indicated by *italic text*.

The FRANC3D main window is shown in Fig 1.1. When you start the tutorial, the first thing to do is set the working directory using the **File** → **Work Directory** menu option, Fig 1.1.

Before or after a model is imported, you can set the FE model units, Fig 1.2. The selected units are applied to plots and figures that FRANC3D displays and are used during crack growth in combination with fatigue crack growth rate data, which also has units.

You can also set your preferences; for example, it is helpful to set the path to the ANSYS executable. You can view preferences using the **Edit** → **Preferences** menu option. The Preferences dialog is described in Section 5.4 of the FRANC3D Reference document. The ANSYS executable and license setting is discussed in Step 5.2 in Section 2.5 of this tutorial.

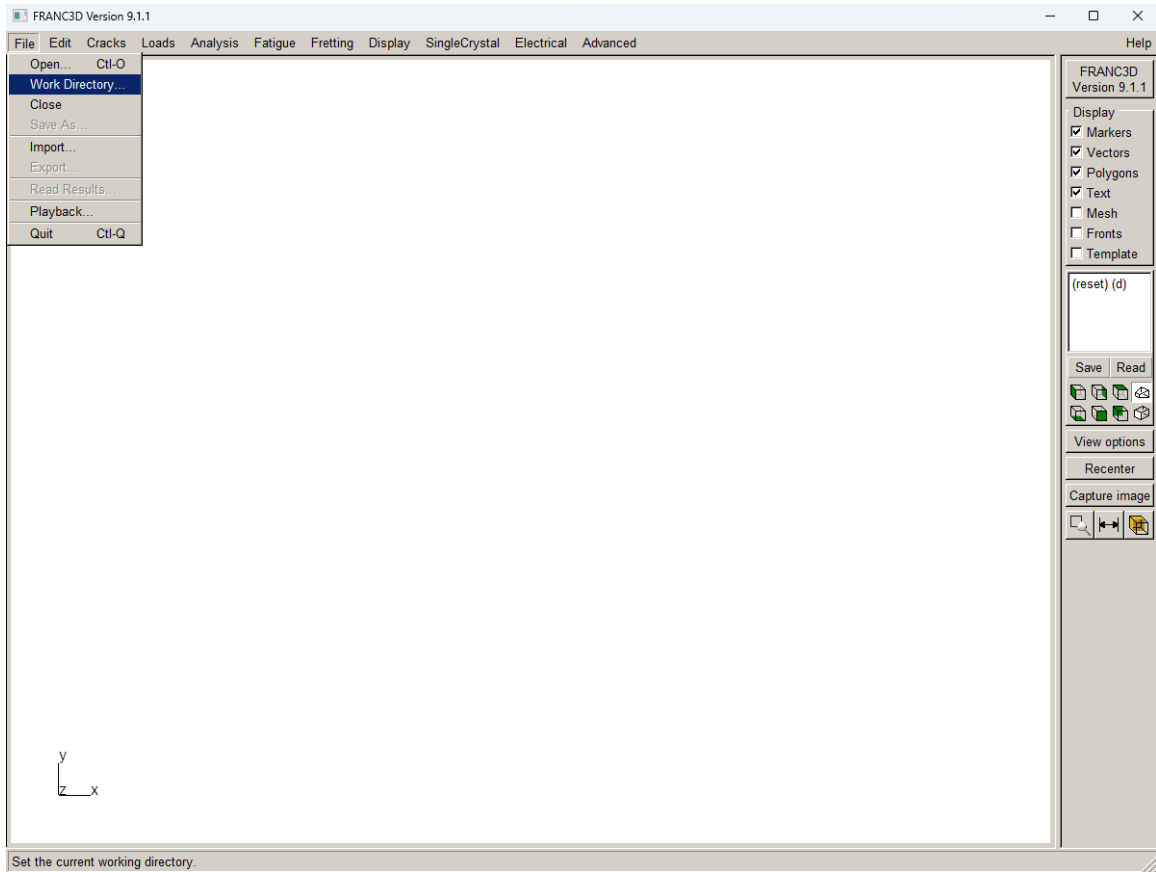


Figure 1.1 FRANC3D main window – Work Directory menu option.

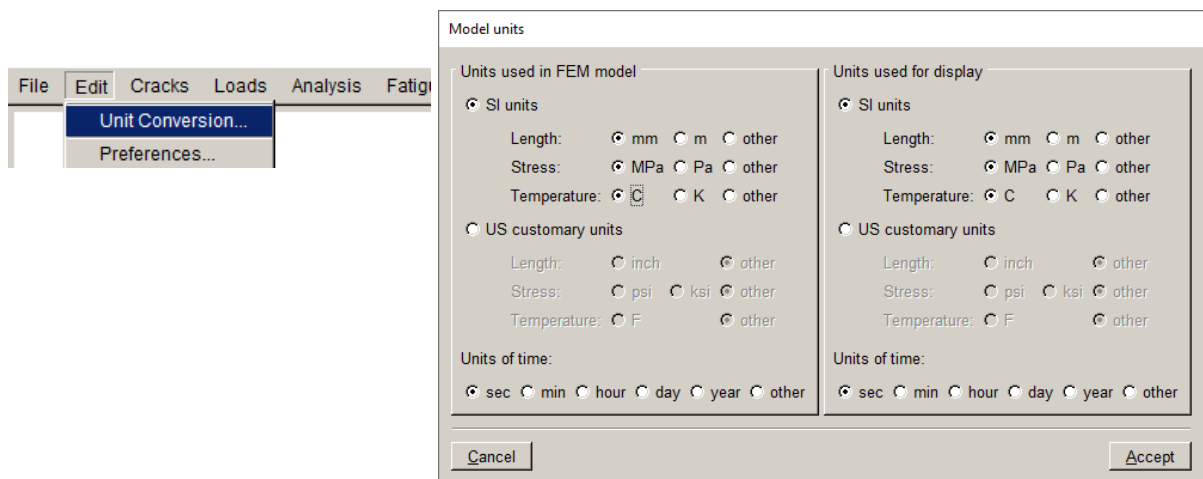


Figure 1.2 FRANC3D FE model unit conversion menu option and dialog.

2.0 Tutorial 1: Crack Insertion and Growth in a Cube

In this tutorial, a single surface crack in a cube under far-field tension is modeled. It is assumed that the user is familiar with a pre-processor for ANSYS; we use ANSYS Mechanical APDL. Once the FE model is created, the FRANC3D steps to: 1) read the model, 2) insert a crack, 3) perform the ANSYS analysis, and 4) compute stress intensity factors, are all described.

2.1 Step 1: Create the ANSYS FE Model

First, you must create a cube model using any pre-processor for ANSYS. We outline the steps to create the FE model with boundary conditions to ensure that you have a model that can be used with FRANC3D.

1. Create a 10x10x10 cube geometry; units of length are mm.
2. Subdivide the edges for meshing, using 10 to 20 subdivisions.
3. Define the element type as quadratic; use brick or tetrahedral elements to mesh the volume.
4. Define the material properties; elastic modulus (E) is 10000 and Poisson's ratio is 0.3; units for E are MPa..
5. Boundary conditions consist of displacement constraints on the bottom surface and uniform traction (a negative pressure) on the top surface. The bottom surface is constrained in the y-direction, the bottom left edge is also constrained in the x-direction, and the point at the origin is also constrained in the z-direction. The top surface traction is 10 MPa.
6. Write the *.cdb* file (*Ansys_Cube.cdb*). When saving the *.cdb* file, you do not need to write the IGES data. ANSYS Classic/APDL and ANSYS Workbench both allow you to write *.cdb* files. See Appendix B for Workbench options.

The resulting model should appear as in Fig 2.1. The symbols for the boundary conditions are displayed attached to the mesh model.

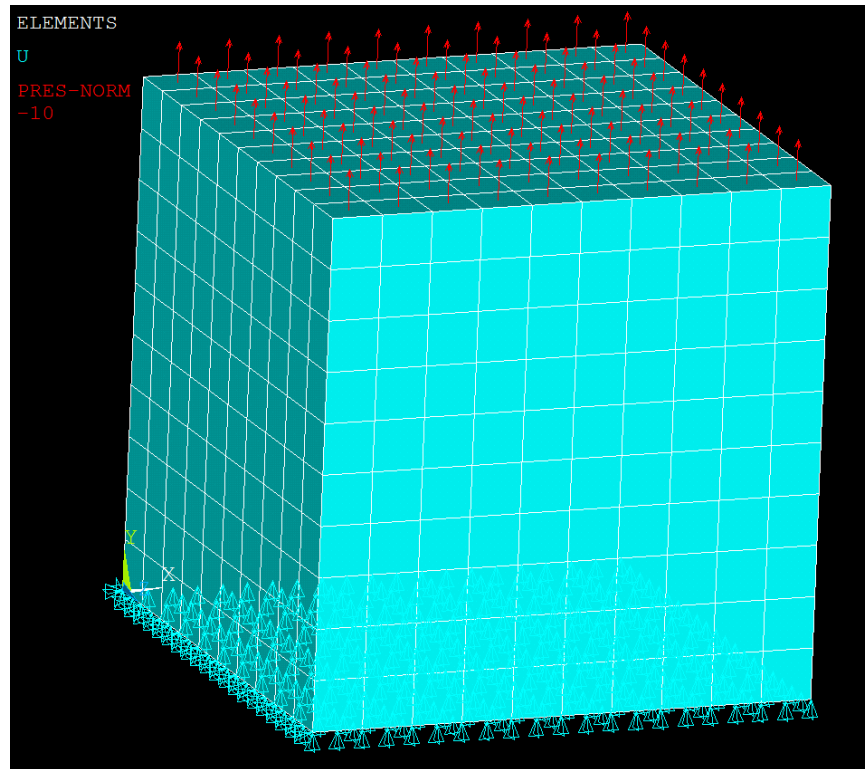


Figure 2.1 ANSYS cube with brick element mesh and boundary conditions.

2.2 Step 2: Reading ANSYS FE Model into FRANC3D

Once the ANSYS FE model is created and the *.cdb* file is saved, we proceed with importing the FE model into FRANC3D. We use the *Ansys_Cube.cdb* file that was created in the previous step.

You can choose to do either Step 2 or Step 3 (see Section 2.3).

Step 2.1: Importing Complete ANSYS FE Model

Start with the FRANC3D graphical user interface (see Fig 1.1) and select **File** → **Import**, Fig 2.2. In the dialog shown in Fig 2.3, choose Complete Model and select **Next**.

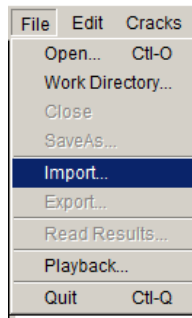


Figure 2.2 File Import menu option

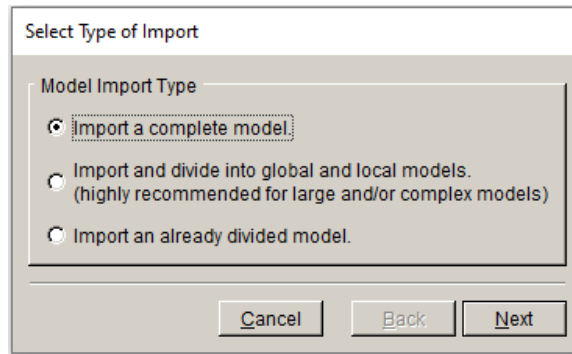


Figure 2.3 Model Import Type dialog

Switch the Mesh File Type radio button in the Select Import Mesh File window, Fig 2.4, to ANSYS and select the file name for the model, *Ansyz_Cube.cdb*, and then select **Next**.

Note that the mesh file type can be set in the Preferences so that you do not have to change the mesh file type radio button if you always use ANSYS.

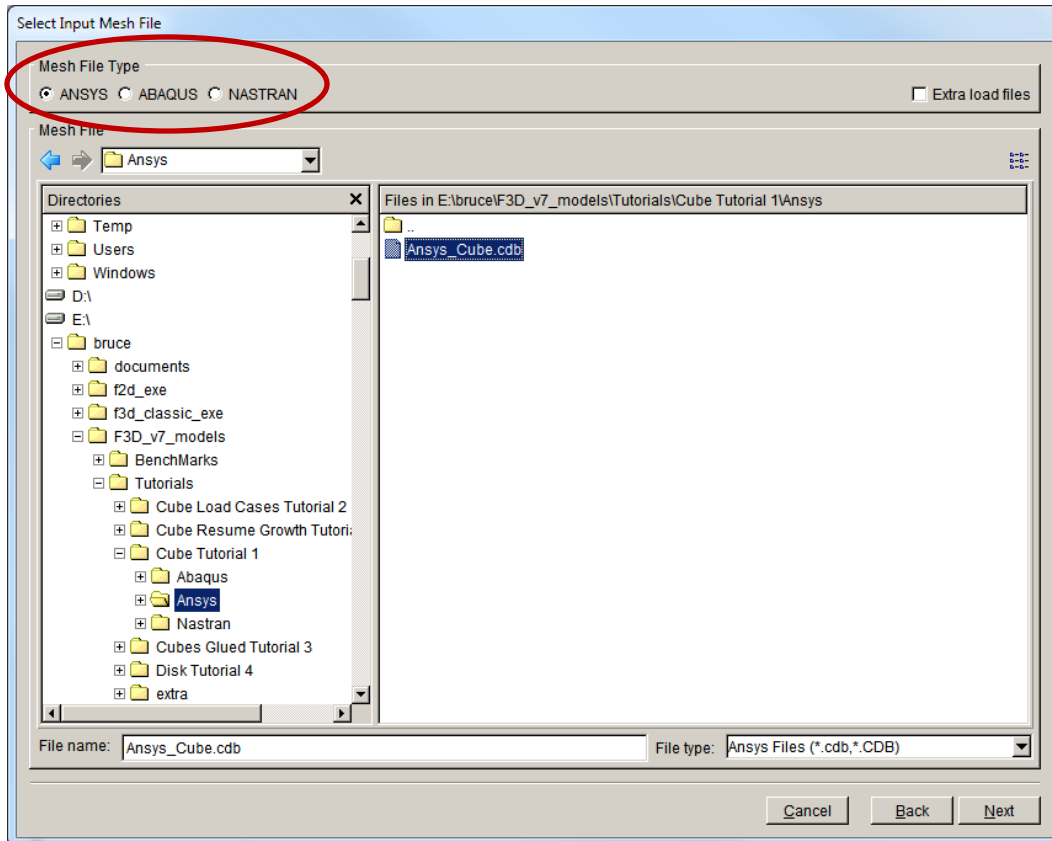


Figure 2.4 Select Input Mesh File dialog box.

Step 2.2: Select the Retained Items in the FE Model

The next panel, Fig 2.5, allows you to choose the mesh surface facets that are retained from the *.cdb* file. Surfaces with boundary conditions are blue, Fig 2.6, and turn red when selected, Fig 2.7. You can rotate the model to view all surfaces. We retain the surfaces with boundary conditions (top and bottom of the cube) by choosing Select all. The boundary conditions are transferred automatically to the new mesh when a crack is inserted.

Select **Finish** to proceed. The Units dialog will be displayed; if you have already set the units using the **Edit** menu, just select **Finish**.

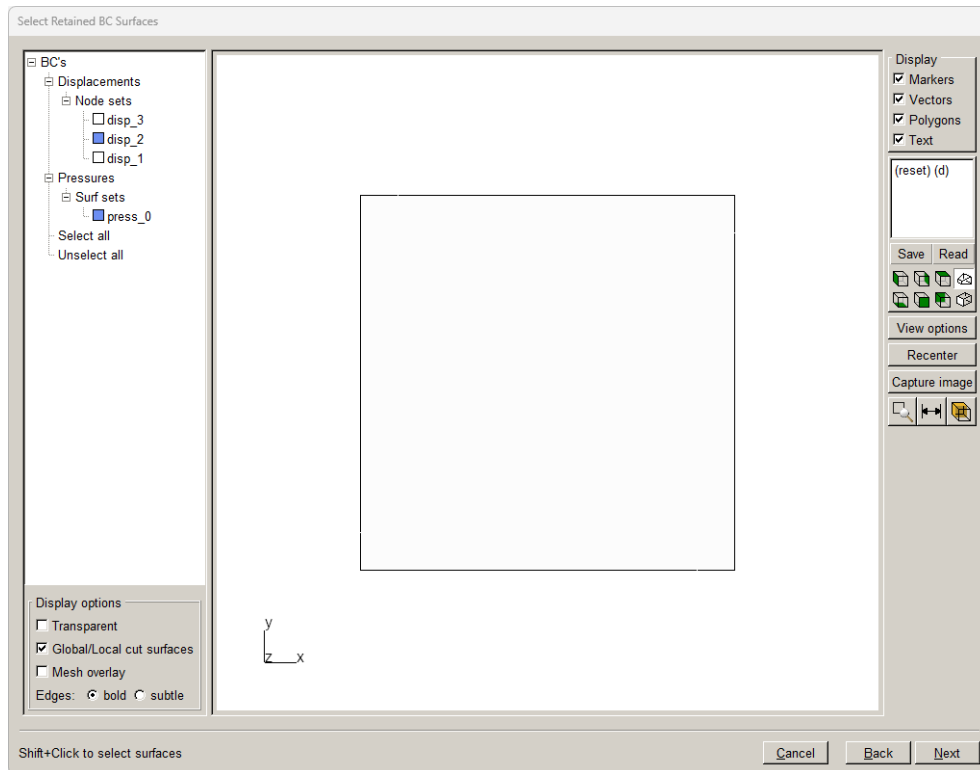


Figure 2.5 Select Retained BC Surfaces wizard panel.

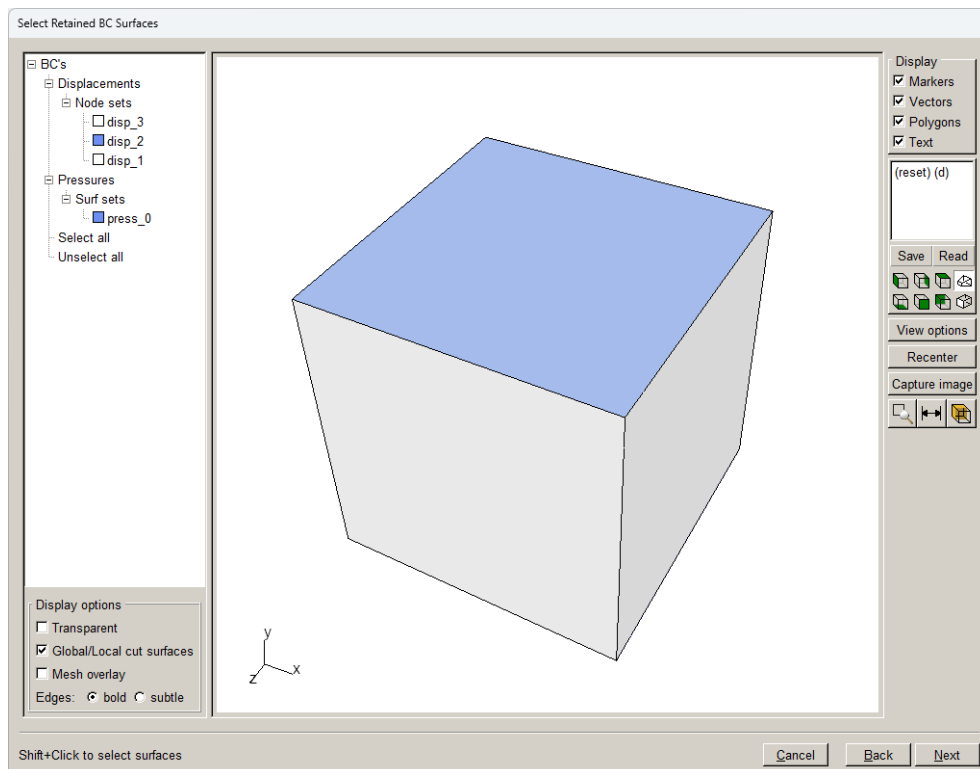


Figure 2.6 ANSYS Model retain BC Surfaces wizard panel - unselected surface.

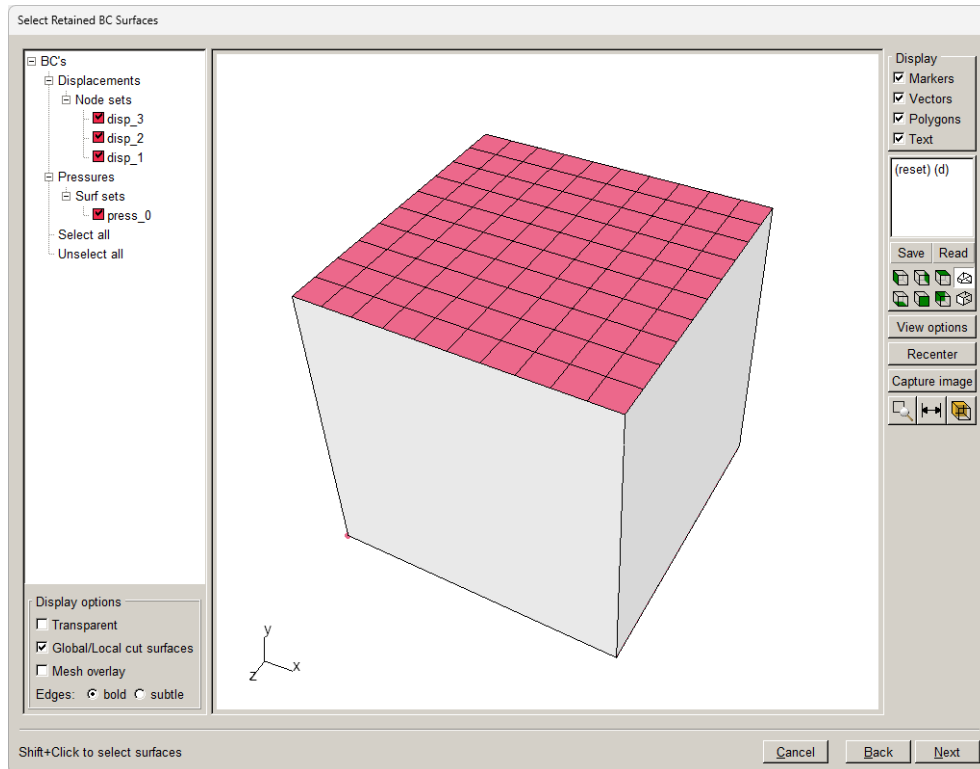


Figure 2.7 ANSYS Model retain BC Surfaces wizard panel – selected surfaces.

Surface Mesh NOT Retained

The surface mesh facets do not have to be retained, and if a crack is inserted into a surface that has boundary conditions, then the surface mesh must not be retained. In such a case, the boundary condition data is mapped to the remeshed surface. In practice, transfer of boundary condition data is simpler and more precise if the surface mesh is retained, but sometimes this is not possible. FRANC3D will automatically transfer the boundary condition data to the new mesh in either case.

The model is displayed in the FRANC3D modeling window, Fig 2.8. You can turn on the surface mesh, or manipulate the model view by rotating, *etc.* Fig 2.8 shows the retained mesh on the top surface (retained bottom surface is not visible) where the boundary conditions are applied. See Section 2 of the Reference document for a description of the view manipulation.

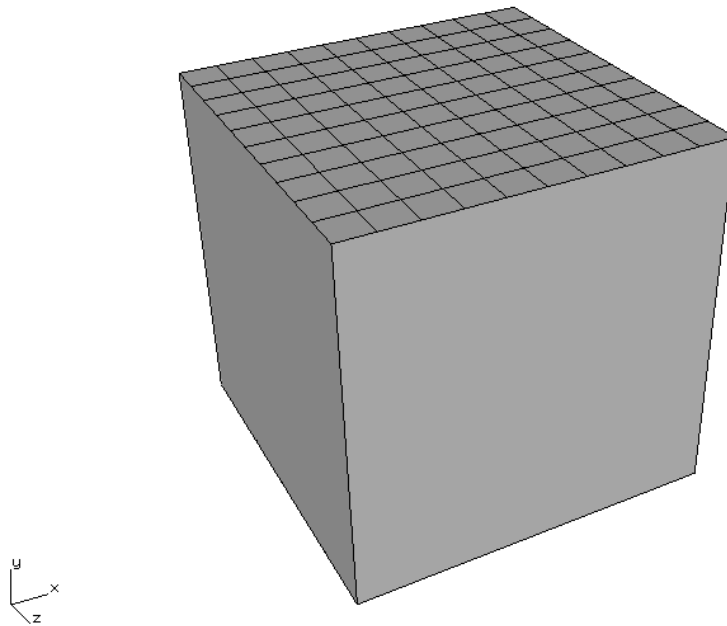


Figure 2.8 ANSYS model imported into FRANC3D with retained facets on top of the cube.

If you have not set a working directory (using the **File** → **Work Directory** menu option), FRANC3D might present the dialog shown in Fig 2.9 prior to displaying the model. Select **Yes** to set the directory as the folder where the *.cdb* file resides. Many files are created during a crack growth simulation, and it is best to keep them together in one folder.

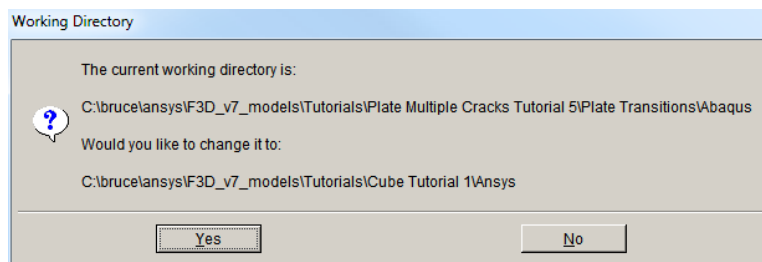


Figure 2.9 Set Working Directory dialog.

2.3 Step 3: Importing and Subdividing the Model

If you did Step 2, close the model using the **File** → **Close** menu option before starting this step.

The ANSYS model can be split into smaller parts before inserting the crack. Go to **File** → **Import** and choose the Import and divide into global and local models radio button, Fig 2.10, and choose the *Ansys_Cube.cdb* model, Fig 2.11. Remember to set the Mesh File Type radio button to ANSYS.

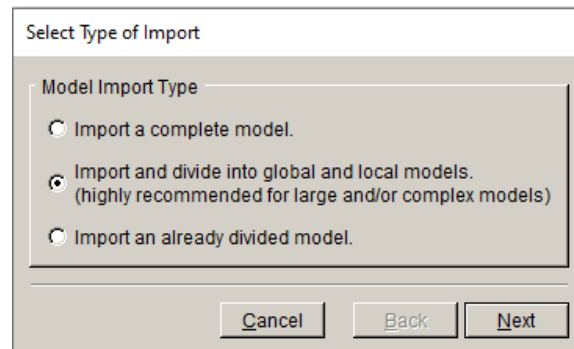


Figure 2.10 Model Import Type dialog with Import and divide selected.

Instead of the retain BC surfaces dialog (see Fig 2.5), the Define a Local Submodel window, Fig 2.12, appears. Figs 2.13 and 2.14 show two of the submodel selection tools. Using the options on the left, a submodel is selected; elements to be retained in the submodel appear red, and the selection is confirmed using the **Crop** button. See Section 4.5 of the Reference document for a more complete description of the submodeling tools.

For this tutorial, the rectangular Rubberband Box is used to create a smaller portion of the cube, Fig 2.15, which has an exposed face for crack insertion.

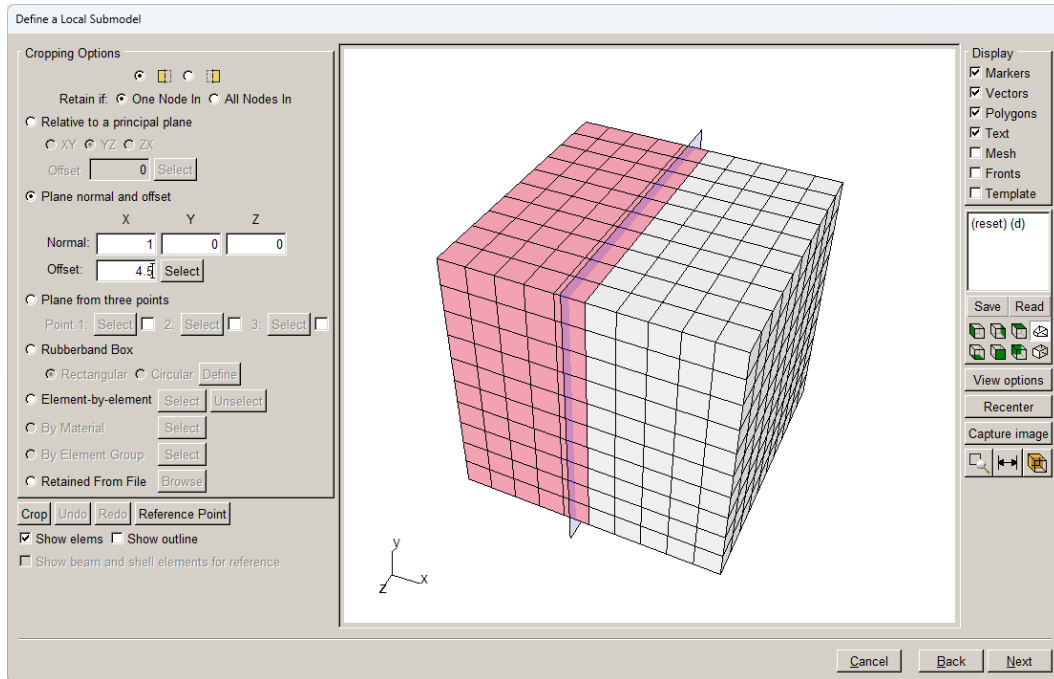


Figure 2.13 Plane Cutting tool – **Crop** will retain the red elements.

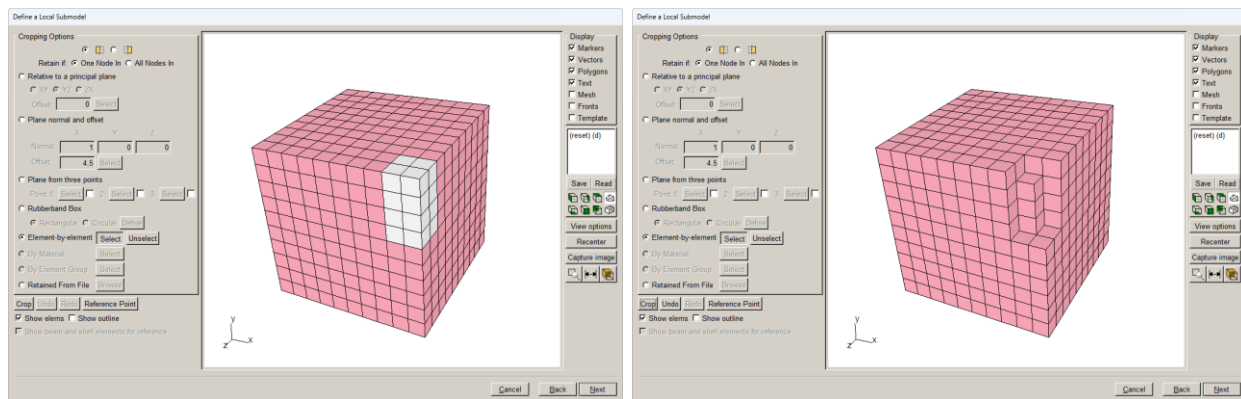


Figure 2.14 Element-by-Element cropping by selecting elements to remove.

We are inserting a surface crack normal to the y-axis (loading direction) and located on the +z surface, and the selection in Fig 2.15 is designed for this. Use the Rubberband Box to select and **Crop** a small portion of the model.

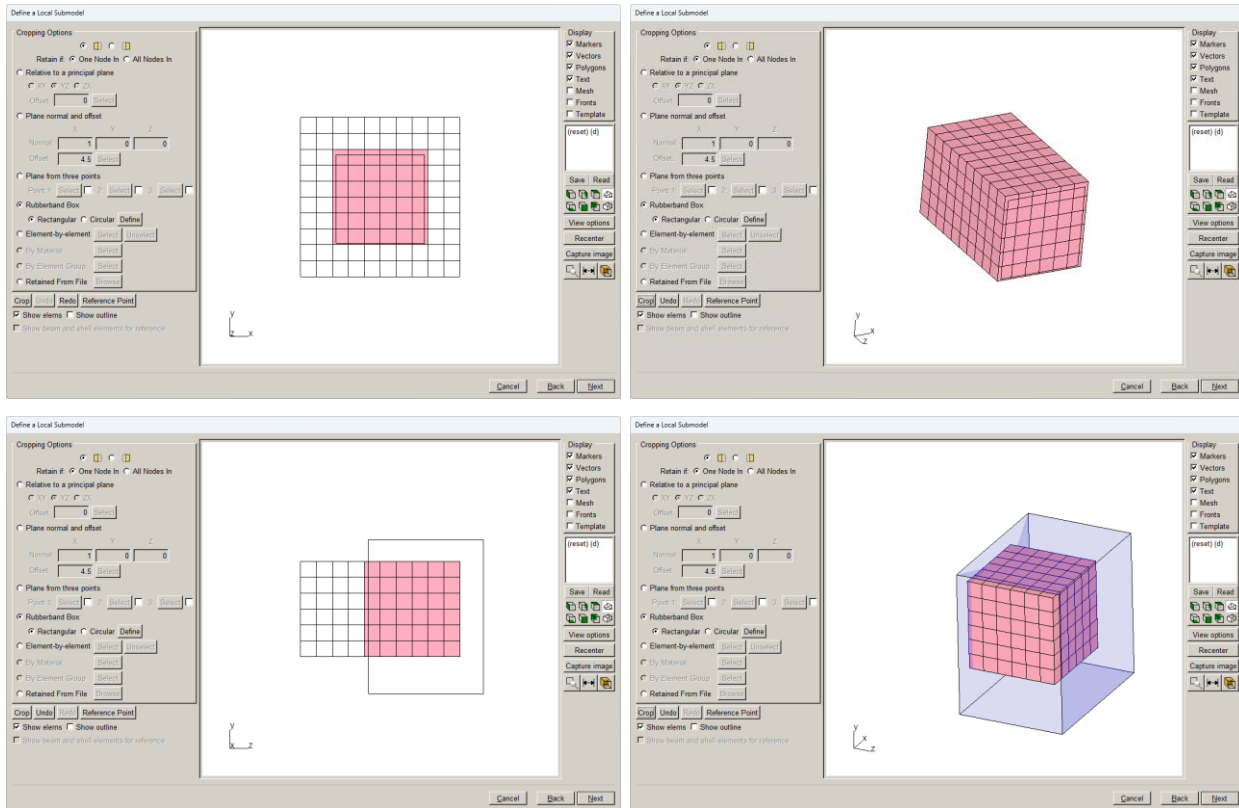


Figure 2.15 Rubberband Box tool used to select a submodel.

Once the elements have been selected and cropped, select **Next**. The Save Files dialog appears, Fig 2.16. Choose the names of the local and global models and their location; we use the default file names and the working directory set earlier. Select **Next**.

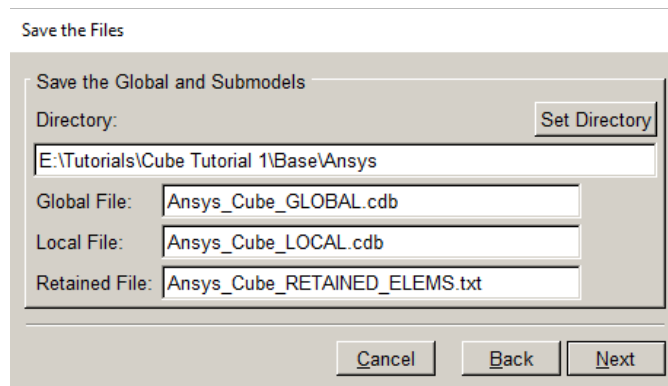


Figure 2.16 Local/Global model save window.

The Select Retained BC Surfaces dialog is displayed next, Fig 2.17, and shows the retained cut-surface. The cropped selection in Fig 2.15 avoids all surfaces with boundary conditions. Click **Next** to proceed.

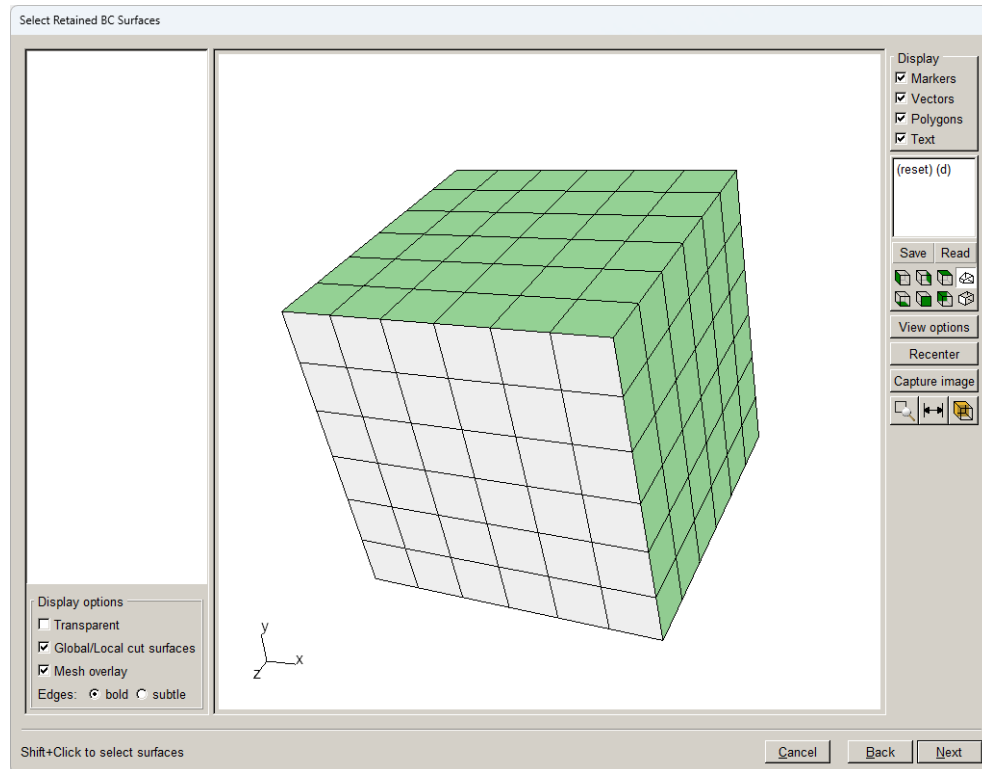


Figure 2.17 Select Retained BC Surfaces wizard panel – showing the retained cut-surface.

The model units, Fig 2.18, must be set. As defined in Step 1, the units for length are mm and the units for stress are MPa. The temperature and time units are not needed for this tutorial.

Select **Finish** and the local model is displayed in the main FRANC3D window, Fig 2.19. The cut-surface mesh facets are retained, while the front surface where the crack will be inserted does not have retained facets.

Specify the units used in FEM model

FEM model units

☒ SI units

Length: ☒ mm ☐ m ☐ other

Stress: ☒ MPa ☐ Pa ☐ other

Temperature: ☒ C ☐ K ☐ other

☐ US customary units

Length: ☐ inch ☒ other

Stress: ☐ psi ☐ ksi ☒ other

Temperature: ☐ F ☒ other

Units of time:

☒ sec ☐ min ☐ hour ☐ day ☐ year ☐ other

Cancel Back Finish

Figure 2.18 FE Model units

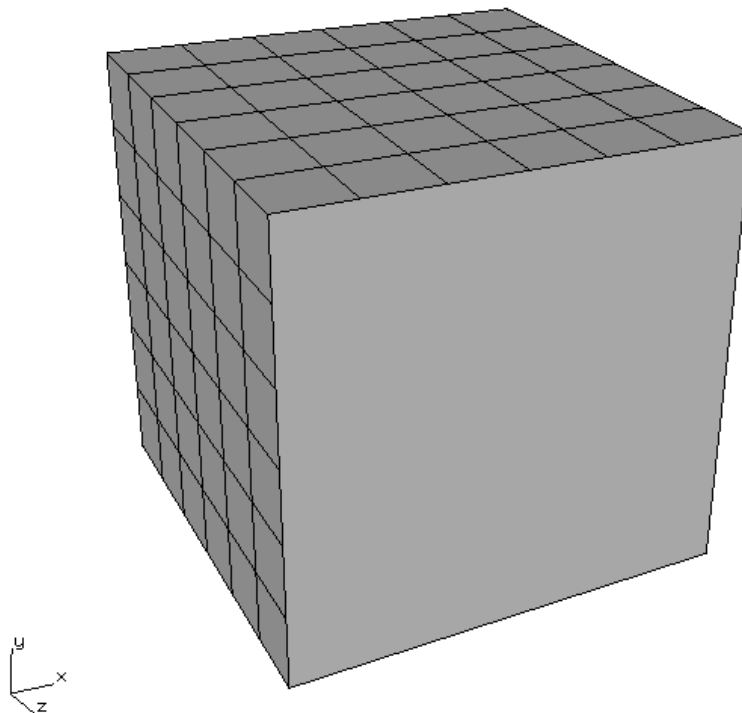


Figure 2.19 Local model with retained mesh facets on cut-surface

2.4 Step 4: Insert a Crack

We now insert a half-penny surface crack into the model. The local submodel from Section 2.3 is used, as opposed to the full model from Section 2.2. In Step 4.1, we describe how to define a new crack, and in Step 4.2, we describe how to insert a flaw from a file.

Note: if you want to try both types of crack insertion, you will have to re-import the uncracked model to insert a crack from a file; inserting a crack from a file will be used in subsequent steps and other tutorials.

Step 4.1: Define a new Crack from FRANC3D Menu

From the FRANC3D menu, select **Cracks** → **New Flaw Wizard**, Fig 2.20. The first panel of the wizard, Fig 2.21, allows you to select the flaw type; a crack with zero volume is the default. We select the Save to file and add flaw radio button so that the *.crk* file is saved for subsequent tutorials. Note that the Symmetry surface crack option is only displayed if symmetry cracks are turned on in the Preferences (see Section 5.2 in the Reference document).

Select **Next**. The second panel, Fig 2.22, lets you choose the crack shape; the default is the ellipse, which is what we want. Select **Next**. The crack is a circle with radius=1, centered on the cube's front (+z) face, and parallel to the xz-plane (normal to y).

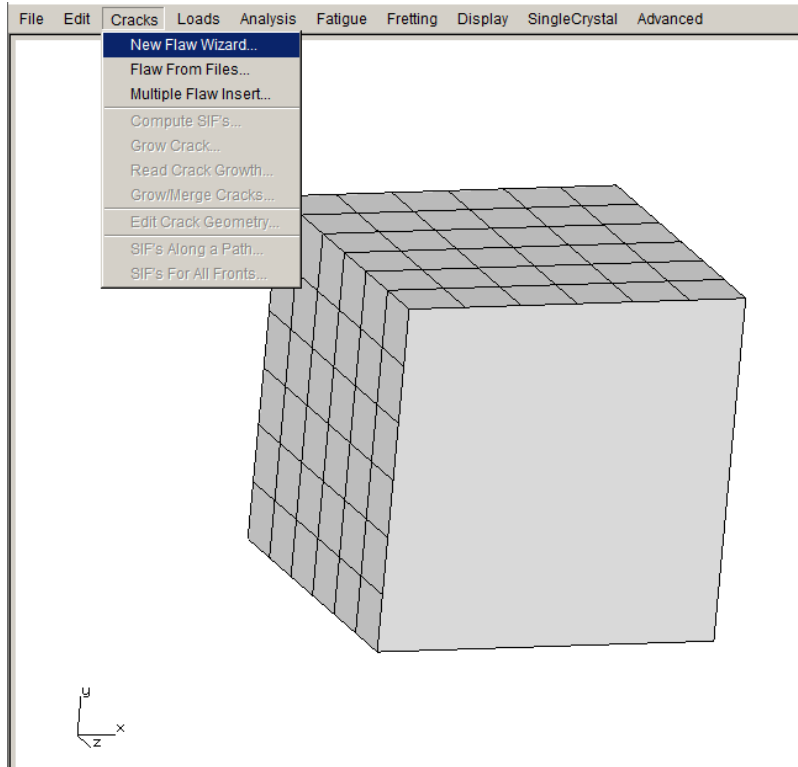


Figure 2.20 New Flaw Wizard menu item selected.

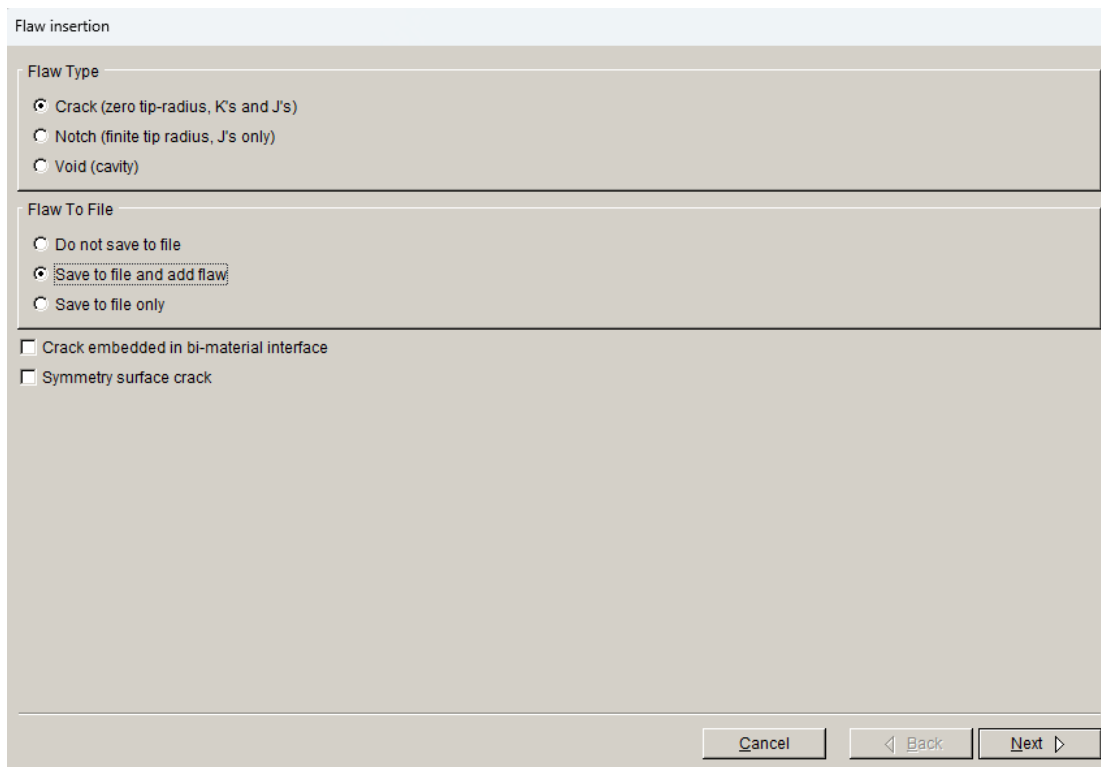


Figure 2.21 New Flaw Wizard first panel – Crack (zero volume flaw) selected.

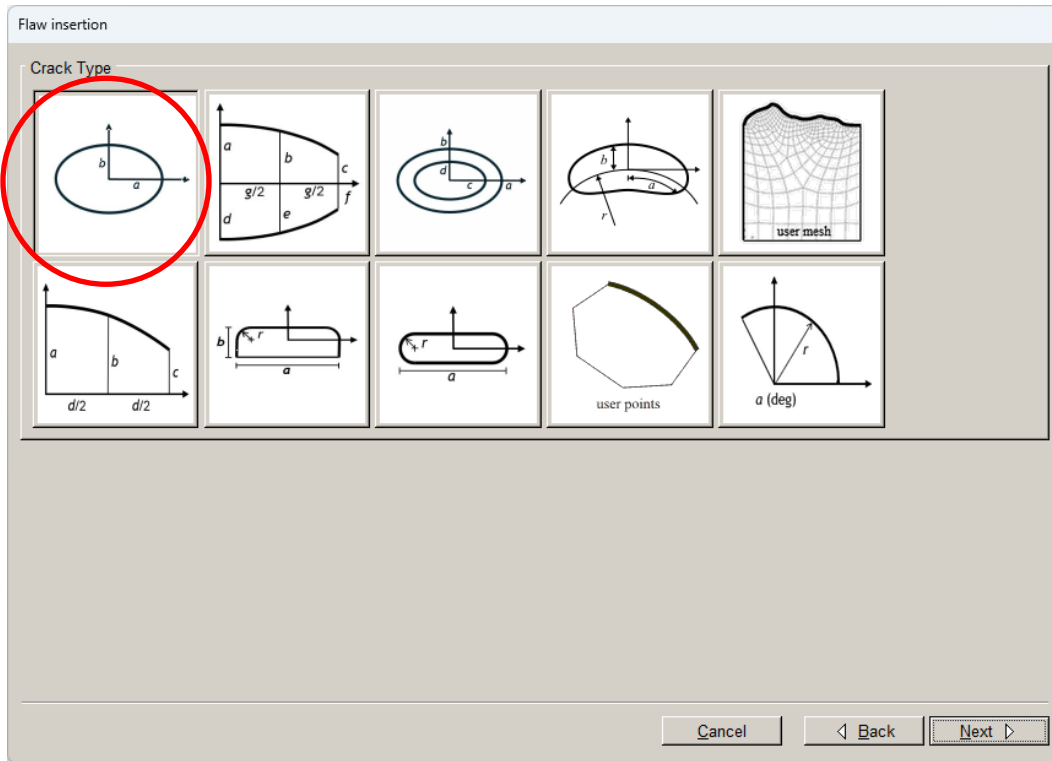


Figure 2.22 New Flaw Wizard second panel – choose the ellipse shape.

The third wizard panel lets you set the dimensions for the ellipse, Fig 2.23; set $\underline{a}=1$ and $\underline{b}=1$. Select **Next**.

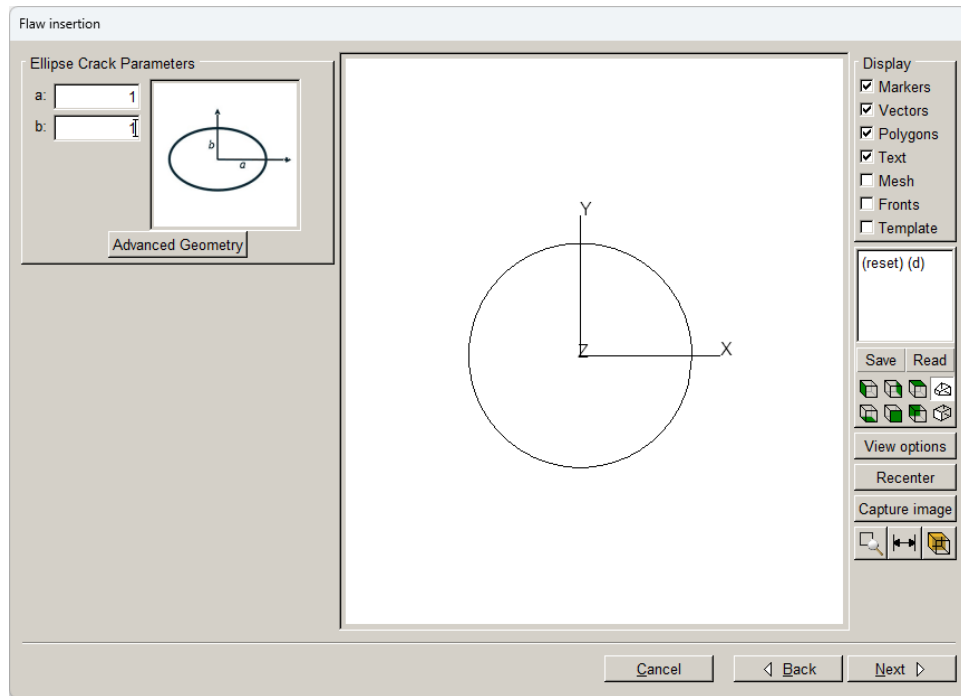


Figure 2.23 New Flaw Wizard third panel – set ellipse dimensions.

The fourth panel lets you set the crack location and orientation, Fig 2.24. We place the crack at the center of the front face at coordinates (5,5,10) and rotate the crack 90 degrees about the Global x-axis. Select **Next**.

The final panel allows you to set the crack front mesh template parameters, Fig 2.25; use the defaults. In practice, the default template radius is based on the crack dimensions and might need to be adjusted depending on the model and crack. Select **Finish**.

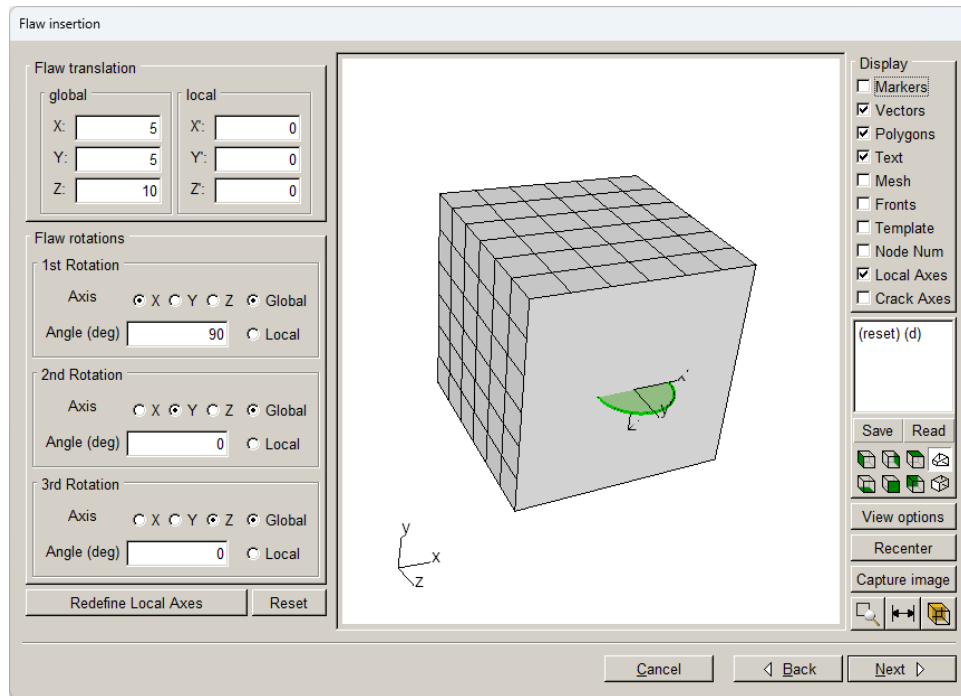


Figure 2.24 New Flaw Wizard fourth panel – set crack location and orientation.

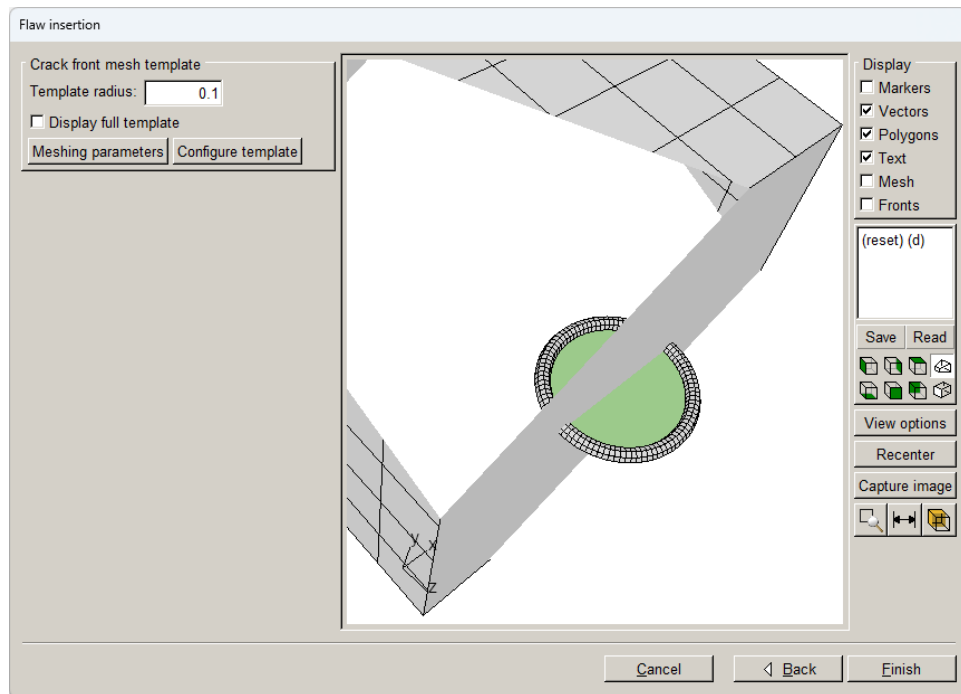


Figure 2.25 New Flaw Wizard final panel – set template mesh parameters.

At this point, you are asked to specify the file name to save the crack information, Fig 2.26; we call it *Cube_Crack.crk*. Select **Accept**.

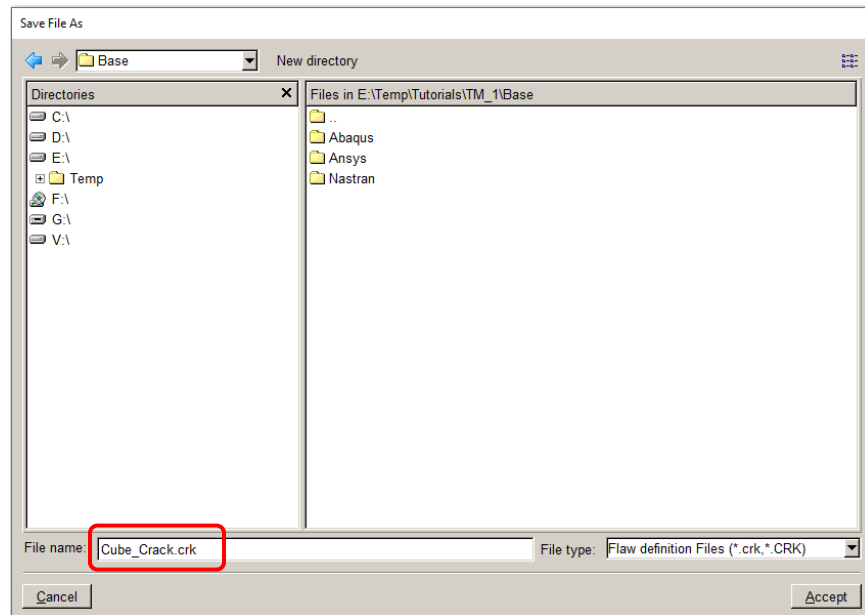


Figure 2.26 New Flaw Wizard final panel – save file as.

The crack is inserted into the model and remeshing occurs. A Flaw Insertion Status dialog is displayed during this process, Fig 2.27. There are three stages: geometric intersection of the crack with the model, surface meshing and volume meshing. The resulting model is displayed in the main FRANC3D window, Fig 2.28; the surface mesh is turned on in this image.

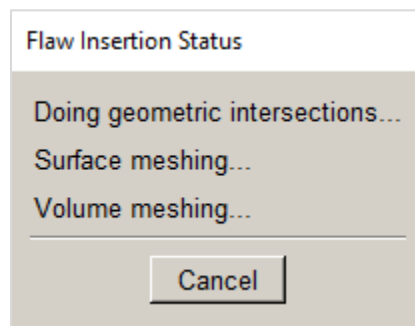


Figure 2.27 Flaw Insertion Status window.

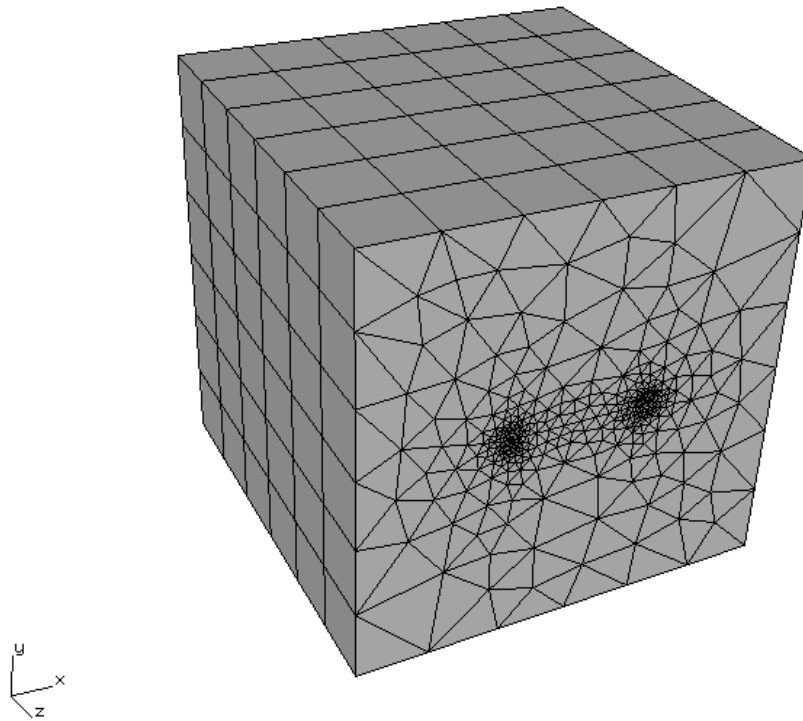


Figure 2.28 Remeshed cracked local model with surface mesh turned on.

Step 4.2: Insert Cracks from Files

If you want to try this step, you must **Close** the model from Step 4.1 and re-import the original uncracked (local) model.

From the FRANC3D menu, select **Cracks** → **Flaw From Files**, Fig 2.29. The first panel, Fig 2.30, lets you select the .crk file. Choose the *Cube_Crack.crk* file name and select **Accept**. This crack was created using the New Flaw Wizard with the save to file option in Step 4.1. The flaw being added is a circle with radius=1, centered on the cube's front face, and parallel to the xz-plane.

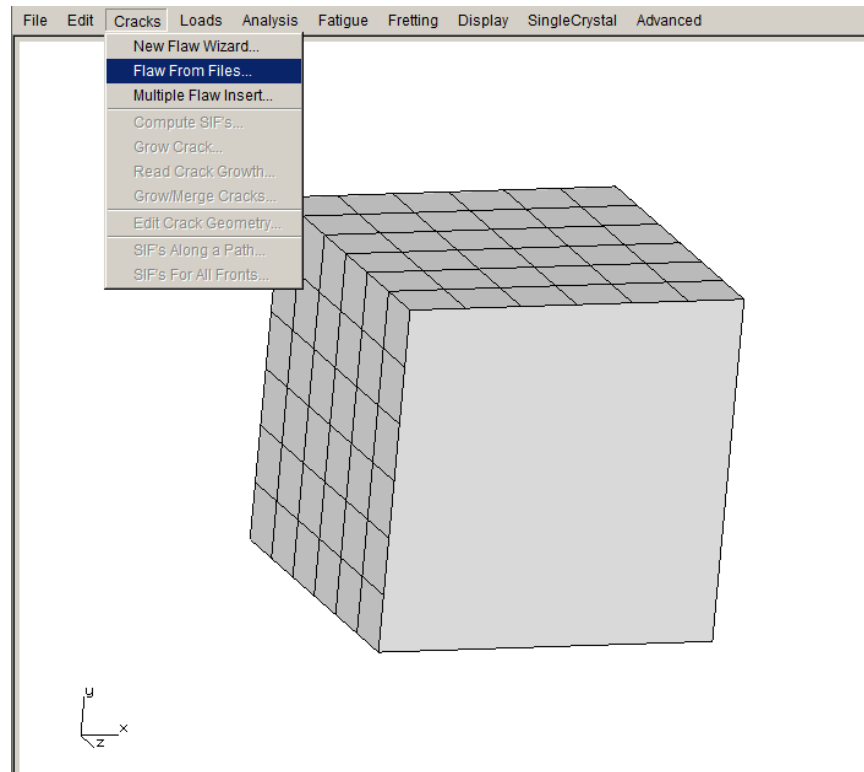


Figure 2.29 Flaw From Files option in Crack menu.

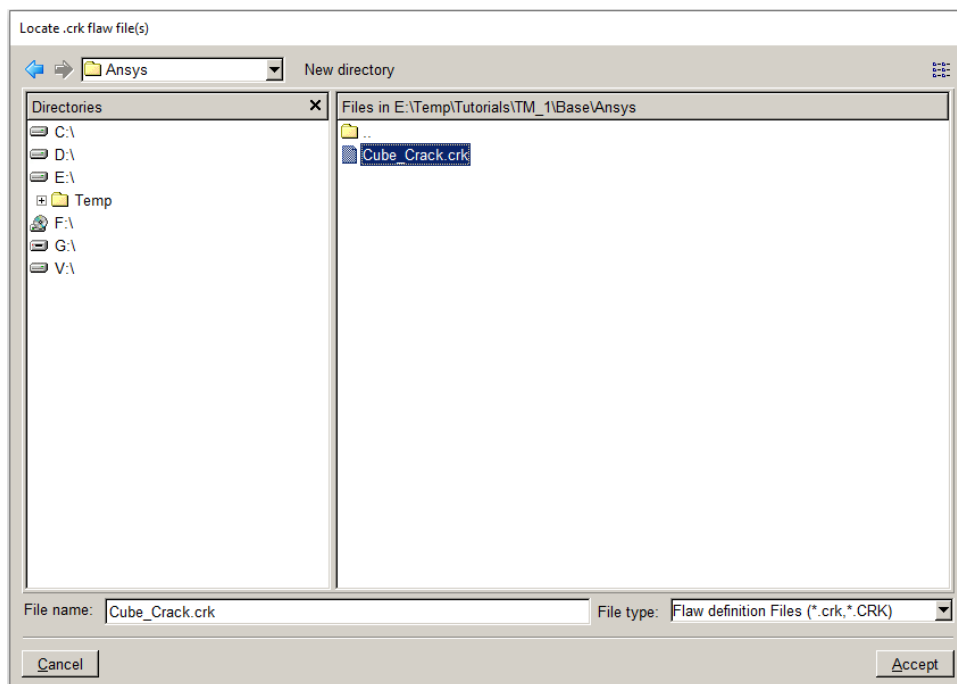


Figure 2.30 Flaw from file dialog to select .crk file.

The next panel of the wizard, Fig 2.31, allows you to adjust the location of the flaw. The orientation cannot be changed. The final panel, Fig 2.32, allows you to set the template mesh parameters. Use the default settings and select **Finish** to complete the insertion and remeshing.

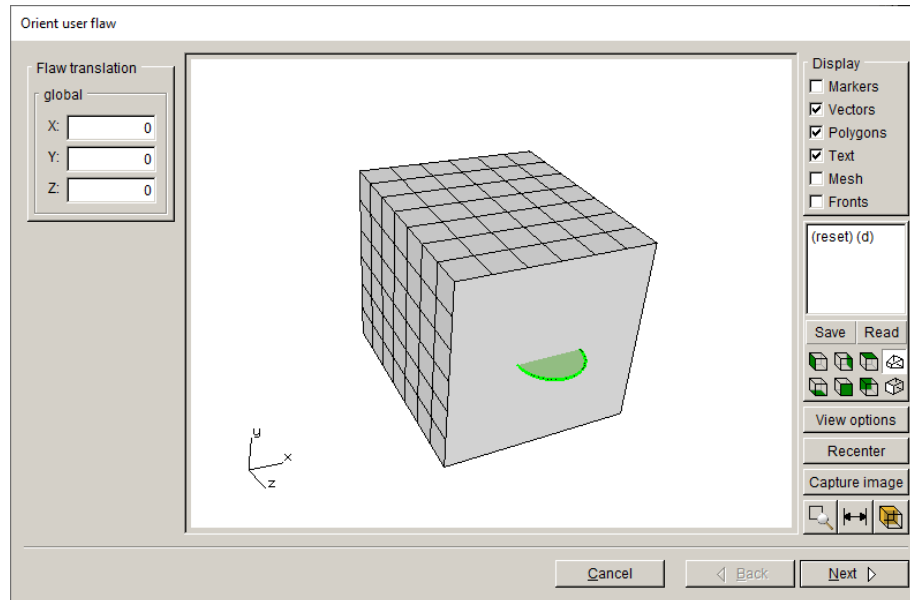


Figure 2.31 Flaw wizard panel to set location.

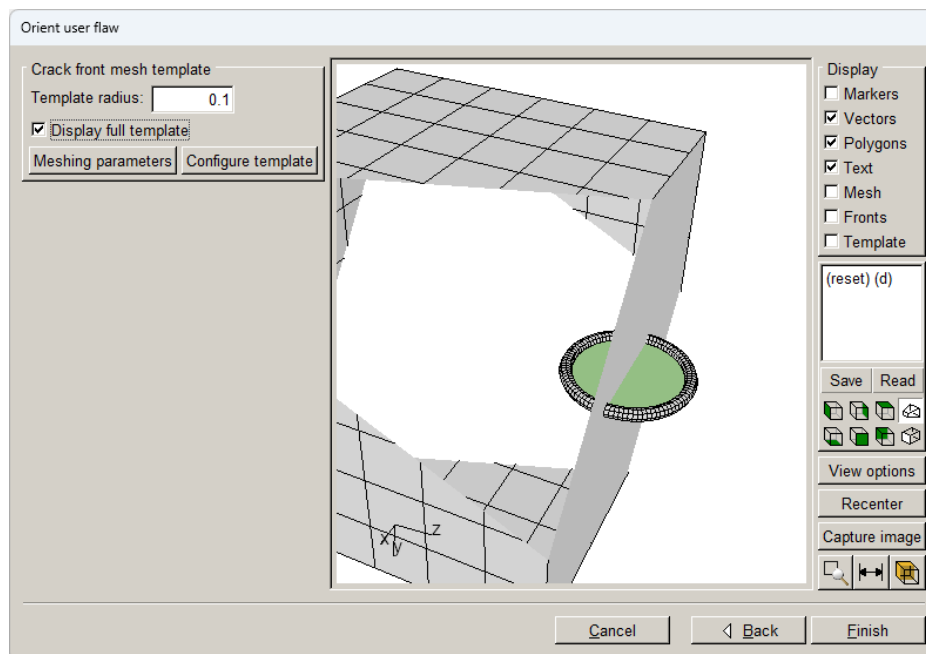


Figure 2.32 Flaw geometry shown in the model with crack template mesh.

2.5 Step 5: Static Crack Analysis

Once the crack is inserted and the model is remeshed, we perform a stress analysis using ANSYS, which provides the displacement results that are needed to compute stress intensity factors (SIFs). Typically, you would run a “static crack analysis” of the initial crack prior to running automated crack growth; this allows you to verify that the crack model is consistent with the uncracked model. You can compare displacement contours using the ANSYS post-processor, for the original uncracked and cracked models to verify that the crack model is valid.

You might also insert the crack with a bigger or smaller template radius to verify that the computed SIFs are accurate; there should be no notable change in SIFs for a reasonable template.

Step 5.1: Select Static Crack Analysis

From the FRANC3D menu, select **Analysis** → **Static Crack Analysis**, Fig 2.33. The first panel, Fig 2.34, requests that you enter the file name for the FRANC3D data. We call it *Ansys_Cube_Subdivide.fdb*; select **Next** once you enter a File Name.

Note that you cannot use the initial uncracked *Ansys_Cube.cdb* or *Ansys_Cube_LOCAL.cdb* file names, because a new *.cdb* file will be created during the analysis. The original uncracked *.cdb* files are reused for each step of crack growth.

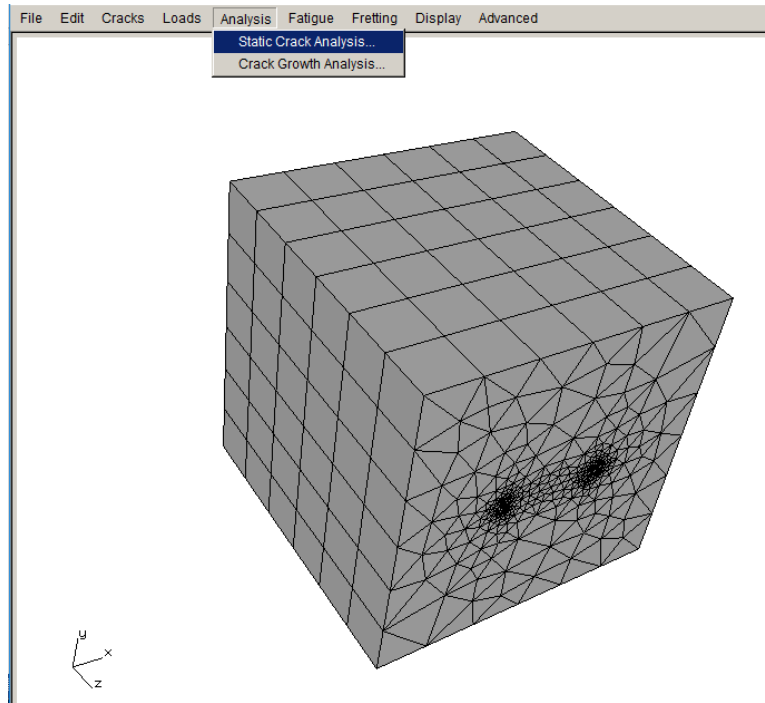


Figure 2.33 Static Crack Analysis menu

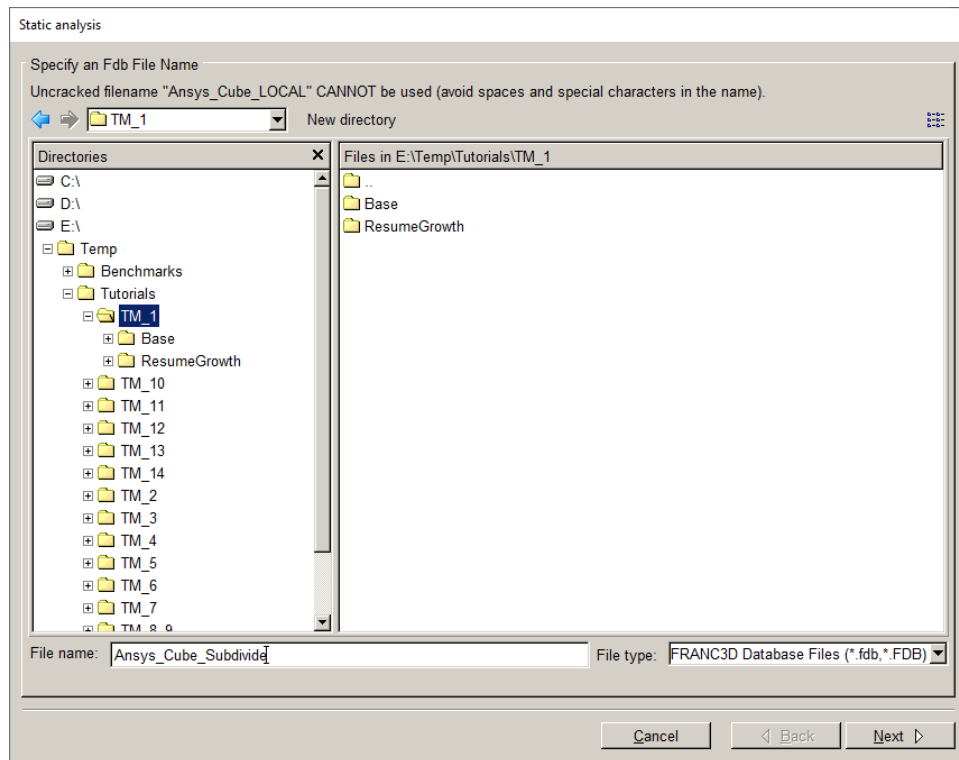


Figure 2.34 Static Analysis wizard first panel – File Name

Step 5.2: Select ANSYS Analysis Options

The next panel of the wizard, Fig 2.35, allows you to specify the ANSYS output and analysis options. The model is a sub-model and must be combined with the global model, so the Connect to global model box is checked automatically if the FRANC3D submodeler tool was used. The next panel, Fig 2.36, allows you to set the local + global model connection.

Note that if the model being analyzed is a full/complete model, the panel in Fig 2.36 is the final panel as there is no global model, and the **Next** button will be a **Finish** button.

Fig 2.36 shows the options for joining the local and global models. AUTO_CUT_SURF and GLOBAL_CONNECT_SURF are the sets/surfaces created by the FRANC3D submodeling tool and should be selected automatically. Click **Finish** to start ANSYS running (in batch/background mode). FRANC3D writes files and then attempts to execute ANSYS based on the ANSYS Executable information (Fig 2.35). You can monitor the FRANC3D CMD window and the ANSYS *.err*, *.out* and *.log* files for errors or messages.

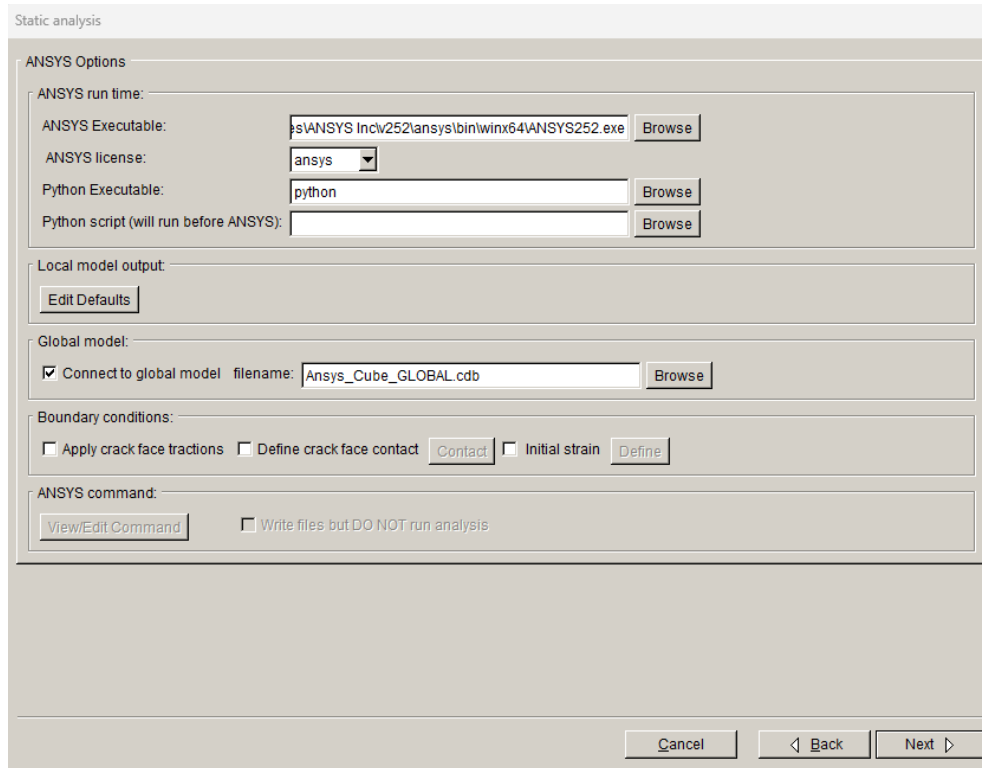


Figure 2.35 Static Analysis wizard third panel – ANSYS output options

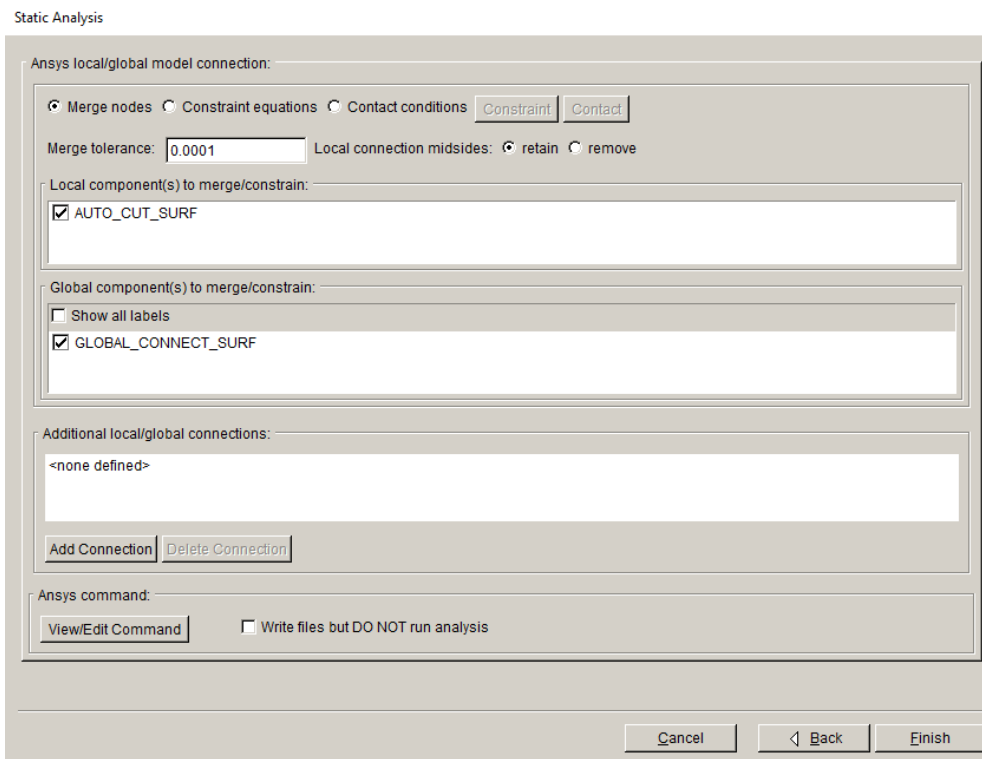


Figure 2.36 Local/Global model connection

Note that if ANSYS does not run, you will see a message in FRANC3D indicating that SIFs cannot be computed because there are no displacements, or you might see a message indicating that FRANC3D cannot find or open the *.dtp* file. This might be due to an incorrect path to the ANSYS executable or an incorrect license string. The ANSYS executable (and the path to it) will depend on the ANSYS installation, ANSYS version, operating system, *etc.* For MS Windows, for a typical installation of ANSYS 2024 R2, the executable will be “ANSYS242.exe” in the folder “C:\Program Files\ANSYS Inc\v242\ansys\bin\winx64”. The default ANSYS license string is “ansys”. You can determine your ANSYS license string by using the Mechanical APDL Product Launcher. Using the top menu bar of the Product Launcher, click on Tools – Display Command Line, and then look for “-p” in the window that is displayed; the license string will follow the -p.

Choosing Write files but DO NOT run analysis in Fig 2.36 will create all the files without running ANSYS, if the analysis needs to be run later or on a different computer.

If you need to run analyses on a different computer, you must bring the results file (*.dtp*) back to the working folder, and then you can use the **File** → **Read Results** menu option to import the results into FRANC3D. The *.dtp* file is created by ANSYS APDL commands, which are part of the *.cdb* file that FRANC3D writes.

2.6 Step 6: Compute Stress Intensity Factors

Step 6.1: Select Compute SIFs

When the ANSYS analysis is done (and the *.dtp* file read back into FRANC3D), we can compute SIFs. From the FRANC3D menu, select **Cracks** → **Compute SIFs**, Fig 2.37. The Stress Intensity Factor dialog is displayed, Fig 2.38. You should use the M-integral, but you can check that the Displacement Correlation results are similar. Select **Accept** and the SIFs Plot window is displayed, Fig 2.39. You can view the three modes of stress intensity factor (SIFs) modes using the upper tabs: KI, KII and KIII; mode II and III values will be negligible compared to mode I.

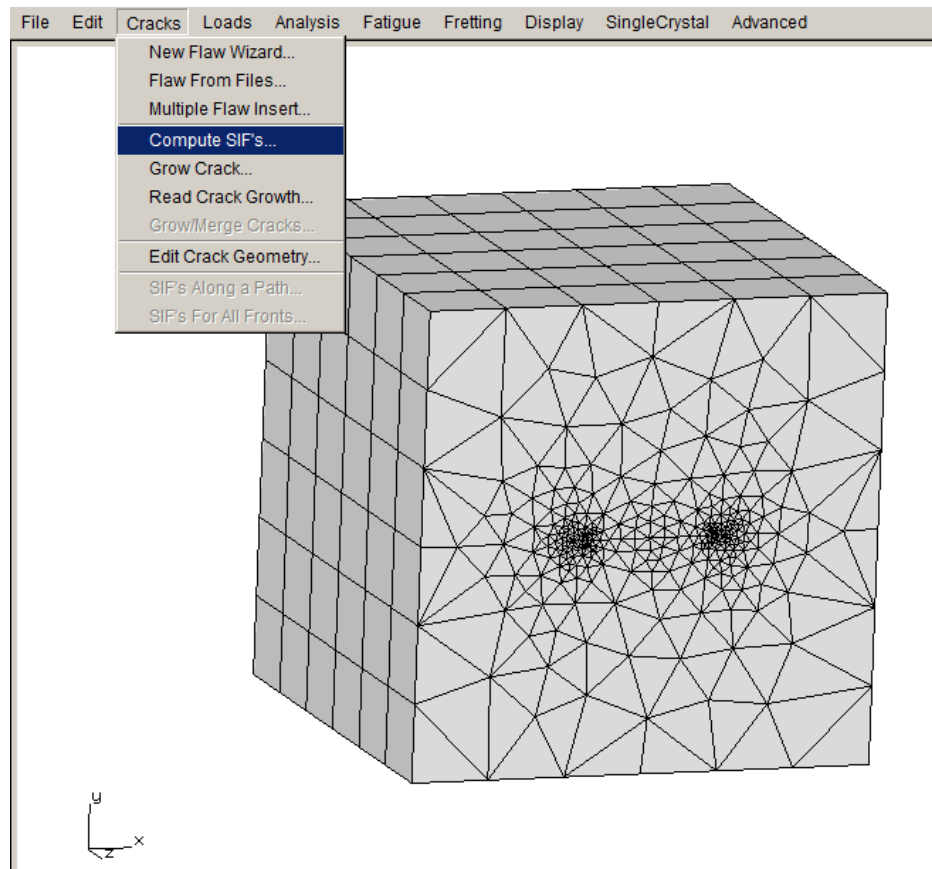


Figure 2.37 Compute SIFs selected from the Cracks menu.

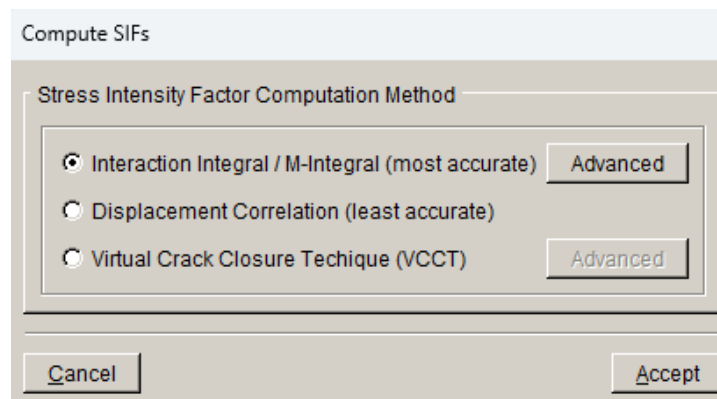


Figure 2.38 Compute SIFs dialog

The units for SIFs are $\text{MPa}\sqrt{\text{mm}}$. The SIFs are plotted versus the normalized crack front length, starting from point **A** – see the image in the left pane of Fig 2.39. SIFs are computed at element midside nodes along the front, so the values do not start at exactly zero.

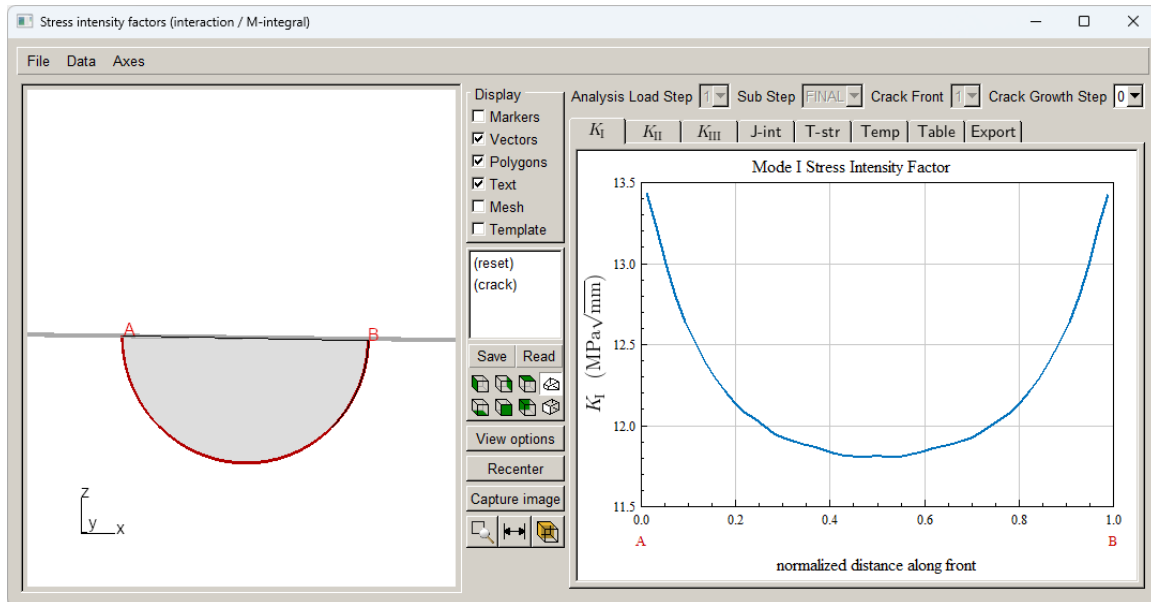


Figure 2.39 Stress Intensity Factor plot

2.7 Step 7: Manual Crack Growth

We will “manually” propagate the crack front; this lets us examine the predicted crack growth to determine suitable parameters for fitting and extrapolation before proceeding with automated crack growth, which is described in Section 2.8.

Step 7.1: Select Grow Crack

From the FRANC3D menu, select **Cracks** → **Grow Crack**, Fig 2.40. In the first panel, Fig 2.41, switch to Quasi-Static crack growth and select **Next**. Use the Max Tensile Stress theory to determine the kink angle, Fig 2.42; in this model, crack growth is mostly planar. Select **Next**. Use a Power Law relationship to determine relative growth among points along the crack front, Fig 2.43; we will use the default power $n=2$. There is only one FEM load step in the model so

select it. Select **Finish** to proceed. You will be asked if you want to save the growth parameters to a file, Fig 2.44, you can select **No**.

This quasi-static growth model can be considered a simplified Paris model; it is simpler to use than the subcritical growth model. Consult the Reference document for descriptions of crack extension models.

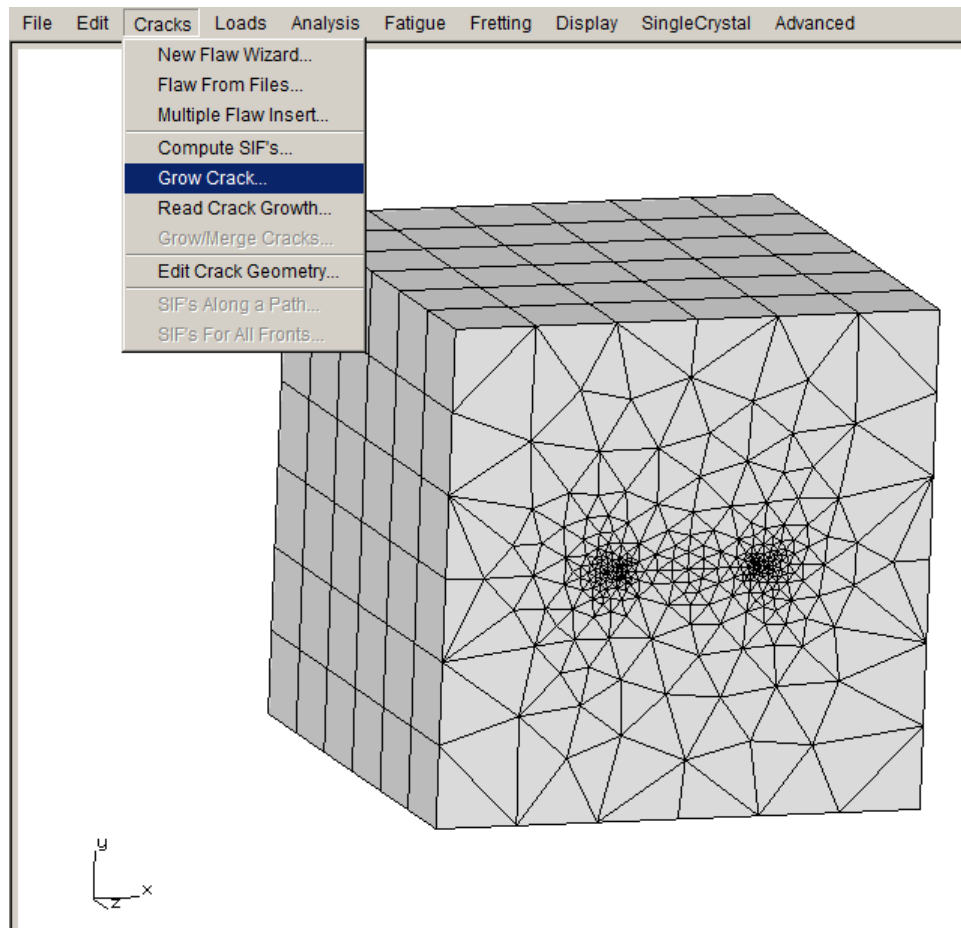


Figure 2.40 Grow Crack menu option.

Growth Type

☐ Subcritical crack growth (fatigue, SCC, creep, etc.)
☒ Quasi-static crack growth
☐ User defined crack growth

Read growth parameters from file:

Figure 2.41 Crack Growth wizard – first panel.

Kink angle model

Kink angle model

☒ Max tensile stress (MTS) $\max(K_I^r(\theta))$
☐ Max shear stress (MSS) $\max\left(\sqrt{(\eta_{II}K_{II}^r(\theta))^2 + (\eta_{III}K_{III}^r(\theta))^2}\right)$
☐ Generalized stress $\max(\text{MTS}, \text{MSS})$
☐ Strain energy release rate $\max(K_I^r(\theta)^2 + (\eta_{II}K_{II}^r(\theta))^2 + (\eta_{III}K_{III}^r(\theta))^2)$
☐ Planar $\theta = 0$
☐ User defined model
☐ Kink angle limit Maximum kink angle (deg):

Non-proportional kink angle parameter: ☒ ΔK ☐ $K_{\max}$

Mixed mode eta factors: η_{II} : η_{III} :

Crack growth resistance

☐ Anisotropic toughness

Figure 2.42 Crack Growth wizard – second panel using quasi-static growth.

Quasi-static growth parameters

Power law growth parameter

$$\Delta a_i = \Delta a_{\text{median}} (K_i / K_{\text{median}})^n$$

n:

Mixed-mode equivalent K

☒ $K^{\text{equiv}} = K_I$ ☐ $K^{\text{equiv}} = \sqrt{K_I^2 + (\gamma_{mII} K_{II})^2 + (\gamma_{mIII} K_{III})^2}$

γ_{II} : γ_{III} :

sign: ☒ from K_I ☐ from K_{II} ☐ from K_{III} ☐ always positive ☐ always negative

FEM Load Steps

Step	Sub	Load mult
1	1	1 Edit

Figure 2.43 Crack Growth wizard – third panel using quasi-static growth.

Save Parameters

? Save growth parameters to a file?

Figure 2.44 Save growth parameters dialog.

Step 7.2: Specify Fitting and Extrapolation

The next panel, Fig 2.45, allows you to specify the crack front point fitting parameters. You can double click on the (crack) view to see the crack surface; it should be the default view.

The median extension is set to 0.15; this value is set automatically based on the initial template radius so your value might be different. We change it to 0.12 and turn on Mesh and Markers (top right checkbox in Fig 2.45). The green boxes represent the computed new front points, and the blue line is the curve-fit, Fig 2.46. Note that the scale node is 0.5; the computed extensions are sorted from minimum to maximum, and the scale node can vary from 0 to 1. For a scale node

of 0.5, the median extension is set to the user-specified value, and all other extensions are scaled accordingly.

A Fixed order polynomial with order set to 3 and extrapolation set to 3% at both ends of the crack front is the default; the fitting parameters might need to be adjusted as the crack grows. A 4th order polynomial provides a slightly better fit in this case; the guidelines for curve-fitting are to use the simplest fit that is reasonable. The extrapolation is automatically adjusted to make sure the end points of the fitted-front fall outside the model surface. We stick with the default settings and select **Next** to proceed.

Note that the fitted (blue) curve must be extrapolated so that both ends fall outside of the model, but it should not be extrapolated too much. Normally, FRANC3D will automatically find reasonable values for extrapolation if Allow fit adjustment is turned on.

The user-specified median extension must be enough to provide finite space between the current and new fronts along the entire front to define new crack geometry. It is best to specify the extension (and have cycles computed) as opposed to specifying the cycles and having extension computed.

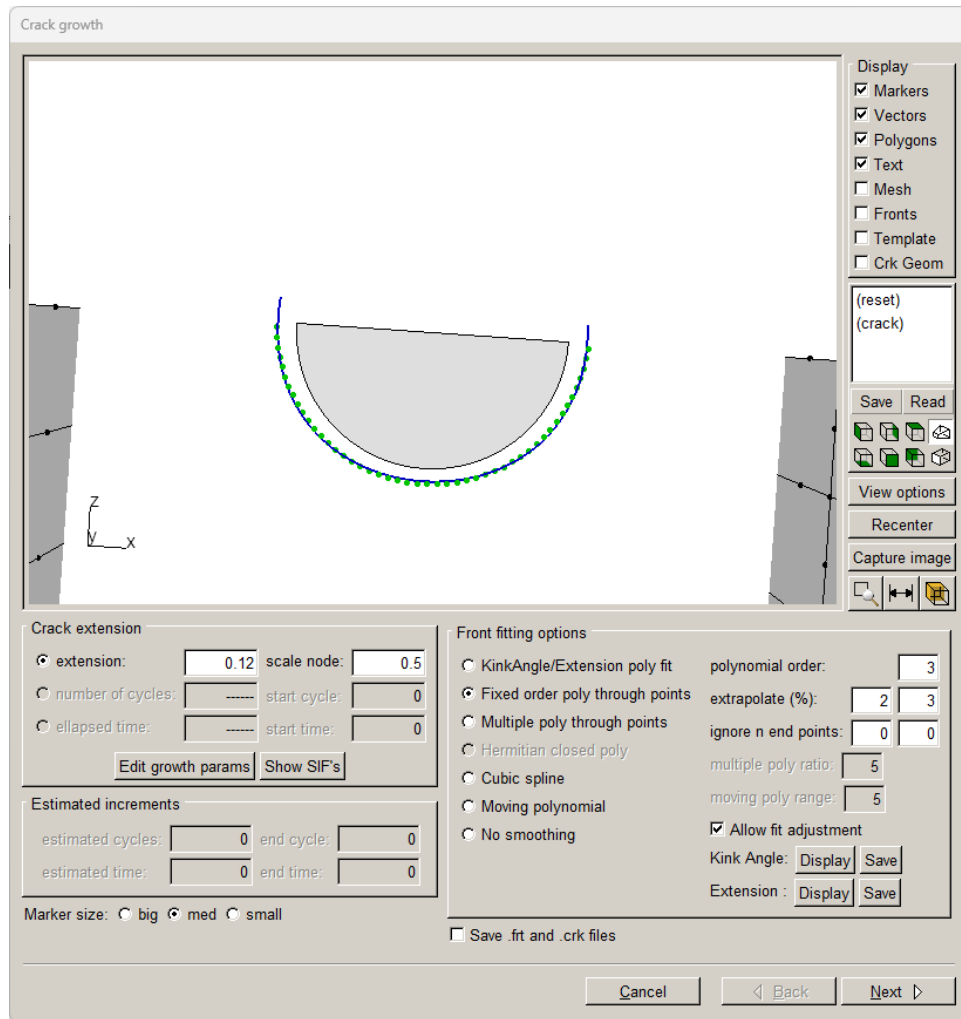


Figure 2.45 Crack growth wizard panel for crack front fitting options

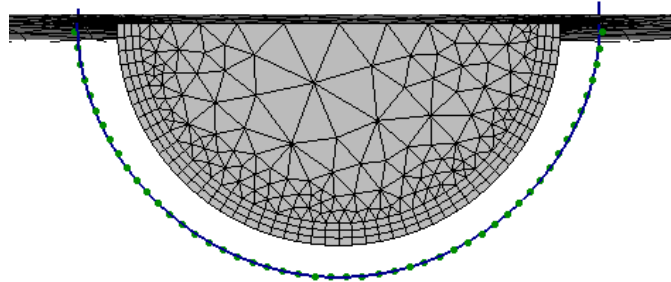


Figure 2.46 Curve-fit through front points.

Step 7.3: Specify Crack Front Template

The final panel, Fig 2.47, allows you to specify the crack front mesh template parameters. The crack front mesh template extends beyond the model surface, which corresponds with the front-fit extrapolation. The template radius can be set as an absolute value or as a % of the median extension; use an absolute value of 0.1, which is the same as the initial crack template radius. Select **Next** when ready to proceed with crack insertion and remeshing.

Note that the crack geometry for the extended crack is inserted into the original uncracked model, so do not overwrite, or remove the original uncracked model files.

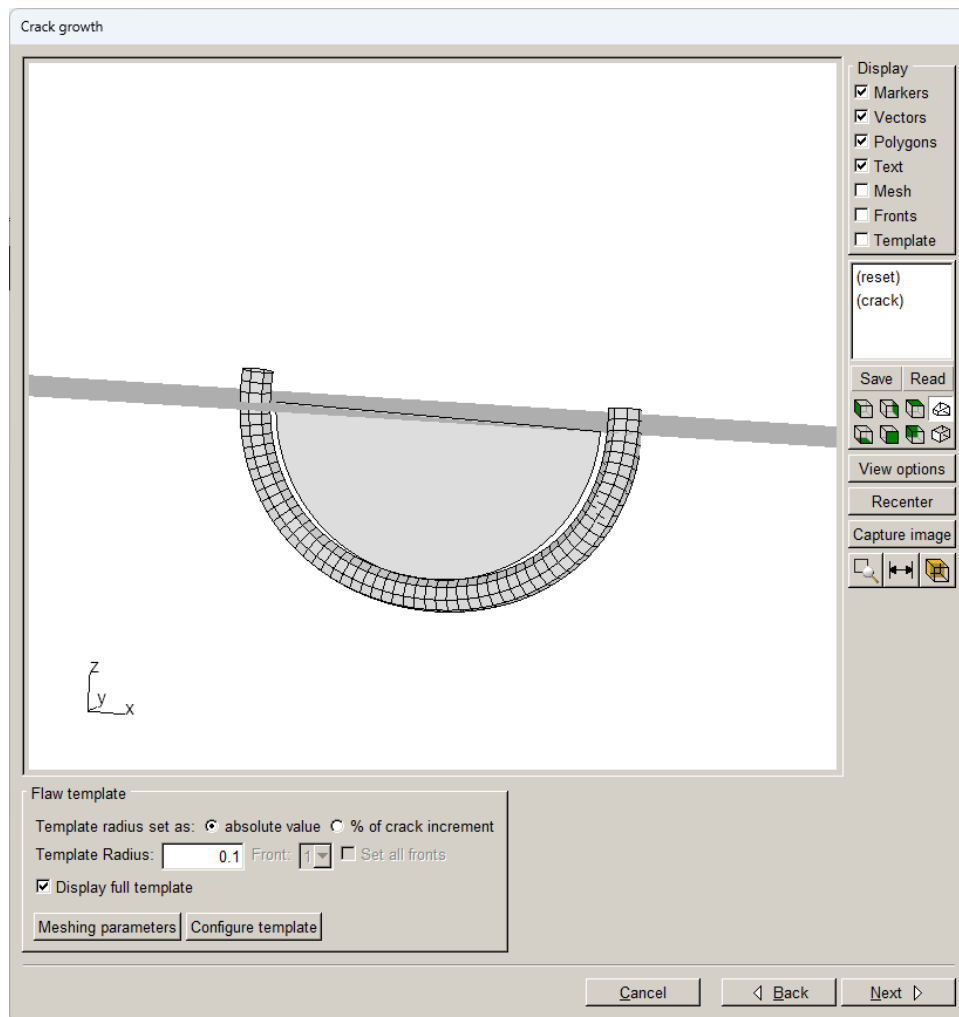


Figure 2.47 Crack growth wizard panel for mesh template options

A template radius that is slightly less than the crack extension, Fig 2.48, works best for cases where there is kinking or turning of the crack front.

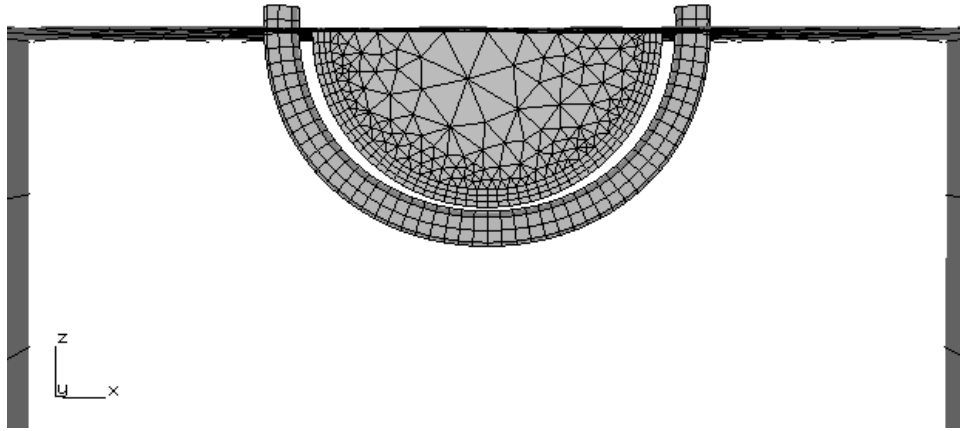


Figure 2.48 Template radius is smaller than extension.

The resulting model can be analyzed as was done for the initial crack (see Step 5 above). Note that you will want to give this model a different name, such as *Ansys_Cube_step_001*, so that you do not overwrite the initial crack files. You could continue with another manual step of growth followed by static analysis, or you can try automated crack growth, which is described next.

2.8 Step 8: Automatic Crack Growth

This section describes the steps for automatic crack growth starting from the initial crack model. We start with the existing FRANC3D model, using the model created in Section 2.4 and 2.5. In most cases, we would simply do automatic growth after the manual step described in Section 2.7, but this will not show all the dialogs, so we restart from the initial crack model.

Step 8.1: Open FRANC3D Restart File

Start with the FRANC3D graphical user interface (see Fig 1.1) and select **File** → **Open**. Select the file name specified in Section 2.5, called *Ansys_Cube_Subdivide.fdb*. Select **Accept**. The

model will be read into FRANC3D (along with the results that were created when running the static analysis). We will ignore the fact that we already propagated the initial crack in the previous step (but we will use the settings). The next section describes the steps for automatic crack growth analysis.

Step 8.2: Select Crack Growth Analysis

From the FRANC3D menu, select **Analysis** → **Crack Growth Analysis**, Fig 2.49. The first panel of the wizard, Fig 2.50, allows you to choose the method for computing SIFs. We use the default M-integral and select **Next**. If this panel does not show up, the SIF method is chosen automatically based on your earlier choice when plotting SIFs.

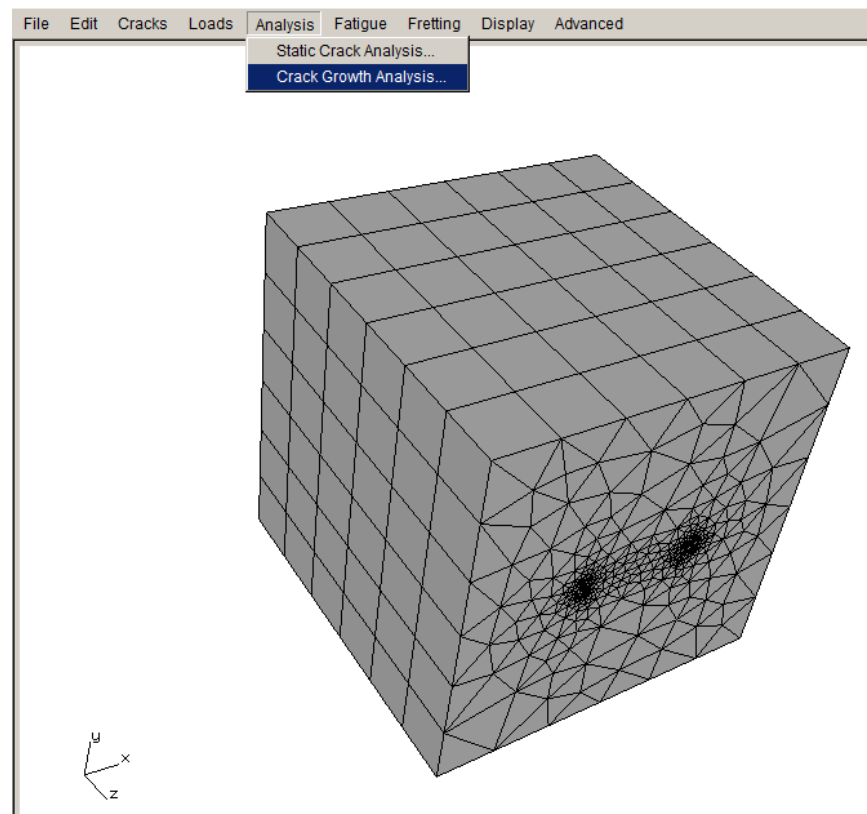


Figure 2.49 Crack Growth Analysis menu

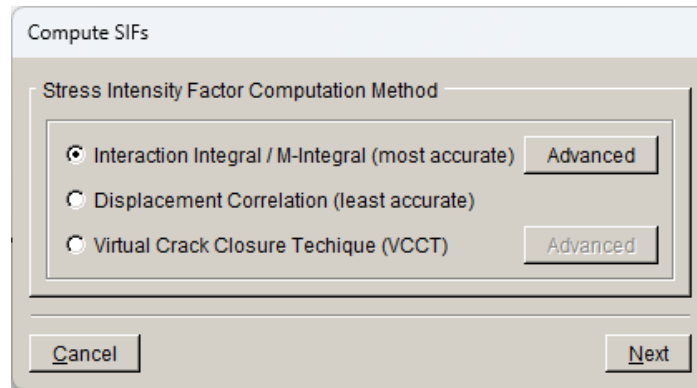


Figure 2.50 Crack Growth Analysis wizard – first panel.

Step 8.3: Specify Growth Rules

The second panel, Fig 2.51, lets us set the crack growth type; we use quasi-static growth. Select **Next** to display the third panel, Fig 2.52; use the default MTS theory to compute the kink angle. Select **Next** to display the fourth panel, Fig 2.53; use the default settings for the quasi-static growth; there is only one FEM load step to select. Select **Next**. You will be asked if you want to save the growth parameters to a file; you can select **No**.

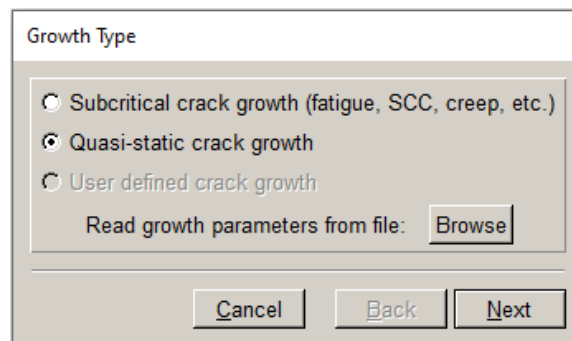


Figure 2.51 Crack Growth Analysis wizard – second panel.

Kink angle model

Kink angle model

☒ Max tensile stress (MTS) $\max(K_I^r(\theta))$

☐ Max shear stress (MSS) $\max\left(\sqrt{(\eta_{II}K_{II}^r(\theta))^2 + (\eta_{III}K_{III}^r(\theta))^2}\right)$

☐ Generalized stress $\max(\text{MTS}, \text{MSS})$

☐ Strain energy release rate $\max\left(K_I^r(\theta)^2 + (\eta_{II}K_{II}^r(\theta))^2 + (\eta_{III}K_{III}^r(\theta))^2\right)$

☐ Planar $\theta = 0$

☐ User defined model

☐ Kink angle limit Maximum kink angle (deg):

Non-proportional kink angle parameter ☒ ΔK ☐ $K_{\max}$

Mixed mode eta factors η_{II} : η_{III} :

Crack growth resistance

☐ Anisotropic toughness

Figure 2.52 Crack Growth Analysis wizard – third panel.

Quasi-static growth parameters

Power law growth parameter

$\Delta a_i = \Delta a_{\text{median}} (K_i / K_{\text{median}})^n$ n:

Mixed-mode equivalent K

☒ $K^{\text{equiv}} = K_I$ ☐ $K^{\text{equiv}} = \sqrt{K_I^2 + (\gamma_{II}K_{II})^2 + (\gamma_{III}K_{III})^2}$

γ_{II} : γ_{III} :

sign: ☒ from K_I ☐ from K_{II} ☐ from K_{III} ☐ always positive ☐ always negative

FEM Load Steps

Step	Sub	Load mult
1	—	1 Edit

Figure 2.53 Crack Growth Analysis wizard – fourth panel.

Step 8.4: Specify Fitting and Template Parameters

The fifth panel, Fig 2.54, lets us set the crack front fitting and mesh template parameters. Use the default 3rd order polynomial and 2% extrapolation. Set the template radius to an absolute value of 0.1. Select **Next**.

Fitting & template parameters

Front fitting options

☐ KinkAngle/Extension poly fit polynomial order:

☒ Fixed order poly through points extrapolate (%):

☐ Multiple poly through points ignore n end points:

☐ Hermitian closed poly multiple poly ratio:

☐ Cubic spline moving poly range:

☐ Moving polynomial ☒ Allow fit adjustment

☐ No smoothing

Flaw template

Template radius set as: ☒ absolute value ☐ % of crack increment

Template Radius: Front: ☐ Set all fronts

Extension

☒ distance extension scale node:

☐ cycles extension

☐ time extension

Figure 2.54 Crack Growth Analysis wizard – fifth panel.

Step 8.5: Specify Growth Plan

The sixth panel, Fig 2.55, lets us define the amount of extension for each step of crack growth. Set the number of steps to five and the constant median extension to 0.2. Select **Next**.

Note that we are taking a fairly large crack growth step; this is to limit the number of steps and time for this tutorial. Large steps can sometimes lead to issues with oscillations in the shape of the crack fronts (see Section 9 of the User's Guide), but for this model 0.2 is acceptable.

Growth plan

Specified crack growth steps

Number of Crack Growth Steps =

☒ Constant crack growth increments

Crack growth step size =

☐ Linearly increasing/decreasing crack growth increments

First crack growth step size =

Change in step size at each step =

☐ User defined crack growth increments

Figure 2.55 Crack Growth Analysis wizard – sixth panel.

Step 8.6: Analysis Code Setting

The seventh panel, Fig 2.56, shows the Base filename and analysis settings. The Current crack growth step defaults to 0 – representing the initial crack. If we “manually” propagated the crack first, we should set this to 1. FRANC3D will extend the initial crack based on the growth rule defined in the previous panels and then name the resulting set of files as *Ansys_Cube_Subdivide_STEP_001.**. Subsequent file names have the step number incremented as the automatic analysis proceeds.

Some options will be specific to your site. You should have already set the ANSYS executable path and license in the FRANC3D Preferences. The Global model settings should be set automatically. The local + global connection is specified in the next panel; select **Next**.

Figure 2.56 Crack Growth Analysis wizard – seventh panel.

The eighth panel, Fig 2.57, displays the local + global connection. The default settings use node-merging of the cut-surfaces between the local and global portions. Select **Finish** to start the automated crack growth process.

FRANC3D extends the initial crack first. It then saves the *.fdb*, *.cdb* and other files for the first crack step model with the name *Ansys_Cube_Subdivide_STEP_001*, and then ANSYS starts in the background. If the analysis stops at any stage, it can be restarted from the last crack step. All the required restart *_STEP_#* files are retained. The model for any step can be read into FRANC3D to view the SIFs or to restart the analysis with a modified crack growth increment (for example).

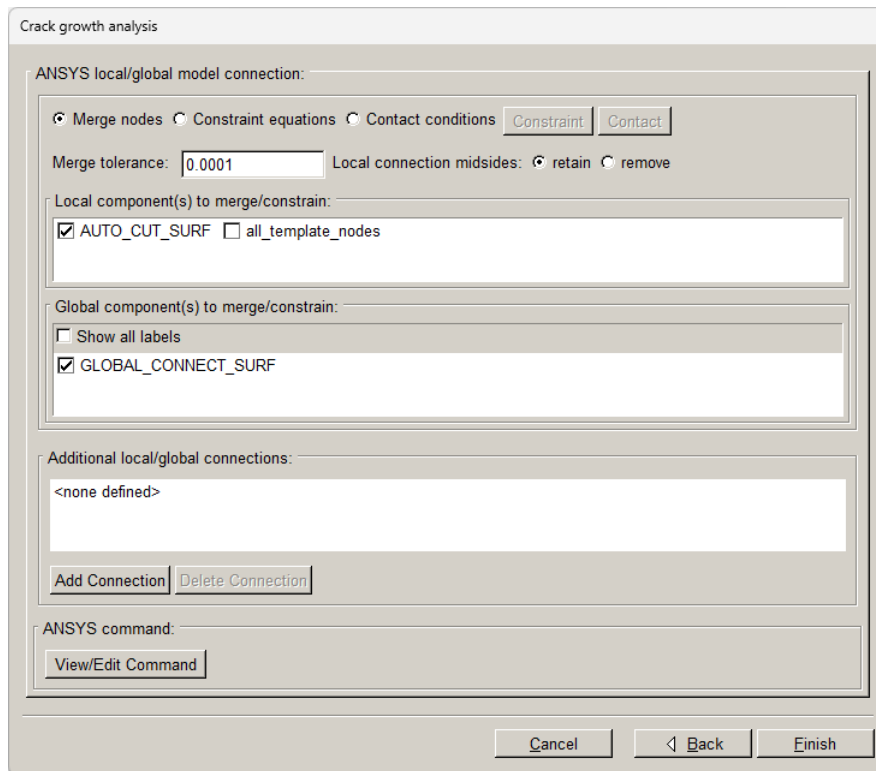


Figure 2.57 Local + Global connection options

A summary of the various files can be found in Section 3 of the User's Guide.

2.9 Step 9: SIF History and Fatigue Life

When the automatic crack growth analysis from Step 8 has finished, the SIF history can be displayed. If you still have FRANC3D open and the crack growth from Step 8 is done, you can proceed with Step 9.1. Otherwise, you can restart FRANC3D and read in the *.fdb* file for the last step, using the **File** → **Open** menu option.

Step 9.1: Select SIFs Along a Path

From the FRANC3D menu, select **Cracks** → **SIFs Along a Path**, Fig 2.58. If the SIFs have not been computed yet, the Compute SIFs dialog (see Fig 2.38) will be displayed; leave the defaults and select **Finish**.

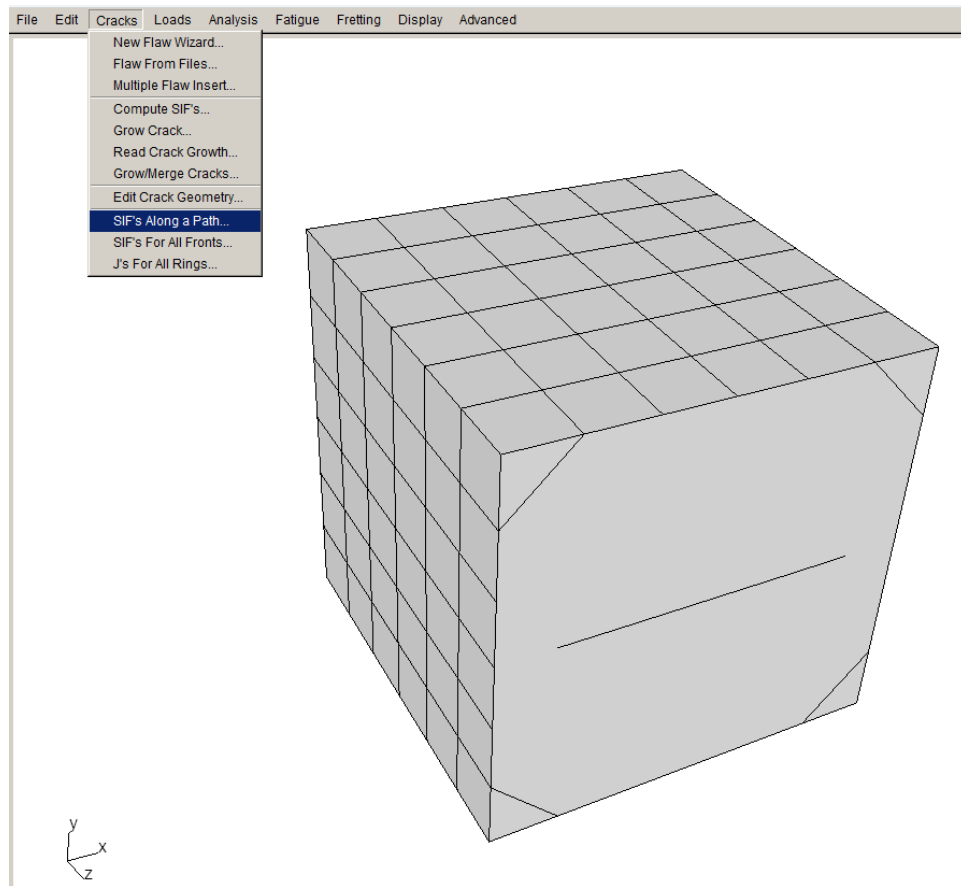


Figure 2.58 SIFs Along a Path menu option

The Stress Intensity Factors (along a path) dialog should appear, Fig 2.59. The crack fronts are displayed on the left along with the path through the fronts; the SIF history along the path is shown in the graph on the right. You can use the tabs above the graph to plot the Mode II and III SIFs as well as the elastic J-integral and T-stress values along the path. You can define a new path, Fig 2.60. You can export the data also; for example, you might need to export Mode I SIF history (K vs a) to compute fatigue cycles using a different program. The exported data files are ASCII txt files.

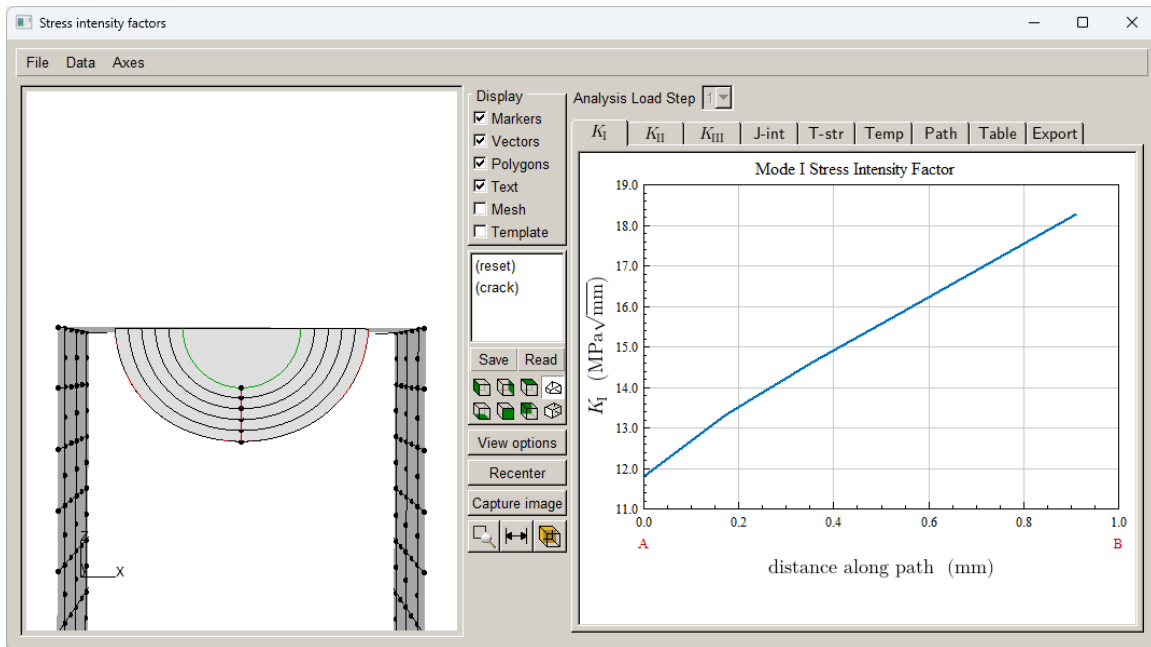


Figure 2.59 SIFs Along a Path dialog – KI plot

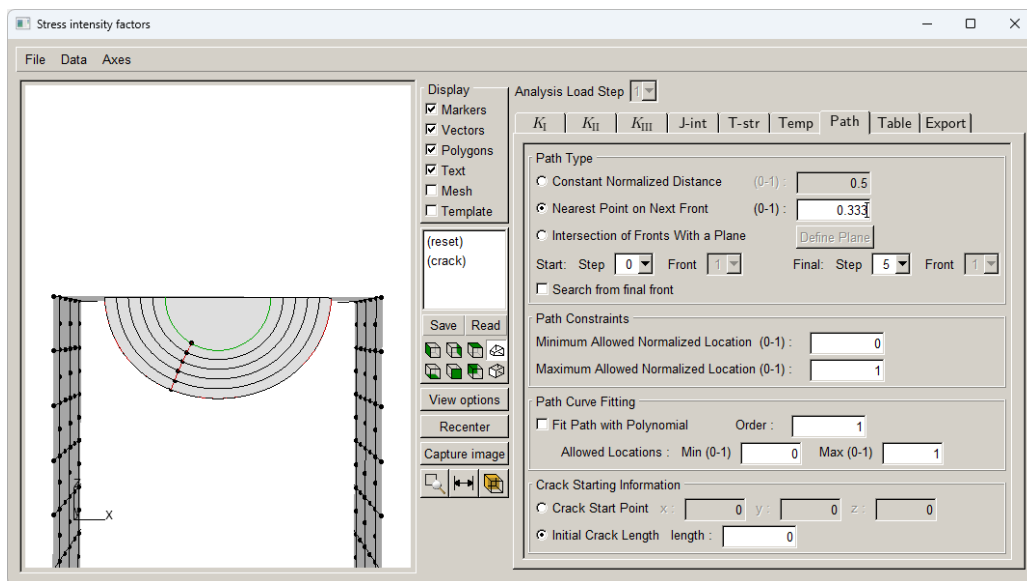


Figure 2.60 SIFs Along a Path dialog – Define Path

Step 9.2: Select SIFs For All Fronts

From the FRANC3D menu, select **Cracks** → **SIFs For All Fronts**. If the SIFs have not been computed yet, the Compute SIFs dialog will be displayed; leave the defaults and select **Finish**.

The Stress Intensity Factors (for all fronts) dialog should appear, Fig 2.61. The crack fronts are displayed on the left and the SIFs for all crack fronts are displayed on the right. You can use the tabs above the plot to display Mode II and III SIFs as well as the elastic J-integral and T-stress values. If there are multiple crack fronts or multiple load steps, these can be selected using the drop-down lists; this will be seen in subsequent tutorials.

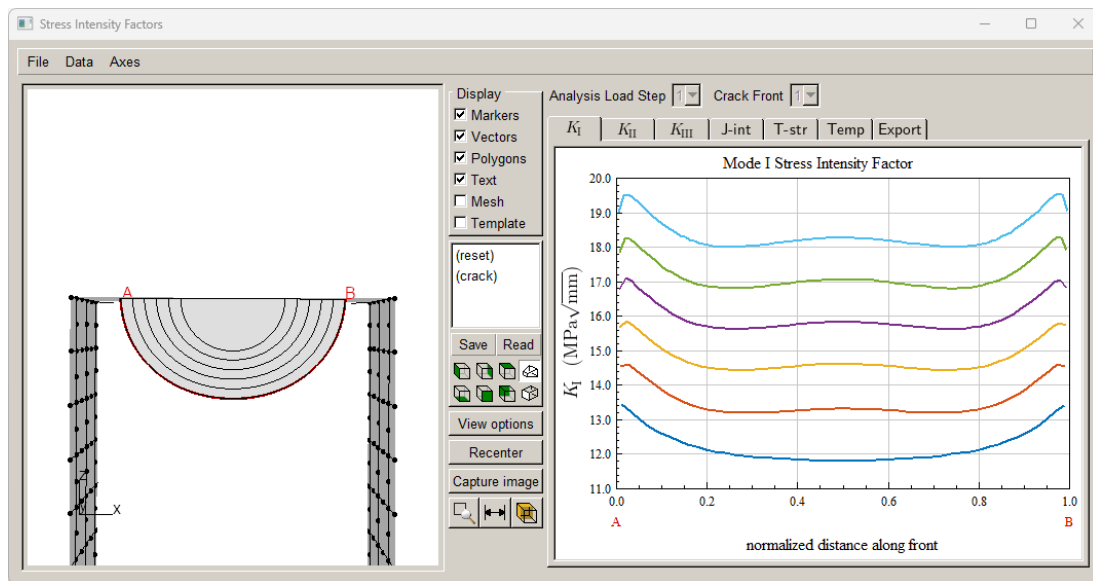


Figure 2.61 SIFs for all Fronts dialog – KI plot

Step 9.3: Select Fatigue Life Predictions

Fatigue life or cycles can be computed next. From the FRANC3D menu, select **Fatigue** → **Fatigue Life Predictions**, Fig 2.62. If the SIFs have not been computed yet, the Compute SIFs dialog will be displayed; leave the defaults and select **Finish**.

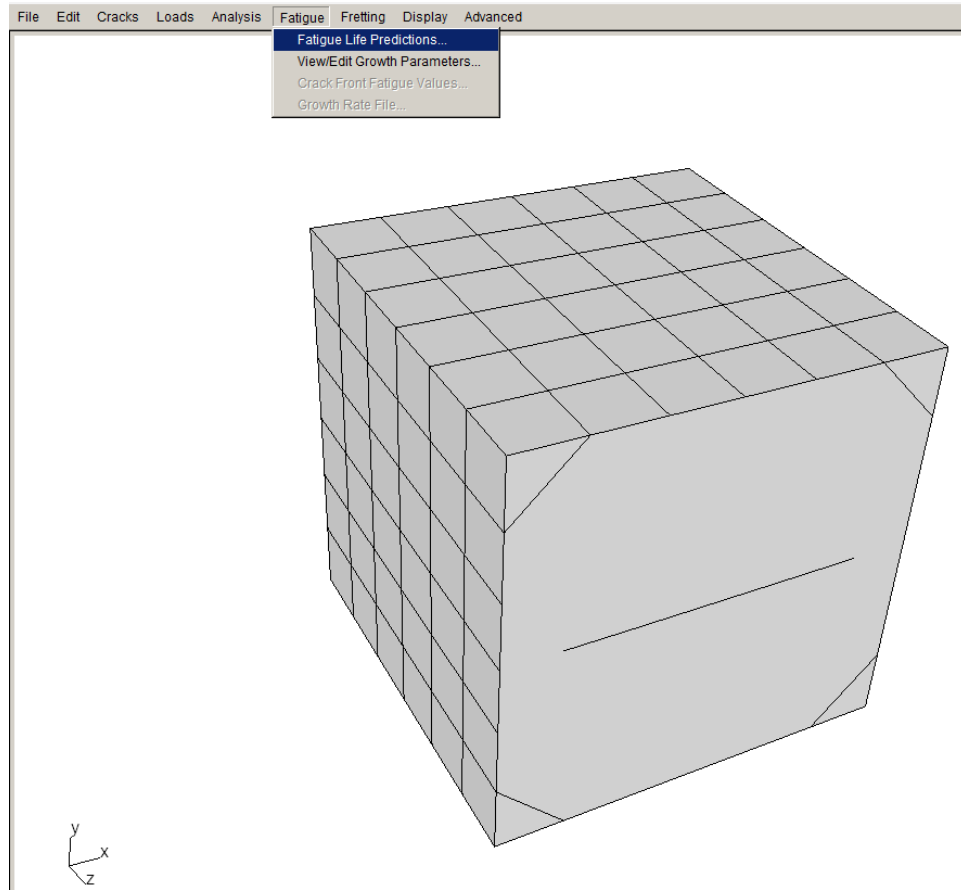


Figure 2.62 Fatigue Life Predictions menu option

The Fatigue Life dialog should appear as in Fig 2.63. The crack fronts are displayed on the left and the window on the right side should be blank (assuming that lifing parameters have not been defined previously). You must **Set (or Read) Parameters**. Selecting the **Set Parameters** button displays the dialog shown in Fig 2.64.

The units are MPa and mm; select the **Change** button to display the dialog in Fig 2.65 and set the units to MPa and mm; select **Accept**.

The FEM units can also be set in the main FRANC3D window using the **Edit** menu.

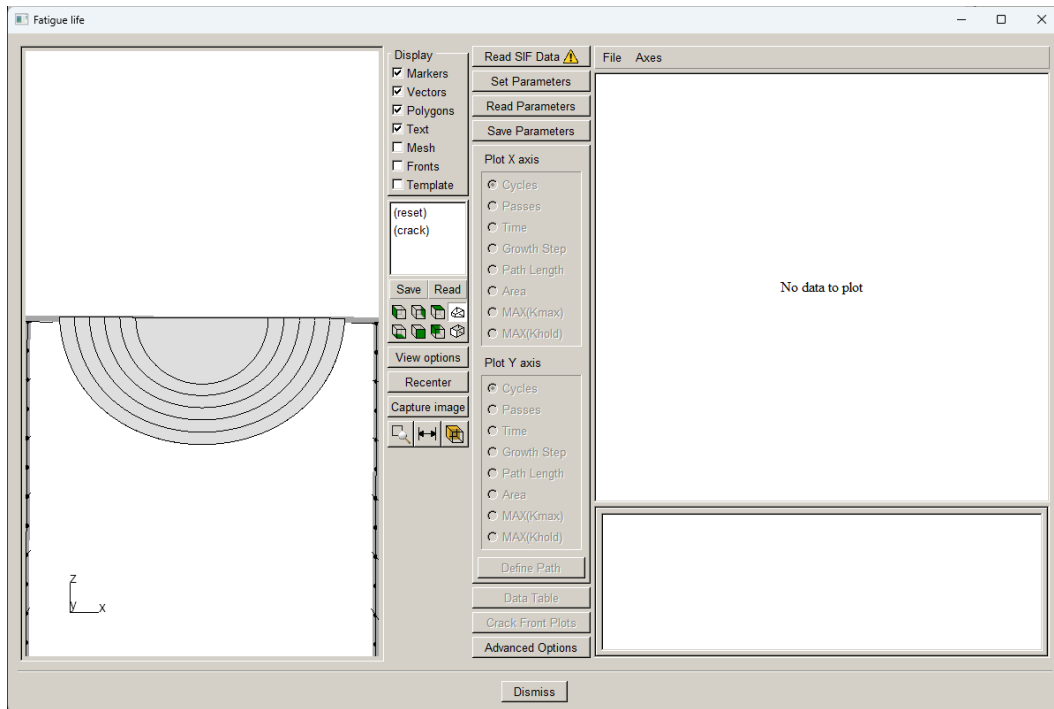


Figure 2.63 Fatigue Life dialog

Growth Parameters

Units used in FE model
 stress: MPa length: mm temp: C time: sec Change

Crack growth load schedule
New schedule Read from file Wizard View/Edit Save to file

Crack growth rate model
New model Read from file View/Edit da/dN View/Edit da/dt Save to file

Mixed-mode equivalent K
☒ $K^{\text{equiv}} = K_I$ ☒ $K^{\text{equiv}} = \sqrt{K_I^2 + (\gamma_{II} K_{II})^2 + (\gamma_{III} K_{III})^2}$
 γ_{II} : γ_{III} :
 sign: ☒ from K_I ☐ from K_{II} ☐ from K_{III} ☐ always positive ☐ always negative

Effective delta K
☒ $\Delta K_{\text{eff}} = K_{\text{max}} - K_{\text{min}}$ ☐ $\Delta K_{\text{eff}} = K_{\text{max}} - \max(K_{\text{min}}, 0)$
☐ Use load interaction model Set/Edit parameters

Integration options
☒ Accelerated counting ☐ Constant K for time integration

Dynamic pairing metric
☒ $\Delta K_{\text{eff}}^{\text{equiv}}$ ☐ da/dN ☐ Simplified HCF/LCF pairing

Cancel Accept

Figure 2.64 Fatigue Life parameters

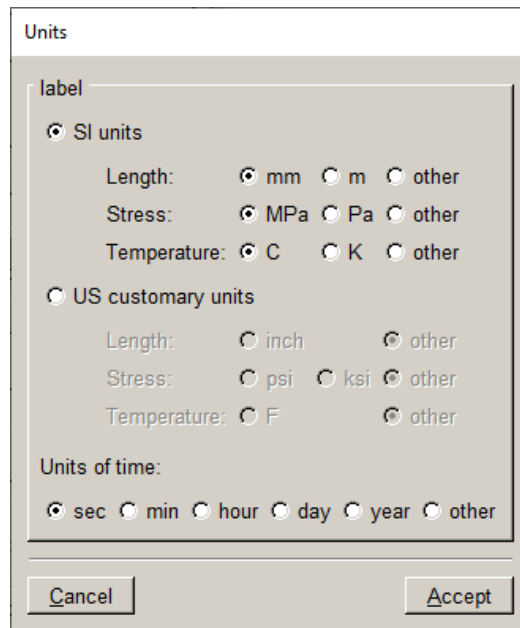


Figure 2.65 FEM units dialog

The crack growth load schedule is defined next. Select the **New Schedule** button (see Fig 2.64) to display the dialog in Fig 2.66. Select the Schedule and then select the **Add** button; see the right side image of Fig 2.66. This will display the dialog shown in Fig 2.67; we use a simple cyclic load schedule where the applied load produces K_{max} , and K_{min} is zero. Select **Accept** to return to the Load Schedule, Fig 2.68.

The stress ratio is set to 0.0. There is only one load case, which represents the K_{max} condition. The Repeat count is set to FOREVER. Select **Accept** to finish the load schedule. There is only one event in the schedule so the repeat forever can be set for either the event or the schedule; it is usually best to set Repeat FOREVER for the schedule.

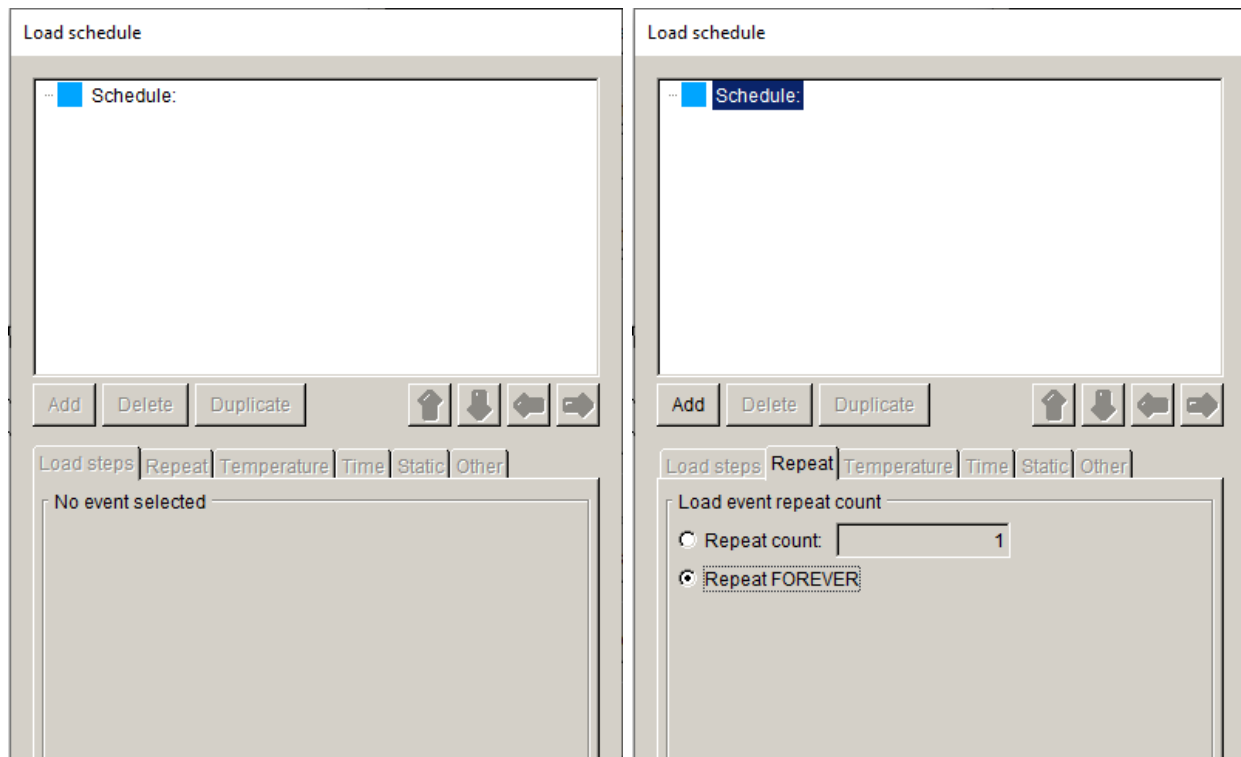


Figure 2.66 New Load Schedule dialog

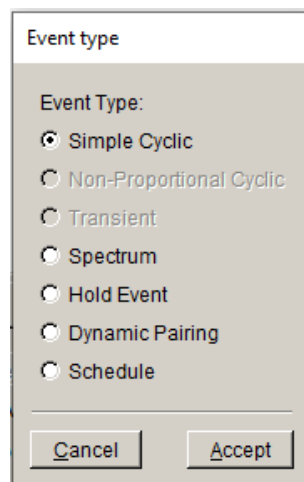


Figure 2.67 Event type dialog

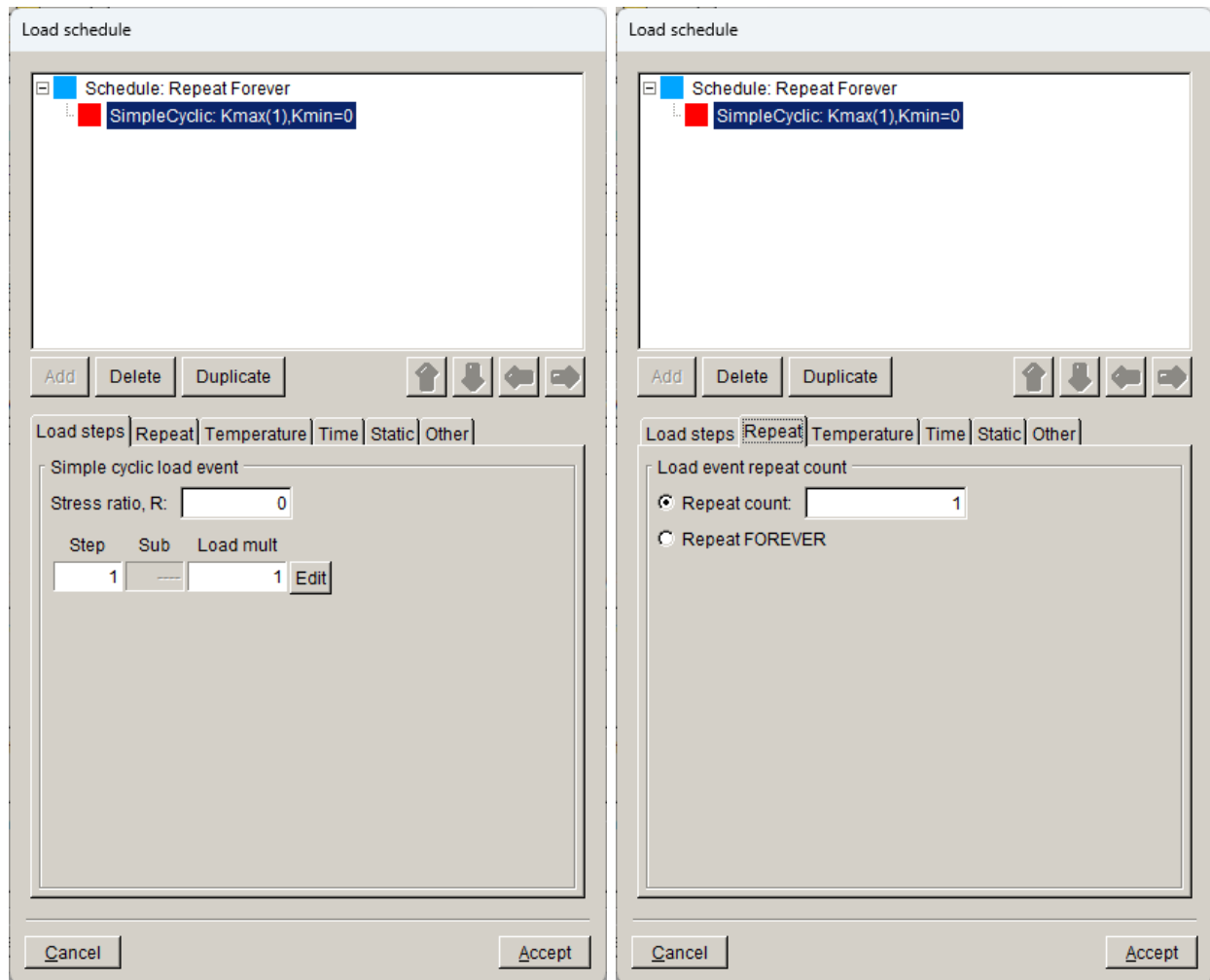


Figure 2.68 Simple cyclic load event repeated once and schedule repeated forever

The crack growth rate model is defined next. Select the **New Model** button (see Fig 2.64) to display the dialog in Fig 2.69. Choose the cyclic loading growth rate model, and select **Next** to display the dialog in Fig 2.70.

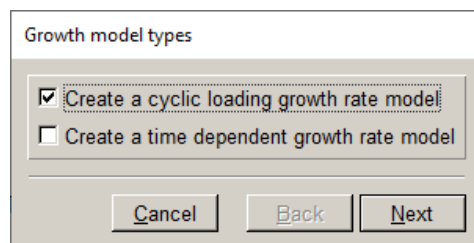


Figure 2.69 Growth model type dialog

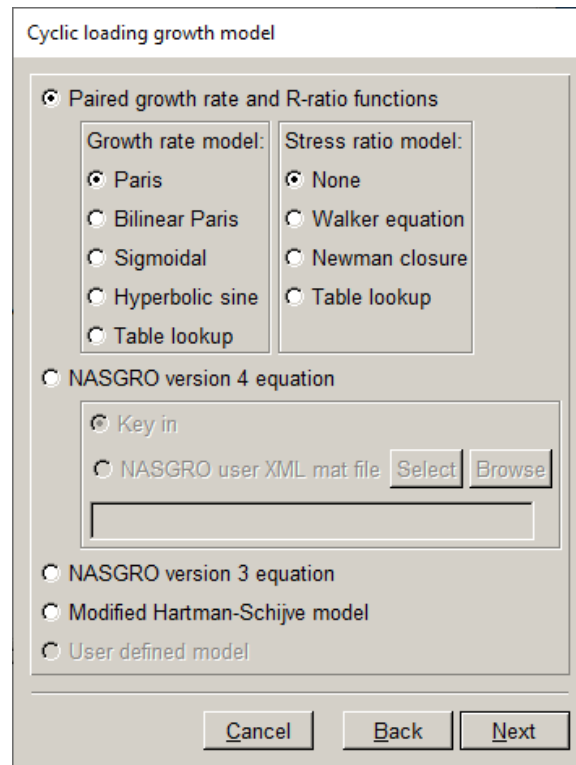


Figure 2.70 Cyclic loading growth model dialog

We use a Paris growth model, which is the default; select **Next** to display the dialog in Fig 2.71. The growth rate model will be temperature independent. Select **Next** to display the Paris growth model dialog, Fig 2.72.

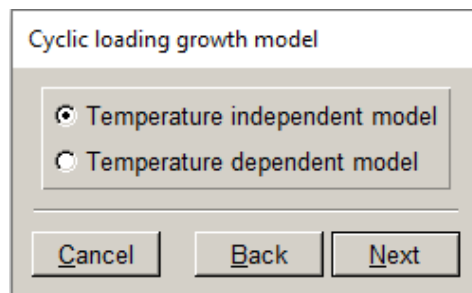


Figure 2.71 Temperature dependent or independent growth rate model dialog

Cyclic loading growth model

Paris growth model:
 $da / dN = C \Delta K^n$

Plot

Model label (optional):

Description (optional):

Units: stress: unset length: unset temp: unset time: unset  Change

	C	n	DKth	Kc
1				

Cancel Back Next

Figure 2.72 Paris growth rate model dialog

First, set the units for the Paris growth model. Select the **Change** button to display the Units dialog, Fig 2.73. Set the units the same as the FEM units that we set earlier, which were MPa and mm. If your growth model units are different from the FEM units, FRANC3D will do the unit conversions automatically. Select **Accept**.

Units

label

☒ SI units

Length: ☒ mm ☐ m ☐ other

Stress: ☒ MPa ☐ Pa ☐ other

Temperature: ☒ C ☐ K ☐ other

☐ US customary units

Length: ☐ inch ☐ other

Stress: ☐ psi ☐ ksi ☐ other

Temperature: ☐ F ☐ other

Units of time:

☒ sec ☐ min ☐ hour ☐ day ☐ year ☐ other

Cancel Accept

Figure 2.73 Paris growth rate model units

In the Paris growth model dialog, click on the fields for C, n, DKth and Kc and enter the appropriate values, Fig 2.74. Select **Next** to return to the main dialog (Fig 2.64). Other settings and options are left at their default values. Select **Accept** (Fig 2.64) to return to the Fatigue Life dialog, Fig 2.75.

The dialog box is titled "Growth rate model". It contains the following elements:

- Paris growth model:** $da / dN = C \Delta K^n$
- Plot** button
- Model label (optional):** A text input field.
- Description (optional):** A text input field with a scroll bar.
- Units:** A text input field showing "stress: MPa length: mm temp: C time: sec" and a **Change** button.
- Table:**

	C	n	DKth	Kc
1	1e-10	3	1	100
- Cancel** and **Accept** buttons at the bottom.

Figure 2.74 Paris growth rate model values

Note that the Paris exponent 'n' is set to 3 whereas we used $n=2$ for the quasi-static growth. In practice, you can use the "correct" fatigue model and material data during the growth, so that growth is consistent with the fatigue life computations. For a single crack front with relatively uniform K_I values along the front, the quasi-static growth option is acceptable. If there are multiple crack fronts with significantly different growth rates, the "correct" fatigue model should always be used.

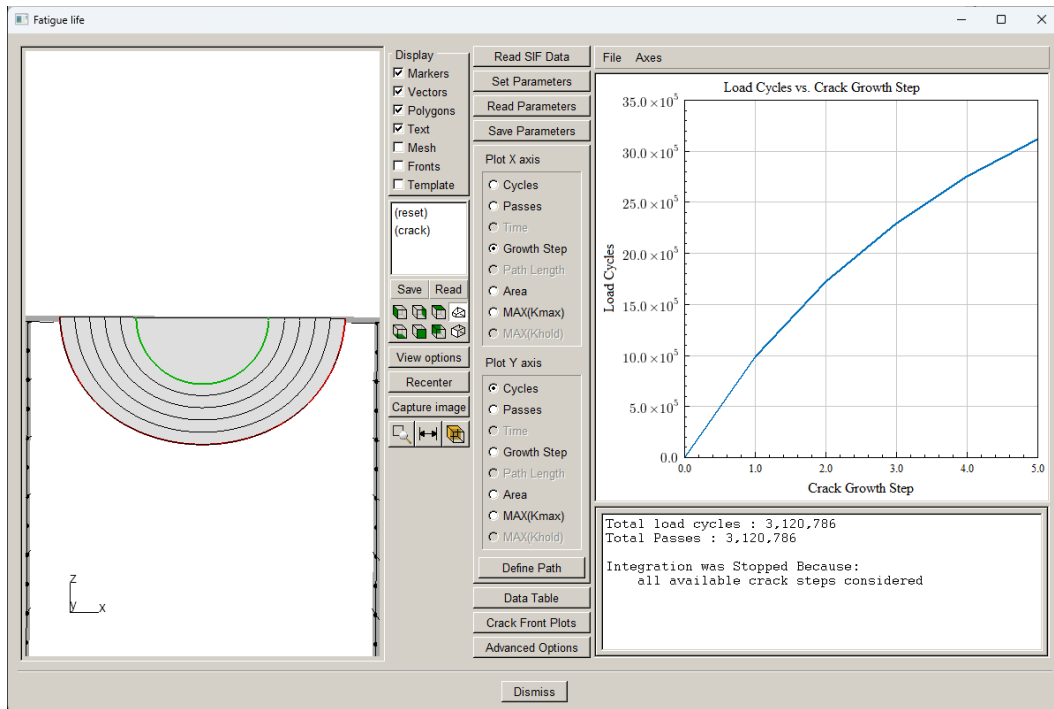


Figure 2.75 Fatigue life dialog showing Cycles vs Crack Growth Step

If we choose to plot cycles based on a path, we must select the Path radio button and then use the **Define** button to create a path, Fig 2.76. The default path is through the crack front midpoints. Once a path is defined, the cycles versus crack path length is displayed, Fig 2.77.

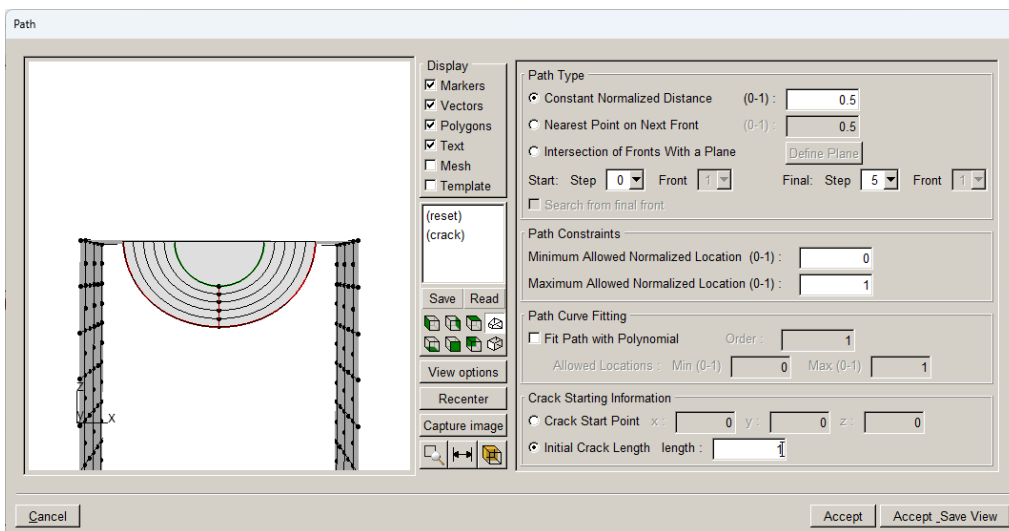


Figure 2.76 Define crack path dialog

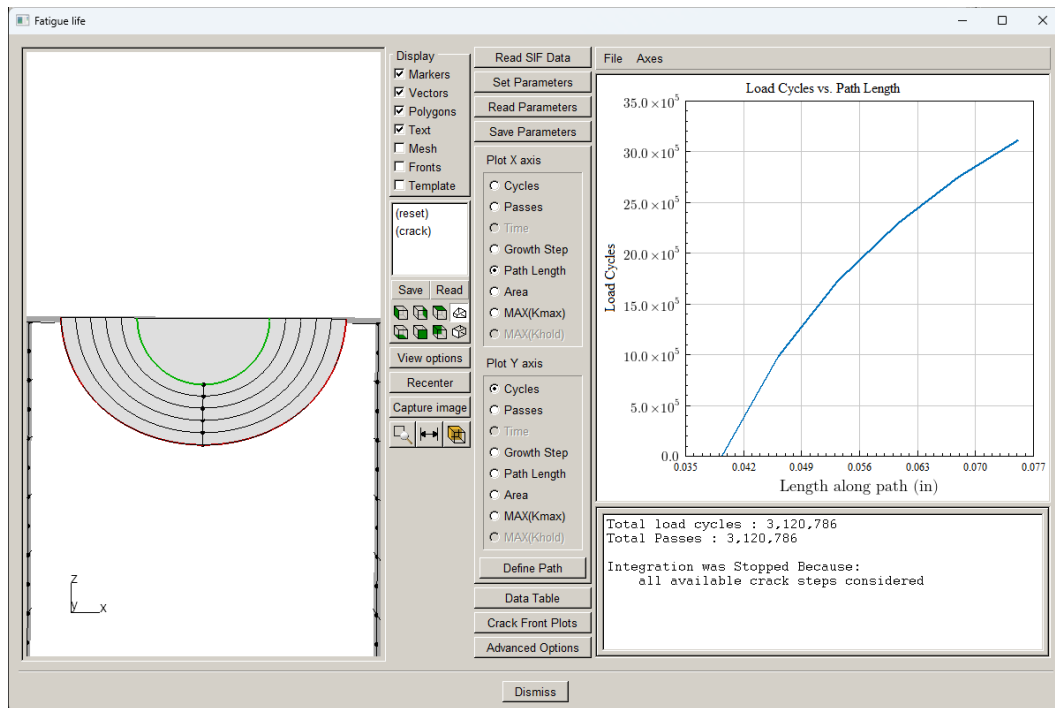


Figure 2.77 Fatigue cycles versus crack length

Appendix A:

ANSYS Workbench

A.1 APDL Commands in Workbench

ANSYS Workbench combined with APDL commands can be used to create *.cdb* files that can be used in FRANC3D. A simple plate model is constructed in Workbench, Fig A-1, with two regions, Fig A-2. The small central region is defined to create a local model for FRANC3D; in practice, the local region can be extracted with FRANC3D without having to define two regions in Workbench. Fig A-3 shows the APDL commands that are added to Workbench to write the *.cdb* files. Three *.cdb* files are created; the *BOXALL.cdb* will be the full model and could be imported and divided in FRANC3D. Alternatively, the *BIG_BOX.cdb* and *SMALL_BOX.cdb* files could be imported as “already divided”.

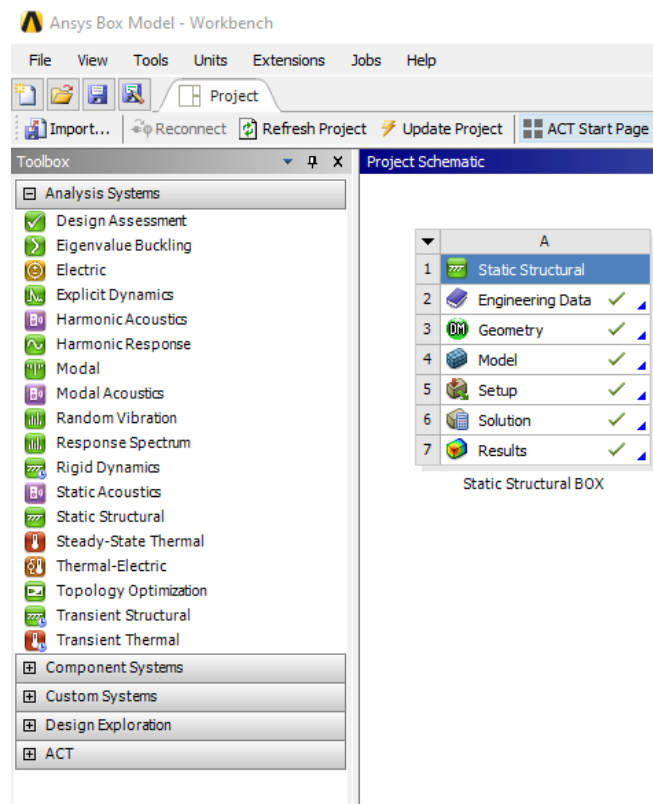


Figure A-1 Workbench Interface.

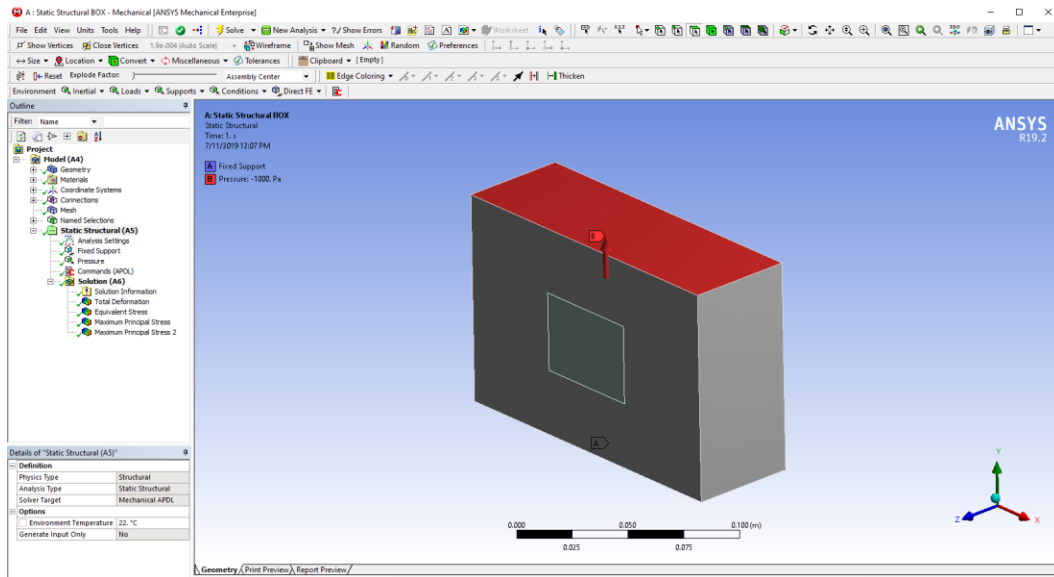


Figure A-2 Workbench model of a simple plate.

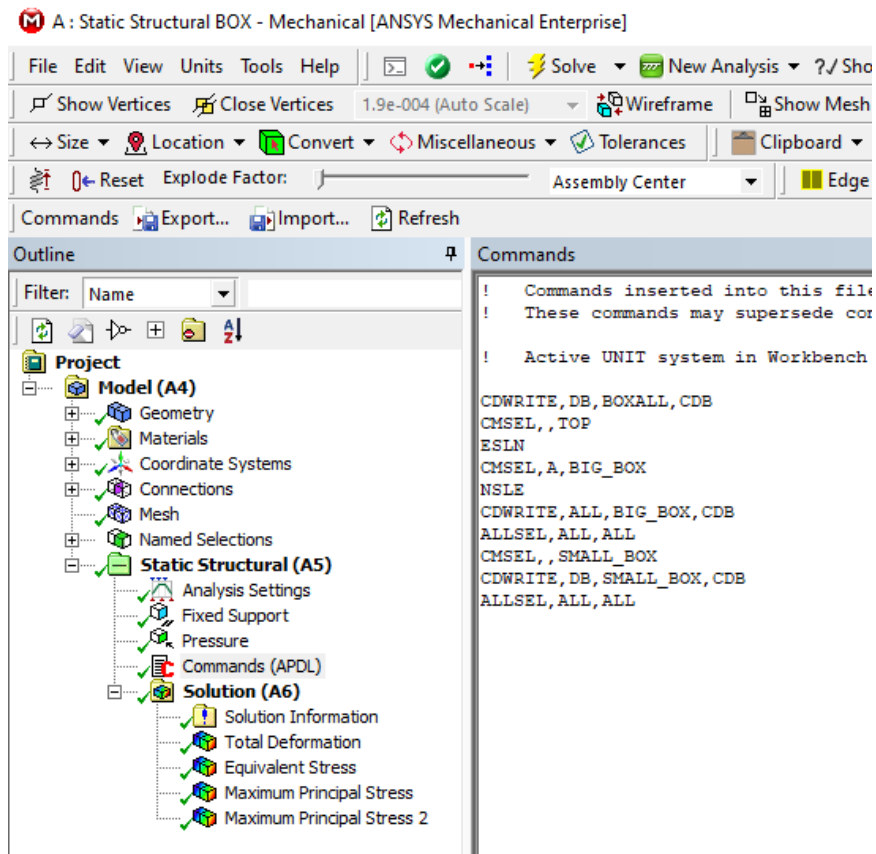


Figure A-3 Workbench APDL commands to write .cdb files.

A slightly more complicated process and model is described next. A model is created in Workbench and a Mechanical APDL is linked to the setup, Figs A4-A6. This is used to edit the model in ANSYS Classic to create the *.cdb* files if commands in Workbench are not used.

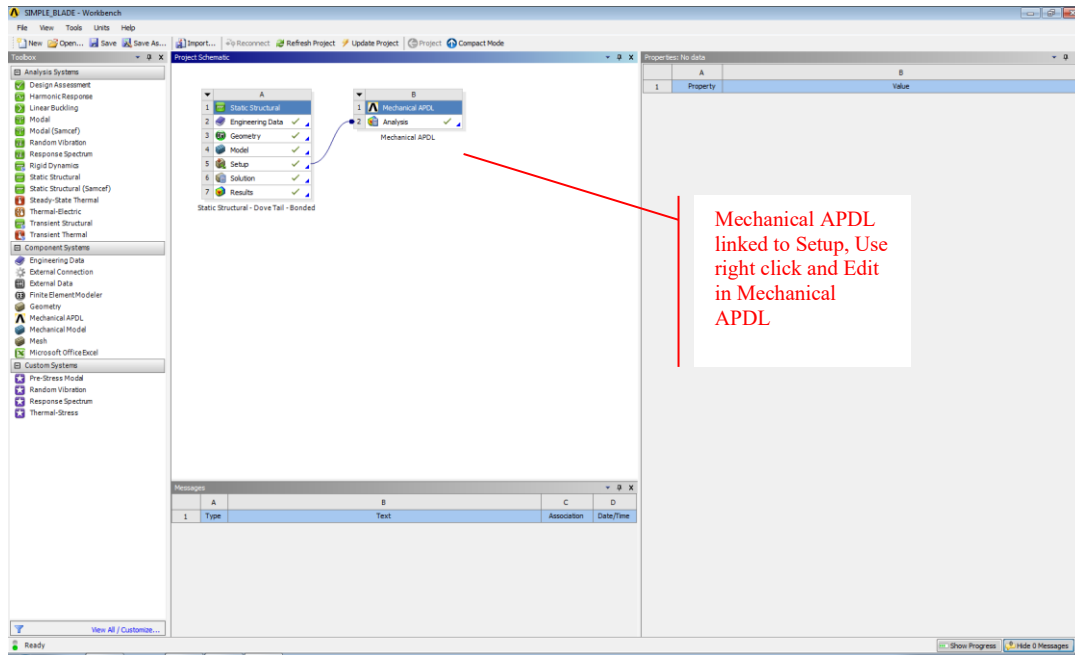


Figure A-4 Workbench interface to ANSYS Classic.

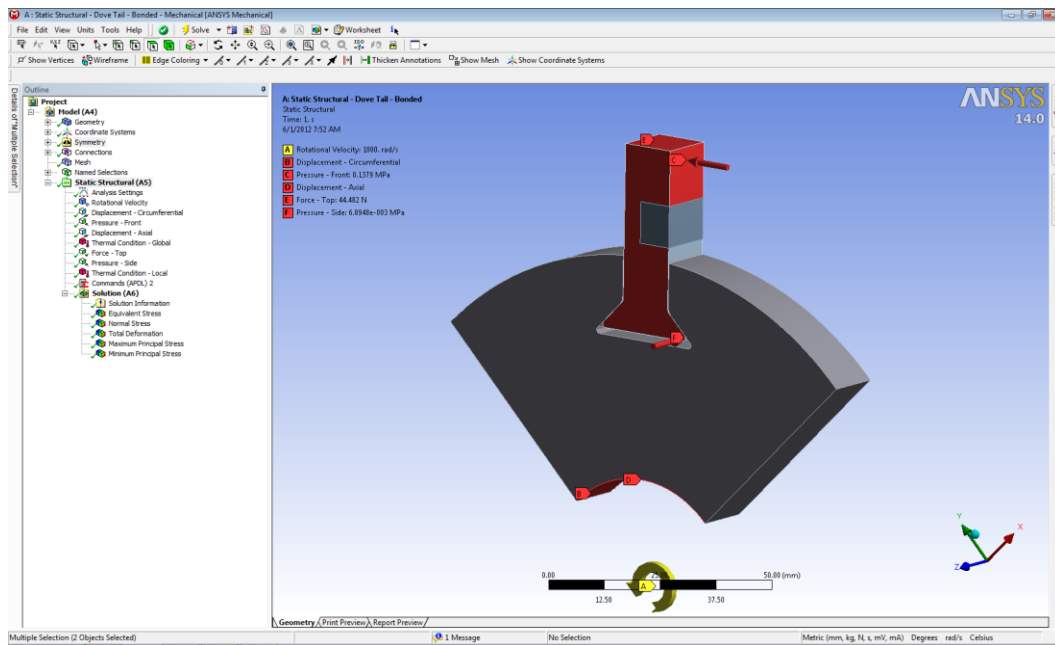


Figure A-5 Example model with temperatures, pressures, forces, contacts, and cyclic symmetry.

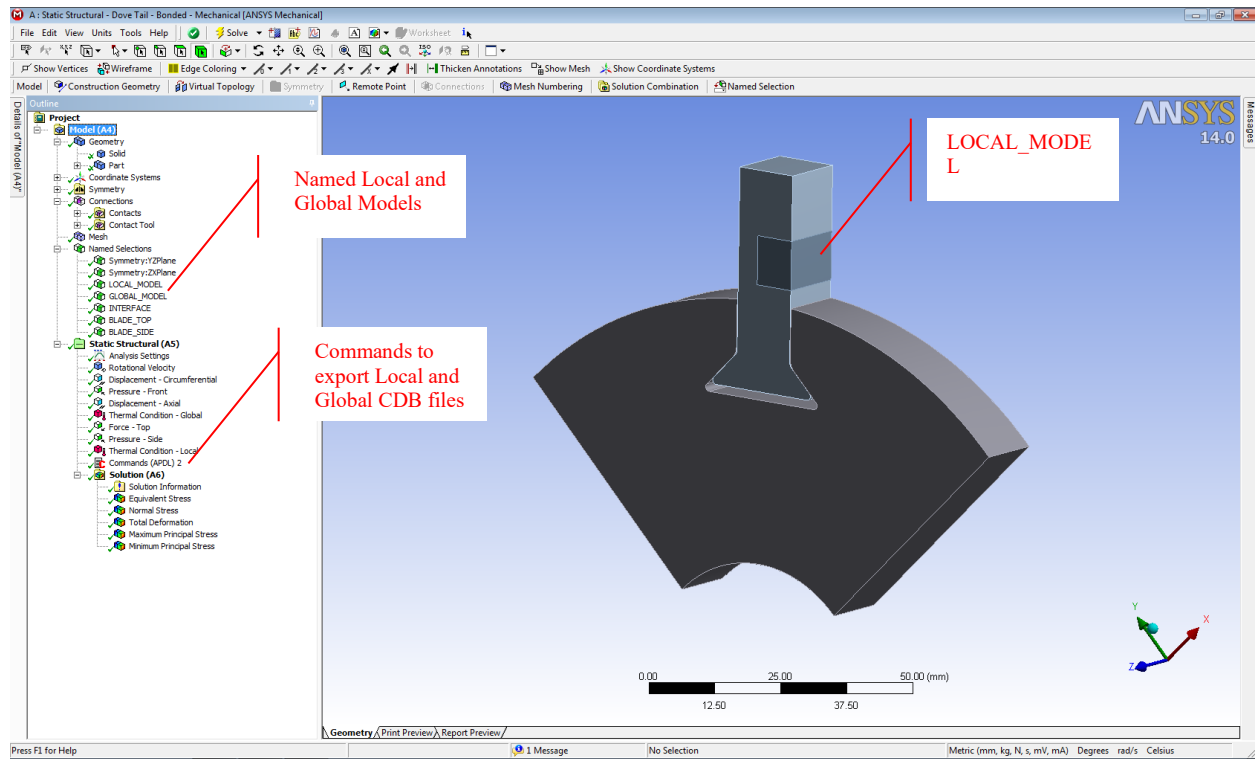
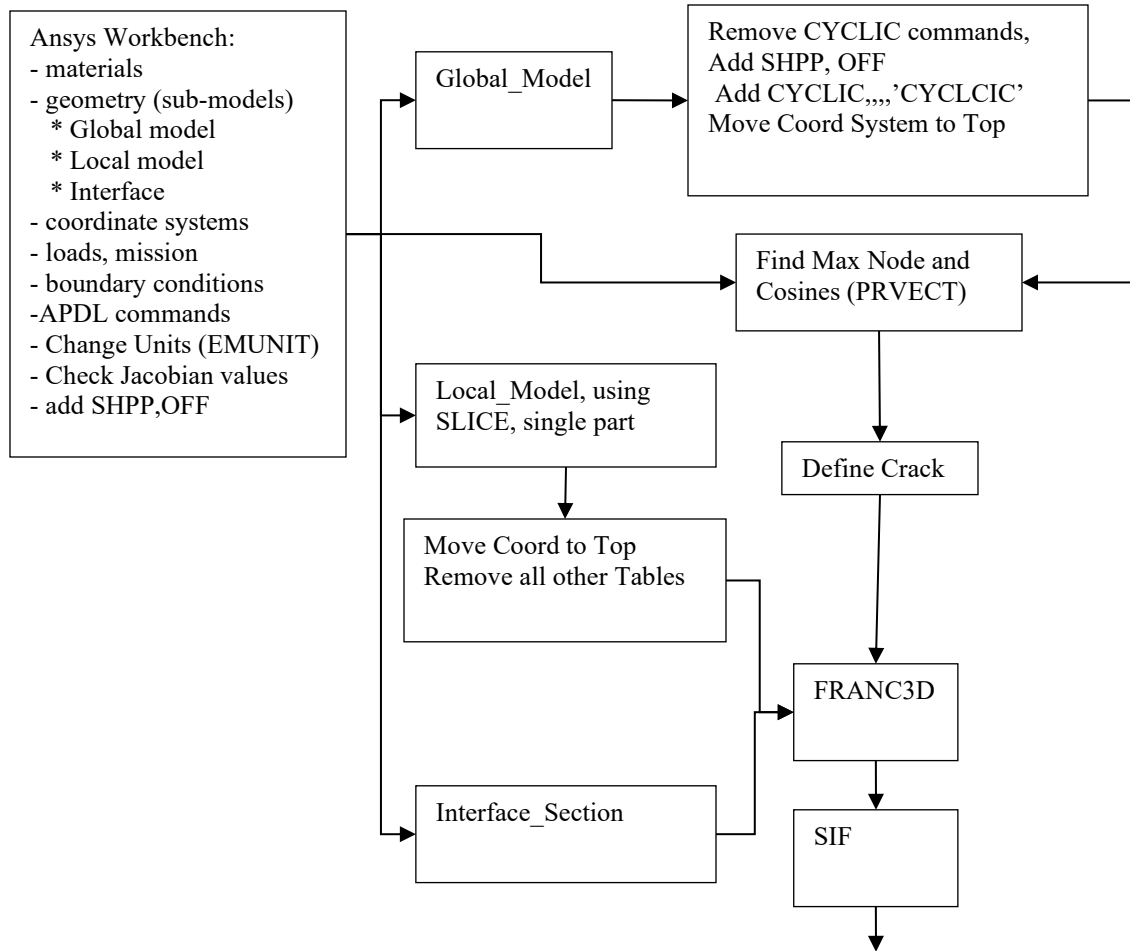


Figure A-6 Model Defining the Local and Global Models

The following method was used to create the local and global *.cdb* files for use in FRANC3D.

ANSYS:

1. Create model in Workbench
 - a. Set up coordinate systems
 - b. Name parts as GLOBAL_MODEL, LOCAL_MODEL, and INTERFACE
 - c. Add Temperatures, Pressures, Forces, Contacts, *etc.*
 - d. Add Boundary Conditions
 - e. Use consistent Units
2. Link APDL to Setup of Structural Model
 - a. Edit Mechanical in APDL (This will launch ANSYS Classic)
3. In Ansys Classic the following commands (with potential modifications) will be needed to create the CDB files or add commands to the Workbench Mechanical Setup:



ANSYS CLASSIC APDL CLICKS/COMMANDS:

→Select→Everything

CDWRITE,DB,F3D_FULL_MODEL,CDB

CMSEL,U,LOCAL_MODEL

NSLE

CDWRITE, DB,F3D_GLOBAL_MODEL,CDB

CMSEL,,LOCAL_MODEL

NSLE

CDWRITE, DB,F3D_LOCAL_MODEL,CDB

ANSYS WORKBENCH APDL COMMNADS:

ALLSEL,ALL,ALL

CDWRITE,DB,D:\F3DMODELS\DOVETAIL\F3D_FULL_MODEL,CDB

ALLSEL,ALL,ALL

CMSEL,U,LOCAL_MODEL

NSLE

CDWRITE,DB,D:\F3DMODELS\DOVETAIL\F3D_GLOBAL_MODEL,CDB

ALLSEL,ALL,ALL

```
CMSEL,,LOCAL_MODEL
NSLE
CDWRITE,DB,D:\F3DMODELS\DOVETAIL\F3D_LOCAL_MODEL,CDB
ALLSEL,ALL,ALL
```

GLOBAL MODEL FILE EDITS:

The following was changed in the GLOBAL_MODEL.CDB file:

1. REMOVED the following lines in the CDB file:
 *DIM,_CYCLICMAP
 *SET,_CYCLICMAP
 CYCL,CDWR
2. ADDED the following Element Shape Checking Command:
 SHPP,OFF
3. ADDED the following lines in the CDB file before the /GO and /FINISH Commands:
 CYCLIC,,,,'CYCLIC'

A.2 Workbench *ds.dat* File

ANSYS Workbench creates a *ds.dat* file when the model is submitted for analysis. This file contains all the model data and is an ASCII file, so it can be edited. A simple method of creating a *.cdb* file is to edit the *ds.dat* file and add a “cdwrite” command. The edited file can be processed by ANSYS and a *.cdb* file will be written.

FRANC3D can process a *.db* or a *.dat* file directly during the GUI import. A working ANSYS version must be accessible as FRANC3D will execute ANSYS in the background to convert these files into *.cdb* files. The resulting *.cdb* file is then imported. The process is briefly described here.

Starting with a *ds.dat* file from WB, use the FRANC3D Import and divide option, Fig A-7. Choose the *ds.dat* file, Fig A-8, after changing the File type filter. Once the file is selected, click **Next**; ANSYS runs in the background to create a *cdb* file along with *.s##* files, Fig A-9.

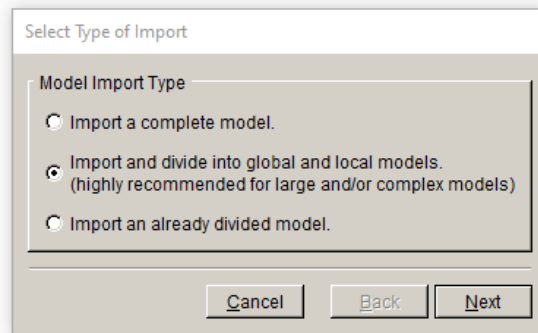


Figure A-7 Import and divide selected.

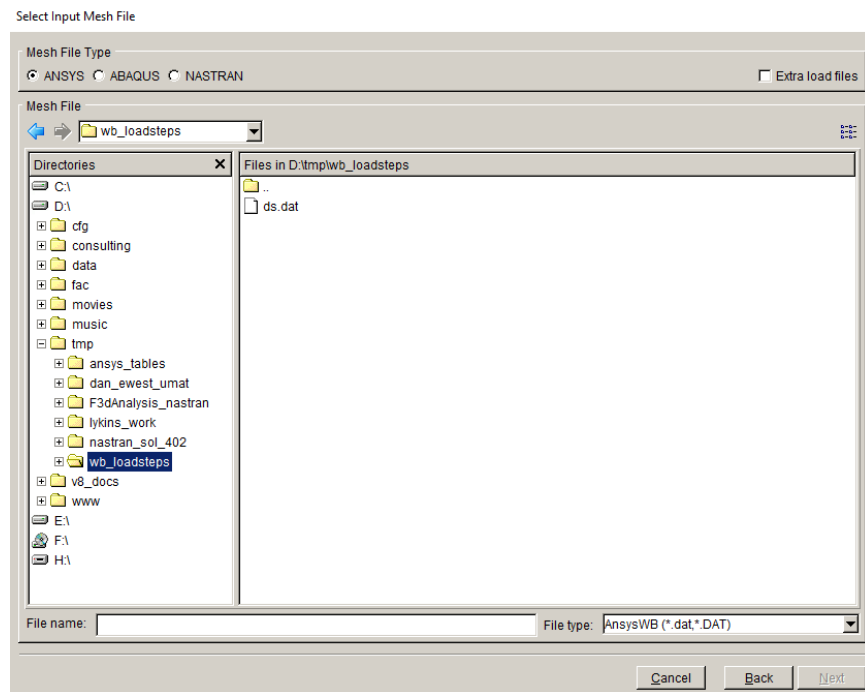


Figure A-8 Change the File type and select the ds.dat file.

```

C:\bruce\v8\franc3d_v8\Drivers\Release\franc3d.exe
Editing .dat file to comment /eof, turn off solve commands, and add lswrite,#
1 file(s) copied.
1 file(s) copied.
run ansys to convert the WB dat file to cdb
executing: "C:\Program Files\ANSYS Inc\v212\ansys\bin\winx64\ANSYS212.exe" -b -p ansys -j ds -i "D:\tmp\wb_loadsteps\ds.
dat" -o "D:\tmp\wb_loadsteps\ds.out"

```

Figure A-9 ANSYS is used to create .cdb and .s## files from the ds.dat file.

When ANSYS finishes, FRANC3D automatically imports the *.cdb* and *.s##* files into the submodeler dialog, Fig A-10. One can then cut out a local model, and save the Local and Global *.cdb* files, Fig A-11. The Local and Global *cdb* files will contain the relevant boundary condition data and load steps.

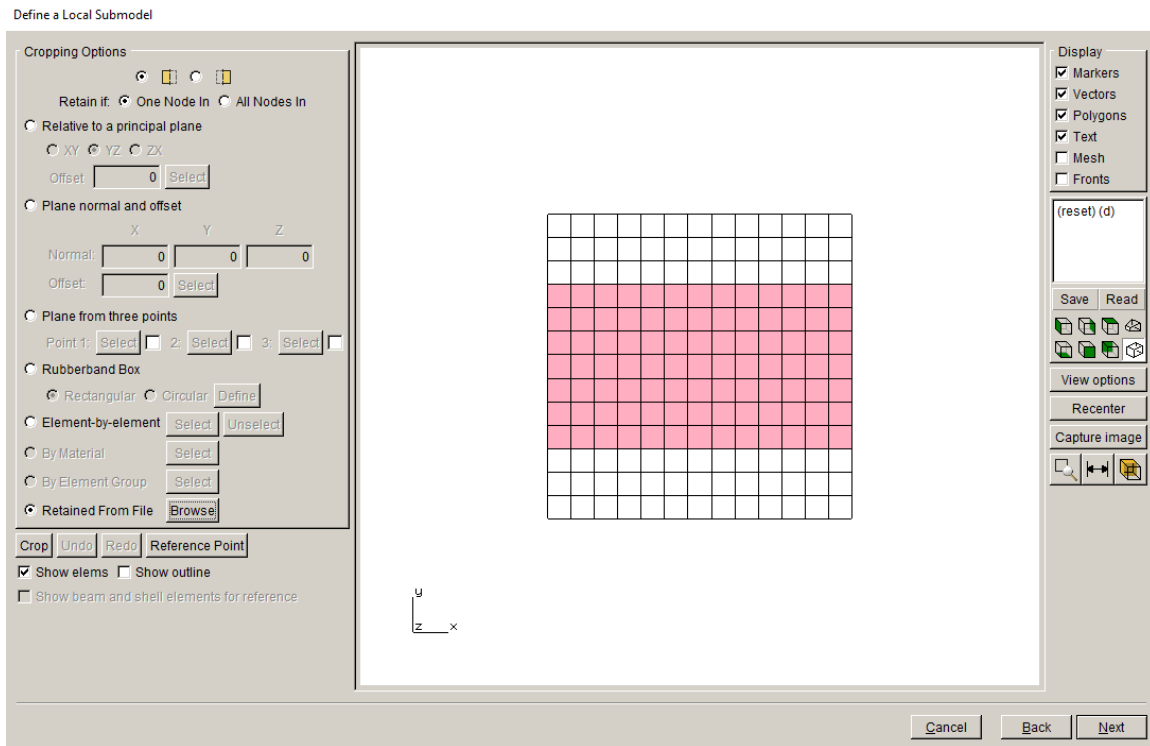


Figure A-10 FRANC3D submodel dialog.

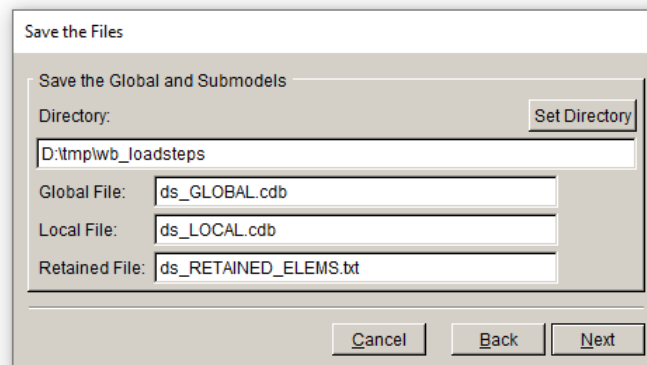


Figure A-11 FRANC3D dialog for saving the LOCAL and GLOBAL cdb files.

Appendix B:

Python Script to Add/Modify ANSYS CP Data

ANSYS CP (coupled degrees of freedom) data is usually passed through by FRANC3D. It is best to have all the nodes with CP data in the “global” model portion, but in some cases, one might need to have some of the nodes with CP data in the “local” model portion. In this case, the surface mesh facets should be retained for those CP nodes. The resulting cracked *_full.cdb* file should be edited to recreate the CP data; this can be done using a simple Python script that will run prior to ANSYS.

We will demonstrate this using a simple disk, Fig B-1. The boundary conditions include CP data on the symmetry surfaces, Fig B-2. The nodes on the two symmetry surfaces are put into a “cp_surfs” component (cmblock), which allows us to retain the nodes and mesh facets when we import the model into FRANC3D.

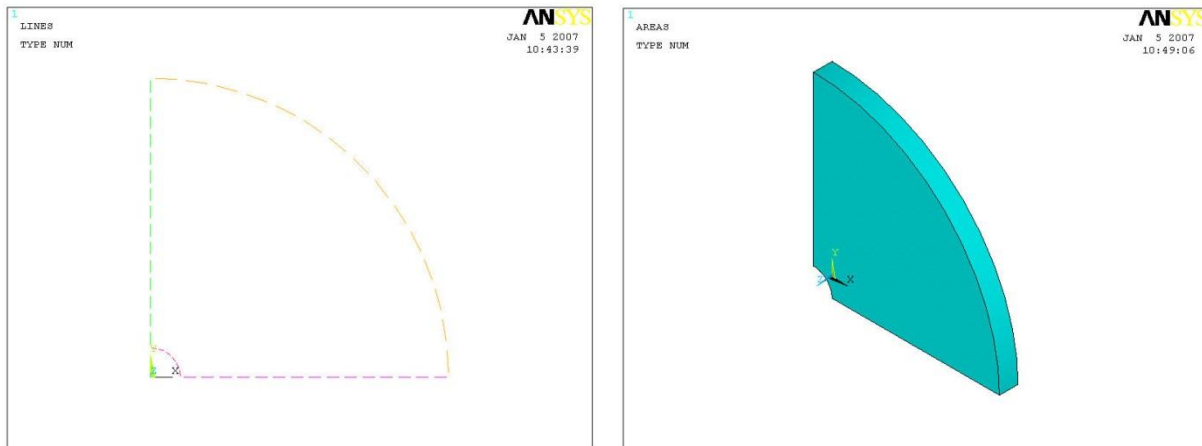


Figure B-1 ANSYS disk geometry.

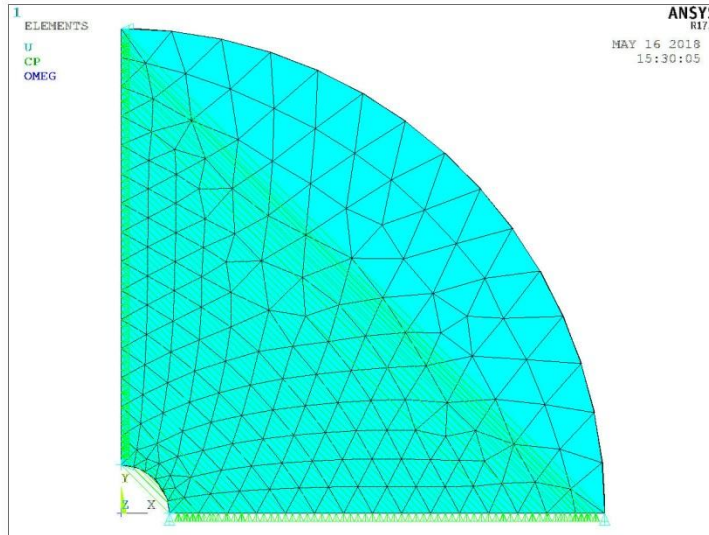


Figure B-2 ANSYS CP on symmetry surface nodes.

B.1 Import Disk into FRANC3D

Start FRANC3D and select **File** and **Import**. In the Select Type of Import panel, choose the Import and divide into global and local models radio button and select **Next**. Switch the Mesh File Type radio button in the Select Import Mesh File window to ANSYS and select the file name for the model, called *disk_CP.cdb*, Fig B-3. Select **Next**.

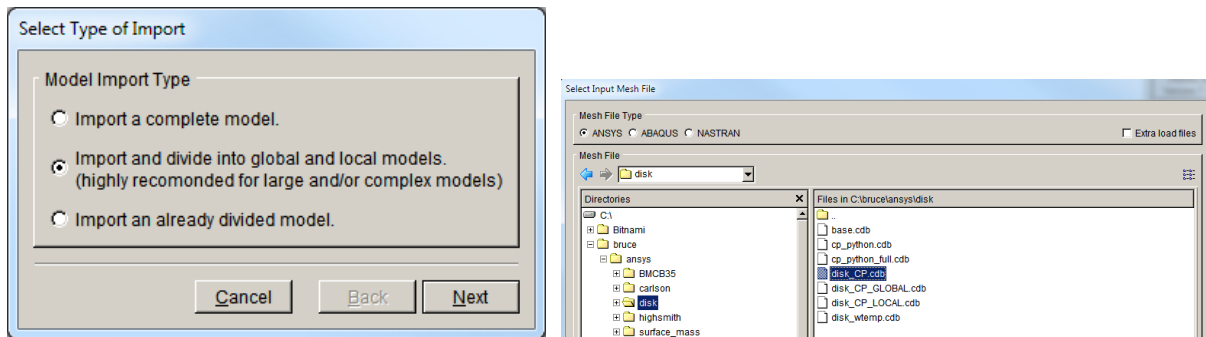


Figure B-3 FRANC3D FE model import.

Create a local model using the Rubberband Box tool in the Define Local Submodel window, Fig B-4. Select **Next** after using the **Crop** button and then save the local and global model files using the default file names. In the Select Retained BC Surfaces window, Fig B-5, click Select

all under BCs and also check the CP_SURFS box. Select **Next** when ready to proceed. The model will be displayed in the FRANC3D main window, and we are ready to insert the crack into the local model.

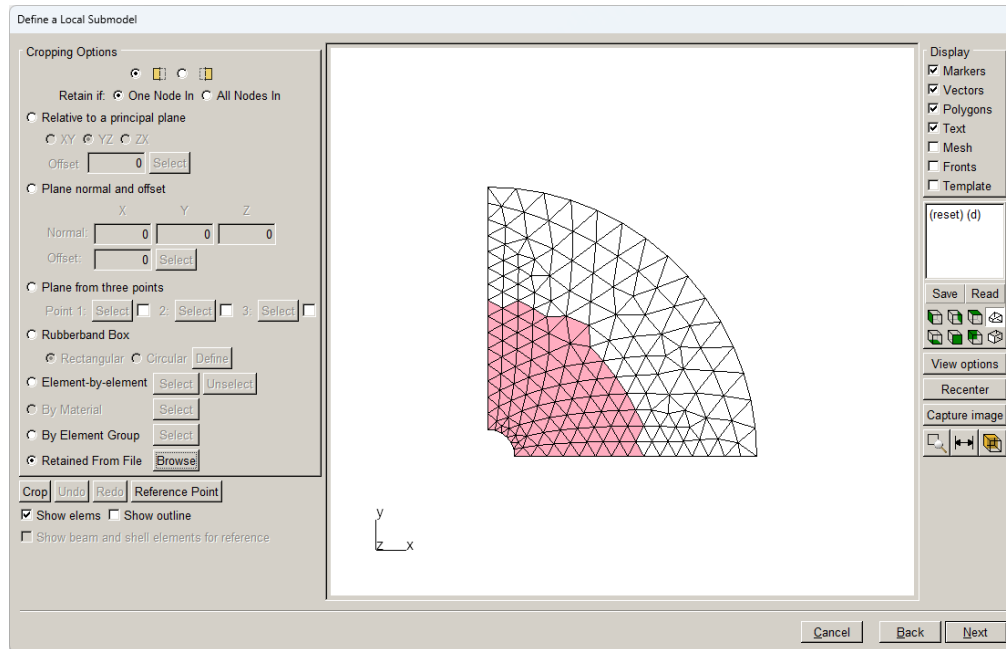


Figure B-4 FRANC3D FE model import.

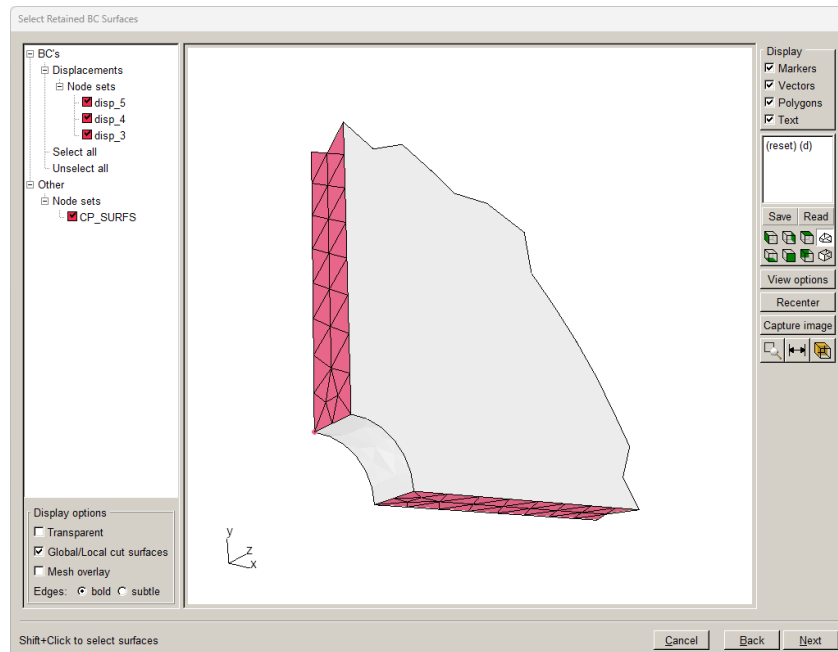


Figure B-5 FRANC3D Select Retained BC Surfaces.

B.2 Insert Initial Crack

We insert an elliptical crack with a radius of 0.1, Fig B-6. Note the surface mesh facets and nodes are retained on the CP symmetry surfaces. This allows us to regenerate the CP data in ANSYS as described in the next two sections.

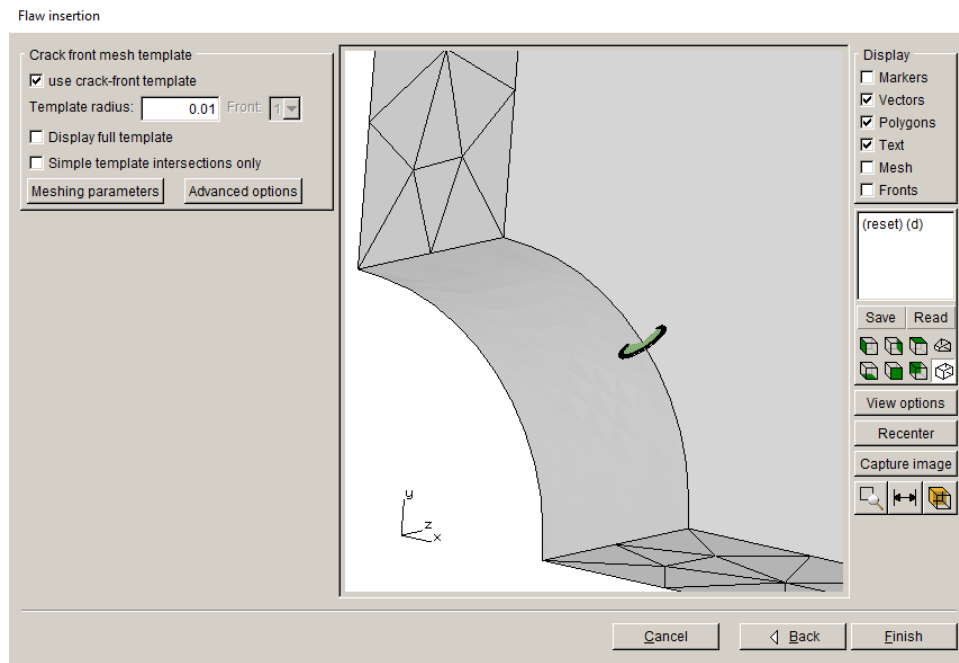


Figure B-6 Crack inserted into local model.

B.3 Create Python Script

We need to create a Python script that will modify the `_full.cdb` file that is written during the next step. A sample script is shown here.

```
#!/usr/bin/python
import sys
import string

if __name__ == '__main__':
```

```

fname = sys.argv[-1]
# read the input file
fdi = open(fname,'r')
buff = fdi.readlines()
output = []

for lb in buff:
    if len(lb) > 0 :
        if lb[0] == '/' or lb[0] == '!' :
            str = lb
            vals = str.split()
            if vals[0].upper() == '/SOLU' :
                output.append(lb)
                cpdel = "CPDEL,all,all\n"
                output.append(cpdel)
                cpcyc = "CPCYC,UY,0.001,1,0,90,0\n"
                output.append(cpcyc)
                continue

        output.append(lb)

fdi.close()
fdo = open(fname,'w')
for lo in output:
    fdo.write("%s"%(lo))
fdo.close()

```

The Python script reads the file and then iterates through the lines looking for the ‘/SOLU’ string. When it finds this string, the CP commands are added, which will cause ANSYS to delete and then recreate the CP data. The file is overwritten with the edited information.

Note that the Python script operates on the *_full.cdb* file prior to running ANSYS; this is described next.

B.4 Static Crack Analysis

We perform a static crack analysis using ANSYS. Choose **Analysis** and **Static Crack Analysis** from the FRANC3D menu. Provide a file name (*disk_cp_crack.fdb*) and choose ANSYS as the solver. The ANSYS options are shown in Fig B-7. The ANSYS executable and license string

should be set for your installation. The Python executable can be found using the **Browse** button if it is not in your PATH environment variable. The Python script can be set using the **Browse** button. We have saved the script from the previous section to a file: *editAnsysCdb.py*.

The Connect to global model option is checked automatically, and the global file name is set. Select **Next** to continue. The next dialog box, Fig B-8, has options for connecting the local and global portions. We use node merging with the automatically selected local and global cut-surfaces.

The image shows a 'Static analysis' dialog box with the following sections and controls:

- ANSYS Options:**
 - ANSYS run time:**
 - ANSYS Executable:
 - ANSYS license:
 - Python Executable:
 - Python script (will run before ANSYS):
- Local model output:**
 -
- Global model:**
 - ☒ Connect to global model filename:
- Boundary conditions:**
 - ☐ Apply crack face tractions ☐ Define crack face contact ☐ Initial strain
- ANSYS command:**
 - ☐ Write files but DO NOT run analysis

At the bottom are buttons for , , and .

Figure B-7 Static analysis options for ANSYS with Python script provided.

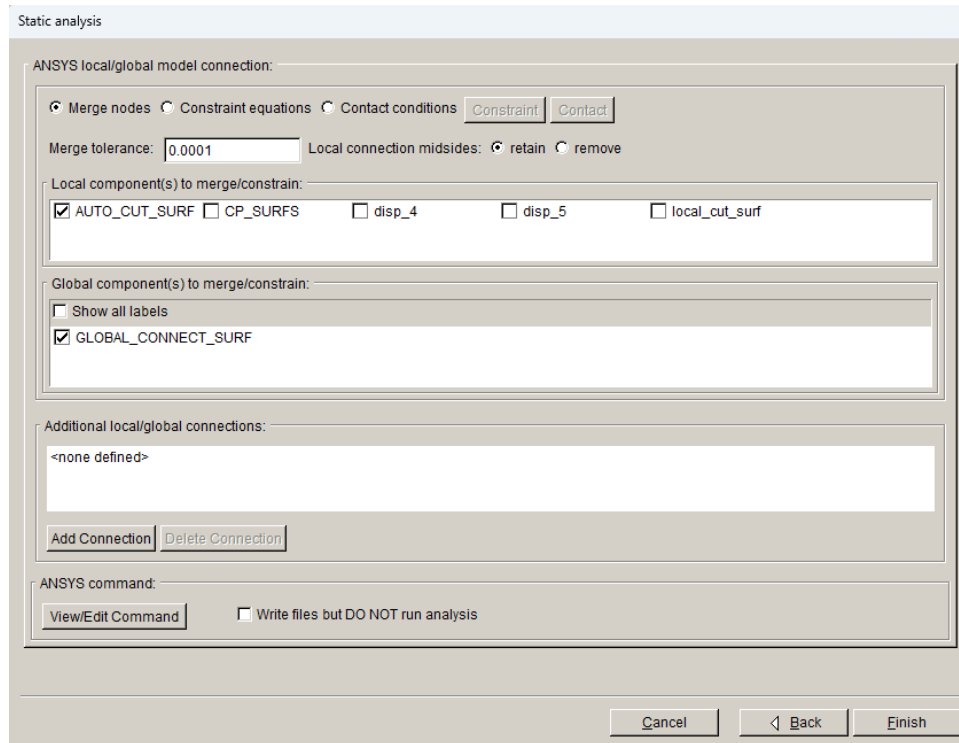


Figure B-8 Static analysis ANSYS local + global model connection dialog.

Click **Finish** to begin the process of running Python followed by the ANSYS analysis.

You can verify that the *_full.cdb* file was edited correctly, Fig B-9. You can also verify that the ANSYS solution is correct, Fig B-10.

```

/COM, select everything and solve
allsel,all,all
/GOPR
/solu
CPDEL,all,all
CPCYC,UY,0.001,1,0,90,0
solve

```

Figure B-9 Portion of the *_full.cdb* file with CP commands added.

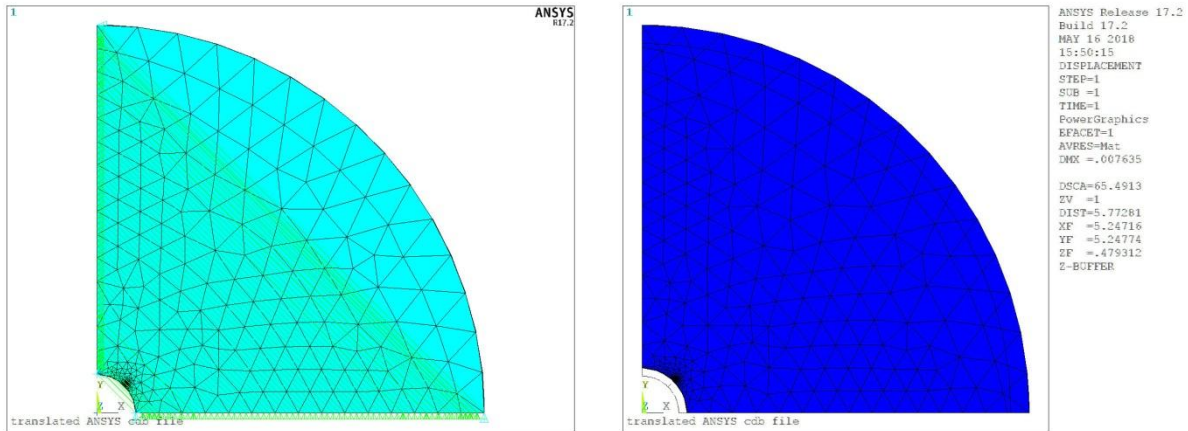


Figure B-10 ANSYS _full.cdb model showing CP data (left) and deformed shape (right).

B.5 Compute SIFs

Once ANSYS has finished running, we compute the SIFs; choose **Cracks** and **Compute SIFs** from the FRANC3D menu. The dialog, Fig B-11, allows you to specify the method for computing SIFs. If you use the M-integral, you can make sure that the Include Thermal Terms option is checked, and that the Reference Temperature is 0 degrees by clicking the **Advanced** button. The SIFs based on the M-integral are shown in Fig B-12. The SIFs should match the SIFs from Tutorial #4, which uses roller boundary conditions on the symmetry-surfaces.

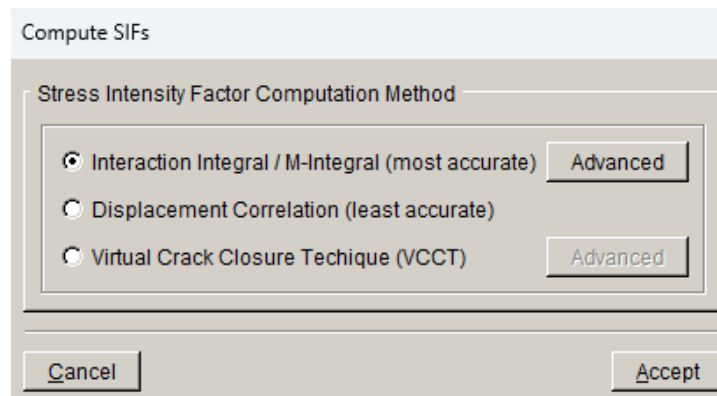


Figure B-11 Compute SIFs dialog.

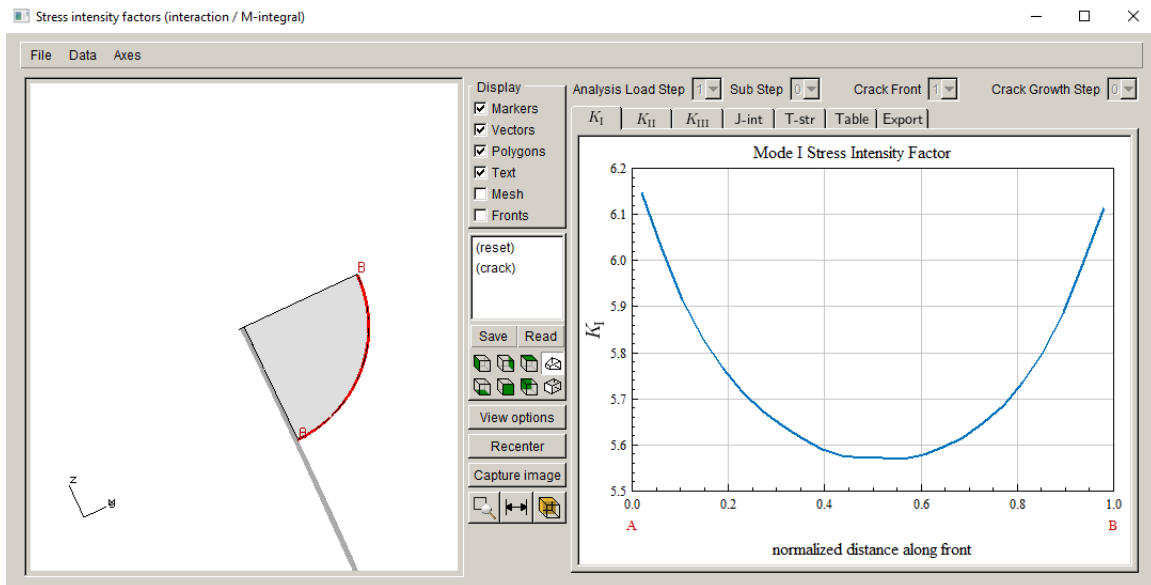
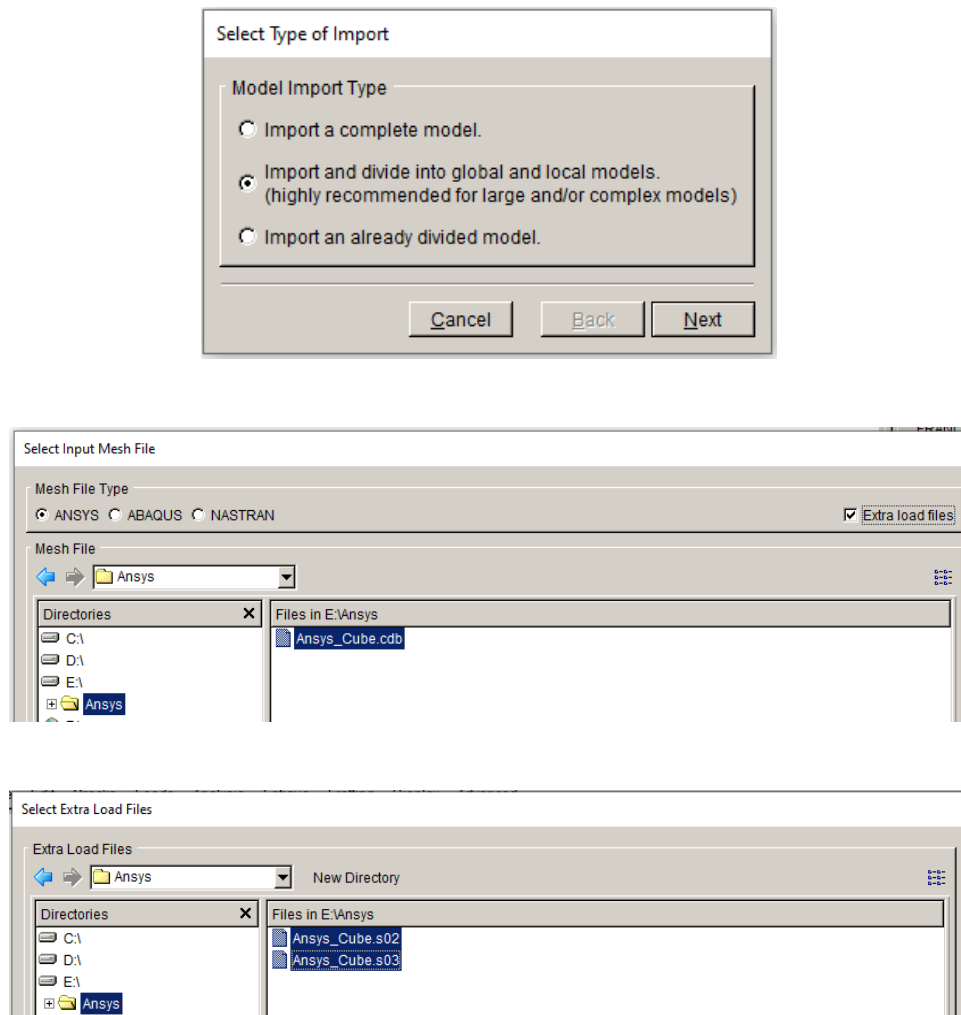


Figure B-12 M-Integral based SIFs.

Appendix C: ANSYS Model with Extra Load Steps

When importing and dividing an ANSYS model with extra load step files, it is important to note that the extra loads will be added to both LOCAL and GLOBAL *.cdb* files. Fig C.1 shows the sequence of import dialogs for an ANSYS model with two extra load step files.



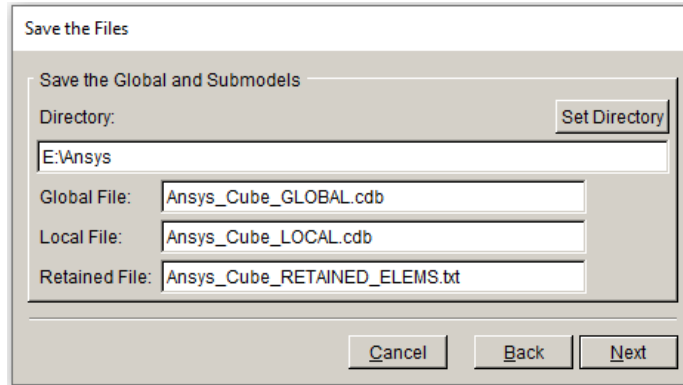


Figure C.1 Sequence of dialogs for importing and dividing an ANSYS model with two .s## files.

To verify that the extra load step information is included, we can open the GLOBAL .cdb file in a text editor and search for “_lnum=” commands. For example:

```
allsel,all,all
/SOLU
_lnum= 1
solve
```

The _lnum provides the FRANC3D load step ID. The original *Ansys_Cube.cdb* file contains one load step and the .s02 and .s03 files each contain one load step; these could correspond with arbitrary ANSYS load steps, but FRANC3D will renumber internally. Older versions of FRANC3D, used ANSYS “lswrite” and “lssolve” commands.

The user should not include the extra .s## files if the local and global model files are imported again (as an already divided model). The Extra load files check box, Fig C.1, is not available when importing using the “already divided” option.

Tutorial #2 uses multiple load steps, and the ANSYS version of this model has two extra load step files. If you decide to create your own local and global portions outside of FRANC3D, then it is important that you put all the load step information for each portion in the appropriate .cdb file. FRANC3D needs to keep track of the load steps and needs to recombine the cracked-local with the global portion; this can only be done if all information is contained in the .cdb files.

Appendix D: ANSYS Keywords

The list of ANSYS keywords that FRANC3D recognizes is provided here. FRANC3D must process some of the data, but a large amount of the *.cdb* file can be passed through to the *_GLOBAL.cdb* and then to the *_full.cdb*, which is the combination of the cracked/remeshed portion of the model and the *_GLOBAL.cdb* portion.

```
dispatch.Store(String("/COM"), &AnsysCdbReader::Comment) ;
dispatch.Store(String("/TITLE"), &AnsysCdbReader::Title) ;
dispatch.Store(String("*SET"), &AnsysCdbReader::sSet) ;
dispatch.Store(String("*DIM"), &AnsysCdbReader::sDim) ;
dispatch.Store(String("*IF"), &AnsysCdbReader::IfEndif) ;

dispatch.Store(String("ET"), &AnsysCdbReader::Et) ;
dispatch.Store(String("KEYOP"), &AnsysCdbReader::Keyop) ;
dispatch.Store(String("ETBLOCK"), &AnsysCdbReader::EtBlock) ;
dispatch.Store(String("TYPE"), &AnsysCdbReader::EType) ;

dispatch.Store(String("LOCAL"), &AnsysCdbReader::LocalCoordSys) ;
dispatch.Store(String("N"), &AnsysCdbReader::SingleNode) ;
dispatch.Store(String("NBLOCK"), &AnsysCdbReader::NodeBlock) ;
dispatch.Store(String("EN"), &AnsysCdbReader::SingleElem) ;
dispatch.Store(String("EBLOCK"), &AnsysCdbReader::ElemBlock) ;
dispatch.Store(String("MPTEMP"), &AnsysCdbReader::MPTemp) ;
dispatch.Store(String("MPDATA"), &AnsysCdbReader::MPData) ;
dispatch.Store(String("MP"), &AnsysCdbReader::MP) ;
dispatch.Store(String("TB"), &AnsysCdbReader::TBData) ;
dispatch.Store(String("_LSNUM"), &AnsysCdbReader::LSNum) ;
dispatch.Store(String("_LSNUM="), &AnsysCdbReader::LSNum) ;
dispatch.Store(String("LSWRITE"), &AnsysCdbReader::LSWrite) ;
dispatch.Store(String("D"), &AnsysCdbReader::DispBC) ;
dispatch.Store(String("F"), &AnsysCdbReader::ForceBC) ;
dispatch.Store(String("SFE"), &AnsysCdbReader::PressBC) ;
dispatch.Store(String("SFEBLOCK"), &AnsysCdbReader::PressBC) ;
dispatch.Store(String("SFGRAD"), &AnsysCdbReader::Sfgrad) ;
dispatch.Store(String("BFBLOCK"), &AnsysCdbReader::BFBlock) ;
dispatch.Store(String("BF"), &AnsysCdbReader::BodyForceBC) ;
dispatch.Store(String("BFE"), &AnsysCdbReader::BodyForceEBC) ;
dispatch.Store(String("BFUNIF"), &AnsysCdbReader::BodyForceUnif) ;
dispatch.Store(String("TREF"), &AnsysCdbReader::TempRef) ;
dispatch.Store(String("TIME"), &AnsysCdbReader::TimeBC) ;
dispatch.Store(String("CMBLOCK"), &AnsysCdbReader::CompBlock) ;
```

```

dispatch.Store(String("DDEL"), &AnsysCdbReader::DelDisp) ;
dispatch.Store(String("DDELE"), &AnsysCdbReader::DelDisp) ;
dispatch.Store(String("SFEDEL"), &AnsysCdbReader::DelSfe) ;
dispatch.Store(String("SFEDELE"),&AnsysCdbReader::DelSfe) ;
dispatch.Store(String("FDEL"), &AnsysCdbReader::DelForce) ;
dispatch.Store(String("FDELE"), &AnsysCdbReader::DelForce) ;
dispatch.Store(String("BFDEL"), &AnsysCdbReader::DelBf) ;
dispatch.Store(String("BFDELE"), &AnsysCdbReader::DelBf) ;
dispatch.Store(String("LSCLEAR"),&AnsysCdbReader::LsClear) ;
dispatch.Store(String("NLDIAG"), &AnsysCdbReader::Nldiag) ;
dispatch.Store(String("ANTYPE"), &AnsysCdbReader::AnType) ;
dispatch.Store(String("NSUBST"), &AnsysCdbReader::NSubSt) ;
dispatch.Store(String("NSUBS"), &AnsysCdbReader::NSubSt) ;
dispatch.Store(String("SOLVE"), &AnsysCdbReader::Solve) ;
dispatch.Store(String("SOLV"), &AnsysCdbReader::Solve) ;
dispatch.Store(String("LSSOLVE"),&AnsysCdbReader::LsSolve) ;
dispatch.Store(String("CDWRITE"),&AnsysCdbReader::CdWrite) ;
dispatch.Store(String("EXTOPT"),&AnsysCdbReader::Ignore) ;

```

```

// this is passed through, but we need to process to count load steps
dispatch.Store(String("ACEL"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("OMEGA"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("DOMEGA"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("CGLOC"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("CGOMEGA"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("DCGOMG"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("CMOM"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("CMOMEGA"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("CMDO"), &AnsysCdbReader::Omega) ;
dispatch.Store(String("CMDOMEGA"),&AnsysCdbReader::Omega) ;
dispatch.Store(String("CMRO"), &AnsysCdbReader::Omega) ;

```

```

dispatch.Store(String("CP"), &AnsysCdbReader::CoupleDof) ;

```

```

// currently this is passed through
dispatch.Store(String("CE"), &AnsysCdbReader::ConstraintEqn) ;
dispatch.Store(String("CERIG"), &AnsysCdbReader::ConstraintEqn) ;
dispatch.Store(String("SECTYPE"), &AnsysCdbReader::Pretension) ;
dispatch.Store(String("SECDATA"), &AnsysCdbReader::Pretension) ;
dispatch.Store(String("SECMODIF"),&AnsysCdbReader::Pretension) ;

```

```

// these are passed through
dispatch.Store(String("CYCLIC"), &AnsysCdbReader::Cyclic) ;
dispatch.Store(String("CYCL"), &AnsysCdbReader::Cyclic) ;

```

```

// we generate ansys commands to create contact

```

```
// we are using cmsel as the dispatch key for this
dispatch.Store(String("CMSEL"), &AnsysCdbReader::Cmsel) ;

dispatch.Store(String("/PREP7"), &AnsysCdbReader::Prep) ;
dispatch.Store(String("/SOLU"), &AnsysCdbReader::Solu) ;
dispatch.Store(String("/SOLUTION"),&AnsysCdbReader::Solu) ;
dispatch.Store(String("/POST1"), &AnsysCdbReader::Post) ;
dispatch.Store(String("FINISH"), &AnsysCdbReader::Finish) ;
dispatch.Store(String("FINI"), &AnsysCdbReader::Finish) ;
```