



# Tutorial

Simulation of a heat source in gas metal arc  
welding by using LS-PrePost<sup>®</sup>

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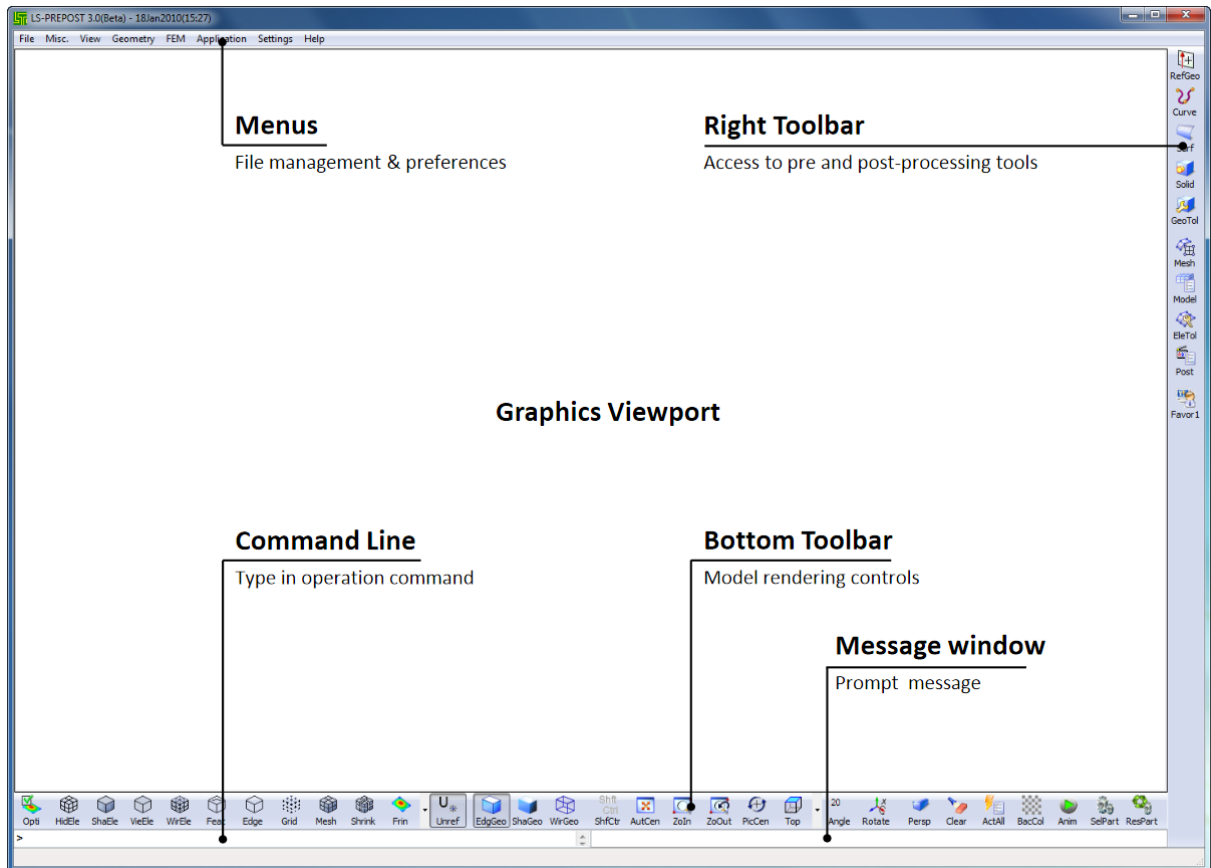
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# 1 Introduction

LS-DYNA<sup>®</sup> is a finite element code released by Livermore Software Technology Corporation (LSTC) for transient dynamic analysis capable of simulating complex real world problems. For solving coupled multi-physics and multi-stage problems, LS-DYNA<sup>®</sup> provides solvers, like arbitrary LAGRANGIAN EULERian (ALE), coupled fluid dynamics (CFD), smooth particle hydrodynamics (SPH), *etc.*

LS-DYNA<sup>®</sup> include a pre-processing and postprocessing environment for the analysis of finite element models. It is used in a wide range of industries like automotive, aerospace *etc.*, and also is extensively used in academic and research institutions due to its capability to address non-linear problems. The LS-PrePost<sup>®</sup> graphical user interface (GUI) is shown in Fig. 1.1. Fig. 1.1 helps the user to get familiar with the icons.



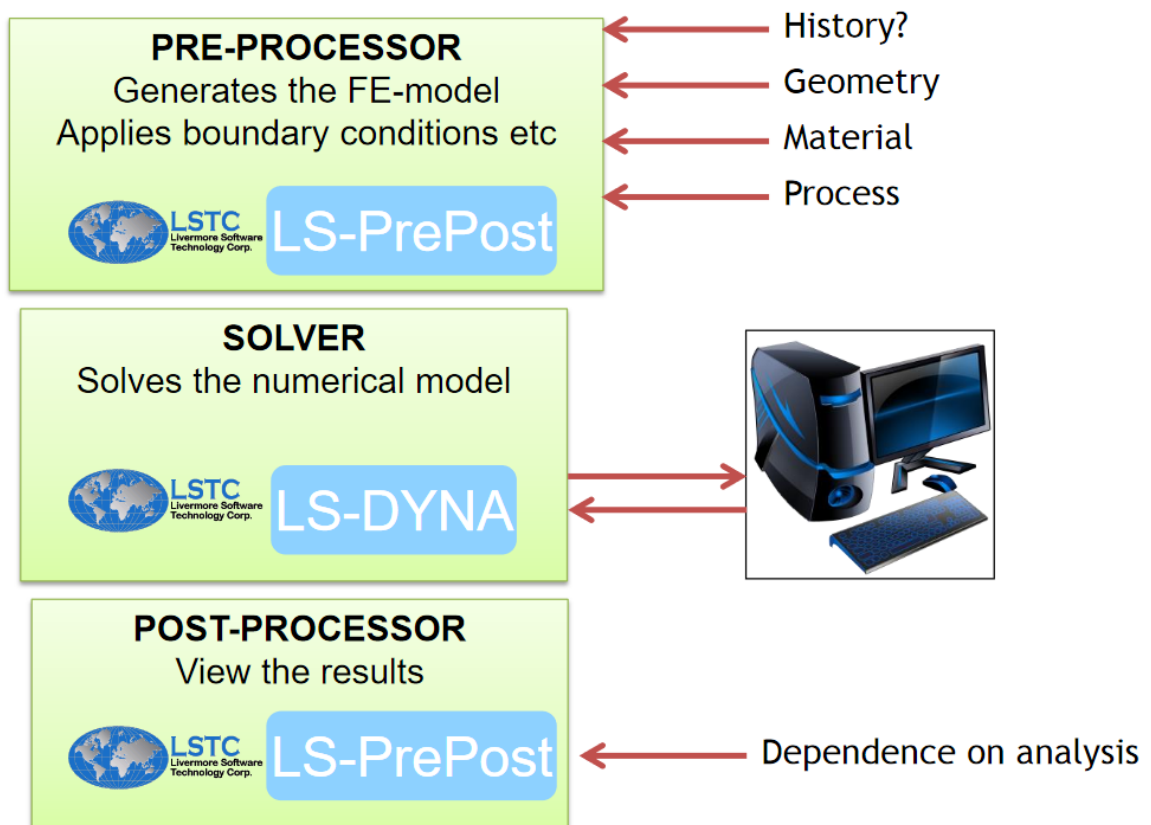
**Figure 1.1:** Graphical user interface (GUI) of LS-PrePost

The complete finite element analysis (FEA) is done in three different stages, namely preprocessing, analysis, and postprocessing, respectively. During preprocessing the given physical problem is modeled to generate a keyword file (.k). In the second stage, *i.e.* in analysis the modeled physical problem is solved numerically. The du-

ration of the analysis depends on the complexity of the given problems. In the last stage the results are evaluated by using a binary output database file (d3plot), see Fig. 1.2.

The general approach involved in solving physical problems by using LS-DYNA<sup>®</sup> are mentioned below:

- Create finite element (FE) Model
- Choose Material Model and Properties
- Assign Material and Property
- Assign loads and boundary condition
- Specify control parameters
- Create ".k" input file
- Solve the .k file in LS-DYNA<sup>®</sup> solver to get "d3plot" output file
- Post process the d3plot file in LS-PrePost<sup>®</sup>



**Figure 1.2:** Different steps of complete LS-DYNA<sup>®</sup> finite element analysis

## 1.1 Aim

The aim of this tutorial is to develop a step-by-step instructions for simulation of a welding process with the FEA-program LS-PrePost<sup>®</sup>. The tutorial contain a simulation of a single-pass gas metal arc welding, but we can use the same method to simulate multi-pass welding.

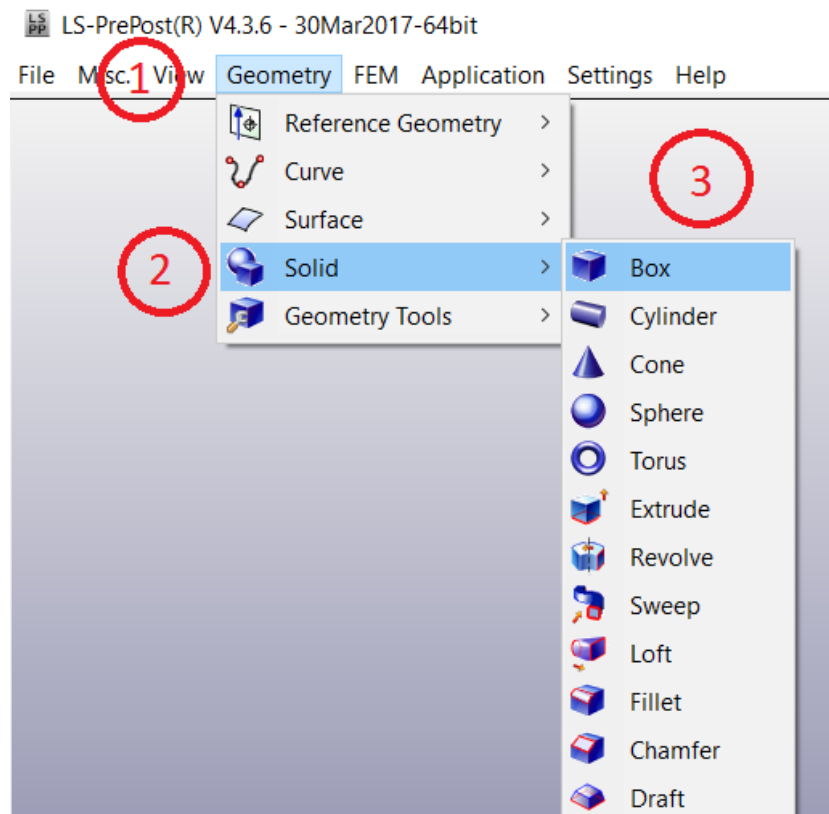
## 2 LS-PrePost

In the preprocessing, we generate a FE model for a physical problem. There are two methods to use LS-PrePost<sup>®</sup>. The first is by means of the graphical user interface (GUI). The second is by means of keyword file. The keyword file is a text file, which builds a bridge for communication between the preprocessor and the analysis. It contains a complete description of the FE-model, like geometry, material properties, boundary condition, mesh, *etc.* This tutorial is designed to describe the user both the GUI and the keyword file approach.

### 2.1 Geometry

The first step in this simulation is to create and give the model the right geometrical dimensions. Geometry module is used to build different parts of the model. For complex structure we can divide the whole model into various parts and create each part using geometry module. Later all the parts can be assembled together. Here, we have a simple rectangular plate of dimension  $300 \times 60 \times 6$  mm. To create the plate geometry go to **Geometry** → **Solid** → **Box**, see Fig. 2.1.

It is also possible to import an already existing model from a CAD program.



**Figure 2.1:** Selection of a solid box geometry



In **create Box** window choose the method parameter and enter value of the coordinate  $(x_1, y_1, z_1)$  and  $(x_2, y_2, z_2)$ . At the bottom click **Apply** → **Done**, see Fig. 2.2. Instead of the menu bar one can also use the shortcut in the right toolbar, see Fig. 1.1.

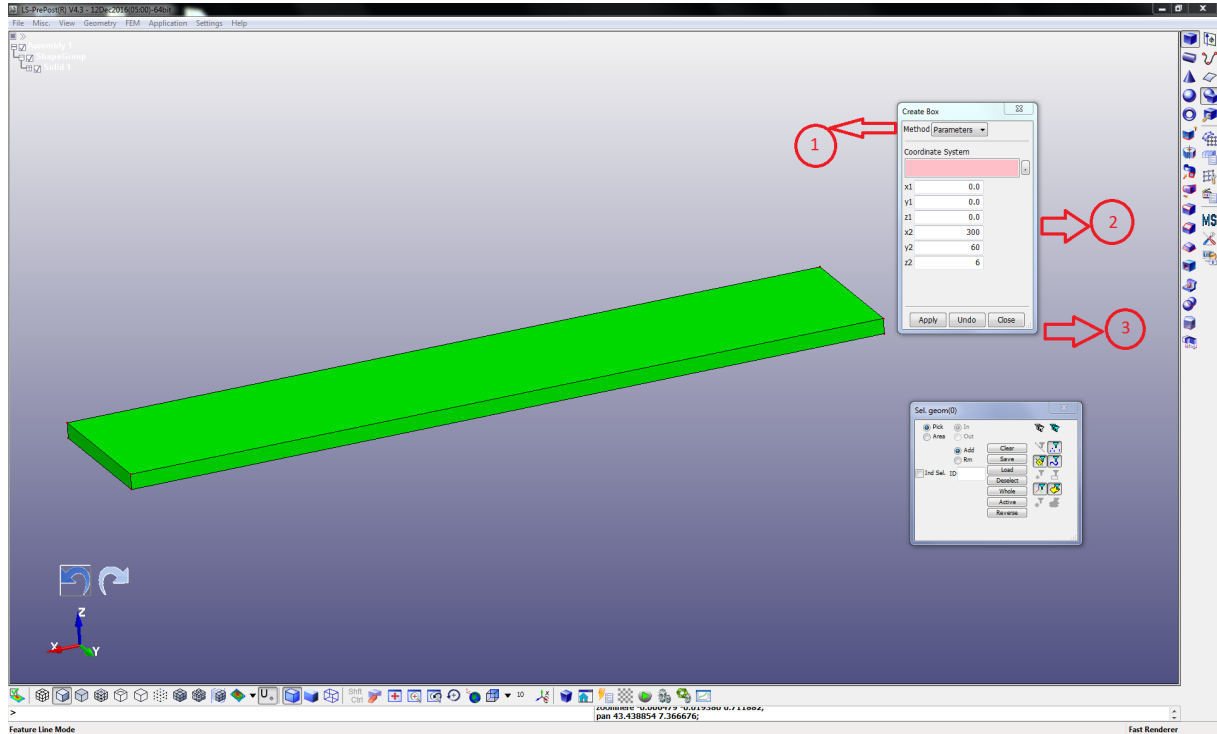


Figure 2.2: Geometry of the plate

## 2.2 Mesh

To create the mesh for the given plate we should use the following steps, see Fig. 2.3 and Fig. 2.4:

1. In the menu bar, click **FEM** → **Element** and **Mesh** → **Solid Mesher**
2. Choose Operation: **Meshing**
3. Click Volume by closed faces
4. Enter element size
5. Click **Try Meshing Automatically**
6. At the bottom click **Accept** → **Done**

This will create the mesh as shown in Fig. 2.4.

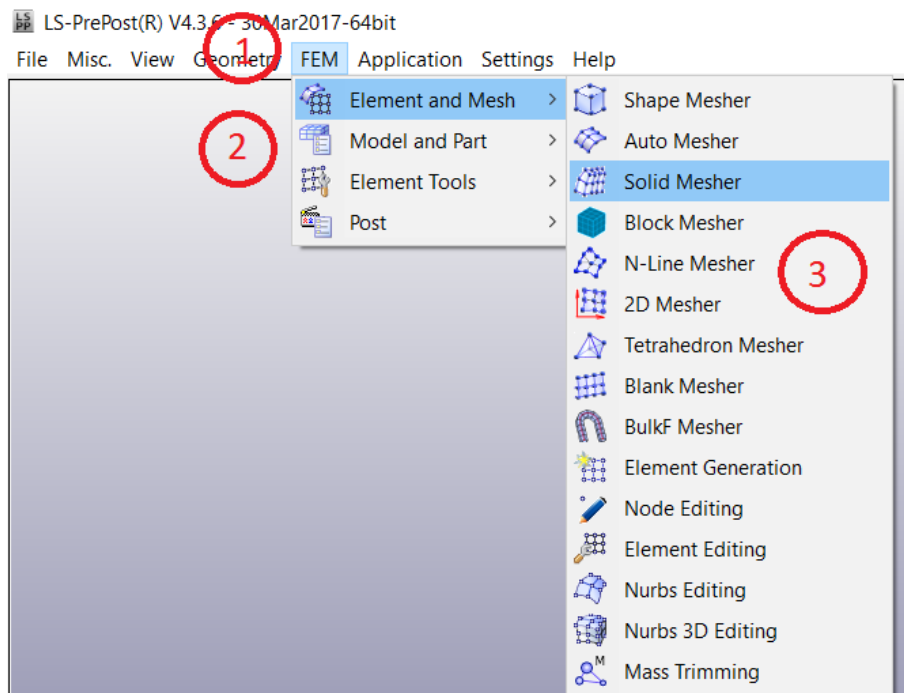


Figure 2.3: Selection of a solid mesher

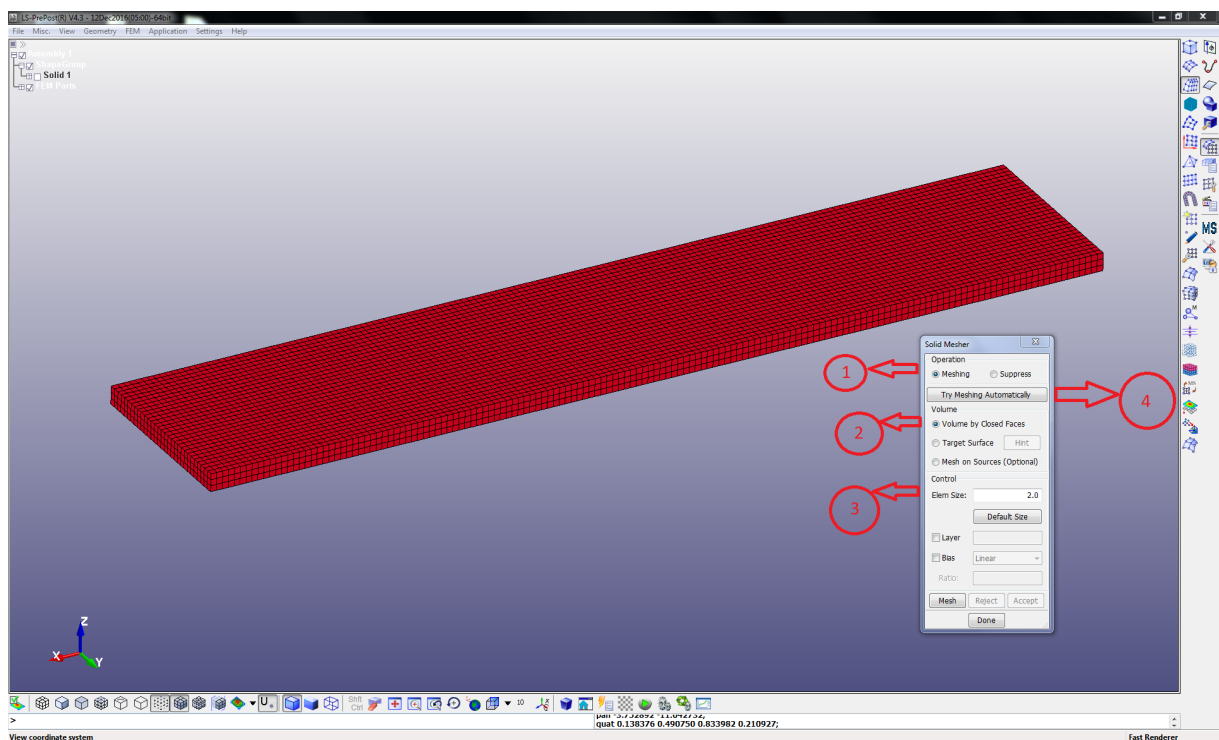
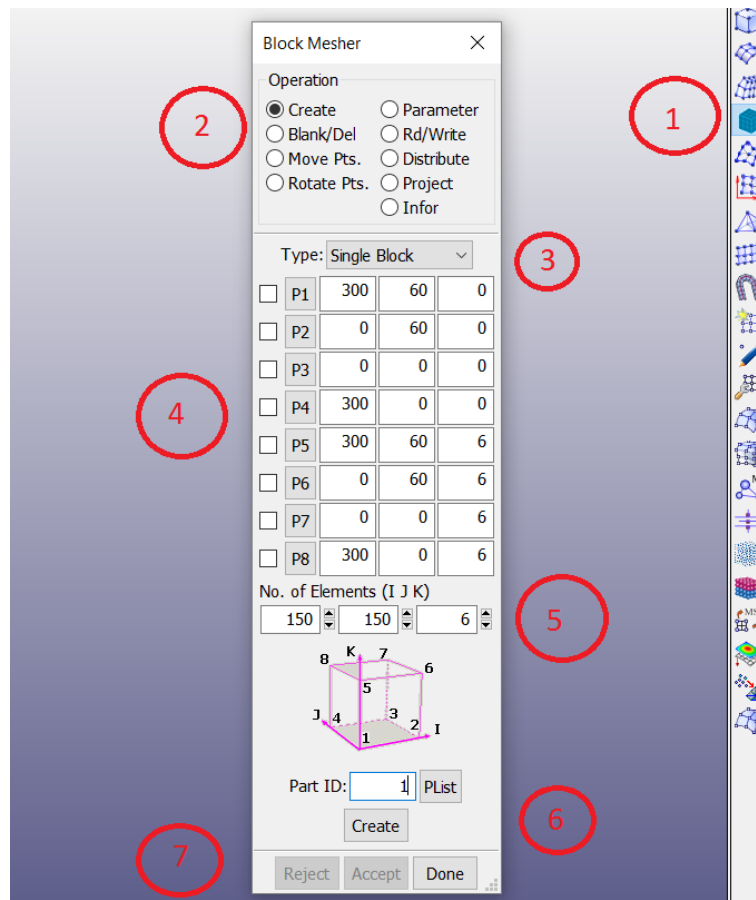


Figure 2.4: Plate mesh

We can also create mesh by using **Block Mesher**. This give us freedom to choose the number of elements in each direction manually. To create mesh by using **Block Mesher**

we can use the following steps, see Fig. 2.5:

1. In the right toolbar select **Element and Mesh** → **Block Mesher**
2. Choose Operation: **Create**
3. Choose Type: **Single Block**
4. Choose eight different corner of the plate
5. Enter the number of elements in I (x-axis), J (y-axis) and K (z-axis) direction, respectively
6. At the bottom click **Create** → **Accept** → **Done**



**Figure 2.5:** Dialog box for a Block mesher in LS-PrePost<sup>®</sup>

### 2.3 Welding simulation setup

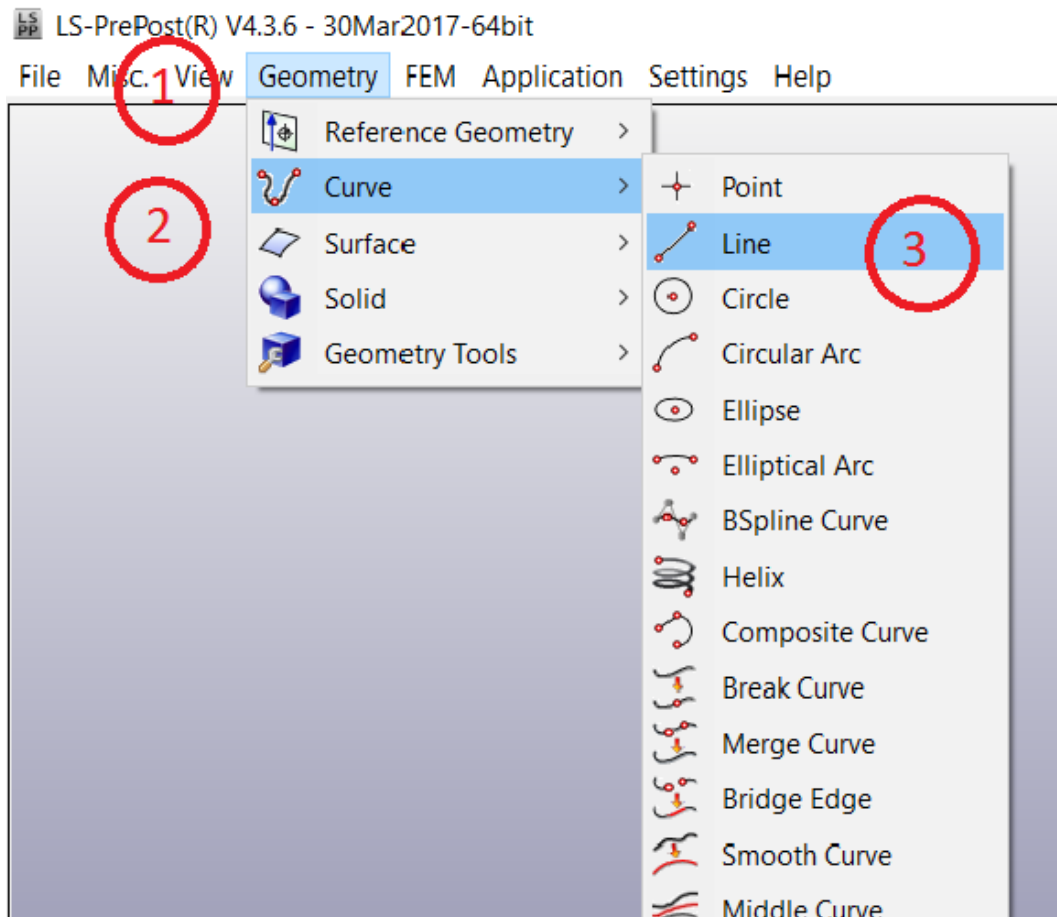
For welding setup in LS-PrePost<sup>®</sup> we need to create weld and orientation paths as a beam elements. The weld path defines the path of the moving heat source represented by a local moving coordinate system. The weld path is composed of geometrical information provided from the trajectories and the time information, initial time and velocity provided from the process plan. To obtain a local coordinate system, two trajectories are required

for each weld path: the weld line which describes the origin of the coordinate system and the reference line which describes its orientation.

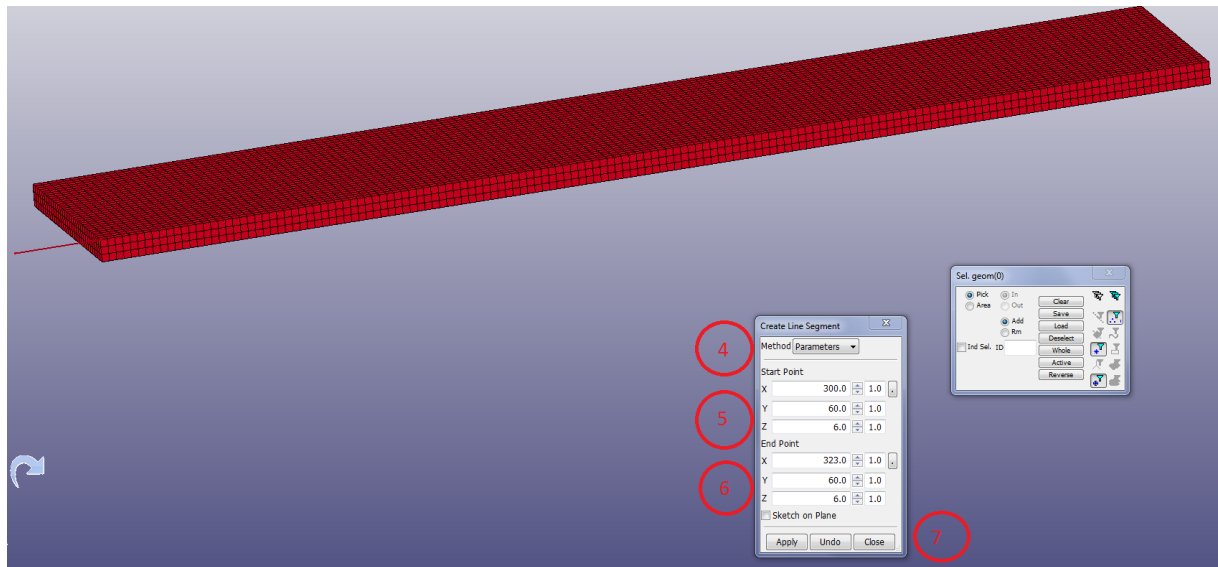
### 2.3.1 Weld line

To create a weld line we need to apply the following steps, see Fig. 2.6:

1. From the LS-PrePost® menu bar, select **Geometry** → **Curve** → **Line**. A sub menu called "Create Line Segment" will pop up.
2. In the Create Line Segment dialog box choose **Parameters** as the Method, see Fig. 2.7
3. Enter the coordinate of the Start Point
4. Enter the coordinate of the End Point
5. At the bottom of the dialog box click **Apply** and **Close**



**Figure 2.6:** Selection of the curve toolbar to create a line



**Figure 2.7:** Geometry of the welding line

### 2.3.2 Beam element

By following the steps below we can create the beam element, see Fig. 2.8:

1. From the right toolbar, select **FEM** → **Element and Mesh** → **Element Generation**. A sub menu "Element Generation" will pop up.
2. In the Element **Element Generation** dialog box, select the option **Beam**
3. **Element ID** and **Part ID** will be automatically generated
4. Choose Beam by: **Curve**
5. Click the **Third Node** option and give the values for x, y and z
6. Choose the option **Num** and enter the value 1 in order to divide the weld line in one beam element.
7. In the next step select the line by using your cursor
8. At Bottom of the dialog box, select **Create** → **Accept** → **Done**
9. In the **Sel. geom** window, choose the options **Pick** and **add**.

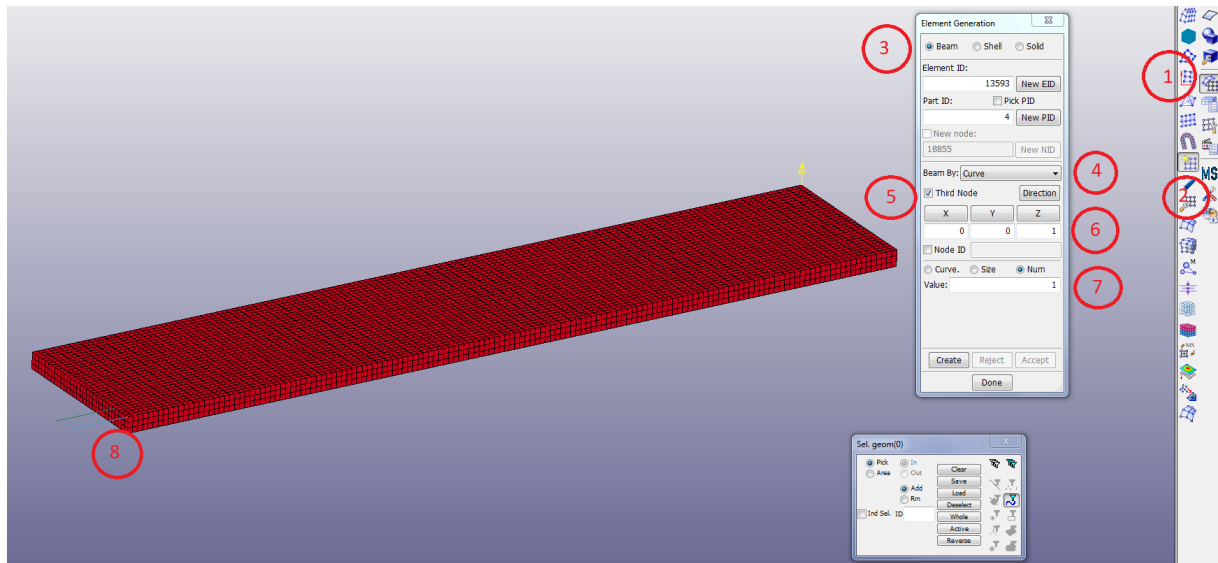


Figure 2.8: Beam element creation

### 2.3.3 Reference line

In LS-PrePost<sup>®</sup> we need to define a orientation path as beam element to model heat source. To create this reference line we can follow the same procedure as mentioned for welding torch, see Fig. 2.9

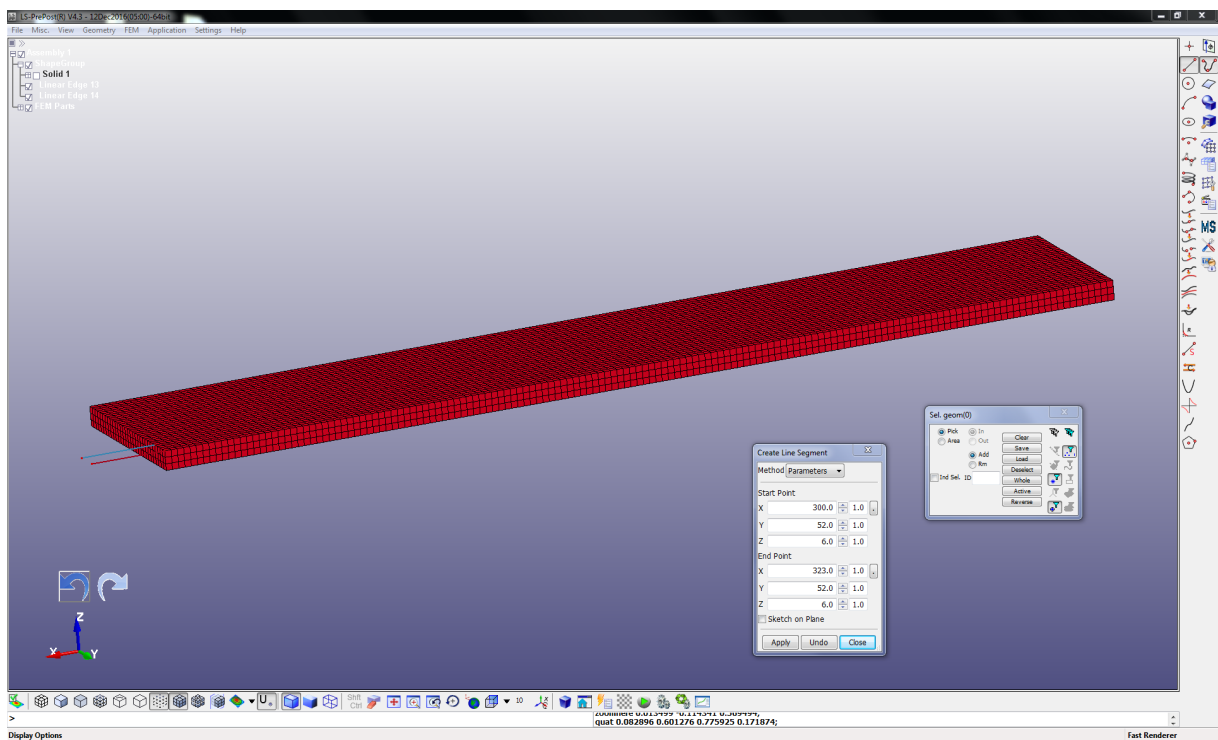


Figure 2.9: To create reference line

## 2.4 Weld and reference line path

To define the path of the moving heat source and its orientation, we can follow the steps below, see Fig. 2.10 und Fig. 2.11.

1. From the right toolbar, select **Model and Part** → **Create Entity**
2. From the **Entity Creation** box, select **Set Data** → **\*SET\_NODE**
3. Choose the option **cre** to create path for the weld line
4. Enter a **title** (This is optional).
5. From the **Sel. Nodes** dialog box, select **pick** and **add**.
6. Choose the nodes of interest along which the line will move
7. In the last step select **Apply** → **Done** from the Entity creation box

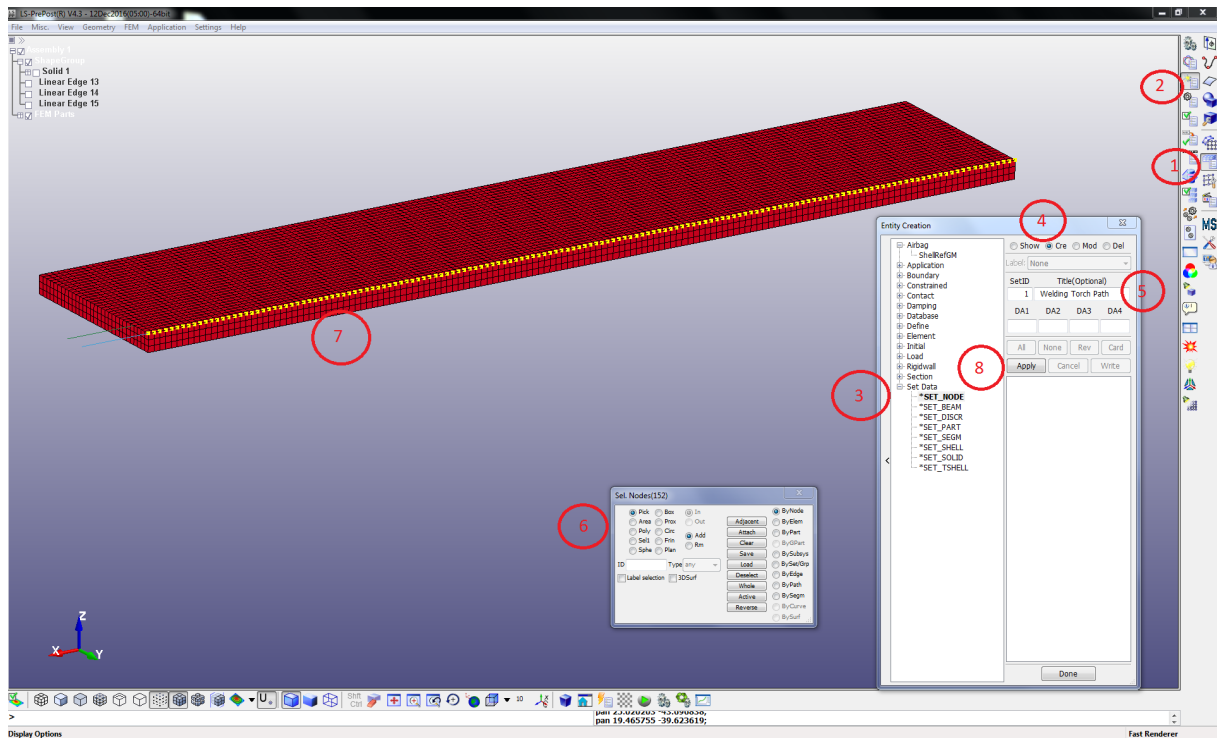


Figure 2.10: Path of the welding line

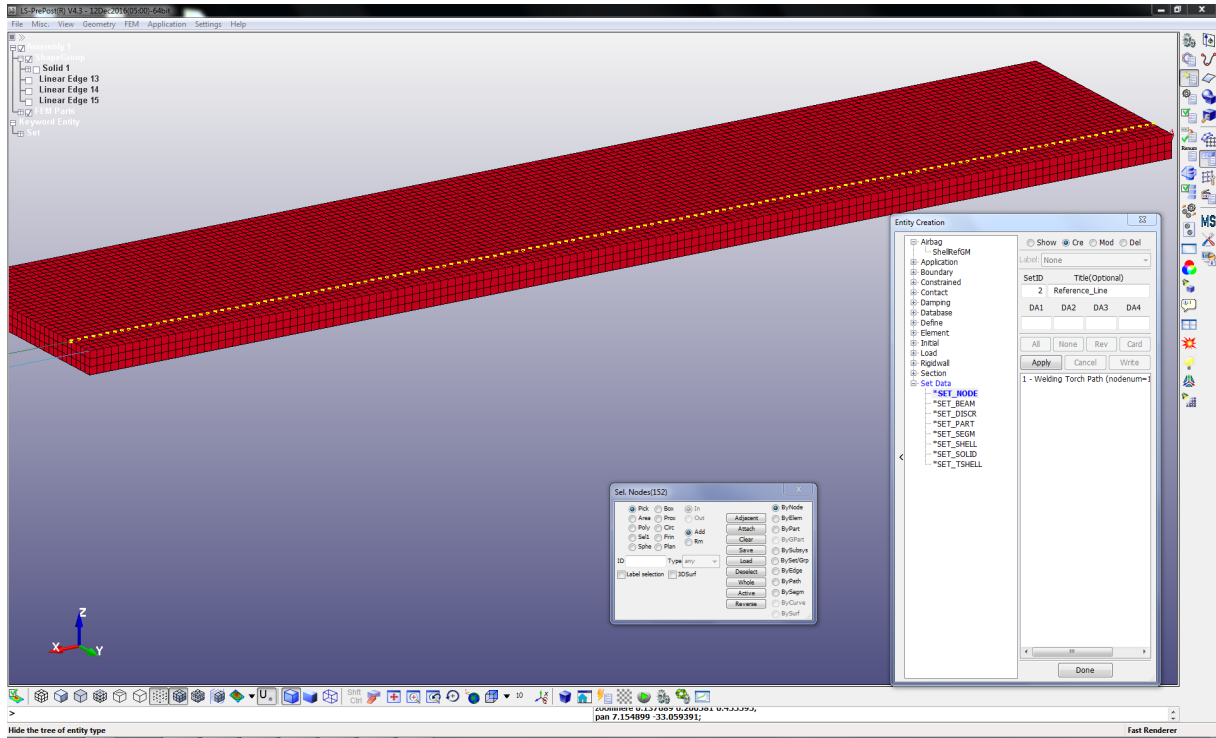


Figure 2.11: Reference line path

## 2.5 Material Models for Welding Simulations

The keyword **\*MAT** is used to define the properties of various materials used in the model. Then, sections are created and materials are assigned to each section. For this simulation we have used the material type 270 (**\*MAT\_CWM**). It is a thermo-elasto-plastic material model based on the VON MISES yield function with isotropic, kinematic or mixed hardening. This material model allows the parameter to be defined as a function of temperature, see Fig. 2.12. This serves as a very simplistic model for the phase change in the microstructure.

The variables, **LCEM**, **LCPR**, **LCAT**, **LCSY** and **LCHR**, in Fig. 2.12 represents the load curve for YOUNG's modulus, POISSON's ratio, thermal expansion coefficient, yield limit and hardening modulus as a function of temperature, respectively, see Fig. 2.13 - 2.17.

The material model **\*MAT\_UHS\_STEEL** and **\*MAT\_PHS\_BMW** are implemented for detailed studies of the material properties in the heat affected zone. This material model contains elaborate phase kinetics models and expect the input of the chemical composition and the thermos-mechanical properties of the individual phases of the microstructure.



```

$=====
$ Mechanical material properties of the plate
$=====
*MAT_CWM_TITLE
Plate
$#      mid      ro      lcem      lcpr      lcsy      lchr      lcat      beta
      17.85000E-9      10001      10002      eghost      pghost      aghost
$# tstart      taend      tlstart      tlend
$# t2phase      t1phase
      800.0      500.0
$=====

```

Figure 2.12: Material property of the plate

```

$=====
$ LCEM: E-Modul in N/mm^2
$=====
*DEFINE_CURVE_TITLE
E-Modul(T)
$#      lcid      sidr      sfa      sfo      offa      offo      dattyp      lcint
      10001      0      1.0      1.0      0.0      0.0      0      0
$#      a1      o1
      24.0      210000.0
      200.0      200000.0
      400.0      175000.0
      600.0      135000.0
      800.0      78000.0
      1000.0      15000.0
      1100.0      7000.0
      1200.0      3000.0
      1300.0      1000.0
      1500.0      1000.0
      100000.0      1000.0
$=====

```

Figure 2.13: YOUNG's modulus as a function of temperature

```

$=====
$ LCPR: Poisson's ratio
$=====
*DEFINE_CURVE_TITLE
Poisson's ratio (T)
$#      lcid      sidr      sfa      sfo      offa      offo      dattyp      lcint
      10002      0      1.0      1.0      0.0      0.0      0      0
$#      a1      o1
      25.0      0.3
      1400.0      0.3
      1500.0      0.3
      100000.0      0.3
$=====

```

Figure 2.14: Poisson ratio as a function of temperature

```

$=====
$ LCAT: Thermal expansion coeff in 1/K
$=====
*DEFINE_CURVE_TITLE
Thermal expansion coeff (T)
$#   lcid   sidr   sfa   sfo   offa   offo   dattyp   lcint
      10005      0   1.01.00000E-6   0.0   0.0      0      0
$#
      a1      o1
      25.0      17.49
      20.0      17.49
      200.0      18.55
      400.0      19.4
      600.0      20.1
      800.0      20.85
      1000.0      21.7
      1200.0      22.35
      1300.0      22.56
      1350.0      19.66
      1500.0      0.0
      1000000      0.0
$=====

```

Figure 2.15: Thermal expansion coefficient as a function of temperature

```

$=====
$ LCSY : Yield point in N/mm^2
$=====
*DEFINE_CURVE_TITLE
Streckgrenze (T)
$#   lcid   sidr   sfa   sfo   offa   offo   dattyp
      10003      0   1.0   1.0   0.000   0.000      0
$#
      a1      o1
      0      240
      100      206
      200      178
      300      160
      400      143
      500      135
      600      125
      700      110
      800      100
      900      85
      1000      60
      1100      30
      1200      15
      1300      5
      1500      5
.

```

Figure 2.16: Yield limit as a function of temperature

```

$=====
$ LCHR : Hardening modul in N/mm^2
$=====
*DEFINE_CURVE_TITLE
Verfestigungsmodulus (T)
$#      lcld      sidr      sfa      sfo      offa      offo      dattyp
      10004      0      1.0      1.0      0.000      0.000      0
$#      a1      o1
      0      688
      100      591
      200      510
      300      459
      400      410
      500      387
      600      358
      700      315
      800      287
      900      243
      1000      172
      1100      86
      1200      5
.

```

Figure 2.17: Hardening modulus as a function of temperature

To define the thermal material property we should use the keyword `*MAT_THERMAL_CWM`. It is a thermal material model with temperature dependent properties, see Fig 2.18.

The variable, **LCHC** and **LCTC**, in fig. 2.18 represents specific heat- and thermal conductivity as function of temperature, respectively, see Fig. 2.19 and Fig. 2.20.

```

$=====
$ Thermal material properties
$=====
*MAT_THERMAL_CWM_TITLE
Plate, Trajectory and reference
$#      tmid      tro      tgrlc      tgmult
      17.85000E-9      0.0      0.0
$#      lchc      lctc      tlstart      tlend      tistart      tiend      hghost      tghost
      10006      10007
$=====

```

Figure 2.18: Thermal property

```

$=====
$ HCLC, LCHC: Specific heat capacity in mJ/(t K) = 1e-06 J/Kg K
$=====
*DEFINE_CURVE_TITLE
Specific heat capacity (T)
$#   lcid   sidr   sfa   sfo   offa   offo   dattyp   lcint
    10006     0     1.0 1000000     0.0     0.0         0         0
$#
    a1         o1
    00.0       430.0
    100.0      500.0
    200.0      550.0
    300.0      580.0
    400.0      610.0
    500.0      650.0
    600.0      710.0
    700.0      790.0
    800.0      865.0
    900.0      965.0
   1440.0     630.0
   2500.0     707.0
$=====

```

Figure 2.19: Specific heat as a function of temperature

```

$=====
$ TCLC, LCTC: Thermal conductivity mW/ mm K = W/ m K
$=====
*DEFINE_CURVE_TITLE
Thermal Conductivity (T)
$#   lcid   sidr   sfa   sfo   offa   offo   dattyp   lcint
    10007     0     1.0     1.0     0.0     0.0         0         0
$#
    a1         o1
    00.0       0.046
    100.0      0.046
    200.0      0.045
    300.0      0.043
    400.0      0.041
    500.0      0.038
    600.0      0.035
    700.0      0.029
    800.0      0.024
   1500.0      0.032
$=====

```

Figure 2.20: Thermal conductivity as a function of temperature

## 2.6 Section

The keyword **\*SECTION** is used to define the element formulation, integration rule, nodal thickness and cross sectional properties.

### 2.6.1 Plate section

After the materials have been created we need to assign these materials to the parts which were previously created. For this purpose we can use GUI as well as keyword file. To assign the plate section with corresponding material we need a section identifier (**SECID**), which must be unique and element formulation (**ELFORM**). Here, we have used the

keyword **\*SECTION\_SOLID**, which define section properties for solid continuum and fluid elements, see Fig. 2.21.

```

$=====
$ Section - Element formulation
$=====
*SECTION_SOLID
$#   secid   elform       aet
      1       1         0
$=====

```

**Figure 2.21:** Section plate

## 2.6.2 Beam section

To define cross sectional properties for beam we should use the keyword **\*SECTION\_BEAM**, see Fig. 2.22.

```

$=====
$ Section Element formulation
$=====
*SECTION_BEAM
$#   secid   elform   shrf   qr/irid   cst   scoor   nsm
      2       1       1.0     2         0     0.0     0.0
$#   ts1     ts2      tt1     tt2     nsloc   ntloc
      1.0     1.0     1.0     1.0     0.0     0.0
$=====

```

**Figure 2.22:** Section beam

## 2.7 Heat source

A weld heat source model by J. GOLDAK is used by LS-DYNA to model the weld heat source. The GOLDAK weld heat source model is based on a Gaussian distribution of power density in space. An important feature of this model is that it uses a double ellipsoidal method of heat deposition such that the size and shape of the energy source can be easily changed to account for different types of welds namely, arc welding, laser and beam welding, and friction stir welding. The energy deposited in the front and rear quadrants are defined by the fractions  $F_f$  and  $F_r$ . The weld shape parameters,  $a$ ,  $b$ ,  $c_f$ ,  $c_r$  are defined using **\*BOUNDARY\_THERMAL\_WELD** keyword in input file, see Fig. 2.23.

```

$=====
$ Trajectory - Heatsource in Ws
$=====
*BOUNDARY_THERMAL_WELD
$#   pid      ptyp      nid      nflag      x0      y0      z0      n2id
$#   1001      2      128956      1      0.0      0.0      0.0      128959
$#   a        b        cf        cr        lcid      q        ff        fr
$#   20      20      15      30      0      5950000      1.086      0.904
$#   pid      ptyp      nid      nflag      x0      y0      z0      n2id
$#   1002      2      128962      1      0.0      0.0      0.0      128965
$#   a        b        cf        cr        lcid      q        ff        fr
$#   20      20      15      30      0      5950000      1.086      0.904
$=====

```

Figure 2.23: The GOLDAK weld heat source

## 2.8 Contact cable

To bring the weld line in contact with plate, we can use the keyword **\*CONTACT\_GUIDE\_CABLE**, see Fig. 2.24.

```

$=====
$ Contact - Cabel: Define a siliding contact that guides beam element through a list of nodes
$=====
*CONTACT_GUIDED_CABLE
$#   nsid      pid      soft      ssfac      fric
$#   1         2         0         1.0         0.0
$#   nsid      pid      soft      ssfac      fric
$#   2         4         0         1.0         0.0
$=====

```

Figure 2.24: Contact

## 2.9 Heat source motion

The heat source movement like, translation and rotation, is defined by using the keyword **\*BOUNDARY\_PRESCRIBED\_MOTION\_RIGID**, see Fig. 2.25. .

```

$=====
$ Provide motion to the heat source
$=====
*BOUNDARY_PRESCRIBED_MOTION_RIGID
$#   pid      dof      vad      lcid      sf      vid      death      birth
$#   2         1         2         4001      1.0      0         0.0         0.0
$#   pid      dof      vad      lcid      sf      vid      death      birth
$#   3         1         2         7001      1.0      0         0.0         0.0
$#   pid      dof      vad      lcid      sf      vid      death      birth
$#   4         1         2         4001      1.0      0         0.0         0.0
$#   pid      dof      vad      lcid      sf      vid      death      birth
$#   5         1         2         7001      1.0      0         0.0         0.0
$=====

```

Figure 2.25: Motion of the weld and reference line

## 2.10 Weld line motion

The weld line is considered to be a rigid body and we can use the keyword **\*MAT\_RIGID** to define its material property, see Fig. 2.26.

```

$=====
$ Rigid bodies properties for the welding torch: Trajectory
$=====
*MAT_RIGID_TITLE
Welding torch
$#      mid      ro      e      pr      n      couple      m      alias
      27.85000E-9  210000.0      0.3      0.0      0.0      0.0
$#      cmo      con1      con2
      1.0      0      7
$#lco or a1      a2      a3      v1      v2      v3
      0.0      0.0      0.0      0.0      0.0      0.0
$=====

```

Figure 2.26: Rigid body property for the weld line

## 2.11 Initial condition

To initialize a particular value we can use the keyword **\*INITIAL**. To define initial nodal point temperature we can use the keyword **\*INITIAL\_TEMPERATURE\_SET**. These initial temperatures are used in a thermal only analysis or a coupled thermal/structural analysis. In Fig. 2.27 we have used the numerical value 0 for **NSID**. Here, 0 represents that we have included all nodes. Instead of that we can also use particular set of nodes.

```

$=====
$ Initial value
$=====
*INITIAL_TEMPERATURE_SET
$#      nsid      temp      loc
      0      20.0      0
$=====

```

Figure 2.27: Initial condition

## 2.12 Thermal boundary condition

For a thermal analysis we need to apply a convection boundary condition on a **SEGMENT** or **SEGMENT\_SET**. For this we can use the keyword **\*BOUNDARY\_CONVECTION\_SET** (This card is used to specify the convection coefficients to a group of element segments), see Fig. 2.28.

The variable **HLCID** represents convection heat transfer coefficient,  $h$ .

```

$=====
$ Boundary conditions - Thermal
$=====
*BOUNDARY_CONVECTION_SET
$#  ssid
    1
$#  hlcid  hmult  tlcid  tmult  loc
    -51    1.0    0    20.0    0
$=====
$
$=====
*DEFINE_CURVE_TITLE
Radiation and Convection
$#  lcid  sidr  sfa  sfo  offa  offo  dattyp
    51    0    1.0  1.0e-03  0.000  0.000  0
$#
    a1    o1
$-10000,20.0
0,20.00
100,20.00
200,20.00
300,20.00
400,20.00
500,20.00
600,20.00
700,20.00
800,20.00
1000,20.00
1250,20.00
1500,20.00
1750,20.00
2000,20.00
50000,20.00
$=====

```

Figure 2.28: Thermal boundary condition

## 2.13 Curve

The keyword **\*DEFINE\_CURVE** is used to define trajectory and reference line travel distance, see Fig. 2.29.

```

$=====
$ Define curve for Trajectory
$=====
*DEFINE_CURVE_TITLE
Trajectory x (t)
$#  lcid  sidr  sfa  sfo  offa  offo  dattyp  lcint
    4001    0    1.0  1.0  0.0  0.0  0  0
$#
    a1    o1
    0.0    0.0
    16.7  -310.0
    10000.0 -310.0
*DEFINE_CURVE_TITLE
Trajectory x (t)
$#  lcid  sidr  sfa  sfo  offa  offo  dattyp  lcint
    7001    0    1.0  1.0  0.0  0.0  0  0
$#
    a1    o1
    0.0    0.00
    16.7  -310.00
$=====

```

Figure 2.29: Curve definition for weld and reference line



## 2.14 Parameter

This option provides a way to specifying numerical values of parameter names that are referenced throughout the input file. The parameter definitions, if used, should be placed at the beginning of the input file following **\*KEYWORD**. Here, we have assign numerical values to time step and heat source velocity, see Fig. 2.30.

```
$=====
$ Parameter: Define the numerical values of parameter names referenced throughout the inout file
$=====
*PARAMETER
R      dt 0.1      R      v 6.0
$=====
```

Figure 2.30: Parameter

## 2.15 Part group

By using the keyword **\*SET\_PART\_LIST** we can create part set ID, which can be used for GOLDAK double ellipsoid heat source, see Fig. 2.31.

```
$=====
$ Part Groups
$=====
*SET_PART_LIST_TITLE
Heat source Trajectory
$#      sid      da1      da2      da3      da4      solver
      1001      0.0      0.0      0.0      0.0MECH
$#      pid1      pid2      pid3      pid4      pid5      pid6      pid7      pid8
      1          2          3          0          0          0          0          0
*SET_PART_LIST_TITLE
Heat source Trajectory
$#      sid      da1      da2      da3      da4      solver
      1002      0.0      0.0      0.0      0.0MECH
$#      pid1      pid2      pid3      pid4      pid5      pid6      pid7      pid8
      1          4          5          0          0          0          0          0
$=====
```

Figure 2.31: Part Group

## 2.16 Implicit analysis

Thermal problems in LS-PrePost<sup>®</sup> are solved using implicit time integration. Therefore, there is no stability condition on the thermal time step. Much larger time steps can be used for the thermal solution as opposed to the mechanical solution, which uses explicit time integration. For implicit analysis we need to use the following keywords, see Fig. 2.32 and Fig. 2.33:

**\*CONTROL\_SOLUTION**: This card is used to specify the analysis solution procedure if thermal only or coupled thermal analysis is performed.

**\*CONTROL\_START:** This keyword define the start time of analysis.

**\*CONTROL\_TERMINATION:** Used to terminate the job.

**\*CONTROL\_IMPLICIT\_GENERAL:** This keyword is necessary to activate implicit analysis and define associated control parameters, like time step. This keyword is required for all implicit analyses.

```

=====
$ Solver
=====
$ Purpose: To specify the analysis solution procedure
$-----x-----x-----x-----
*CONTROL_SOLUTION
$#   soln      nlq      isnan      lcint
      2         0         0         100
$-----x-----x-----x-----
$Purpose:Define the start time of analysis
$-----x-----x-----x-----
*CONTROL_START
$#  BEGTIM
      0.00
$-----x-----x-----x-----
$ Purpose: Stop the job
$-----x-----x-----x-----
*CONTROL_TERMINATION
$#  endtim  endcyc      dtmin      endeng      endmas
      30.0      0      0.000      0.000      0.000
.

```

Figure 2.32: Solver for thermal analysis

**\*CONTROL\_IMPLICIT\_SOLUTION:** This keyword is used to specify whether a linear or nonlinear solution is desired.

**\*CONTROL\_IMPLICIT\_DYNAMICS:** The purpose of this keyword is to activate implicit dynamic analysis and define time integration constants.

**\*CONTROL\_IMPLICIT\_AUTO:** This card is used to define parameters for automatic time step control during implicit analysis.

**\*CONTROL\_THERMAL\_NONLINEAR:** By using this keyword we can set parameters for a nonlinear thermal or coupled structural/thermal analysis.

**\*CONTROL\_THERMAL\_TIMESTEP:** This card is used to define the timestep parameters for the thermal only analysis.

```

$-----x-----x-----x-----x-----
$ Purpose: Activate implicit analysis and is required for all implicit analysis
$-----x-----x-----x-----x-----
*CONTROL_IMPLICIT_GENERAL
$# imflag 1 dt0 2 inform 2 nsbs 1 igs 2 cnstn 0 form 0 zero_u 1
$-----x-----x-----x-----x-----
$ Purpose: To Specify whether a linear or nonlinear solution is desired
$-----x-----x-----x-----x-----
*CONTROL_IMPLICIT_SOLUTION
$# nsolvr 12 ilimit 1 maxref 15 dctol 0.05 ectol 100.0 rctol 0.1 lstol 0.9 abstol
$# dnorm 2 diverg 2 istif 1 nlprint 1 nlnorm 1 d3itctl 1
$# arcctl 0 arcdir 0 arcclen 0.0 arccth 1 arcdmp 2 ARCPST ARCALF ARCTIM
$-----x-----x-----x-----x-----
$ Purpose: Activate implicit analysis
$-----x-----x-----x-----x-----
*CONTROL_IMPLICIT_DYNAMICS
$# IMASS 0 GAMMA BETA TDBIR TDYDTH TDYBUR IRATE
$-----x-----x-----x-----x-----
$ Purpose : Define parameters for automatic time step control during implicit analysis
$-----x-----x-----x-----x-----
*CONTROL_IMPLICIT_AUTO
$# IAUTO 1 ITEOPT 16 ITEWIN 10 DTHIN 0.001 DTHMAX -111 DTEXP KFAIL KCYCLE
$-----x-----x-----x-----x-----
$ Purpose: Set parameters for a nonlinear thermal or coupled structural/thermal analysis
$-----x-----x-----x-----x-----
*CONTROL_THERMAL_NONLINEAR
$# refmax 60 tol 0.0 dcp 0.5 lumpbc 0 thlstl 0.1 nlthpr 1 phchpn 0.0
$-----x-----x-----x-----x-----
$ Purpose: Set options for the thermal solution in a thermal only or coupled structural thermal analysis
$-----x-----x-----x-----x-----
*CONTROL_THERMAL_SOLVER
$# atype 1 ptype 1 solver 11 cgltol 1.0E-4 gpt 8 eqheat 1.0 fwork 1.0 sbc 0.0
$# msglvl 0 maxitr 50 abstol 0 reltol 0 omega 1.0 unused 0 tsf 1.0
$-----x-----x-----x-----x-----
$ Purpose: Set time step control for the thermal solution
$-----x-----x-----x-----x-----
*CONTROL_THERMAL_TIMESTEP
$# ts 0 tip 1.00 its 1 tmin 0.000 tmax 0.000 dtemp 5000 tscp 0.50 lcts 112
$=====

```

Figure 2.33: Solver for implicit thermal analysis

## 2.17 Output file

The keyword **\*DATABASE** with a combination of options is used for controlling the output of **ASCII** databases and binary files output. With this keyword the frequency of writing the various databases can be determined.

The keyword **\*DATABASE\_BINARY** is necessary to obtain output files containing results information, see Fig. 2.34. The keyword **\*DATABASE\_TPRINT** gives thermal output from a coupled structural/thermal or thermal only analysis.

The keyword **\*DATABASE\_BINARY** is used for binary output option and the attribute **D3PLOT** represent the database for entire model. If **\*DATABASE\_BINARY\_D3PLOT** is not specified in the keyword file then the

output interval for d3plot is automatically set to  $1/20^{\text{th}}$  the termination time.

```

$=====
$ Postprocessing
$=====
*DATABASE_IPRINT
$#      dt      BINARY      LCUR      IOOPT      OPTION1      OPTION2      OPTION3      OPTION4
      &dt      1      0      0      0
*DATABASE_BINARY_D3PLOT
$#      dt      lcdt      beam      npltc      psetid
      &dt      0      0      0      0
$#      ioopt
      0
$=====

```

Figure 2.34: Binary output file

By using the keyword **\*DATABASE\_EXTENT\_BINARY** we may control to some extent the content of binary output databases.

```

$=====
$ Postprocessing: Structural mechanics
$=====
*DATABASE_EXTENT_BINARY
$#      neiph      neips      maxint      strflg      sigflg      epsflg      rltflg      engflg
      11      11      3      1      1      1      1      1
$#      cmpflg      ieverp      beamip      dcomp      shge      stssz      n3thdt      ialemat
      0      0      0      1      1      1      2      1
$#      nintsld      pkp_sen      sclp      unused      msscl      therm      intout      nodout
      0      0      1.000000      0      0      2ALL      ALL
$#      dtdt      resplt
      1      1
.

```

Figure 2.35: Extended binary output file

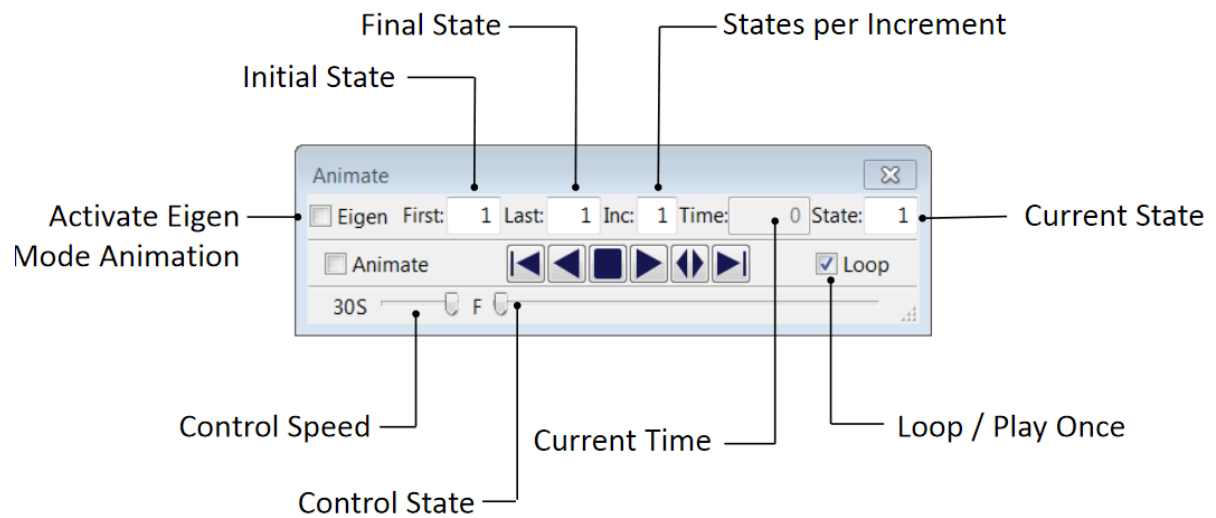
### 3 Postprocessing

The results generated from the analysis can be enormous so it requires additional processing which is termed postprocessing. LS-PrePost<sup>®</sup> GUI is used for postprocessor. It supports the latest Open-GL standards to provide fast rendering for fringe plots and animation results. It also handles the **ASCII** output data and links it to the input files and animations.

LS-PrePost<sup>®</sup> also include features like animations, contour plots, X-Y graphs, overlay plots, vector plots, ASCII plotting, mesh manipulation, *etc.*.

#### 3.1 Animation interface

The Animation controls are displayed when d3plots are loaded. If the animation interface is closed, it can be restored by clicking the **Anim** render button, see Fig. 3.1.



**Figure 3.1:** Animation interface

The results generated from the analysis can be viewed in LS-PrePost<sup>®</sup>, see Fig. 3.5:

1. From the right toolbar, select the **Fringe component** option
2. In the **Fringe component** dialog box, choose the option **Misc**
3. Choose the variable of your interest, here it is temperature
4. Use the animation interface, see Fig. 3.1, to animate the temperature field

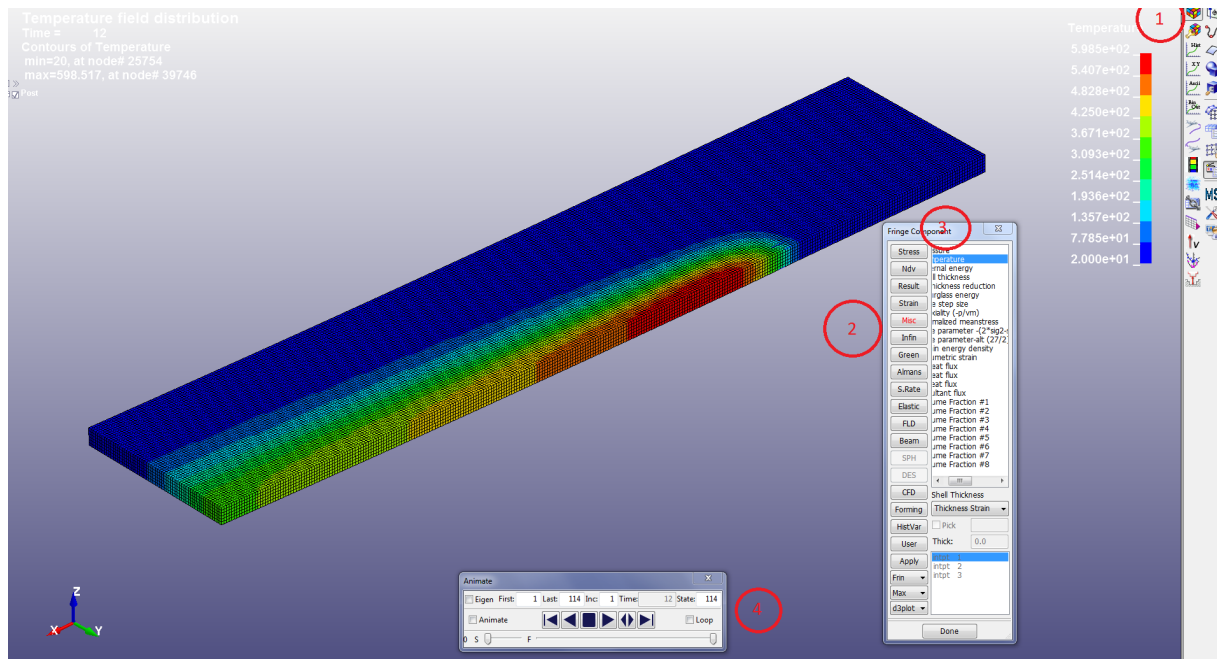


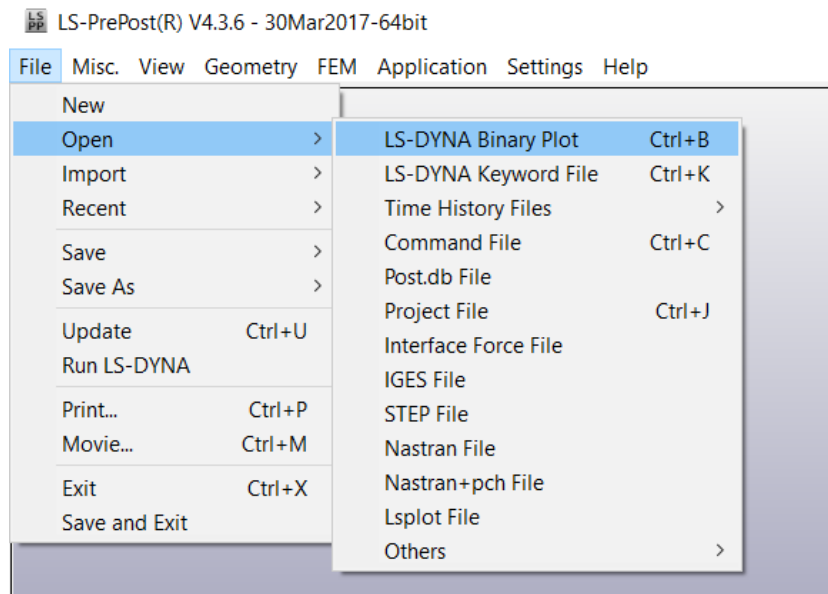
Figure 3.2: Animation interface

### 3.2 Plot

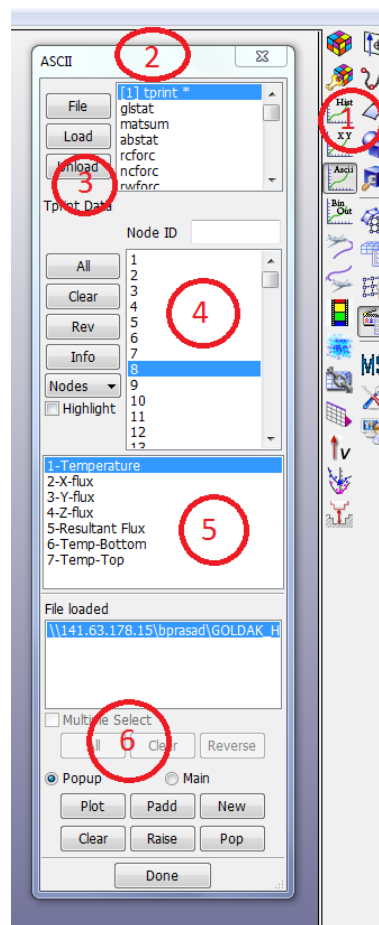
The temperature vs time plot can be created as shown in Fig. 3.4 - 3.5:

1. In the menu bar, select **File** → **Open** → **LS – DYNA Binary Plot**, see Fig. 3.3
2. From the right toolbar, click the **ASCII** option
3. Choose the option **tprint\*** for temperature data
4. In the next step load the computed temperature data by clicking the option **Load**
5. Select the node at which you want to know the temperature distribution
6. Click the variable of interest, here temperature
7. At the bottom of the dialog box click the option **Plot** and this will generate the temperature field vs time graph as shown in Fig. 3.5.

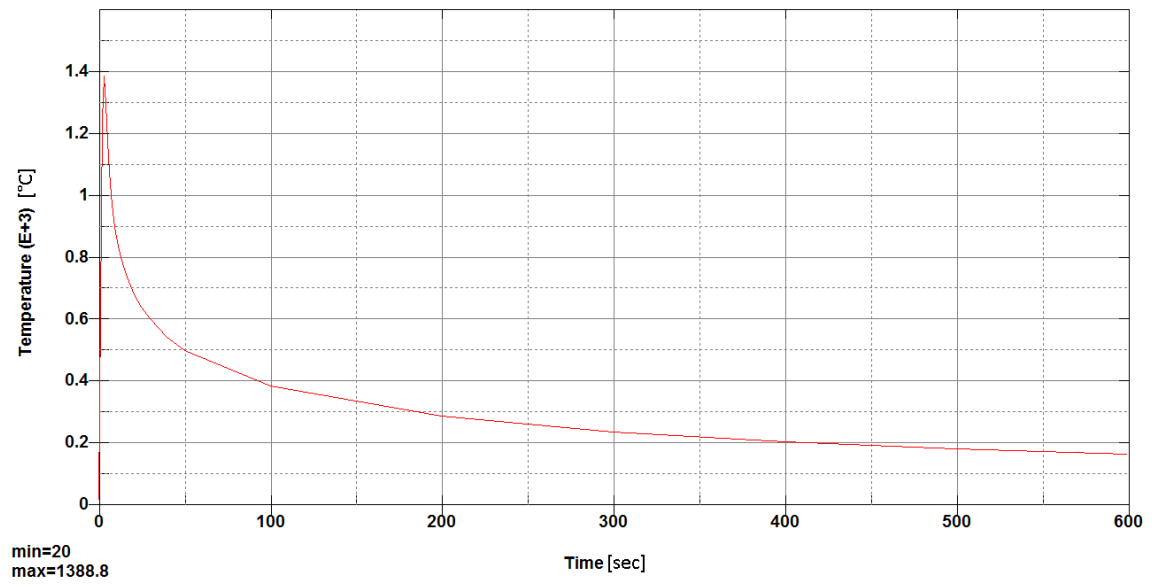
Here, not all the modules of postprocessing are explained. One can explore different modules and try to generate plots for different parameters and get meaningful results out of them.



**Figure 3.3:** Sub menu to open the binary file



**Figure 3.4:** ASCII dialog box



**Figure 3.5:** Temperature vs time graph