True*Grid*® User's Manual

A Guide and a Reference

by

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VOLUME 2:

Geometry, Assembly, Global Properties, and Output

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I. Geometry Commands

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1. 2D Curves

These commands define and modify 2D curves. All 2D curves use a 2 dimensional local coordinate system where **x′** is the abscissa (horizontal axis) and **z′** is the ordinate. A 2D curve can be rotated around the **z′**-axis of the curve's local coordinate system to form a surface of rotation. You can choose how to orient this local **z′**-axis in the global coordinate system (see the **sd** command with the option **crx**, **cry**, **crz**, or **cr)**. A 2D curve can be extruded or lofted in the third direction to make an infinite surface (see the **sd** command using the **cp** option). Two 2D curves can be placed into the global 3D coordinate system and a ruled surface formed between them (see the **sd** command using the **rule2d** option). 2D curves can form the cross sections of a surface (see the **sd** command using the **swept** option). For more details, see the **sd** command, and the surface dictionary. You can embed a 2D curve in the global 3D coordinate system or turn the 2D curve into a 3D curve (see the **curd** command using the **ld2d3d** option).

Some 2D curves are special purpose and cannot be used to form 3D geometry. For example, load curves are used, when defining boundary conditions, to regulate a time dependent variable.

ld initialize a 2D curve

ld *2D_curve_# segments*

where a

2D_curve_#	number of the curve to be initialized
segment	2D curve segment from the 2D Curve Segment Dictionary

Remarks

You may use as many segments as you like; each is appended to the end of the previous one until the next 2D curve is initialized. Use the automatic parameter %nextld for the next available 2D curve number.

Using this command you construct a curve by appending curve segments to each other, end-to-end. The **ld** command initializes a new 2D curve definition. All subsequent 2D curve segments append to the end of this 2D curve. Once a new 2D curve is created, the previous 2D curve can no longer be appended. Many kinds of curve segments are available, and you may combine them in almost any order. For details, see the 2D Curve Segment Dictionary. This method of constructing 2D curves is fashioned after the way a draftsman might draw a complex 2D curve.

These 2D curves can be used to form 3D curves and surfaces. For example, a 2D curve can be rotated about any axis of symmetry, extruded indefinitely in any direction, or combined with another 2D curve to produce a ruled surface between the two 2D curves. See the **sd** command.

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Example

```
ld 1 lp2 10.7 -2 11.4 -2;
ld 2 lp2 1 0;
lfil 0 10.7 -2 90 1;
lp2 10.7 -2;
ld 3 lp2 1 0 0 2.7 3.05 5.75;
ltp 3.75 7.5 2.55;
lp2 3.75 8.75;ltp 2.8 9.9 1.45;
lp2 0 11.2 0 14 8 13
  8.35 12.75;
ld 4 lp2 11.4 -2;
lfil 90 7.3 .6 0 1;
lfil 180 8.4 2.3 90 1.45;
lp2 8.4 2.3;
ld 5 lep .5 .35 6.3525 5.4475
  -180 -90 45;
lep .5 .8 8.75 2.65 90 180 45;
ld 6 lp2 6 5.09393 5.3 5.8;
ltp 4.65 7.5 2.6;
lp2 4.65 8.7;ltp 6.9 9.95 1.5;
lp2 7.45 9.6;ltp 8.8 9.4 2.5;
ltp 9.45 10.05 .65;
lp2 9.45 10.6 9.3 11.05 9.2
  11.05 9.05 10.6 9.05 10.2;
ltp 8.7 10.15 .25;lp2 7.55 10.85;
ltp 7.55 12.15 .8;lp2 8.35 12.75;
```

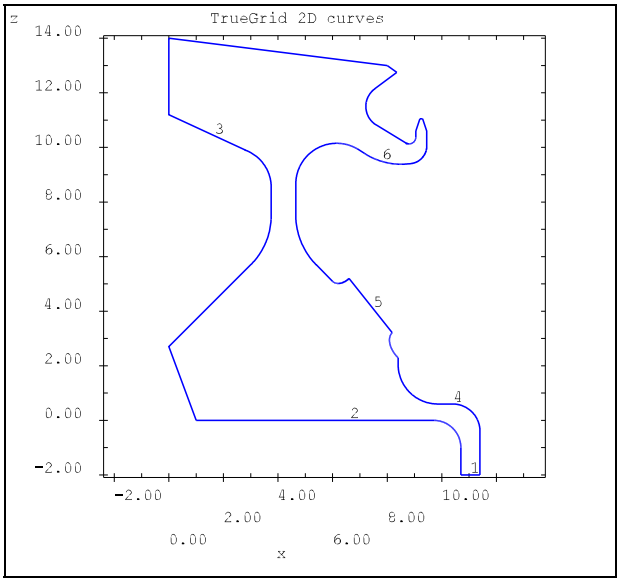

Figure 1 Complex 2D Geometry

apld append curve segments

apld *segments*

where a *segment* any one of the curve segments described in the 2D Curve Segments Dictionary

Remarks

A 2D curve is initialized with an **ld** command with options. Then this command can be used to append other 2D curve segments to the end of it. This command is used primarily by the GUI (graphical user interface) so that you have a command to click on in the menus when you want to interactively append another 2D curve segment to the previously initiated 2D curve. Otherwise, the **apld** command does nothing.

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Examples

The following two curve definitions are equivalent:

```
ld 1 lp2 1.6 5 1.6 4;
lat 1.6 4 5 4 .2;
lat 2.8 4 3 5 .2;
```

and

```
ld 1 lp2 1.6 5 1.6 4;
apld lat 1.6 4 5 4 .2;
apld lat 2.8 4 3 5 .2;
```

lcc concentric arcs

lcc x' z' θ_{begin} θ_{end} $radius_1$ $radius_2$ ... $radius_n$;

where

(x',z')	coordinates of the center of the circle
θ_{begin}	angle which defines the beginning of the arc
θ_{end}	angle which defines the end of the arc
$radius_1$ . . . $radius_n$	radii of the various arcs

Remarks

The 2D curves are numbered sequentially, beginning with the last-defined 2D curve number, plus one.

Example

```
lcc 1 2 45 135 1 4 9 16 25
36;
```

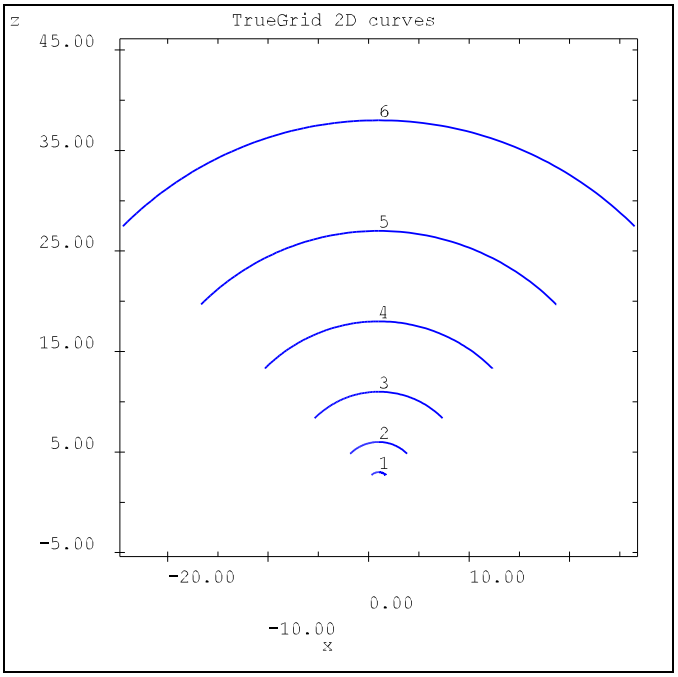

Figure 2 Concentric Arcs

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ckl check curvature

ckl *2D_curve$_1$ 2D_curve$_2$ angle*

where

2D_curve$_1$	number of the first curve to be checked
2D_curve$_2$	number of the last 2D curve to be checked
angle	is the greatest permissible angular deviation from 180 degrees.

Remarks

This command will check each defined 2D curve whose number lies in the specified range. The angle at each point of the polygonal line approximation to the curve is calculated and. reported when this angle differs from 180 degrees by more than the specified angular deviation, *angle*. Then the specified curves are displayed..

Example

```
ld 1 lp2 0 0 1 1 2 1 2 0;
ckl 1 1 60
```

The above example will cause the following warning message:

warning - curve 1 at node 3 has an angle of 90.00 degrees

gset set 2D Curves window

gset *xc′ zc′ size*

where

(*xc′, zc′*)	coordinates of the center of the 2D Curves window
size	length of the square 2D Curves window

Remarks

The default center and size of the 2D Curves window is based on the 2D curves being drawn. This window is made just large enough to contain the selected curves for display while remaining a square. In order to return to this default, use a size of 0.

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mazt intersection tolerance

mazt *tolerance*

Remarks

The **mazt** command controls the tolerance for the intersection of the 2D curves. The default is .001. This tolerance is used in the **lpil** 2D curve segment type.

lcd load curve definition

lcd *ld_curve_# options curve ;*

where

ld_curve_#	load curve number

where an *option* can be

sidr *type*	
where *type* can be:	
0	transient analysis or other
1	stress initialization only
2	stress initialization and transient analysis
sfa *factor*	time (abscissa) scale factor
sfo *factor*	amplitude (ordinate) scale factor
offa *offset*	time (abscissa) offset
offo *offset*	amplitude (ordinate) offset
dattyp *type*	
where *type* can be	
0	monotonic time (abscissa)
1	non-monotonic time (abscissa)
g	modal damping viscous dampers
crit	modal critical damping
q	modal damping quality factor
d1	frequency/time dependency dynamic load
d2 *offset*	frequency/time dependency dynamic load
d3 *offset scale*	frequency/time dependency dynamic load
d4 *offset scale min*	frequency/time dependency dynamic load
m1	temperature dependence
m2 *offset*	temperature dependence
m3 *offset scale*	temperature dependence
m4 *offset scale min*	temperature dependence

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where

offset	offset for amplitude
scale	scale factor for amplitude
min	lower clamp for the amplitude

s1 stress dependent

st material property temperature dependent

rnd1 power spectral density

where a *curve* must be one of

$t_1 f_1 \; t_2 f_2 \; \ldots \; t_n f_n$

lp $t_1 f_1 \; t_2 f_2 \; \ldots \; t_n f_n$

where (t_i, f_i) are the pairs of time and load amplitudes

func *npnts* u_0 u_x *load_expression*

where

npnts	number of points in curve
u_0	minimum value of the independent variable u
u_x	maximum value of the independent variable u
load_expression	algebraic expression with independent variable u

Remarks

The load curve number is used for identification and is unique for each load curve. The automatic parameter %nextlc has the next available load curve number.

Load curves are usually for dynamic simulations, variable properties, or tables. They are used differently by different simulation codes. Sometimes a load command, such as **fc**, will require a load curve for completeness. Depending on your output selection, the load curve number you select in these types of commands will be used as a set id and will not require that you define a corresponding load curve with **lcd**.

The easiest way to make a load curve is:

```
lcd 1 0 0 .001 1;
```

which defines load curve 1 with two points. At time 0 the amplitude of the associated load will be 0. Then the amplitude is ramped up linearly to full amplitude at time .001. This is equivalent to:

```
lcd 1 lp 0 0 .001 1;
```

for backward compatibility.

The options **sidr**, **sfa**, **sfo**, **offa**, **offo**, and **dettyp** are for LS-DYNA load curves.

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The options **s1**, **st**, **rnd1**, **g**, **crit**, **q**, **d1**, **d2**, **d3**, **d3**, **m1**, **m2**, **m3**, and **m4** are for NASTRAN and NE/NASTRAN tables.

Example

```
lcd 3 0 0 1 1 2 1 2.00001 0 3 0;
```

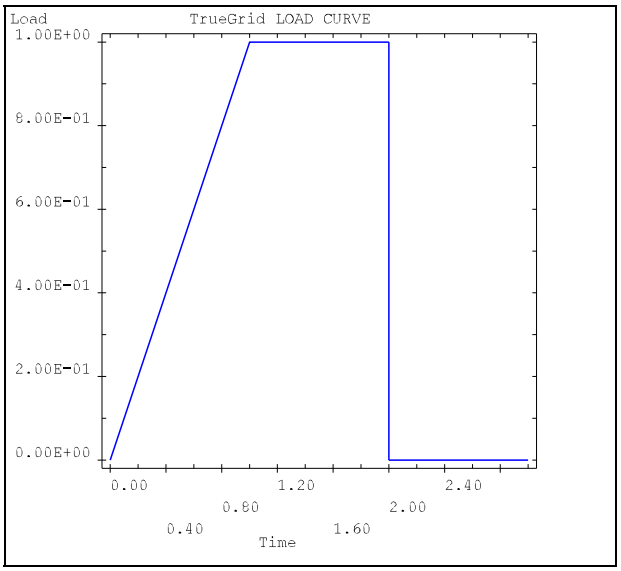

Figure 3 Load Curve Drawn Using **LCV**

Example

```
lcd 2 sidr 0 func 100 0 .001
sin(u*180000);
```

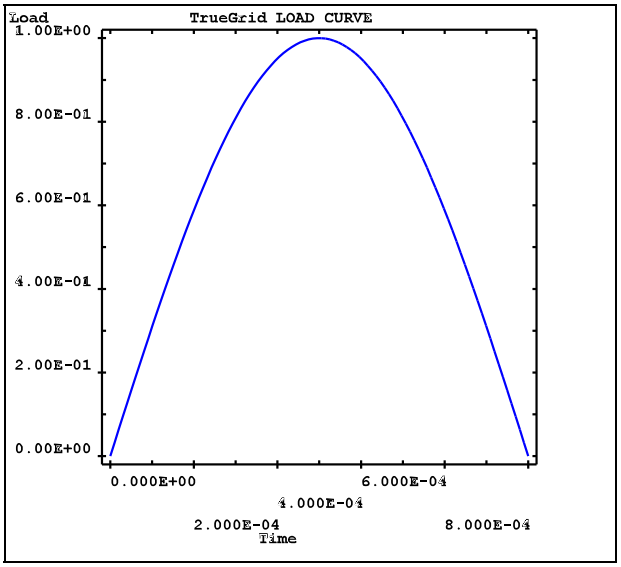

Figure 4 Function load curve

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lcinfo information about load curves

lcinfo *<no arguments>*

Remarks

Prints the number of points of each load curve.

Example

```
flcd 1 phase 180 cosine 30 5 .1 lp .1 1; tinit .1 finit 1
fsca .1 foff 1 phase 0 sine 60 60 .1;
lcd 2 0 0 .1 .04 .4 .08 .9 .12 1.6 .16 2.5 .2;
lcinfo
```

The above example produces 2 load curves. The table below is written to the tsave file and to the text window if the command is issued interactively.

```
  LOAD CURVE DEFINITIONS
load curve number      1 has       91 points
load curve number      2 has        6 points
```

ld3d2d convert 3D curves to 2D curves

ld3d2d *reduct_coord first_2D_curve_# 3D_curve$_1$ 3D_curve$_2$ ... ;*

where

reduct_coord	**x**, **y**, or **z** coordinate by which to reduce the 3D curve to a 2D curve
first_2D_curve_#	is the number to be assigned to the first 2D curve
3D_curve$_i$	are the numbers of the 3D curves

Remarks

The symbols **x**, **y**, or **z** specify the coordinate to be eliminated. The specified 3D curve is projected onto the coordinate plane defined by the remaining two coordinates.

Example

One of the edges of a surface is converted into a 3D curve. The **ld3d2d** command converts the 3D curve into a 2D curve by eliminating the z-coordinate.

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```
sd 1 function 1000 2000 100 2000 u*(cos(v)+v*sin(v));
u*(sin(v)-v*cos(v)); 1000*(u+v*.25);;
curd 1 sdedge 1.2
ld3d2d z 1 1;
```

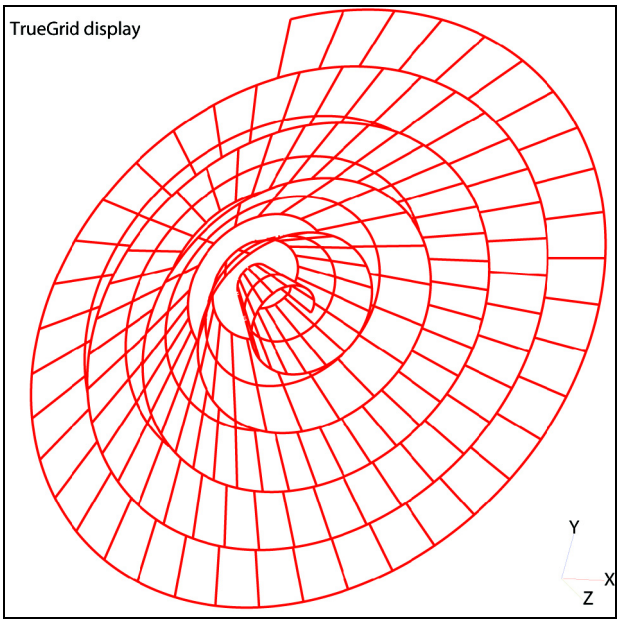

Figure 5 Spiral Surface

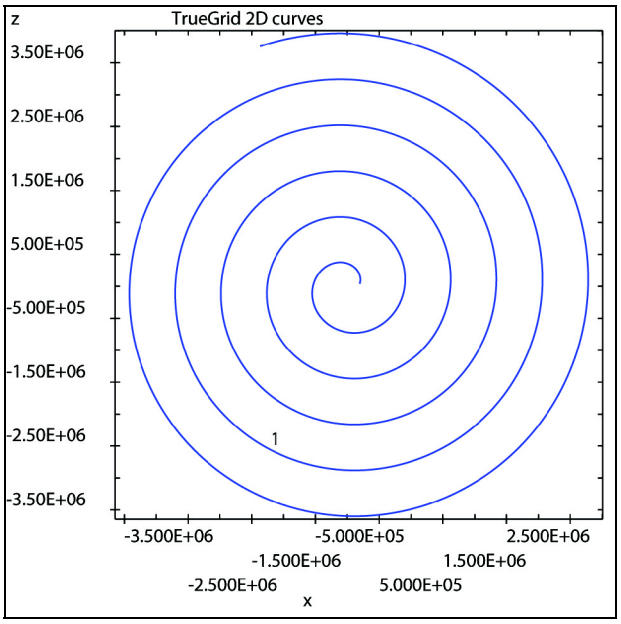

Figure 6 Curve From Spiral Surface

ldinfo information about 2D curves

ldinfo *<no arguments>*

Remarks

A table is printed listing each defined 2D curve. This data is always written to the tsave file. When in the interactive mode, the data is also written to the text window. This table contains a list of 2D curves that have been defined, the number of points used in their definition, and the length of the curve.

Example

For a model with several curves, this command was issued:

```
ldinfo
```

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```
2D CURVE DEFINITIONS
  2D curve definition number   1 has  73 points & length 2.13012
  2D curve definition number   2 has   8 points & length 17.1542
  2D curve definition number   4 has   3 points & length 0.01230
  2D curve definition number   9 has  13 points & length 9.00000
  2D curve definition number  14 has  16 points & length 16.0001
  2D curve definition number  30 has   4 points & length 1.31400
last 2D curve defined is number     14
```

ldprnt print coordinates

ldprnt *2D_curve*

where *2D_curve* is the number of a defined 2D curve.

Remarks

The **x**′ and **z**′-coordinates forming the polygonal approximation to a 2D curve are printed to the tsave file, and, when the command is issued interactively, to the text window.

Example

```
ldprnt 30

       2D Curve Definition       30
            2.812500E+00   3.843750E+00
            2.812500E+00   3.875000E+00
            2.937500E+00   3.875000E+00
            2.937500E+00   3.843750E+00
```

lrl rays

lrl x'_{center} z'_{center} *length* $angle_1$ $angle_2$ ... $angle_n$;

where

(x'_{center}, z'_{center})	is the point that the rays emanate from,
length	is the length of each ray, and
$angle_i$	is the angle formed by the ***i***-th ray and the **x**′-axis.

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Remarks

The 2D rays are numbered sequentially, beginning with one more than the number of the last 2D curve that was defined.

Example

```
lrl 1 2 3 0 5 15 30 50 70 85 95
100;
```

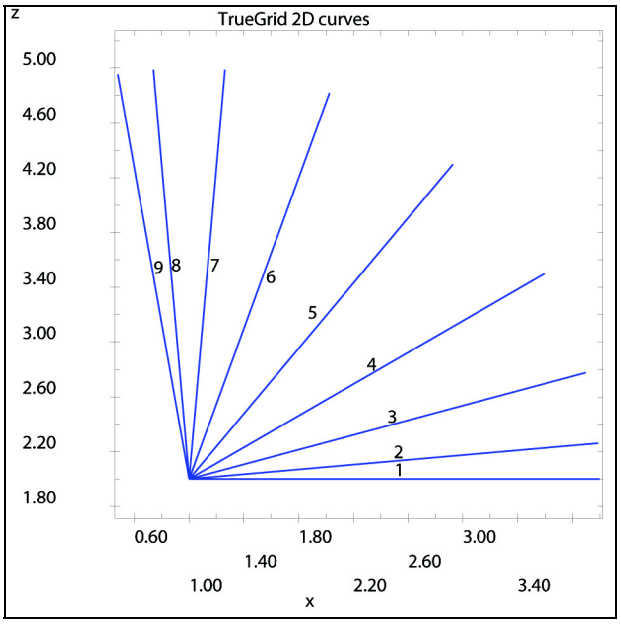

Figure 7 Radial Lines

lrot rotate

lrot *2D_curve angle*

where

2D_curve is the number of the 2D curve and

angle is the amount of rotation in degrees.

Remarks

This function rotates an existing 2D curve where a positive angle is counter-clockwise. See the following page for an example.

Example

```
ld 1 lp2 1 0;lap 3 -.1 2 2;lp2 5
-.1 5 .1 3 .1;
lap 1 .1 2 2.1;lp2 1 0;
ld 2 lp2 1 0;lap 3 -.1 2 2;lp2 5
-.1 5 .1 3 .1;
lap 1 .1 2 2.1;lp2 1 0;
lrot 2 45
```

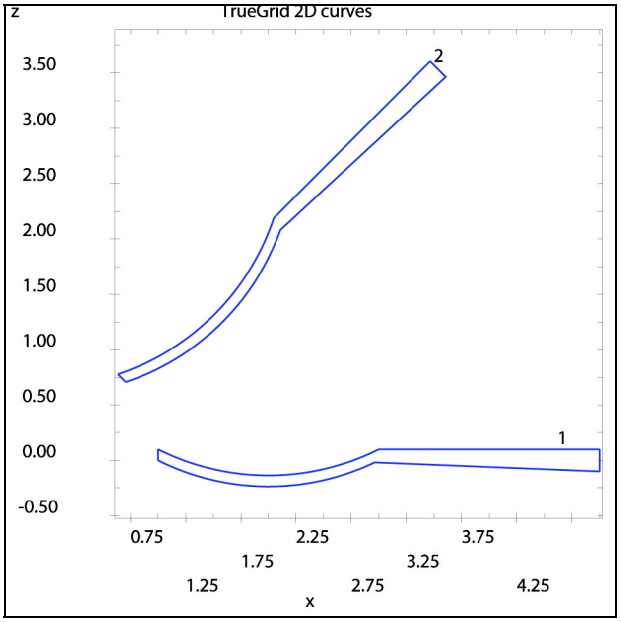

Figure 8 Rotated 2D curve

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lsca scale

lsca *2D_curve scale*

where

2D_curve	number of the 2D curve
scale	scale factor

Remarks

Each coordinate of the curve are multiplied by the value of *scale*.

Example

```
ld 1 lp2 1.5 0;lfil 0 2 1 -75 .2;lfil
-75 [cos(30)] [sin(30)] 30 .2 lfil 30
[cos(15)] [sin(15)] 105 .2 ;lfil 105
1.5 0 0 .2;lp2 1.5 0;
ld 2 lstl 1 1.525 .385;lsca 2 .5
```

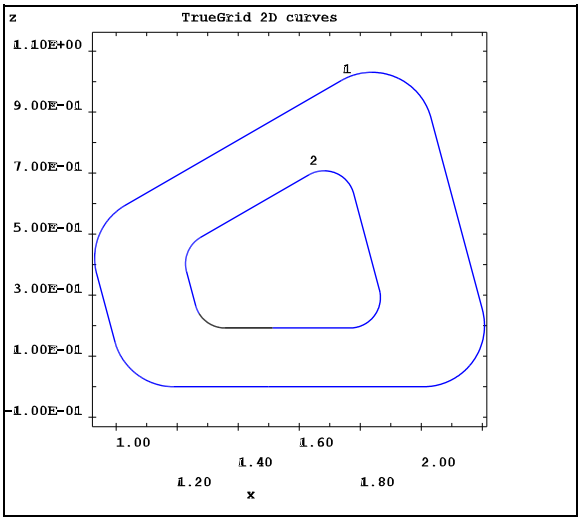

Figure 9 Scale Both Coordinates

lscx scale first coordinate

lscx 2D_curve *x'_scale*

where

2D_curve	number of the 2D curve, and
x'_scale	scale factor

Remarks

Each first coordinate on the curve is multiplied by *x'_scale*.

Example

```
ld 1 lp2 0 0;lad 1 0 360 ;
ld 2 lp2 0 0;lad 1 0 360;lscx 2 3
```

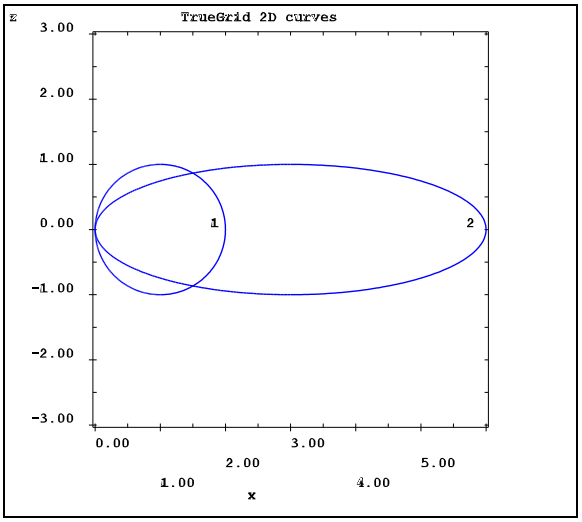

Figure 10 Scale First Coordinate Only

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lscz scale second coordinate

lscz *2D_curve z′_scale*

where

2D_curve number of the 2D curve

z′_scale scale factor.

Remarks

Each second coordinate on the curve is multiplied by *z′_scale*.

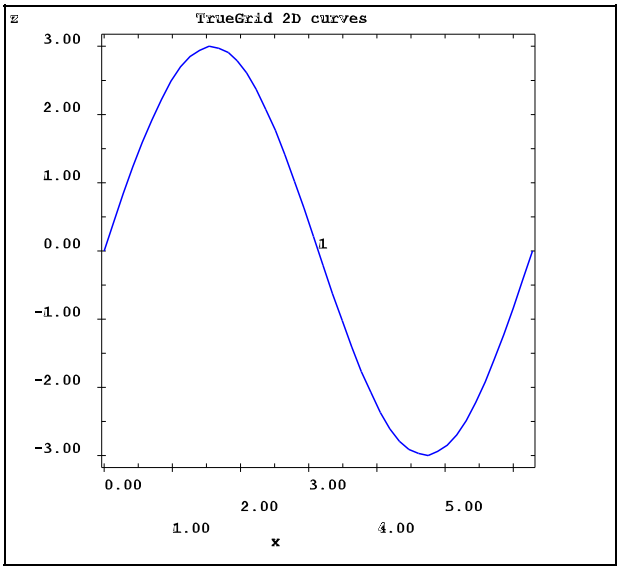

Figure 11 Scale Second Coordinate Only

Example

```
ld 1 lp2 0 0 .14 .14 .28 .28 .42
.41 .56 .53 .7 .64 .84 .74
.98 .83 1.12 .9 1.26 .95 1.4 .98
1.54 1 1.68 .99 1.82 .97  1.95.93
2.09 .87 2.23 .79 2.37 .69 2.51 .59
2.65 .47  2.79 .34 2.93 .21 3.07
.07 3.21 -.07 3.35 -.21 3.49 -.34
3.63 -.47 3.77 -.59 3.91 -.69 4.05
-.79 4.19 -.87 4.33 -.93 4.47 -.97
4.61 -.99 4.75 -1 4.89 -.98 5.03
-.95 5.17 -.9 5.31 -.83 5.45 -.74
5.59 -.64 5.72 -.53 5.86 -.41 6 -.28 6.14 -.14 6.28 0;lscz 1 3
```

lt translate

lt *2D_curve Δx′ Δz′*

where

2D_curve is the number of a previously defined 2D curve, and

(*Δx′*,*Δz′*) is the translation to be applied.

Remarks

This is different from the **lstl** option of the **ld** command because **lt** modifies an existing 2D curve. See the example on the following page.

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Example

```
ld 1
lp2 0 0 .14 .14 .28 .28 .42 .41 .56
.53 .7 .64 .84 .74
.98 .83 1.12 .9 1.26 .95 1.4 .98
1.54 1 1.68 .99 1.82 .97
1.95 .93 2.09 .87 2.23 .79 2.37 .69
2.51 .59 2.65 .47
2.79 .34 2.93 .21 3.07 .07 3.21
-.07 3.35 -.21 3.49 -.34
3.63 -.47 3.77 -.59 3.91 -.69 4.05
-.79 4.19 -.87 4.33
-.93 4.47 -.97 4.61 -.99 4.75 -1
4.89 -.98 5.03 -.95 5.17
-.9 5.31 -.83 5.45 -.74 5.59 -.64
5.72 -.53 5.86 -.41 6
-.28 6.14 -.14 6.28 0;
lscz 1 3
lt 1 1.570796 0
```

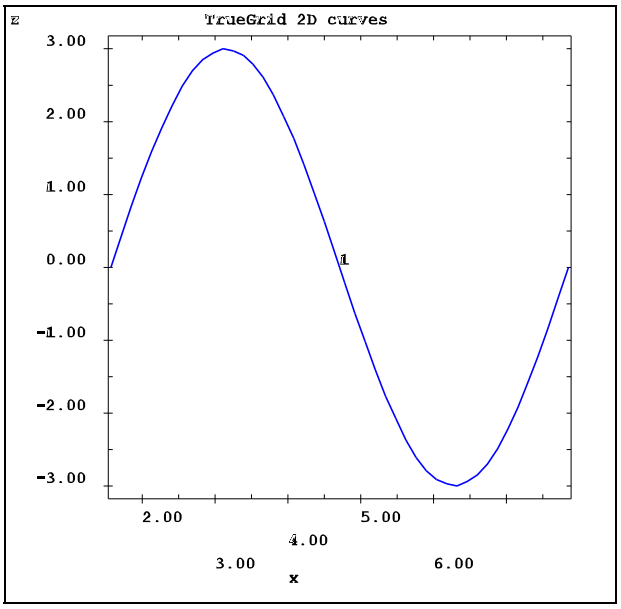

Figure 12 Translate Existing 2D Curve

flcd piece-wise trigonometric load curve

flcd *ld_curve_# options_1 options_2 ;*

where

ld_curve_# load curve number

where an option_1 can be any set of

SIDR type

where type can be:

0 for transient analysis or other

1 for stress initialization only

2 for stress initialization and transient analysis

SFA abscissa_scale_factor

SFO ordinate_scale_factor

OFFA abscissa_offset

OFFO ordinate_offset

DATTYP type

where type can be

0 for monotonic abscissa

1 for non-monotonic absciss

where an *option*_2 can be:

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tinit *initial_time*	specify the initial time
tsca *time_scale*	scale the time variable
toff time_offset	translate time variable
finit initial_load	specify the initial load
fsca load_scale	scale the load variable
foff load_offset	translate the resulting curve after scaling
phase angle	specify the phase for a **sine** or **cosine** curve
lp t_1 f_1 t_2 f_2 ... t_n f_n ;	append a polygonal line or curve
sine #_points #_cycles_per_time time	append a sine curve, computed as the sine of the scaled parameter plus the phase
cosine #_points #_cycles_per_time time	append a cosine curve, computed as the cosine of the scaled parameter plus the phase.

Remarks

You specify a number of curve segments, with the **lp**, **sine**, and **cosine** options. The whole load curve is constructed by appending them; each curve begins where the last one ended.

Use the automatic parameter %nextlc for the next available load curve number.

The initial time, **tinit**, and the initial load, **finit**, are global values. Their last values are the offsets applied to the entire curve. The variables **toff**, **tsca**, **foff**, and **fsca** can be set for each segment. The time variable of a segment is relative to that segment and is assumed to start at zero for each segment.

In the following discussion, the sequences of time and function values are denoted by T_i and F_i respectively. The index i is incremented by 1 for each new value in the sequence. For the polygonal line, you specify the time and function values t_i and f_i. Any transformations you specified would change these to the actual time and function values T_i and F_i. The curve is linear between them. For each trigonometric curve, you specify the number of points N, the number of cycles per unit time *cps*, and the time interval Δt covered. For each new i value, t_i is incremented by $\delta t = \Delta t/N$. The value of the cosine or sine function gives f_i. In the equations below, **T0** is the untransformed time variable at the end of the last segment and is initially 0.

Thus, after taking into account the other parameters you may specify – **tinit**, **tsca**, **toff**, **finit**, **fsca**, **foff**, and **phase** – the time and function values are:

For a polygonal line (lp):

T_i = tsca * (t_i + T0) + toff + tinit

F_i = fsca * f_i + foff + finit

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For a sine curve (sine):

T_i = tsca * (t_i + T0) + toff + tinit

F_i = fsca * sin(*cps* * (t_i + T0) + phase) + foff + finit

For a cosine curve (cosine):

T_i = tsca * (*ti* + T0) + toff + tinit

F_i = fsca * cos(*cps* * (t_i + T0) + phase) + foff + finit

Example

```
flcd 1 phase 180 cosine 30 5 .1
lp .1 1;tinit .1 finit 1
fsca .1 foff 1 phase 0 sine 60 60
.1;
```

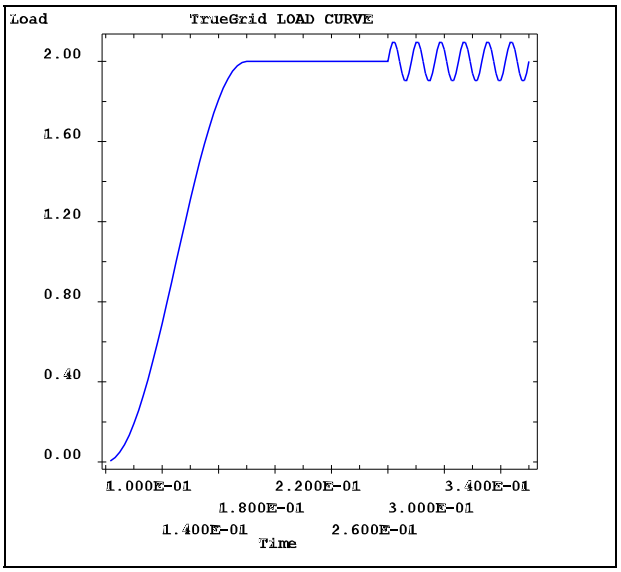

Figure 13 Oscillating Load Curve

it2dc intersect two 2D curves

it2dc *curve$_1$ curve$_2$*

where *curve$_1$ curve$_2$* are the two 2D-curve numbers

Remarks

This command creates the list of intersections of two 2D curves. All points of intersection will be determined and stored in automatic parameters.

%n2dint will be set to the number of intersections.
%x2dci will be set to the x-coordinate for every point of intersection.
%y2dci will be set to the y-coordinate for evey point of intersection.
%rarcl1 will contain the relative position of the corresponding intersection point along the first curve.
%rarcl2 will contain the relative position of the corresponding intersection point along the second curve.

The first 2D curve in the argument list is searched from start to finish for intersections with the second curve. This order is preserved in all four arrays.

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edgefile identify file used for edge data

edgefile *file_name*

Remarks

The **rln**, **rlns**, and **rseg** commands retrieve data from this file in order to define 2D curves. This feature is archaic but may still be useful. The file format consists of curve segment data in sets of coordinates as follows. Each set of coordinates begins with a header record with a fixed format:

The first field (I5) is the number of coordinate pairs to follow.

The second field (I5) is the coordinate system flag: 0 for Cartesian coordinates, 1 for the $\mathbf{x}'$ and $\mathbf{z}'$-coordinates to be switched, and 2 for polar coordinates.

The third field (I5) is the reflection flag: 0 for no reflection, 1 to reflect the 2D curve, or 2 to reflect the 2D curve and also keep the unreflected 2D curve.

The fourth field (E10.0) is the Cartesian coordinate scale factor.

The fifth field (E10.0) is the Cartesian coordinate translation in the second coordinate.

The sixth field (E10.0) is the constant second Cartesian coordinate to be used to reflect the 2D curve.

The seventh field (E10.0) is the Cartesian coordinate translation in the first coordinate.

Thereafter are two fields (2E10.0) per record giving each coordinate pair. For Cartesian coordinates, they are interpreted as $\mathbf{x}'$ and $\mathbf{z}'$-coordinates in the $\mathbf{x}'\mathbf{z}'$-plane. In polar, they are the angle (in degrees) and radius, respectively. Translations, scaling, and reflections are all done in Cartesian coordinates.

The E10.0 fields may be in either floating-point or exponential Fortran format.

This file may contain any number of these 2D curve definitions.

Example

The following edge file example has 5 curve segments. It is referred to in the examples for the **rln**, **rlns**, and **rseg** commands.

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```
   4    0    0 1.0        0.0        0.0        0.0
1.0        1.0
1.5        0.9
2.0        1.0
2.1        1.7
   5    0    0 1.0        0.0        0.0        0.0
2.0        2.0
1.7        2.2
1.0        2.0
0.8        1.7
1.0        1.0
   9    0    0 1.0        2.0        0.0        2.0
1.0        1.0
1.5        0.9
2.0        1.0
2.1        1.7
2.0        2.0
1.7        2.2
1.0        2.0
0.8        1.7
1.0        1.0
   4    0    2 1.0        2.0        3.8        0.0
1.0        1.0
1.5        0.9
2.0        1.0
2.1        1.7
   4    2    1 1.0        0.0        1.0        2.0
135.0      1.5
140.0      1.5
140.0      2.0
145.0      2.1
```

Line number 22 begins the definition of another curve segment with 4 points.

```
   4    0    2 1.0        2.0        3.8        0.0
```

The second number, 0, indicates that the following 4 records will contain (x', z') Cartesian coordinates. The next integer, 2, says the curve segment will be reflected across the line at $z=3.8$, specified in the second to last field. It also means that the unreflected curve segment will be included in the segment definition. The next number is 1.0 which means that the coordinates will not be scaled. The fifth field specifies that each $\mathbf{z}'$-coordinate will be translated by 2.0. The last field, 0.0, specifies that the $\mathbf{x}'$-coordinate will not be translated.

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rln import an edge file 2D curve

rln *(no arguments)*

Remarks

Prior to using this command, you must use the **edgefile** command to specify the file which contains the curve data. This command can then be used repeatedly. Each time this command is used, the next 2D curve number is assigned to the next 2D curve segment data found in this file. When all of the data in this file has been used, then the file is automatically closed.

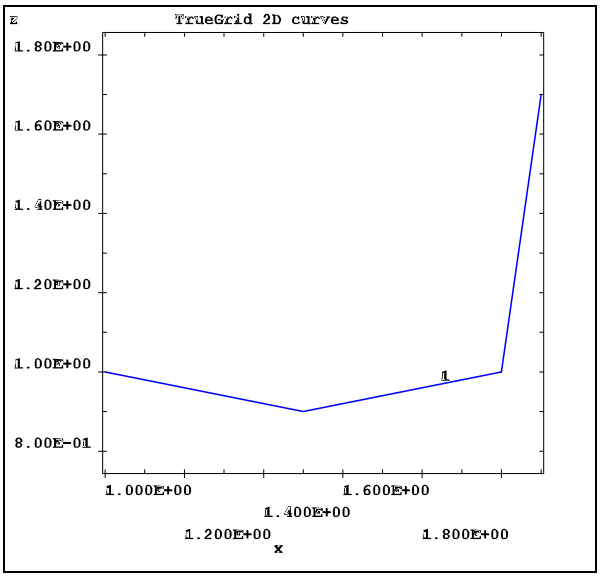

Figure 14 First curve segment

Example

The example edge file in the description on the **edgefile** command above can be converted to 2D curve definitions using the **rln** command. For example, the first curve can be defined with the command:

```
rln.
```

rlns import all edge file 2D curves

rlns *(no arguments)*

Remarks

Prior to using this command, you must use the **edgefile** command to specify the file which contains the curve data. The next 2D curve number is assigned to the next 2D curve segment data found in this file. This is repeated until all of the 2D curve segments have be converted to numbered 2D curve definitions. The edge file is then automatically closed.

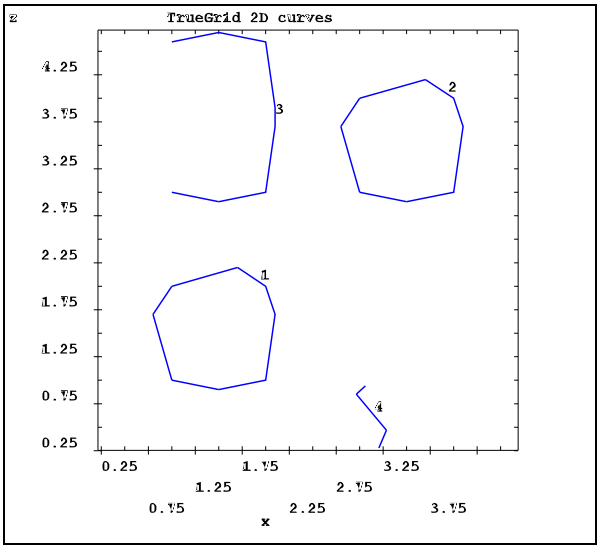

Figure 15 Remaining segments

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Example

The example edge file in the description on the **edgefile** command above can be converted to 2D curve definitions using the **rlns** command. For example,

```
rln rseg rlns
```

2. 2D Curve Segment Dictionary

This section describes all the 2D curve segment types. You can use them to construct composite 2D curves with such commands as **ld** and **apld**.

lp2 polygonal line - pairs

lp2 $x'_1\ z'_1\ x'_2\ z'_2 \ldots x'_n\ z'_n\ ;$

Remarks

lp is the same command, for historical reasons.

Example

```
ld 1 lp2
0 [21/32] .0625 [20.8/32] .125
[20.4/32] .1875 [20/32]
.25 [19.3/32] .3125 [18/32] .375
[16.5/32] .4375 [15/32]
.5 [12.5/32] .5625 [9.6/32] .625
[6/32] .6875 [3.2/32]
.75 [1.8/32] .8125 [1/32] .875
[.2/32] .9375 0;
```

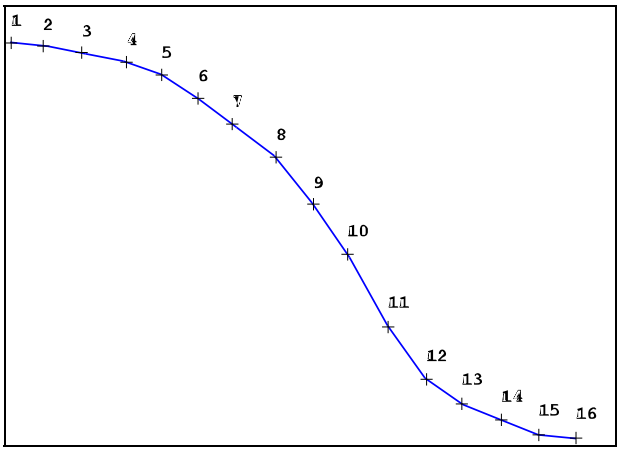

Figure 16 Polygonal line with 16 points

lq polygonal line - lists

lq $x'_1\ x'_2 \ldots x'_n\ ;\ z'_1\ z'_2 \ldots z'_m\ ;$

where

$x'_1\ x'_2 \ldots x'_n\ ;\ z'_1\ z'_2 \ldots z'_m\ ;$ are the two coordinate lists.

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Remarks

This command takes a list of first coordinates, followed by a list of second coordinates. If one list is shorter than the other, it is automatically extended. See the picture above for the example that follows.

Example

```
ld 1 lq 0.0000 0.0625 0.1250 0.1875 0.2500 0.3125 0.3750 0.4375
0.5000 0.5625 0.6250 0.6875 0.7500 0.8125 0.8750 0.9375;
[21.0/32.0]  [20.8/32.0]  [20.4/32.0]  [20.0/32.0]  [19.3/32.0]
[18.0/32.0]  [16.5/32.0]  [15.0/32.0]  [12.5/32.0]  [9.6/32.0]
[6.0/32.0] [3.2/32.0] [1.8/32.0] [1.0/32.0]
[0.2/32.0] [0.0/32.0];
```

lpil intersect 2 curves

lpil *curve$_1$ curve$_2$*

where *curve$_1$ curve$_2$* are the 2 2D-curve numbers

Remarks

This command is used to append the intersection point of two previously defined 2D curves. The 2D curve number 3, shown in the example on the next page, uses this feature. The **mazt** is used to specify the tolerance for this intersection. The coordinates are stored in the automatic parameters %xprj and %zprj.

This also creates the list of intersections of the two 2D curves. All points of intersection will be determined and stored in automatic parameters.

%n2dint will be set to the number of intersections.
%x2dci will be set to the x-coordinate for every point of intersection.
%y2dci will be set to the y-coordinate for evey point of intersection.
%rarcl1 will contain the relative position of the corresponding intersection point along the first curve.
%rarcl2 will contain the relative position of the corresponding intersection point along the second curve.

The first 2D curve in the argument list is searched from start to finish for intersections with the second

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curve. This order is preserved in all four arrays.

Example

```
ld 1 lp2
0 [21/32] .0625 [20.8/32] .125
[20.4/32] .1875 [20/32]
.25 [19.3/32] .3125 [18.0/32.0]
.375 [16.5/32] .4375 [15/32]
.5 [12.5/32] .5625 [9.6/32] .625
[6/32] .6875 [3.2/32]
.75 [1.8/32] .8125 [1/32] .875
[.2/32] .9375 [0/32] ;
ld 2 lstl 1 0 0;lrot 2 45 lt 2
.6 -.4
ld 3 lp2 1 1;lpil 1 2;
```

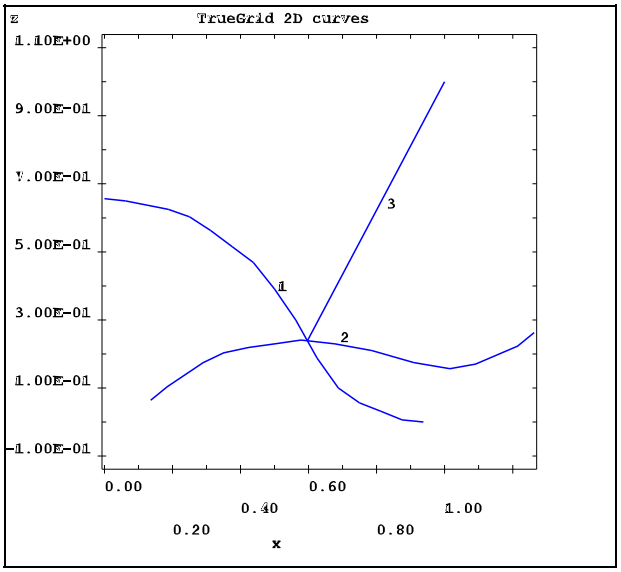

Figure 17 Intersection of 2 curves

lpta tangent to a circle

lpta $x'_0\ z'_0\ radius$

where

(x'_0, z'_0)	is the center of the circle, and
radius	is the circle's radius, with the sign indicating which side of the circle is to be used.

Remarks

If the *radius* > 0, then the point of tangency is found by moving counterclockwise around the circle from the point formed by the intersection of the circle with the line segment extending from its center to the curves previous point. If the *radius* < 0, then the point is found by moving clockwise. The current 2D curve must have at least one point before this option can be used.

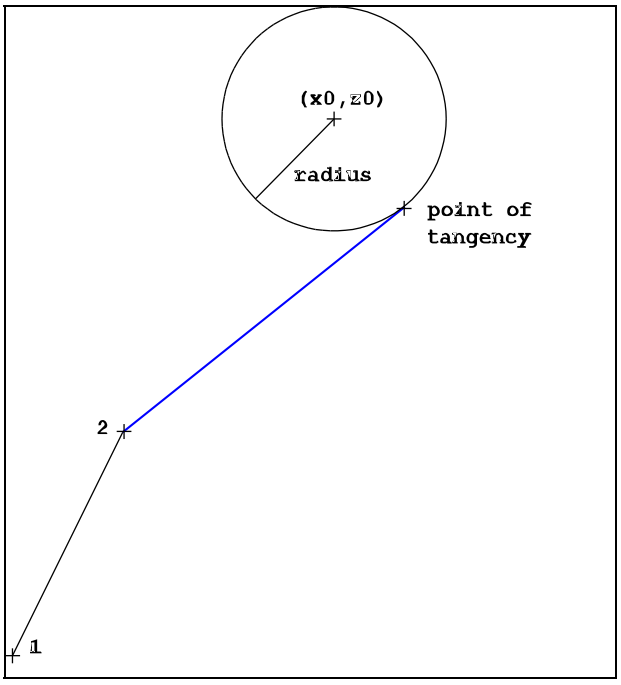

Figure 18 Counterclockwise rotation

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Examples

Compare the examples on the following 2 pages. The first has a positive radius, the second has a negative radius.

```
ld 1 lp2 .1 .2 1.1 2.2; lpta 3 5 1;
```

Example

```
ld 1 lp2 .1 .2 1.1 2.2; lpta 3 5
-1;
```

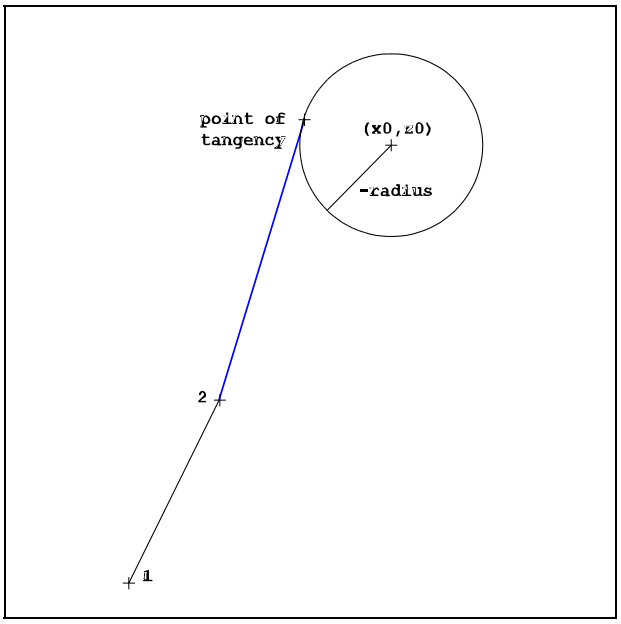

Figure 19 Clockwise rotation

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ltas tangent to 2 circles

ltas $x'_0\ z'_0\ flag\ x'_1\ z'_1\ radius$

where

(x'_0, z'_0)	is the center of the first circle,
flag	is 1 for a clockwise arc and -1 for a counter-clockwise arc,
(x'_1, z'_1)	is the center of the second circle, and
radius	is the radius of the second circle.

Remarks

This command appends an arc of a circle joined to a line tangent to two circles. For the first circle, you only specify the center (x'_0, z'_0); the radius is determined by the requirement that it pass through the last point of the current 2D curve. First an arc of the first circle is appended from the last point of the current 2D curve to a line tangent to both circles. This arc travels around the first circle clockwise if *flag* is -1, counterclockwise if 1. Then the tangent line segment from the first circle to the second circle is appended. You specify the center (x'_1, z'_1) and radius of the second circle; the magnitude of *radius* is the radius, and the sign determines which of the possible tangent lines is chosen, where the same convention is used as in the previous command, **lpta**.

The current 2D curve must have at least one point.

There are generally four lines tangent to two circles. The line is chosen that ensures that the 2D curve's direction will be continuous as we go from an arc of the first circle to the line and then around the second circle (although no arc of the second circle is actually included in the current 2D curve). So *flag*, by determining the direction on the first circle, selects two possible lines. The sign of *radius* determines a direction on the second circle, hence selects one line from those two. For *radius*, a negative number chooses counter-clockwise, and a positive number clockwise.
The four examples that follow show each of the four cases.

Examples

```
ld 1 lp2 1.75 .875 2 1;ltas 2.5 1.3 %r 4 1.9 %s;
```

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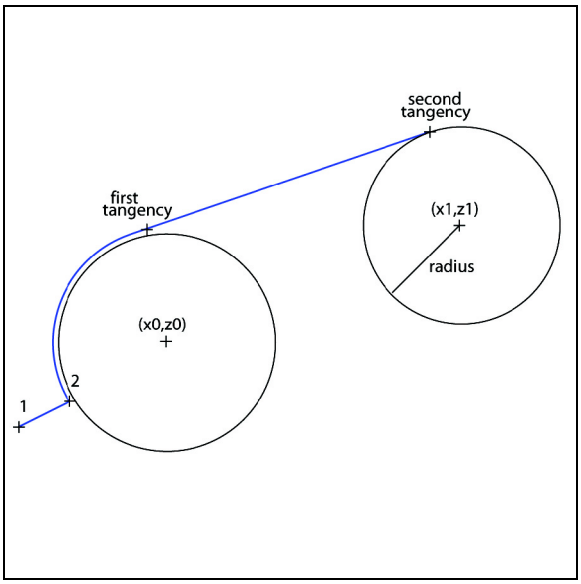

Figure 20 r = -1 , s = .5

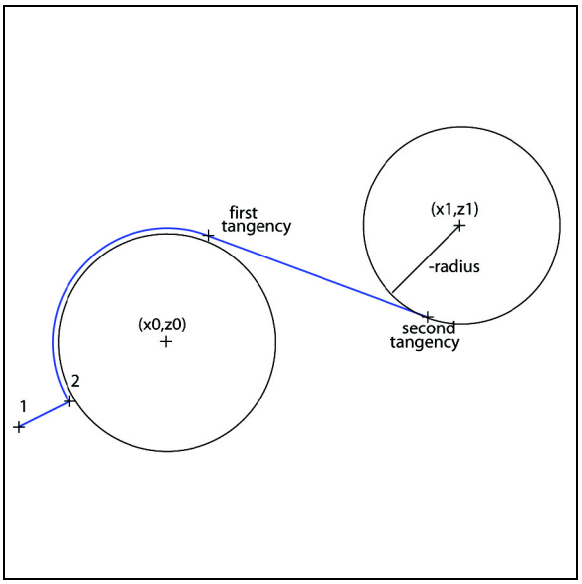

Figure 21 r = -1 , s = -.5

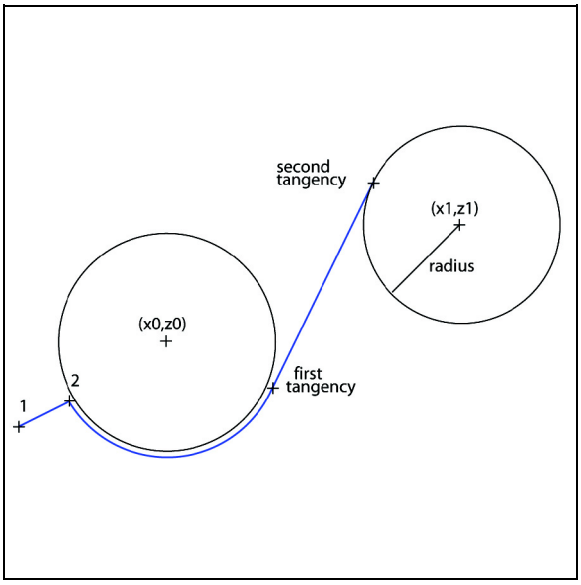

Figure 22 r = 1 , s = .5

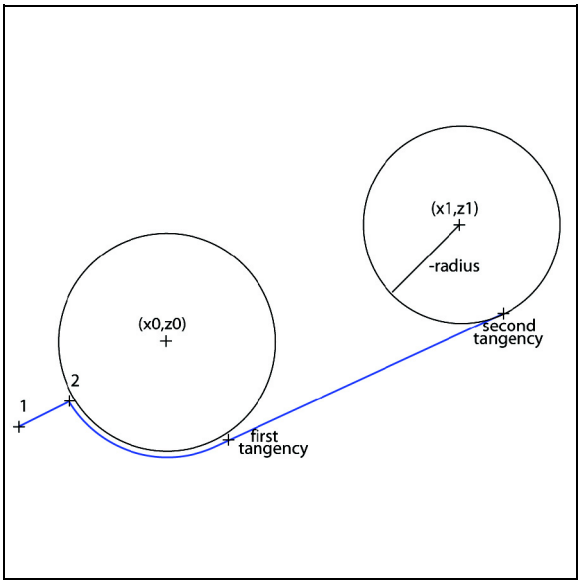

Figure 23 r = 1 , s = -.5

lep elliptic arc

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lep $radius_1$ $radius_2$ x'_0 z'_0 θ_{begin} θ_{end} φ

where

$radius_1$	is the length of the semi-major axis,
$radius_2$	is the length of the semi-minor axis,
(x'_0, z'_0)	is the center of the ellipse,
θ_{begin}	is the beginning angle of the elliptical arc,
θ_{end}	is the ending angle of the elliptical arc, and
φ	is the angle between the major axis and the positive **x**-axis.

Remarks

The ellipse is centered at (x'_0, z'_0). The first radius is measured along the axis forming an angle of φ degrees with the x-axis of the 2 dimensional local coordinate system, and is referred to as the major axis. The second radius is measured along the axis orthogonal to the first axis, and is referred to as the minor axis. θ_{begin} and θ_{end} are the beginning and ending angles of the arc. θ_{begin} is the angle between the major axis and the line from the beginning point to the center of the ellipse. θ_{end} is the angle between the major axis and the line from the ending point to the center of the ellipse. (The same side of the major axis is used for all three angles). See the example on the following page.

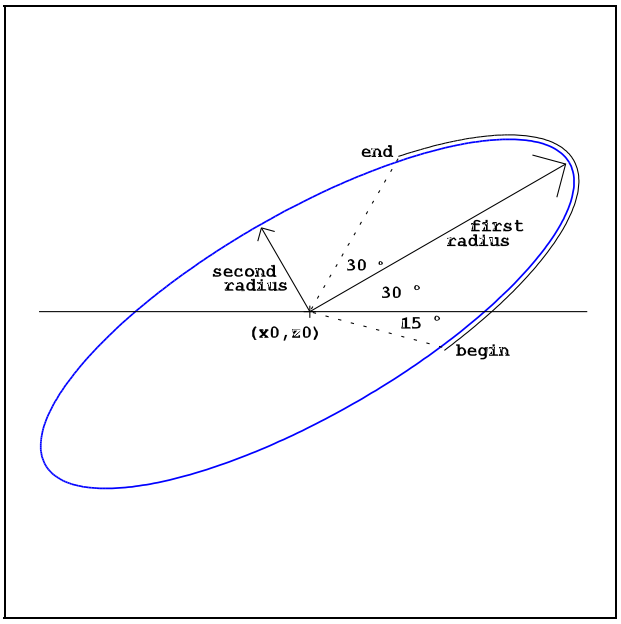

Figure 24 Elliptic Arc

Example

```
ld 2 lep 3 1 1 1 -45 30 30;
```

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lod normal offset

lod *curve offset*

where

curve	is the ID of a previously-defined 2D curve, and
offset	is the distance that this curve is offset in a normal direction.

Remarks

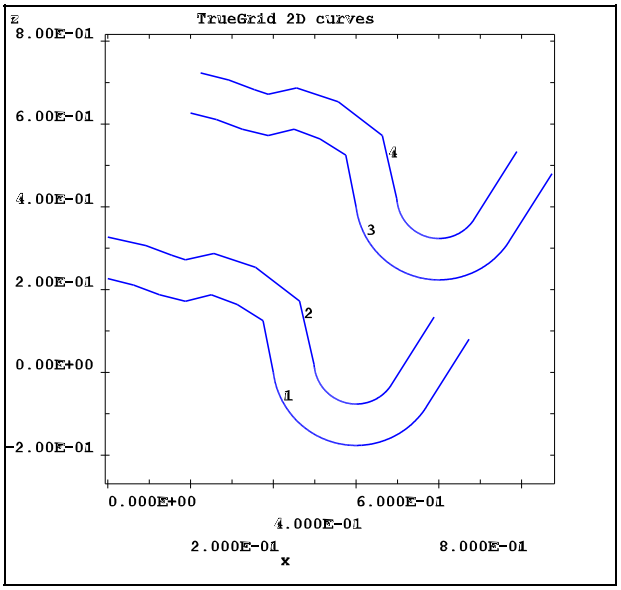

Figure 25 Normal Offset Of 2D Curves

Use this option to take a previously defined 2D *curve* and *offset* it in a normal direction. Positive normals are on the left side of the curve as you move from the beginning to the end. The normals at the endpoints are calculated independently. If an endpoint of the previously defined 2D curve meets the z′-axis, then the z′-axis is used as the normal direction at that endpoint, causing the offset curve to meet the z′-axis as well. The *offset* is used with the normal directions to determine the endpoints of the new segment. The central difference is used for the normal direction for the interior points. See the next page for an example. In this example, the first curve is on the z′-axis, causing the offset curve 2 to also be on the z′-axis. Since curve 3 is not on the z′-axis, the offset curve 4 is treated differently than curve 2 at the endpoints.

Example

```
ld 1 lp2 0 .2266 .0625 .2109 .125 .1875 .1875 .1719 .25 .1875 .3125
.1641 .375 .125 0.4 0; ltas .6 .025 1 1 0 .15;
ld 2 lod 1 .1 ;
ld 3 lstl 1 .2 .4 ;
ld 4 lod 3 .1 ;
```

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lnof normal averaging offset

lnof *curve offset*

where

curve	is a the ID number of a previously-defined 2D curve, and
offset	is the offset distance.

Remarks

Use this option to take a previously defined 2D curve and offset in a direction using the normal averaging method. For each point in the original curve, both the left and right sided normals are calculated. Then they are averaged. This average normal is used to calculate the direction of the offset. In the example below, compare curves 2 and 3 offset from curve 1 using the **lnof** and **lod** options, respectively.

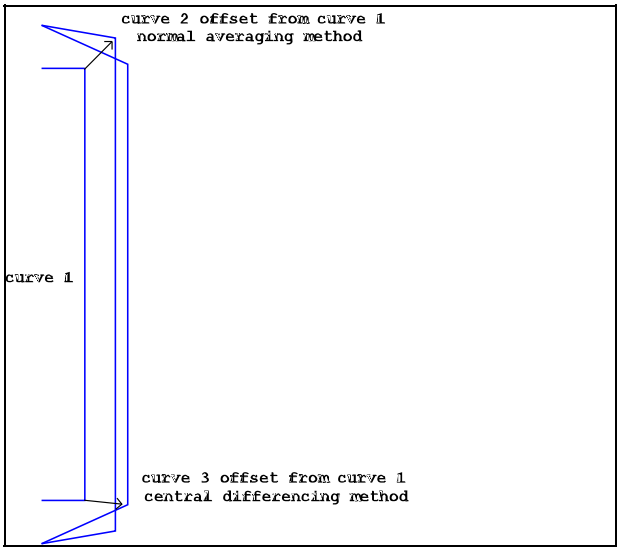

Figure 26 Comparison of two offset methods

Example

```
ld 1 lp2 1 0 2 0 2 10 1 10;
ld 2 lnof 1 -1 ;
ld 3 lod 1 -1 ;
```

lfil fillet

lfil $\Theta\ x'_{end}\ z'_{end}\ \Phi\ radius$

where

Θ	is the angle between the positive **x** axis and the line segment which joins the last point of the current curve and the fillet arc,
$(x'_{end}\ z'_{end})$	is the endpoint of this 2D curve segment,
Φ	is the angle between the positive **x** axis and the line segment which joins the fillet arc and the endpoint (x'_{end}, z'_{end}), and
radius	is the radius of the fillet.

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Remarks

Use this option to add a fillet made from 2 points and 2 angles to form an arc of a circle of the specified *radius*. This arc is met by a tangent line segment originating at the last defined point in the curve, having the slope specified by the angle Θ. That is, Θ is the angle between the positive x-axis and the line segment. The arc terminates at a point such that if a line is drawn from that point to the point (x'_{end}, z'_{end}), it would be tangent to the circle with the slope specified by the angle Φ. The picture which accompanies the following example should explain everything. The current 2D curve must already have at least one point.

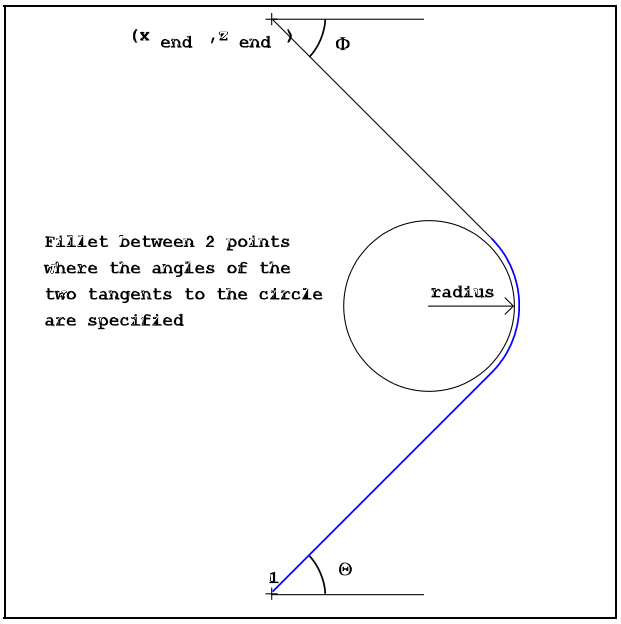

Figure 27 Fillet

Example

```
ld 1 lp2 1 1; lfil 45 1 2 -45 .25 lp2 1 2;
```

lap arc by center and end point

lap $x'_{end}\ z'_{end}\ x'_0\ z'_0$

where

(x'_{end}, z'_{end})	is the end of the line segment which defines the end of the arc, and
(x'_0, z'_0)	is both the center of the circular arc and the beginning of the line segment which defines the end of the arc.

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Remarks

Use this option to create a circular arc passing through the end of the current curve and terminating at the point of the circle intersecting with the ray containing the point (x_{end}, z_{end}). The circle's center is (x_0, z_0). Thus, the radius of the circle is the distance between (x_0, z_0) and the end of the current curve. The direction (clockwise or counter clockwise) is determined so as to produce the shortest arc between the two endpoints. The current 2D curve must already have at least one point. See the picture on the next page.

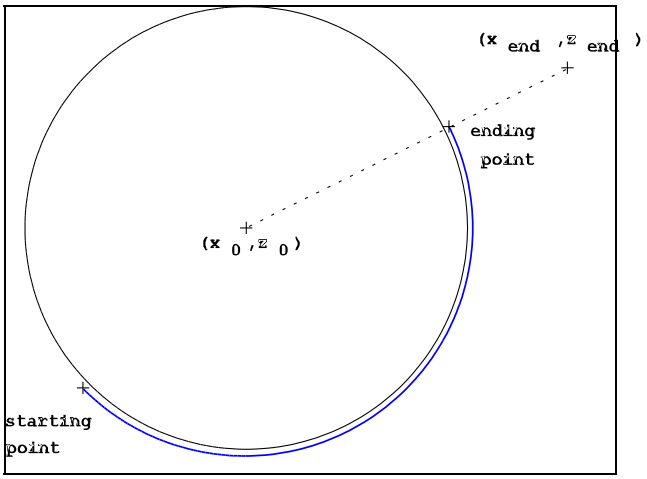

Figure 28 Arc of a circle by point and center

Example

```
ld 1 lp2 1 2;lap 4 4 2 3;
```

lar arc by radius and 2 points

lar x'_{end} z'_{end} *radius*

where

(x'_{end}, z'_{end})	is the end of the arc which is to be created, and
radius	is the radius of this circular arc.

Remarks

Use this option to create a circular arc passing through the end of the current 2D curve and the point (x'_{end}, z'_{end}), with the circle's *radius* also specified. There are two ways to do this. If *radius* is positive, the arc will be counter-clockwise; if negative, the arc will be clockwise. See the example on the following page.

This option requires that the current 2D curve have at least one point.

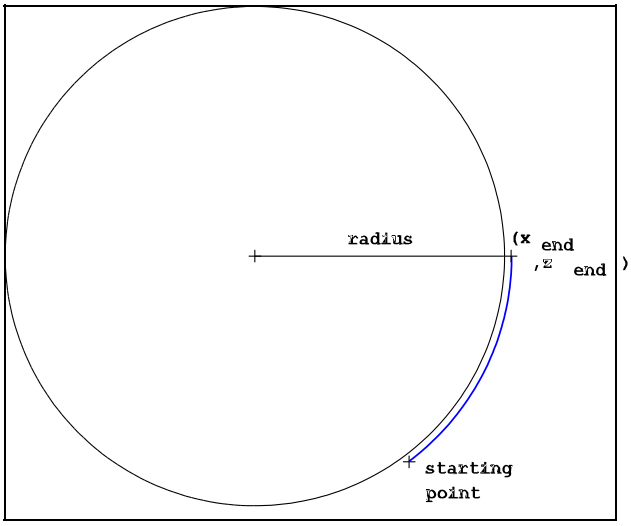

Figure 29 Arc of a circle by radius

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Example

```
ld 1 lp2 1 1 ;lar 2 3 2.5;
```

ltp tangent arc by radius and end point

ltp x'_{end} z'_{end} *radius*

where

(x'_{end}, z'_{end})	is the point at which the arc will end, and
radius	is the radius of the arc.

Remarks

Use this option to create a circular arc tangent to the tangent extension of the current curve, ending at a point (**t**angent-to-**p**oint), with the given *radius*. The current 2D curve must have at least two points. The endpoint of the current 2D curve will be extended or cut back to tangentially meet the circular arc. The circular arc will end at the point (x'_{end}, z'_{end}). See the example and picture on the next page.

Example

```
ld 1 lp2 1 1;
lar 2 3 2.5 ltp 3 6 3 ;
```

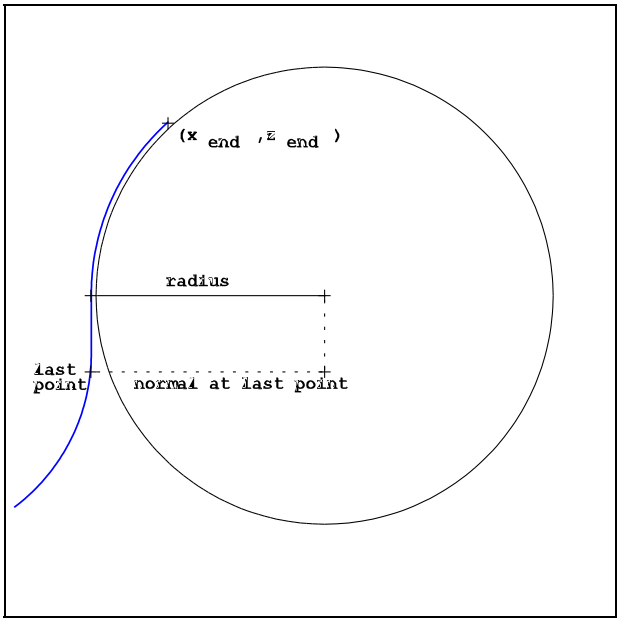

Figure 30 Tangent arc

lpt arc by radius and end tangent line

lpt x'_{begin} z'_{begin} x'_{end} z'_{end} *radius*

where

(x'_{begin}, z'_{begin})	is the end of the specified arc and beginning of the line segment,
(x_{end}, z_{end})	is the end of the line segment, and
radius	is the radius of the circular arc.

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Remarks

Use this option in order to create a circular arc extended by a tangent line segment (point-to-tangent), where the endpoints of the line segment, (x'_{begin}, z'_{begin}) and (x'_{end}, z'_{end}) respectively, and the radius are specified. The point (x'_{begin}, z'_{begin}) will be modified so that it is the point of tangency. This point is only used to specify the slope of the tangent line segment and it does not have to be at the point of tangency. The arc will begin at the end of the current 2D curve and end at the tangent line segment. The current 2D curve must have at least one point. See the example and the picture on the following page.

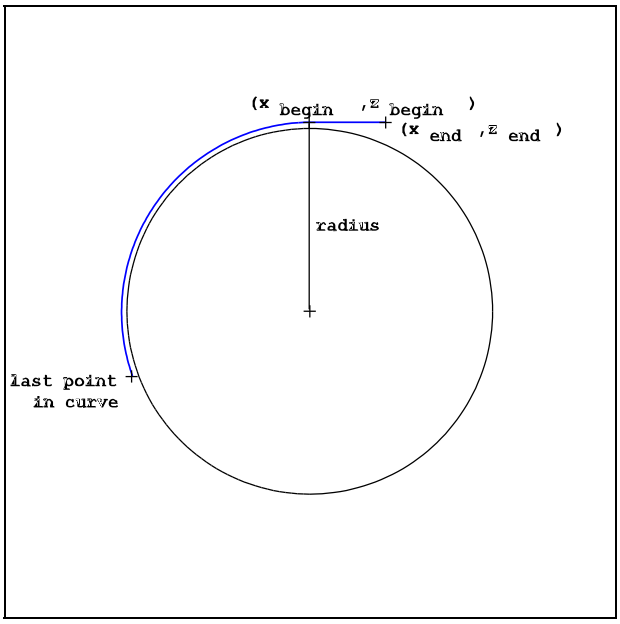

Figure 31 Append arc and tangent line

Example

```
ld 1 lp2 1 1;lpt 2 3 3 3 1.5;
```

lat fillet by radius and end tangent line

lat $x'_{begin}\ z'_{begin}\ x'_{end}\ z'_{end}$ *radius*

where

(x'_{begin}, z'_{begin})	is the point at the beginning of the tangent extension,
(x'_{end}, z'_{end})	is the point at the end of the tangent extension, and
radius	is the radius of the fillet.

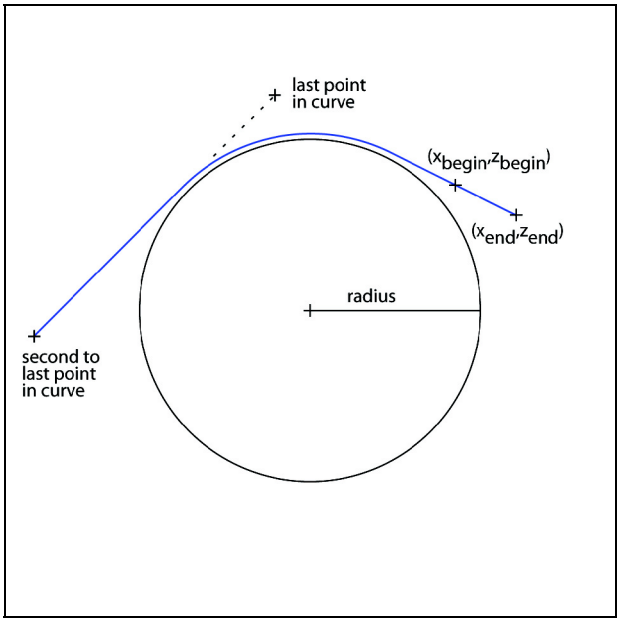

Figure 32 Fillet between two lines

Remarks

Use this option in order to append a fillet and a tangent extension by specifying the tangent extension between the points (x'_{begin}, z'_{begin}) and (x'_{end}, z'_{end}) and the *radius* of the fillet (**a**ll

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tangents). It may extend or truncate the last segment of the current 2D curve. The tangent extension may also be extended or truncated. Notice the difference between this and the previous option. This option guarantees a smooth transition. The trade-off is that the current 2D curve must have at least two points.

Example

```
ld 1 lp2 1 1 2 2;
lat 2.75 1.625 3 1.5 .75;
```

lad arc by center and angle

lad $x'_0\ z'_0\ \theta$

where

(x'_0, z'_0)	center of the arc
θ	subtended angle which defines the arc

Remarks

Use this option in order to append an arc of a circle with the specified center (x'_0, z'_0), starting at the last point in the current curve and rotating counterclockwise about a circle, by the specified positive angle θ. A negative angle rotates in a clockwise direction. The current 2D curve must have at least one point. The last point in the current 2D curve is used (with the arc's center) to determine the radius of the arc.

Example

```
ld 1 lp2 2 1 2 2;lad 0 3 45
```

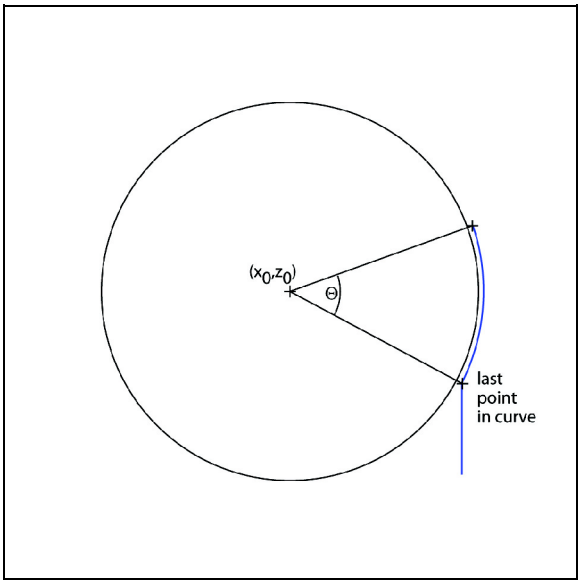

Figure 33 Circular arc

lvc point in polar coordinates

lvc θ *radius*

where

θ	is the angle along which the line segment will be drawn, and
radius	is the length of the defined segment.

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Remarks

Use this option in order to append a line segment by end point described as an offset from the end of the existing curve. The offset is in polar coordinates. The angle, θ, is measured with respect to the positive x-axis and counter clockwise is taken as positive.

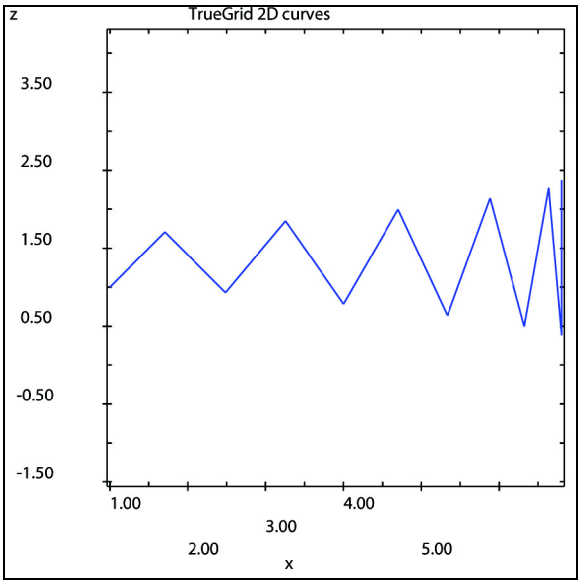

Figure 34 Saw Tooth using polar offsets

Example

```
ld 1 lp2 1 1;
lvc 45 1 lvc -45 1.1
lvc 50 1.2 lvc -55 1.3
lvc 60 1.4 lvc -65 1.5
lvc 70 1.6 lvc -75 1.7
lvc 80 1.8 lvc -85 1.9
lvc 90 2
```

lstl translate

lstl *curve* $\Delta x'$ $\Delta z'$

where

curve	ID number of a previously-defined 2D curve
$\Delta x'$	translation distance in the **x**′ (horizontal) direction
$\Delta z'$	translation distance in the **z**′ (vertical) direction

Remarks

Use this command in order to append a translation of a previously defined and numbered 2D curve to the current 2D curve. See the example and picture on the following page.

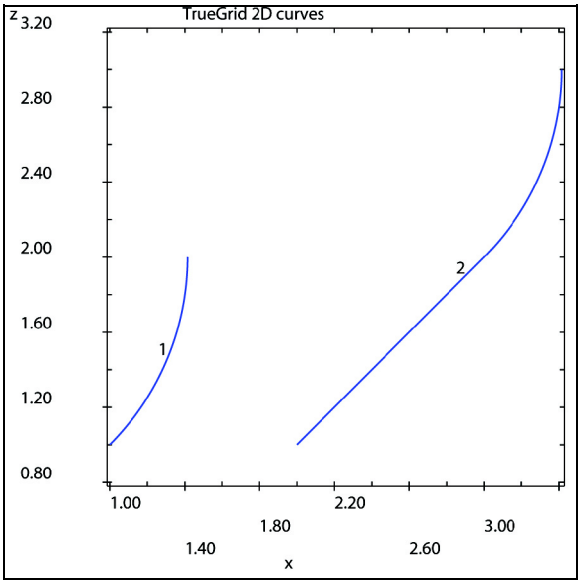

Figure 35 Append a 2D curve

Example

```
ld 1 lp2 1 1; lad 0 2 45
```

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```
ld 2 lp2 2 1; lstl 1 2 1
```

ltbc list of radial coordinates

ltbc θ_{begin} $\Delta\theta$ *scale radii*

where

θ_{begin}	is the first angular coordinate,
$\Delta\theta$	is the angular increment,
scale	is a radial multiplier, and
radii	is the list of radii which will be scaled.

Remarks

Use this option to append points defined in polar coordinates. *Radii* is a list of their radial coordinates. The angular coordinate of the first point is θ_{begin}, and the angular coordinate increases by $\Delta\theta$ for each successive point. Each *radii* is scaled by *scale*. Use the **ltbo** to modify these radii.

Example

The picture which corresponds to the following command file excerpt is shown on the next page.

```
ld 1 ltbc 0 1 1
1.00000 1.02161 1.04201 1.06119
1.07916 1.09592 1.11148 1.12585
1.13904 1.15106 1.16191 1.17163
1.18021 1.18768 1.19406 1.19936
1.20360 1.20682 1.20902 1.21023
1.21049 1.20981 1.20823 1.20577
1.20246 1.19834 1.19344 1.18779
1.18142 1.17436 1.16667 1.15836
1.14948 1.14007 1.13016 1.11980
1.10902 1.09786 1.08637 1.07458
1.06254 1.05028 1.03786 1.02531
1.01268 .999999 .987322 .974687
.962138 .949715 .937462 .925421
.913633 .902140 .890982 .880201
.869837 .859929 .850517 .841639
.833332 .825635 .818584 .812214
.806560 .801656 .797535 .794230
.791770 .790188 .789510 .789766
.790982 .793184 .796396 .800640
.805941 .812316 .819787 .828371
```

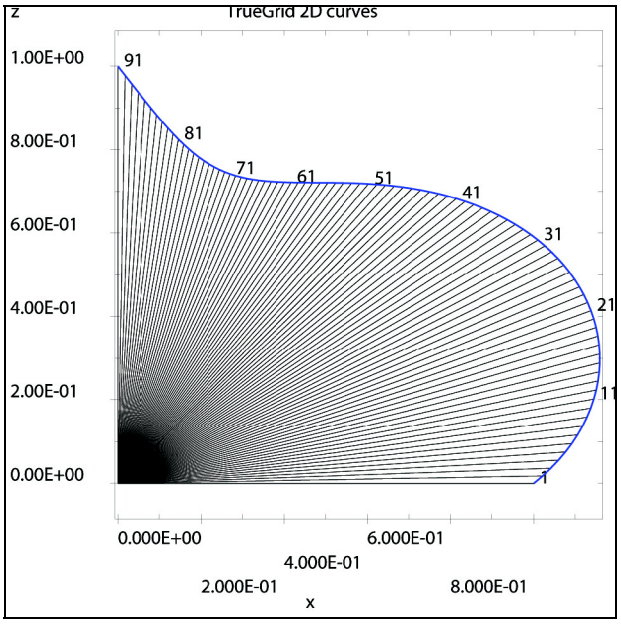

Figure 36 Curve by radial coordinates

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```
.838085  .848943  .860960  .874148  .888519  .904081  .920843  .938811
.957991  .978387  1.00000;
```

ltbo changes to the list of radial coordinates

ltbo $\Delta radius_1\ \Delta radius_2 \ldots ;$

where

$\Delta radius_i$ are the changes for each of the radii.

Remarks

The radial and angular coordinates are initialized (prior to using this command) using the **ltbc** command. The **ltbo** command is subsequently used to change the radial coordinates. These changes are cumulative.

Example

This example is based on the previous example using the **ltbc** command.

```
ld 2 ltbo
0     3.09017E-02     5.87785E-02
8.09017E-02     9.51056E-02
1.00000E-01     9.51057E-02
8.09018E-02     5.87787E-02
3.09019E-02 0 0 0 0 0 0 0 0 0 0
0    3.09017E-02      5.87785E-02
8.09017E-02     9.51056E-02
1.00000E-01     9.51057E-02
8.09018E-02     5.87787E-02
3.09019E-02 0 0 0 0 0 0 0 0 0 0
0     3.09017E-02     5.87785E-02
8.09017E-02     9.51056E-02
1.00000E-01     9.51057E-02
8.09018E-02     5.87787E-02
3.09019E-02 0 0 0 0 0 0 0 0 0 0
0     3.09017E-02     5.87785E-02
8.09017E-02     9.51056E-02
1.00000E-01     9.51057E-02
8.09018E-02       5.87787E-02
3.09019E-02  0  0  0  0  0  0  0  0  0  0  0  3.09017E-02  5.87785E-02
8.09017E-02  9.51056E-02  .1  9.51057E-02   8.09018E-02  5.87787E-02
3.09019E-02 0;
```

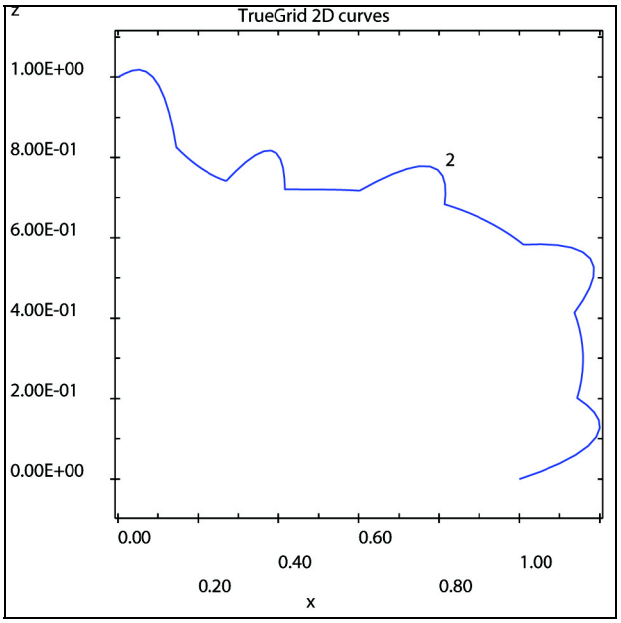

Figure 37 Modified **ltbc** curve

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lint interpolation

lint *curve*$_1$ *curve*$_2$ *interpolation*

where

curve$_1$	ID number of the first 2D curve
curve$_2$	ID number of the second curve
interpolation	interpolation parameter

Remarks

Interpolation is an weight factor for the first curve. (1-*interpolation*) is the weight factor for the second curve. In other words, specifying an *interpolation* of 1 will reproduce the first curve and specifying an interpolation of 0 will reproduce the second curve. The interpolation is done by relative arc length. A parameter from 0 to 1, based on the arc length, is used to traverse both curves uniformly. For each value of this parameter, a point is interpolated along the line between the two points associated with the parameter on the two curves. This is done for every point to form the interpolated third curve. See the example on the next page.

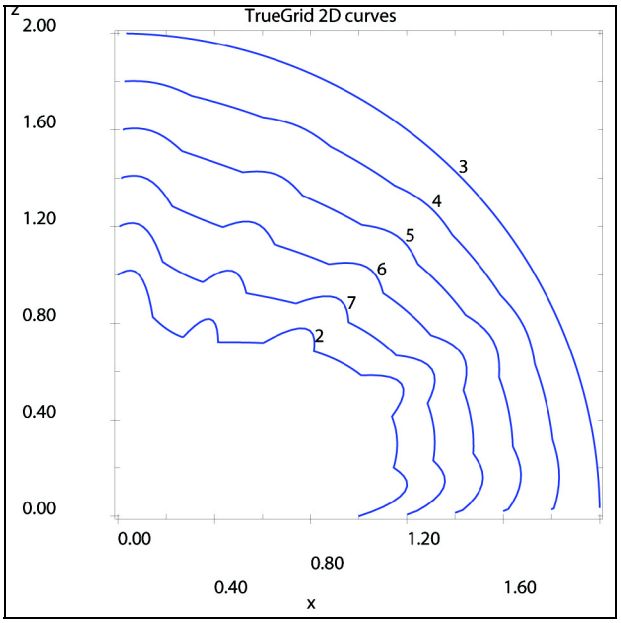

Figure 38 Interpolation Between 2 2D curves

Example

This example is based on the curve defined in the previous example using **ltbo**.

```
ld 3 lep 2 2 0 0 1 89 0;
ld 4 lint 2 3 .2
ld 5 lint 2 3 .4
ld 6 lint 2 3 .6
ld 7 lint 2 3 .8
```

csp2 cubic spline

csp2 *option* $x'_1\ z'_1\ \ldots\ x'_n\ z'_n$

where

option can be

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loop	close curve into a smooth loop
First endpoint with zero 2nd derivative	
00	last with zero 2nd derivative
01 *last_dx′ last_dz′*	last derivative
02 *last_curve_#*	last match first of other curve
03 *last_curve_#*	last match last of other curve
First endpoint derivative	
10 *first_dx′ first_dz′*	last with zero 2nd derivative
11 *first_dx′ first_dz′ last_dx′ last_dz′*	last derivative
12 first_dx′ first_dz′ *last_curve_#*	last match first of other curve
13 first_dx′ first_dz′ last_curve_#	last match last of other curve
First endpoint match first endpoint of another curve	
20 *first_curve_#*	last with zero 2nd derivative
21 *first_curve_# last_dx′ last_dz′*	last derivative
22 *first_curve_# last_curve_#*	last match first of other curve
23 *first_curve_# last_curve_#*	last match last of other curve
First endpoint match last endpoint of another curve	
30 *first_curve_#*	last with zero 2nd derivative
31 *first_curve_# last_dx′ last_dz′*	last derivative
32 *first_curve_# last_curve_#*	last match first of other curve
33 *first_curve_# last_curve_#*	last match last of other curve
$x'_1\ z'_1 \ldots x'_n\ z'_n$	pairs of control points

Example

The polygonal 2D curve number 1 and the spline curve number 2 with natural end derivatives are defined. The spline curve number 3 with loop condition is defined. The loop condition is based on equal coordinates and first derivatives of the start and the end point. Therefore the curve number 3 is smoother than the curve number 2. The start and end point coordinates are (-1,-1). The spline curve number 4 with specified end (-1,-1) and start derivatives (1,1) is defined. The spline curve number 5 is defined by the sane points as curve number 4. It uses start and end derivatives of curve number 4 to form a smooth loop with curve number 4. The spline curve number 6 also uses start and end derivatives of the spline curve number 4, but with different magnitudes. The command file follows:

```
ld 1
```

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```
 lp2 -1 -1 0 -1.5 1 -1 1.5 0 1 1 0 1.5 -1 1 -1.5 0 -1 -1;
    c polygon from control points

ld 2
 csp2 00
   -1 -1 0 -1.5 1 -1 1.5 0 1 1 0 1.5 -1 1 -1.5 0 -1 -1;;
    c spline with natural derivatives - 00

ld 3
 csp2 loop
  -1 -1 0 -1.5 1 -1 1.5 0 1 1 0 1.5 -1 1 -1.5 0 -1 -1;;
    c spline with loop condition - loop

ld 4
 csp2 11 -1 -1 1 1
  -1 -1 0 -1.5 1 -1 1.5 0 1 1 0 1.5 -1 1 -1.5 0 -1 -1;;
    c spline with specified start and end derivatives
    c  -1 -1 and 1 1

ld 5
 csp2 32 4 1 4 1
  -1 -1 0 -1.5 1 -1 1.5 0 1 1 0 1.5 -1 1 -1.5 0 -1 -1;;
    c spline with specified start and end derivatives
    c form curve 4 with magnitude 1      4 1 4 1

ld 5 csp2 32 4 1 4 1
  -1 -1 0 -1.5 1 -1 1.5 0 1 1 0 1.5 -1 1 -1.5 0 -1 -1;;
    c spline with specified start and end derivatives
    c form curve 4 with magnitudes 2 and 4      4 2 4 4
```

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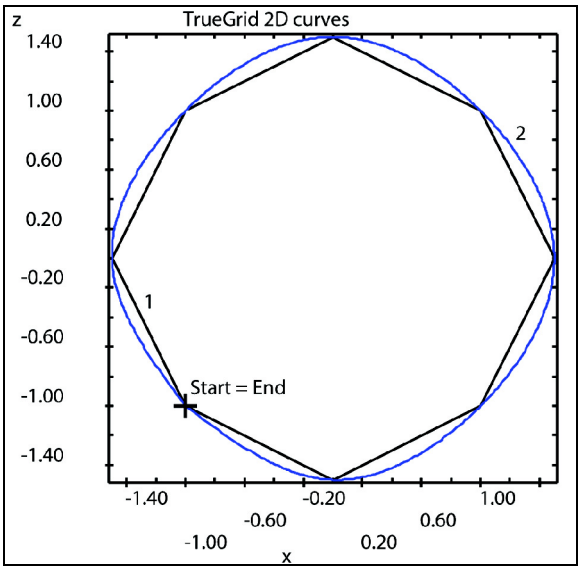

Figure 39 spline with natural end derivatives

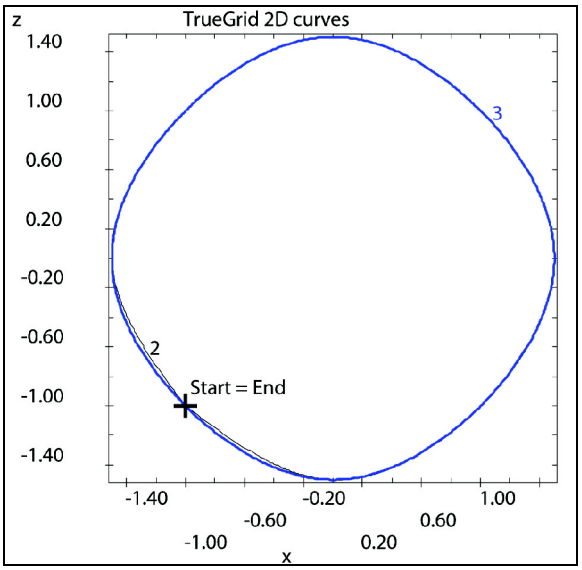

Figure 40 spline with loop condition

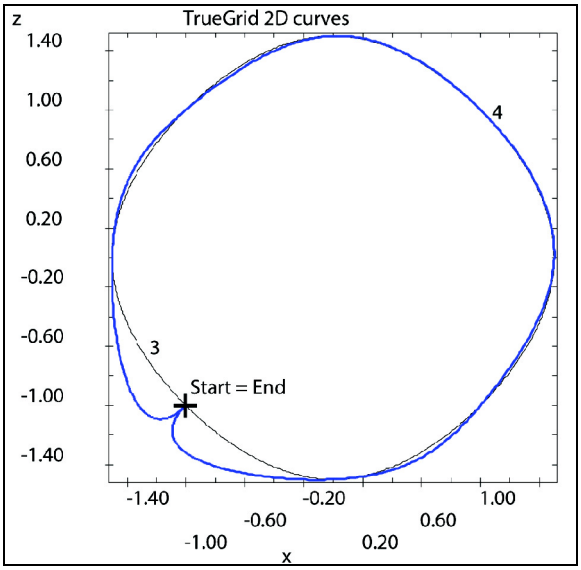

Figure 41 spline with end derivatives

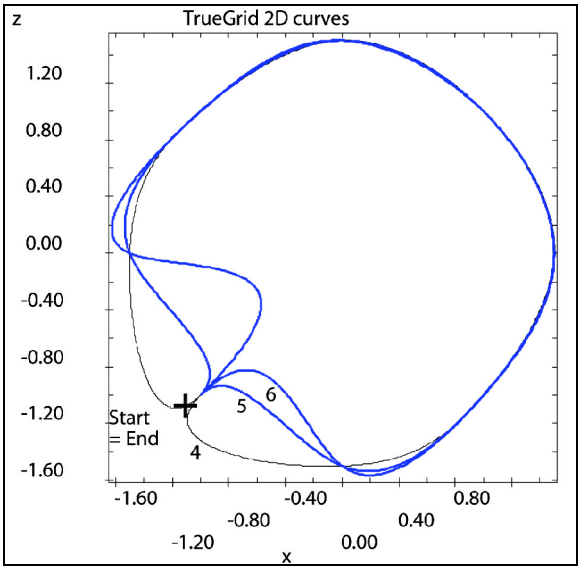

Figure 42 spline with various end derivatives

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fws2 Fowler-Wilson cubic spline

fws2 *option convergence mxiniter mxouiter* $x'_1\ z'_1\ \ldots\ x'_n\ z'_n$;

where

option can be

loop close curve into a smooth loop

First endpoint with zero 2nd derivative

00	last with zero 2nd derivative
01 *last_dx′ last_dz′*	last derivative
02 *last_curve_#*	last match first of other curve
03 *last_curve_#*	last match last of other curve

First endpoint derivative

10 *first_dx′ first_dz′*	last with zero 2nd derivative
11 *first_dx′ first_dz′ last_dx′ last_dz′*	last derivative
12 first_dx′ first_dz′ *last_curve_#*	last match first of other curve
13 first_dx′ first_dz′ last_curve_#	last match last of other curve

First endpoint match first endpoint of another curve

20 *first_curve_#*	last with zero 2nd derivative
21 *first_curve_# last_dx′ last_dz′*	last derivative
22 *first_curve_# last_curve_#*	last match first of other curve
23 *first_curve_# last_curve_#*	last match last of other curve

First endpoint match last endpoint of another curve

30 *first_curve_#*	last with zero 2nd derivative
31 *first_curve_# last_dx′ last_dz′*	last derivative
32 *first_curve_# last_curve_#*	last match first of other curve
33 *first_curve_# last_curve_#*	last match last of other curve

convergence	convergence tolerance
mxiniter	maximum of inner iterations
mxouiter	maximum of outer iterations
$x'_1\ z'_1\ \ldots\ x'_n\ z'_n$	pairs of control points

Remarks

The Wilson-Fowler algorithm fits cubic spline arcs through the set of given points. The basic idea

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was to come as close as possible to the true spline (elastica), which minimizes the energy of the curve, with a curve composed of cubic segments which are joined together to achieve continuous tangent and curvature. This was accomplished by introducing a local u,v-coordinate system for each segment, with the independent variable running along the chord joining (x'_i, z'_i) with (x'_{i+1}, z'_{i+1}). The curvature-matching conditions lead to a tri-diagonal nonlinear system of equations, which must be solved iteratively. The difference between derivatives for the *i* and *i+1* point is minimized until the maximal difference is less than *convergence* or until the *maximum number of iterations* is reached.

The recommended value for *convergence* is from 10^{-3} to 10^{-6}. The recommended value for *mxiniter* is 4. The recommended value for *mxouiter* is 8 times the total number of points.

The process does not necessarily converge for every case. When this method does not converge, an informative message is issued with the largest curvature difference. It is useful for simulation of production processes.

The theoretical description of the Wilson-Fowler Spline can be found in:

> Cubic Spline, a Curve Fitting Routine. Fowler, Wilson.
> Union Carbide Corp., Nuclear Div., Oak Ridge, Tennessee, 1966.

Example

Two Fowler-Wilson spline curves are defined by sets of points. The points are marked and numbered. The **11** option is used to specify first and second endpoint derivatives, both of the same size `0.001` and `1.` Curve number 2 is randomly perturbed with the normal distribution. The command file and pictures follow:

```
para dx 0.001 mp [sqrt(2)/2]
     d 0.02;
 c parameters dx,mp,d
 c are defined
ld 1 fws2 11 %dx 1 %dx 1
.000001 4 40 1 0 0 1 -1 0 0 -1 1
0;;
 c curve 1
 c - Fowler wilson spline
 c option 11 -
 c specified endpoint deriv.
 c first_dx(%dx) first_dz(1)
 c second_dx(%dx) second_dz(1)
```

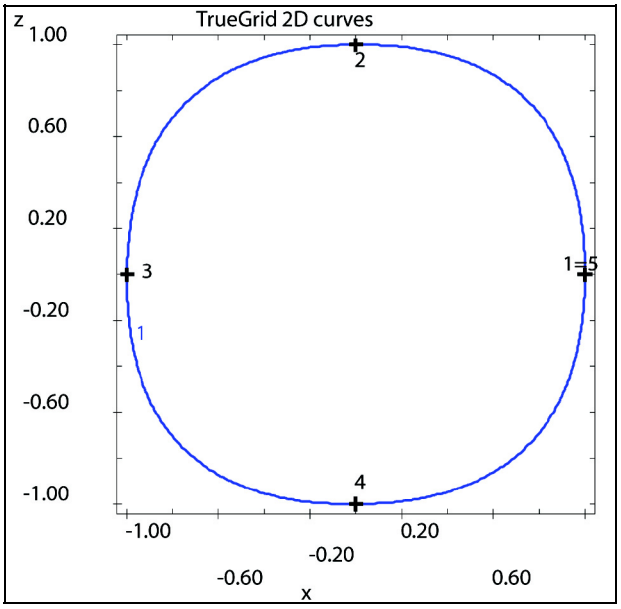

Figure 43 Fowler-Wilson Spline

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```
 c convergence  .0000001
 c max.of inner iterations 4
 c max.of outer iterations 40
 c x and z coordinates
 c   of 5 points
ld 2 fws2 11 %dx 1 %dx 1
  .000001 4 264
 c curve 2
 c - Fowler wilson spline
 c option 11 -
 c specified endpoint deriv.
 c first_dx(%dx) first_dz(1)
 c second_dx(%dx) second_dz(1)
 c convergence  .0000001
 c max.of inner iterations 4
 c max.of outer iterations 264
 c x and z coordinates
 c   of 33 points
```

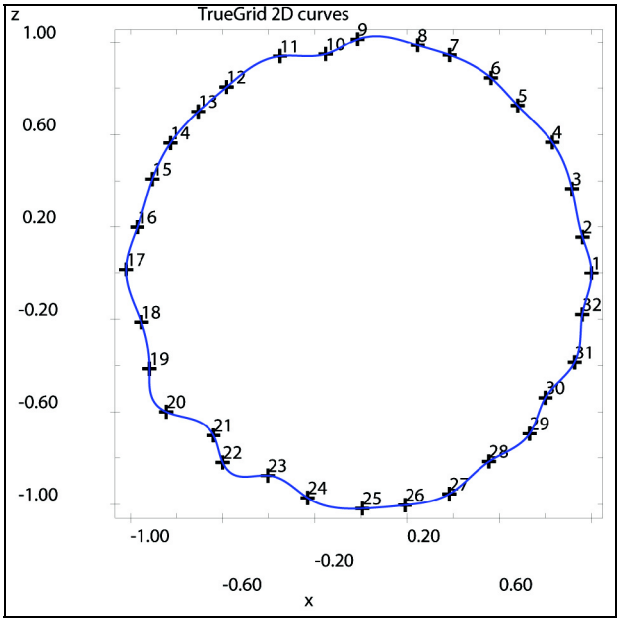

Figure 44 Fowler-Wilson Spline

```
[      1.0          ] [      0.0          ]
[ 0.9772311+%d*norm(0.2138127652)] [ 0.2121777+%d*norm]
[ 0.9238796+%d*norm] [ 0.3826835+%d*norm]
[ 0.8216469+%d*norm] [ 0.5699968+%d*norm]
[ 0.7071068+%d*norm] [ 0.7071068+%d*norm]
[ 0.5699968+%d*norm] [ 0.8216469+%d*norm]
[ 0.3826835+%d*norm] [ 0.9238796+%d*norm]
[ 0.2121777+%d*norm] [ 0.9772311+%d*norm]
[       0.0+%d*norm] [       1.0+%d*norm]
[-0.2121777+%d*norm] [ 0.9772311+%d*norm]
[-0.3826835+%d*norm] [ 0.9238796+%d*norm]
[-0.5699968+%d*norm] [ 0.8216469+%d*norm]
[-0.7071068+%d*norm] [ 0.7071068+%d*norm]
[-0.8216469+%d*norm] [ 0.5699968+%d*norm]
[-0.9238796+%d*norm] [ 0.3826835+%d*norm]
[-0.9772311+%d*norm] [ 0.2121777+%d*norm]
[      -1.0+%d*norm] [       0.0+%d*norm]
[-0.9772311+%d*norm] [-0.2121777+%d*norm]
[-0.9238796+%d*norm] [-0.3826835+%d*norm]
[-0.8216469+%d*norm] [-0.5699968+%d*norm]
[-0.7071068+%d*norm] [-0.7071068+%d*norm]
[-0.5699968+%d*norm] [-0.8216469+%d*norm]
```

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```
[-0.3826835+%d*norm] [-0.9238796+%d*norm]
[-0.2121777+%d*norm] [-0.9772311+%d*norm]
[       0.0+%d*norm] [      -1.0+%d*norm]
[ 0.2121777+%d*norm] [-0.9772311+%d*norm]
[ 0.3826835+%d*norm] [-0.9238796+%d*norm]
[ 0.5699968+%d*norm] [-0.8216469+%d*norm]
[ 0.7071068+%d*norm] [-0.7071068+%d*norm]
[ 0.8216469+%d*norm] [-0.5699968+%d*norm]
[ 0.9238796+%d*norm] [-0.3826835+%d*norm]
[ 0.9772311+%d*norm] [-0.2121777+%d*norm]
[       1.0        ] [       0.0        ];;
```

ctbc polar cubic spline

ctbc *option convergence* θ_{begin} $\Delta\theta$ *scale radii*;

where

option can be

First endpoint with zero 2nd derivative	
00	last with zero 2nd derivative
01 *last_dx′ last_dz′*	last derivative
First endpoint derivative	
10 *first_dx′ first_dz′*	last with zero 2nd derivative
11 *first_dx′ first_dz′ last_dx′ last_dz′*	last derivative

θ_{begin}	first angular coordinate
$\Delta\theta$	angular increment
scale	radial multiplier
radii	list of radii

Remarks

This option is similar to the **csp2** option. The difference is that the spline is calculated in polar coordinates. Points are defined in polar coordinates. *Radii* is a list of their radial coordinates. The angular coordinate of the first point is θ_{begin}, and the angular coordinate increases by $\Delta\theta$ for each successive point. Each *radius* is scaled by *scale*. The derivatives *first_dx′, first_dz′, second_dx′, second_dz′* must be specified in the Cartesian (not polar) coordinate space. Only the direction of the derivative vector is used. The magnitude of the derivative vector is ignored.

Use the **ctbo** option to modify the radii.

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ctbo Modify a Polar Cubic Spline

ctbo *option* $\Delta radius_1$ $\Delta radius_2$ ... ;

where

option can be

First endpoint with zero 2nd derivative	
00	last with zero 2nd derivative
01 *last_dx′ last_dz′*	last derivative
First endpoint derivative	
10 *first_dx′ first_dz′*	last with zero 2nd derivative
11 *first_dx′ first_dz′ last_dx′ last_dz′*	last derivative

$\Delta radius_i$ additive changes to the radii.

Remarks

The spline curve is constructed in the polar coordinate system. All interpolation is done in polar coordinates. The radial and angular coordinates are initialized (prior to using this command) using the **ltbc**, **ctbc**, or **ftbc** option. The **ctbo** option is subsequently used to change the radial coordinates. These changes are cumulative. Only the direction of the derivative vector is used. The magnitude of the derivative vector is ignored.

ftbc Fowler-Wilson polar cubic spline

ftbc *option convergence mxiniter mxouiter* θ_{begin} $\Delta\theta$ *scale radii*;

where

option can be

First endpoint with zero 2nd derivative	
00	last with zero 2nd derivative
01 *last_dx′ last_dz′*	last derivative
First endpoint derivative	
10 *first_dx′ first_dz′*	last with zero 2nd derivative
11 *first_dx′ first_dz′ last_dx′ last_dz′*	last derivative

convergence convergence tolerance

mxiniter maximum of inner iterations

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mxouiter	maximum of outer iterations
θ_{begin}	first angular coordinate
$\Delta\theta$	angular increment
scale	radial multiplier
radii	list of radii

Remarks

This option is similar to the **fws2** option. The difference is that the spline is calculated in polar coordinates. Points are defined in polar coordinates. *Radii* is a list of their radial coordinates. The angular coordinate of the first point is θ_{begin}, and the angular coordinate increases by $\Delta\theta$ for each successive point. Each *radius* is scaled by *scale*. The derivatives *first_dx', first_dz', second_dx', second_dz'* must be specified in the Cartesian (not polar) coordinate space. Only the direction of the derivative vector is used. The magnitude of the derivative vector is ignored.

The recommended value for *convergence* is from 10^{-3} to 10^{-6}. The recommended value for *mxiniter* i 4. The recommended value for *mxouiter* is 8 times the number of points.

Example

A Fowler-Wilson polar spline curve is defined by sets of points (see 45). The points are marked and numbered. The **11** option is used to specify first and second endpoint derivatives, both of the same components `0.001` and `1.` The command file and pictures follow:

```
ld 1 ftbc 11 .001 1 .001 1 .001
4 72 10.  10 1 1 .2 .3 .4 .5 .6
.7 .8 .9
```

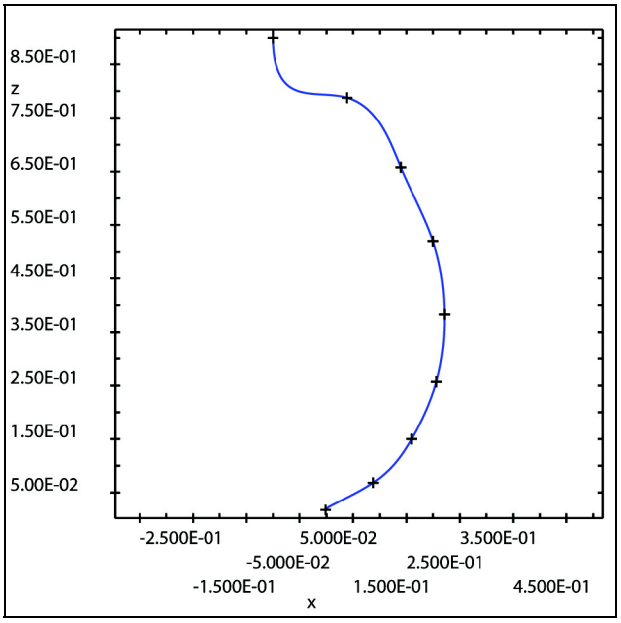

Figure 45 Polar Fowler-Wilson cubic spline

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ftbo modify Wilson-Fowler polar cubic spline

ftbo *option convergence mxiniter mxouiter* $\Delta radius_1 \Delta radius_2 \ldots$;

where

option can be

First endpoint with zero 2nd derivative	
00	last with zero 2nd derivative
01 *last_dx' last_dz'*	last derivative
First endpoint derivative	
10 *first_dx' first_dz'*	last with zero 2nd derivative
11 *first_dx' first_dz' last_dx' last_dz'*	last derivative
convergence	convergence tolerance
mxiniter	maximum of inner iterations
mxouiter	maximum of outer iterations
$\Delta radius_1 \Delta radius_2 \ldots$	list of changes in the radii

Remarks

The spline curve is constructed in the polar coordinate system. All interpolation is done in polar coordinates. The radial and angular coordinates are initialized (prior to using this command) using the **ltbc**, **ctbc**, or **ftbc** command. The **ftbo** command is subsequently used to change the radial coordinates. These changes are cumulative. Only the direction of the derivative vector is used. The magnitude of the derivative vector is ignored.

The recommended value for *convergence* is from 10^{-3} to 10^{-6}. The recommended value for *mxiniter* i 4. The recommended value for *mxouiter* is 8 times the number of points.

Example

The Fowler-Wilson polar spline curve created with the **ftbc** command is modified with the **ftbo** command. The points are marked and numbered.

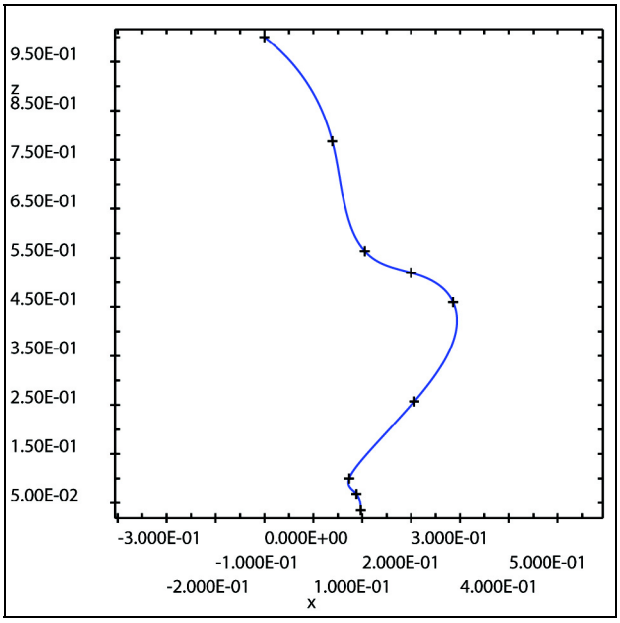

Figure 46 Modified Fowler-Wilson polar

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The **10** option is used to specify the first endpoint derivatives, `0.001` and `1.` The command file and pictures follow:

```
ld 1 ftbc 11 .001 1. .001 1 .001 4 72 10 10 1 .1 .2 .3 .4 .5 .6 .7 .8
.9
ld 1 ftbo 10 .001 1. .001 4 72 .1 0 -.1 0 .1 0 -.1 0 .1
```

rseg import from edge file

rseg <*no arguments*>

Remarks

Use this option to append the result of reading the next segment of a 2D curve definition in an edge file; see the **edgefile** command. It is required that the **edgefile** command was used first to specified the file containing the curve data. This command can be used repeatedly. Each time this command is used, the next 2D curve segment data found in the edge file will be appended to the present 2D curve. When all of the data in this file has been used, then the file is automatically closed.

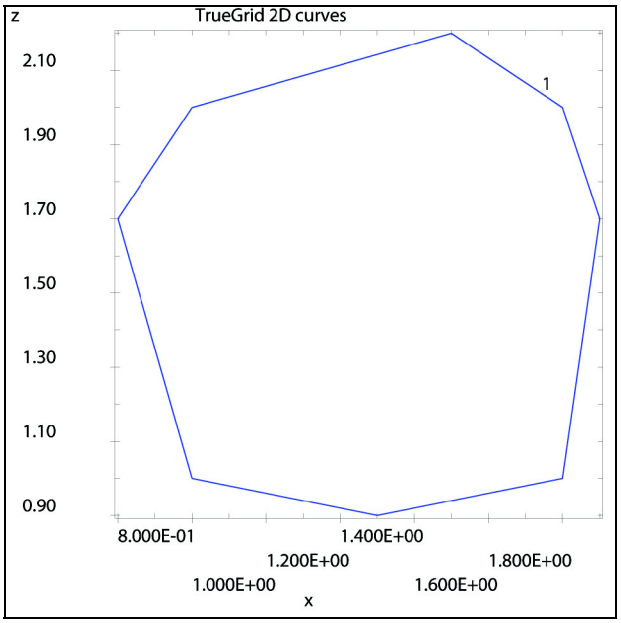

Figure 47 2 segments from edge file

Example

The example edge file in the description on the **edgefile** command above can be converted to 2D curve definitions using the **rseg** command. For example,

```
ld 1 rseg rseg
```

3. 2D Curve Display Commands

These commands control which 2D curves are displayed. Since they are so obvious in what they do, no examples are included in this section.

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lcv display a load curve

lcv *load_curve_number*

where

load_curve_number is the number of the load curve to be displayed.

Remarks

If this is the first invocation of a load curve or 2D curve display command, then a new window for the curves is drawn.

lv display all 2D curves

lv *(no arguments)*

Remarks

The command displays all previously defined 2D curves. If this is the first invocation of a curve display command, then a new window is drawn for the curves.

lvi display list of 2D curves

lvi $curve_1$ $curve_2$ ... ;

where

$curve_i$ is the number of a(the) previously defined 2D curve(s).

Remarks

This command is similar to **dcds**, in that this command allows you to display a list of 2D curves. Other defined 2D curves will not be displayed or will be removed from the display.

lvs display a sequence of 2D curves

lvs *first_line last_line*

where

first_line is the first 2D curve to be displayed, and

last_line is the number of the last curve to be displayed.

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Remarks

This command will display every 2D curve whose ID number lies between the 2 specified curve IDs.

4. Create 3D Curves

These commands define, modify, and display information about 3D curves. There are many uses for 3D curves. Every 3D curve is identified by a number. Use the automatic parameter %nextcrv for the next available 3D curve number.

You may wish to project a face of the mesh to a surface so that the entire surface is covered. In most cases, you will need to attach the edges of the face to the edges of the surface or to 3D curves that form the shape of the surface boundaries. It is typical to import an IGES model with many surfaces that need to be combined for meshing purposes. This is done with the **sds** option of the **sd** command. Then a face of the mesh can be projected to this composite surface. Once again, to cover the entire surface, attach the edges of the mesh to the boundary of the composite surface which can be formed using the **sdedge** option of the **curd** command or the **coedge** interactive equivalent.

When a surface has more than 180 degrees of curvature, the projection to that surface must be helped by attaching the edge of the mesh to the boundaries or a 3D curve. When there is an interior cusp or feature in the surface that requires a mesh line to follow, a 3D curve can be formed. Then an interior edge of the mesh can be attached to this curve.

If two faces, with a common edge, are projected to two different surfaces, then the common edge will be placed along the intersection of the two surfaces. You can accomplish a similar affect along the edge of the mesh by attaching it to a 3D curve.

When two neighboring faces of the mesh are projected to the same surface, you may wish to control the way the common edge is projected onto the surface by initializing that edge along a 3D curve. This is a common procedure when making fine adjustments to the mesh to improve the mesh quality.

3D curves are also useful when no surface information is available. It was noted above that these curves can be used to define surfaces. It is also possible to place the edges of the mesh along the 3D curves and interpolate the faces of the mesh. In fact, this will be done by default, when no surfaces are specified. However, you may wish to specify the type of interpolation between the edges using the **lin** or **tf** commands.

The **curd** command can be used to begin the definition of a 3D curve. A 3D curve can also be imported from an IGES file by using the **IGES** or **IGESCD** commands. A 3D curve can be used to create 2D curves (**ld3d2d**), 3D curves (**contour**), and surfaces (**sd**). Vertices of the mesh can be

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placed at points on a 3D curve (**pbs**). Edges of the mesh can be placed along a piece of a curve either initially (**cur**) or permanently (**curf**) or placed along the entire curve (**cure**). **curs** is useful when you want to be able to control the position of the intermediate vertices along a multiple edge.

A 3D curve is built with 3D curve segments. For example, a composite curve from an IGES file is a segmented 3D curve. In many cases, a single 3D curve segment using only one of the options in this chapter is all that is needed for a 3D curve definition. This segmentation feature is something to keep in mind for later when you may wish to make modifications to the geometry by extending a 3D curve definition. Each 3D curve option appends a curve segment to the 3D curve initialized by the last **curd** command. A segment is appended either to the beginning or ending of the existing curve, and the direction of the curve may be switched. This is done to minimize the distance between the endpoint of the existing curve and the endpoint of the appending segment to form one composite curve. The direction of the curve can be checked with the **labels crvpt** command.

You have to specify curves in the correct order, when you are appending or copying curves (**Figure 48** and **Figure 49**)

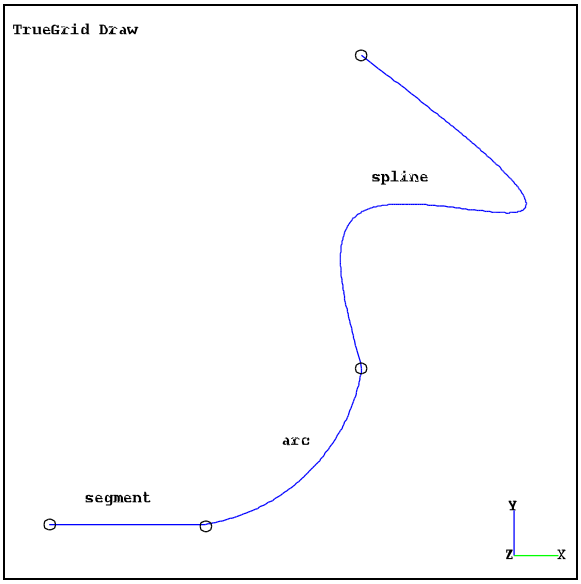

Figure 48 correct order
1 - segment, 2 - arc, 3 - spline

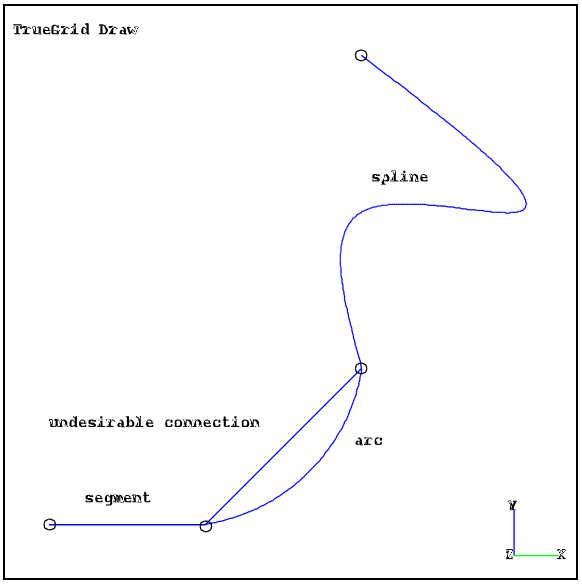

Figure 49 wrong order
1 - segment, 2 - spline , 3 - arc

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curd 3D curve definition

curd 3d_curve_# type_of_curve curve_data_list

where

type_of_curve and *curve_data_list* can be

igc *iges_curve_#*
sdedge *surface_#.edge_#*
lp3 *x1 y1 z1 ... xn yn zn* ; *trans* ;
contour *surface_point_id_1 surface_point_id_2*
csp3 *option x1 y1 z1 ... xn yn zn* ; *trans* ;
bsp3 *knots x1 y1 z1 ... xn yn zn* ;
nrb3 *knots weights x1 y1 z1 ... xn yn zn* ;
ld2d3d *2d_curve_# system coordinate start end trans* ;
intcur *3d_curve_1 3d_curve_2 interpolate trans* ;
lp3pt *point_id_1 ... point_id_n ; trans* ;
3dfunc *min_u max_u x_expres ; y-expres ; z-expres ; trans ;*
projcur *3d_curve_# surface_# trans ;*
pscur *3d_curve_# surface_# #_iterations tol trans ;*
arc3 *option system point system point system point*
twsurf *surface_1 surface_2 x1 y1 z1 ... xn yn zn ; trans ;*
cpcd *curve_# trans* ;
cpcds *list_curve_# trans* ;
rmseg, and

trans is a product from left to right of the following operators

mx *x_offset*
my *y_offset*
mz *z_offset*
v *x_offset y_offset z_offset*
rx *theta*
ry *theta*
rz *theta*
raxis *angle x0 y0 z0 xn yn zn*
rxy
ryz
rzx
tf *origin x-axis y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:

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rt *x y z* (cartesian coordinates)
cy *rho theta z* (cylindrical coordinates)
sp *rho theta phi* (spherical coordinates)
pt *c.i* (label of a labeled point from a 3D curve)
pt *s.i.j* (label of a labeled point from a surface)

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* (cartesian coordinates)
cy *rho theta z* (cylindrical coordinates)
sp *rho theta phi* (spherical coordinates)
pt *c.i* (label of a labeled point from a 3D curve)
pt *s.i.j* (label of a labeled point from a surface)

inv to invert the present transformation
csca scale all coordinates
xsca scale the x-coordinates
ysca scale the y-coordinates
zsca scale the z-coordinates

Remarks

To create a 3D curve, begin by selecting the identification number with the **curd** command. Then select the curve type and associated parameters. Optionally, additional component 3D curve segments can be appended. This is done indefinitely, until a new **curd** command is executed. When a new segment is appended to a 3D curve definition, the data is ordered to best match the endpoint of the last segment appended to the 3D curve definition. The following options are used, within the **curd** command, at any time to append a segment:

Use **igc** to create/append an IGES 3D curve by its sequence number.

Use **sdedge** or **se** to create/append an edge of a defined surface. Each edge is assigned a number. The edge number m of surface number n is identified with the symbol n.m . To view the surface with the numbered edge, use the graphics **label** command with the option **sdedge** or **se**.

Use **arc3** to create/append a 3D arc of a circle passing through 3 points. There are three options. An arc can be created passing from the first to the second to the third point. The second option is the complement of the first option above. That is, the other portion of the circle is used. The third option creates the entire circle.

Use **lp3** to create/append a sequence of points forming a polygonal 3D curve.

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Use **lp3pt** to create/append a sequence of surface/curve points forming a polygonal 3D curve.

Use **cpcd** to create/append a copy of a previously defined 3D curve.

Use **cpcds** to create/append a copy of a sequence of previously defined 3D curves.

Use **contour** to create/append a contour segment from an existing surface or curve. The *surface_point_id* is formed from 3 numbers separated by periods. For example 3.10.12 references a point on surface 3, at position i=10, j=12. The *curve_point_id* is formed from 2 numbers separated by a period. For example, 4.1 references the 1st point in 3D curve 4.

Use **csp3** to create/append a 3D cubic spline through a set of ordered points. The end derivatives can be specified or matched.

Use **bsp3** to create/append a b-spline curve formed from an explicit list of knots and control points.

Use **nrb3** to create/append a nurbs curve formed from an explicit list of knots, weights and control points.

Use **intcur** to create/append a curve interpolated between two other 3D curves.

Use **twsurf** to create/append the curve at the intersection of two surfaces. Initial points are required to determine the endpoints and some intermediate points for surfaces with large curvature (greater than 120 degrees). These points need only be approximate.

Use **ld2d3d** to create/append a 2D curve converted to 3D.

Use **3dfunc** to create/append a parameterized function curve with a specified domain.

Use **projcur** to create/append another 3D curve projected onto a surface.

Use **pscur** to create/append another 3D curve projected onto a surface and smoothed.

Remarks with Examples for Spline Curve Automatic Parameters

There are a number of automatic parameters associated with spline curves. In order to be consistent in the use of these automatic parameters with respect to Cubic Splines, B-Splines, and NURBS, the knots will refer to the monotonic increasing sequence of numbers in the 1D parameter space. Knot points will refer to the 3D coordinates that result when mapping the knots to the 3D curve. Control Points, which are not found in Cubic Splines, refer to the 3D coefficients that scale the basis functions. The 1D weights, which are only found in NURBS, scale the basis functions to form the polynomial in

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the denominator of the rational function.

If the last 3D spline curve was formed using the **cps3** (Cubic Spline) option of the **curd** command, then the following parameters will be automatically set:

splks - 1 denemsion array containing all of the knots
splnk - scalar number of knots in the array splks
splkps - 2 dimensional array (3,*) of knot points on the 3D curve
splord - the order of the polynomials - always 4
splcn - **True*Grid*®** curve number that was created with this data
spltyp - type of spline curve - always 1

For example, if the following command were executed:

```
curd 11 csp3 00
 -3.3954249e+00 1.2127594e-01 -3.5484806e-01
 -3.3571637e+00 1.5977994e-01 -3.3923153e-01
 -3.3138983e+00 2.0643232e-01 -3.1305539e-01
 -3.2234428e+00 2.8636580e-01 -2.4210249e-01
 -3.1053569e+00 3.5468611e-01 -1.2244769e-01
 -3.0454067e+00 3.7062020e-01 -5.3555769e-02
 -3.0135252e+00 3.7547471e-01 -1.6162919e-02
 -3.0040656e+00 3.7634343e-01 -4.9400860e-03
 -2.9906838e+00 3.7713331e-01 1.1124370e-02
 -2.9789010e+00 3.9086247e-01 3.7436009e-02
 -2.9678857e+00 4.1399955e-01 7.1353959e-02
 -2.9503219e+00 4.4951856e-01 1.2442068e-01
 -2.9295484e+00 4.8749987e-01 1.8401165e-01
 -2.9095490e+00 5.1844508e-01 2.3694231e-01
 -2.8932302e+00 5.3794400e-01 2.7557711e-01
 -2.8758707e+00 5.5043960e-01 3.1012778e-01
 -2.8553295e+00 5.4789609e-01 3.3721066e-01
 -2.8470099e+00 5.3850897e-01 3.4150743e-01
 -2.8389274e+00 5.2246647e-01 3.4014181e-01
 -2.8261104e+00 4.7641905e-01 3.2144514e-01
 -2.8157659e+00 4.1792813e-01 2.8925255e-01
 -2.7706413e+00 2.9269938e-01 2.5559440e-01
 -2.7520236e+00 2.6949270e-01 2.6488436e-01
 -2.7330521e+00 2.5119662e-01 2.7857573e-01
 -2.7025447e+00 2.2323597e-01 3.0129292e-01
 -2.5951468e+00 1.1034914e-01 3.5839118e-01
 -2.4999798e+00 0.0000000e+00 3.7500000e-01 ;;;
```

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Then the following automatic parameters would be set to have the values below:

```
%spltyp     =    1
%splcn      =   11
%splord     =    4
%splnk      =   27
%splks(*) =    0.000000D+00
               3.260380D-02
               7.231759D-02
               1.531419D-01
               2.578884D-01
               3.113998D-01
               3.399025D-01
               3.483897D-01
               3.604670D-01
               3.788989D-01
               4.034368D-01
               4.416657D-01
               4.841821D-01
               5.214089D-01
               5.481064D-01
               5.715624D-01
               5.912381D-01
               5.988915D-01
               6.092905D-01
               6.389166D-01
               6.779151D-01
               7.571689D-01
               7.751602D-01
               7.923042D-01
               8.195539D-01
               9.153427D-01
               1.000000D+00
%splkps(3,*)= -3.395425D+00  1.212759D-01 -3.548481D-01
              -3.357164D+00  1.597799D-01 -3.392315D-01
              -3.313898D+00  2.064323D-01 -3.130554D-01
              -3.223443D+00  2.863658D-01 -2.421025D-01
              -3.105357D+00  3.546861D-01 -1.224477D-01
              -3.045407D+00  3.706202D-01 -5.355577D-02
              -3.013525D+00  3.754747D-01 -1.616292D-02
              -3.004066D+00  3.763434D-01 -4.940086D-03
              -2.990684D+00  3.771333D-01  1.112437D-02
              -2.978901D+00  3.908625D-01  3.743601D-02
```

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```
-2.967886D+00  4.139995D-01  7.135396D-02
-2.950322D+00  4.495186D-01  1.244207D-01
-2.929548D+00  4.874999D-01  1.840116D-01
-2.909549D+00  5.184451D-01  2.369423D-01
-2.893230D+00  5.379440D-01  2.755771D-01
-2.875871D+00  5.504396D-01  3.101278D-01
-2.855329D+00  5.478961D-01  3.372107D-01
-2.847010D+00  5.385090D-01  3.415074D-01
-2.838927D+00  5.224665D-01  3.401418D-01
-2.826110D+00  4.764191D-01  3.214451D-01
-2.815766D+00  4.179281D-01  2.892525D-01
-2.770641D+00  2.926994D-01  2.555944D-01
-2.752024D+00  2.694927D-01  2.648844D-01
-2.733052D+00  2.511966D-01  2.785757D-01
-2.702545D+00  2.232360D-01  3.012929D-01
-2.595147D+00  1.103491D-01  3.583912D-01
-2.499980D+00  0.000000D+00  3.750000D-01
```

If the last 3D spline curve was formed using the **bps3** (B-Spline) option of the **curd** command, then the following parameters will be automatically set:

splks - 1 denemsion array containing all of the knots
splnk - scalar number of knots in the array splks
splkps - 2 dimensional array (3,*) of knot points on the 3D curve
splcps - 2 dimensional array (3,*) of control points that form the curve
splord - the order of the polynomials
splcn - **True*Grid*®** curve number that was created with this data
spltyp - type of spline curve - always 2

For example, if the following command were executed:

```
curd 12 bsp3
c knots
0.    0.    0.   0.   0.0625 0.0625 0.125 0.125 0.25  0.25
0.375 .4875 .5 .5 0.625  0.625 0.75  0.75  0.875 0.875
1. 1. 1. 1.;
c control points
-3.395424947 0.121275936 -0.354848062
-3.374553416 0.140465722 -0.348289595
-3.355863722 0.161100044 -0.339381434
-3.319725144 0.200588894 -0.317643003
-3.302032345 0.219324773 -0.304963004
-3.246632986 0.2703958   -0.263503838
```

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```
-3.208025634 0.2994159   -0.229995018
-3.133785624 0.344398736 -0.154716192
-3.096576553 0.360293935 -0.112937594
-3.02172923  0.376659885 -0.026342943
-2.985091236 0.377181147  0.017778912
-2.880315    6.026919D-01 3.450537D-01
-2.843035    5.878149D-01 3.868439D-01
-2.798396013 0.304679856  0.222954109
-2.760616573 0.276025848  0.257583401
-2.686456564 0.209309161  0.314211955
-2.648954363 0.170263866  0.336990097
-2.574723376 0.087828443  0.367203314
-2.537563881 0.044069575  0.375
-2.499979847 0.           0.375;;
```

Then the following automatic parameters would be set to have the values below:

```
%spltyp     =       2
%splcn      =      12
%splord     =       4
%splnk      =      24
%splkns     =  0.000000D+00
               0.000000D+00
               0.000000D+00
               0.000000D+00
               6.250000D-02
               6.250000D-02
               1.250000D-01
               1.250000D-01
               2.500000D-01
               2.500000D-01
               3.750000D-01
               4.875000D-01
               5.000000D-01
               5.000000D-01
               6.250000D-01
               6.250000D-01
               7.500000D-01
               7.500000D-01
               8.750000D-01
               8.750000D-01
               1.000000D+00
               1.000000D+00
```

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```
                1.000000D+00
                1.000000D+00
%splkps(3,*)= -3.395425D+00  1.212759D-01 -3.548481D-01
              -3.337794D+00  1.808445D-01 -3.285122D-01
              -3.337794D+00  1.808445D-01 -3.285122D-01
              -3.283566D+00  2.363484D-01 -2.911433D-01
              -3.283566D+00  2.363484D-01 -2.911433D-01
              -3.170906D+00  3.219073D-01 -1.923556D-01
              -3.170906D+00  3.219073D-01 -1.923556D-01
              -3.085229D+00  3.610343D-01 -9.952367D-02
              -2.992427D+00  3.770003D-01  8.962786D-03
              -2.975566D+00  3.976821D-01  4.753117D-02
              -2.975566D+00  3.976821D-01  4.753117D-02
              -2.820715D+00  4.462474D-01  3.048990D-01
              -2.820715D+00  4.462474D-01  3.048990D-01
              -2.723537D+00  2.426675D-01  2.858977D-01
              -2.723537D+00  2.426675D-01  2.858977D-01
              -2.611839D+00  1.290462D-01  3.520967D-01
              -2.611839D+00  1.290462D-01  3.520967D-01
              -2.499980D+00  0.000000D+00  3.750000D-01
%splcps(3,*)= -3.395425D+00  1.212759D-01 -3.548481D-01
              -3.374553D+00  1.404657D-01 -3.482896D-01
              -3.355864D+00  1.611000D-01 -3.393814D-01
              -3.319725D+00  2.005889D-01 -3.176430D-01
              -3.302032D+00  2.193248D-01 -3.049630D-01
              -3.246633D+00  2.703958D-01 -2.635038D-01
              -3.208026D+00  2.994159D-01 -2.299950D-01
              -3.133786D+00  3.443987D-01 -1.547162D-01
              -3.096577D+00  3.602939D-01 -1.129376D-01
              -3.021729D+00  3.766599D-01 -2.634294D-02
              -2.985091D+00  3.771811D-01  1.777891D-02
              -2.880315D+00  6.026919D-01  3.450537D-01
              -2.843035D+00  5.878149D-01  3.868439D-01
              -2.798396D+00  3.046799D-01  2.229541D-01
              -2.760617D+00  2.760258D-01  2.575834D-01
              -2.686457D+00  2.093092D-01  3.142120D-01
              -2.648954D+00  1.702639D-01  3.369901D-01
              -2.574723D+00  8.782844D-02  3.672033D-01
              -2.537564D+00  4.406957D-02  3.750000D-01
              -2.499980D+00  0.000000D+00  3.750000D-01
```

If the last 3D spline curve was formed using the **nrb3** (NURBS) option of the **curd** command, then the following parameters will be automatically set:

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splks - 1 denemsion array containing all of the knots
splnk - scalar number of knots in the array splks
splkps - 2 dimensional array (3,*) of knot points on the 3D curve
splcps - 2 dimensional array (3,*) of control points that form the curve
splwats - 1 dimensional array of weights
splord - the order of the polynomials
splcn - **True*Grid*®** curve number that was created with this data
spltyp - type of spline curve - always 3

For example, if the following command were executed:

```
curd 13 nrb3
c knots
0.    0.    0.   0.   0.0625 0.0625 0.125 0.125 0.25  0.25
0.375 .4875 .5 .5 0.625  0.625 0.75  0.75  0.875 0.875
1. 1. 1. 1.;
c weights
1. 1. 1. 1. 1. 1. 1. 1. 1. 1.
1. 1. 1. 1. 1. 1. 1. 1. 1. 1. ;
c control points
-3.395424947 0.121275936 -0.354848062
-3.374553416 0.140465722 -0.348289595
-3.355863722 0.161100044 -0.339381434
-3.319725144 0.200588894 -0.317643003
-3.302032345 0.219324773 -0.304963004
-3.246632986 0.2703958   -0.263503838
-3.208025634 0.2994159   -0.229995018
-3.133785624 0.344398736 -0.154716192
-3.096576553 0.360293935 -0.112937594
-3.02172923  0.376659885 -0.026342943
-2.985091236 0.377181147  0.017778912
-2.880315    6.026919D-01 3.450537D-01
-2.843035    5.878149D-01 3.868439D-01
-2.798396013 0.304679856  0.222954109
-2.760616573 0.276025848  0.257583401
-2.686456564 0.209309161  0.314211955
-2.648954363 0.170263866  0.336990097
-2.574723376 0.087828443  0.367203314
-2.537563881 0.044069575  0.375
-2.499979847 0.           0.375;;
```

Then the following automatic parameters would be set to have the values below:

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```
%spltyp     =       3
%splcn      =      13
%splord     =       4
%splnk      =      24
%splkns     =  0.000000D+00
               0.000000D+00
               0.000000D+00
               0.000000D+00
               6.250000D-02
               6.250000D-02
               1.250000D-01
               1.250000D-01
               2.500000D-01
               2.500000D-01
               3.750000D-01
               4.875000D-01
               5.000000D-01
               5.000000D-01
               6.250000D-01
               6.250000D-01
               7.500000D-01
               7.500000D-01
               8.750000D-01
               8.750000D-01
               1.000000D+00
               1.000000D+00
               1.000000D+00
               1.000000D+00
%splkps(3,*)= -3.395425D+00  1.212759D-01 -3.548481D-01
              -3.337794D+00  1.808445D-01 -3.285122D-01
              -3.337794D+00  1.808445D-01 -3.285122D-01
              -3.283566D+00  2.363484D-01 -2.911433D-01
              -3.283566D+00  2.363484D-01 -2.911433D-01
              -3.170906D+00  3.219073D-01 -1.923556D-01
              -3.170906D+00  3.219073D-01 -1.923556D-01
              -3.085229D+00  3.610343D-01 -9.952367D-02
              -2.992427D+00  3.770003D-01  8.962786D-03
              -2.975566D+00  3.976821D-01  4.753117D-02
              -2.975566D+00  3.976821D-01  4.753117D-02
              -2.820715D+00  4.462474D-01  3.048990D-01
              -2.820715D+00  4.462474D-01  3.048990D-01
              -2.723537D+00  2.426675D-01  2.858977D-01
              -2.723537D+00  2.426675D-01  2.858977D-01
```

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```
                -2.611839D+00  1.290462D-01  3.520967D-01
                -2.611839D+00  1.290462D-01  3.520967D-01
                -2.499980D+00  0.000000D+00  3.750000D-01
%splcps(3,*)= -3.395425D+00  1.212759D-01 -3.548481D-01
                -3.374553D+00  1.404657D-01 -3.482896D-01
                -3.355864D+00  1.611000D-01 -3.393814D-01
                -3.319725D+00  2.005889D-01 -3.176430D-01
                -3.302032D+00  2.193248D-01 -3.049630D-01
                -3.246633D+00  2.703958D-01 -2.635038D-01
                -3.208026D+00  2.994159D-01 -2.299950D-01
                -3.133786D+00  3.443987D-01 -1.547162D-01
                -3.096577D+00  3.602939D-01 -1.129376D-01
                -3.021729D+00  3.766599D-01 -2.634294D-02
                -2.985091D+00  3.771811D-01  1.777891D-02
                -2.880315D+00  6.026919D-01  3.450537D-01
                -2.843035D+00  5.878149D-01  3.868439D-01
                -2.798396D+00  3.046799D-01  2.229541D-01
                -2.760617D+00  2.760258D-01  2.575834D-01
                -2.686457D+00  2.093092D-01  3.142120D-01
                -2.648954D+00  1.702639D-01  3.369901D-01
                -2.574723D+00  8.782844D-02  3.672033D-01
                -2.537564D+00  4.406957D-02  3.750000D-01
                -2.499980D+00  0.000000D+00  3.750000D-01
%splwats(*) =  1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
                1.000000D+00
```

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igc create/append an IGES curve to a 3D curve

igc *IGES_curve transformations;*

where

IGES_curve is the sequence number of a curve in the current IGES file, and

trans is a transformation

Remarks

A composite IGES curve is treated as a single curve. After the 3D curve is defined, the **saveiges** and **useiges** commands should be used for fast retrieval of the IGES curve in subsequent sessions. See their command descriptions for more detail. A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

sdedge create/append curve by surface edge

sdedge *surface_#.edge_#*

where

surface_# is the surface ID number of a previously defined surface, and

edge_# is the number of the desired edge. The edge numbers range from 1 to 4.

Remarks

Use **sdedge** or **se** to create/append a 3D curve from an edge of a defined surface. Each edge is assigned a number. To view the surface with the numbered edge, use the graphics **labels** command with the option **sdedge** or **se**. It is a good idea to zoom in tight on the desired edge before turning on these labels, however, particularly if you are using read-in IGES surfaces. You cannot select the edge of an infinite surface such as a cone.

There is an advantage to converting an edge of a surface to a 3D curve and attaching an edge of the mesh to that 3D curve, instead of attaching the same edge of the mesh directly to the edge of the surface using the **edge** command. When using the **edge** command, all interior vertices of the edge of the mesh will be interpolated as a function of the two end vertices. This implies that you have no freedom to position these interior vertices. In contrast, when attaching a multiple edge to a 3D curve, you have freedom to position each vertex, independently, along the 3D curve. If you are attaching a simple edge, one with no interior vertices, then there is no obvious preference, unless, in the future reuse of this model you should choose to insert a partition (**insprt**). If the edge was attached to an edge of the surface using the **edge** command, you will not be able to position the new interior vertex.

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The interactive feature called **coedg** is an easy way to sequence many surface edges together to form a composite curve. The tsave file will record the **coedg** command as a sequence of **se** options in a **curd** command.

The order of the edges in a sequence of **sdeges** options determines the way the edges are sewn together. If there is a gap between two consecutive edges, the gap will be filled with a line segment.

Example

```
sd 1 function 0 10 0 10 u*u ; u*v ; v*v ; ;    c  surface 1
sd 2 function 2 10 10 15 u*u ; u*v ; v*v ; ;    c  surface 2
curd 1 sdedge 1.2 sdedge 2.2   c curve 1 definition from 2 edges
```

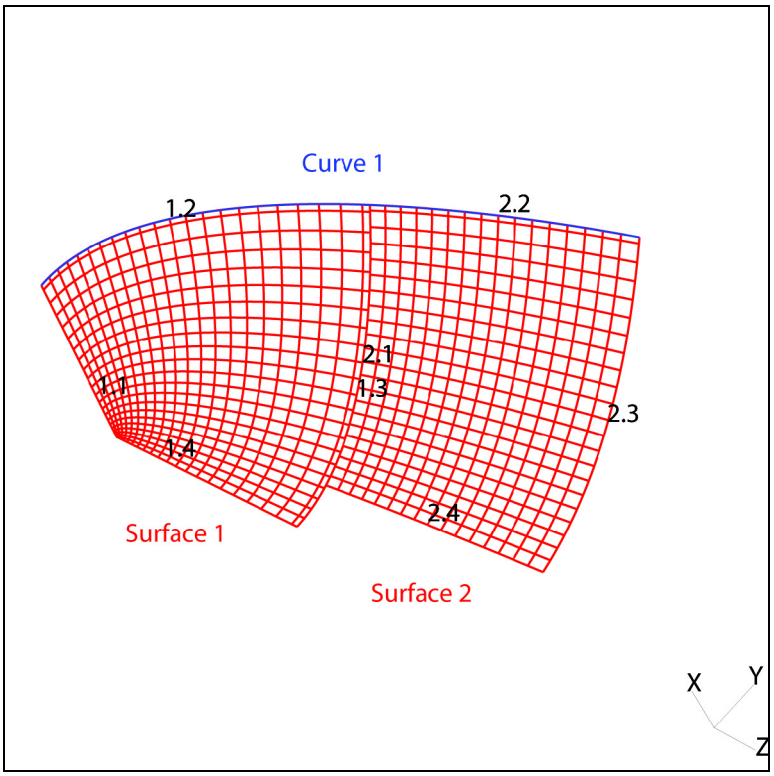

Figure 50 sdedge example

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lp3 create/append polygon of segments

lp3 *x1 y1 z1 ... xn yn zn* ; *trans* ;

where

x1 y1 z1 ... xn yn zn are triplets of coordinates of points forming a polygon, and

trans is a transformation.

Remarks

Use the **lp3** option to create/append a 3D curve by defining a series of points which are used to define line segments. A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

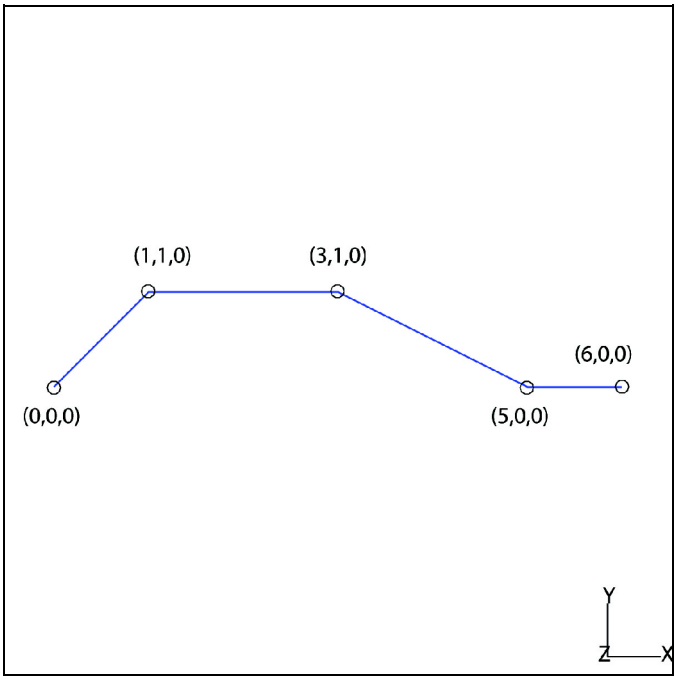

Figure 51 lp3 example

Example

```
curd 1 lp3 0 0 0 1 1 0
     3 1 0 5 0 0 6 0 0;;
```

contour create/append a contour line to a 3D curve

contour $point_1$ $point_2$

where

$point_1$ and $point_2$ identify points of a surface or 3D curve.

Remarks

The term contour is used to refer to a curve connecting two points of a surface or 3D curve. These are points used in the tessellation of the surface or curve and assigned point identifiers. To find the point identifiers, label the points of the surface or curve: either issue **labels sdpt** or **labels crvpt** or use the **Labels** button in the Environment Window. Then you will see point identifiers of the form *s.i.j* on a surface, or *c.i* on a curve. These points along a 3D curve are numbered sequentially. A similar

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scheme is used in two directions for a set of points on a surface. The surface is drawn as a grid of contour lines connecting some of these points on the surfaces (see **sdint**). When choosing a contour from a surface, the two points, $point_1$ and $point_2$, should differ only in their *i* values or only in their *j* values to extract an existing surface contour. *I* or *j* equal to **zero** means the maximum value of *i* or *j*, respectively. If both *i* and *j* are different in the two points, a smooth curve along the surface will be selected to connect the two points.

All of the coordinates of the contour line are written to the **tsave** file.

Example

The contour curve is created on the ruled surface.

```
ld 1 lp2 1 1 -1 1 -1 -1 1 -1 1 1;
ld 2 lp2 1 1;
lad .5 .5 360 ;
sd 1 rule2d 1 1 2 2 ;     c ruled surface definition
curd 1
contour 1.4.1 1.4.0 ;     c curve defined by first point 1.4.1 and
                          c last point 1.4.0 on contour 4 of surface 1
```

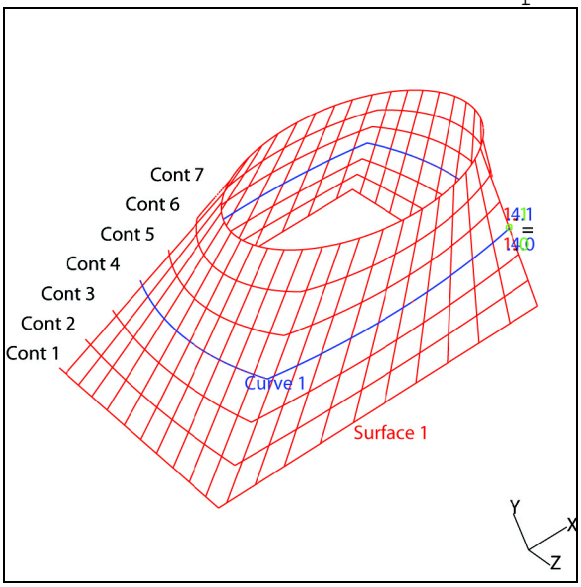

Figure 52 Contour Curve on Ruled Surface

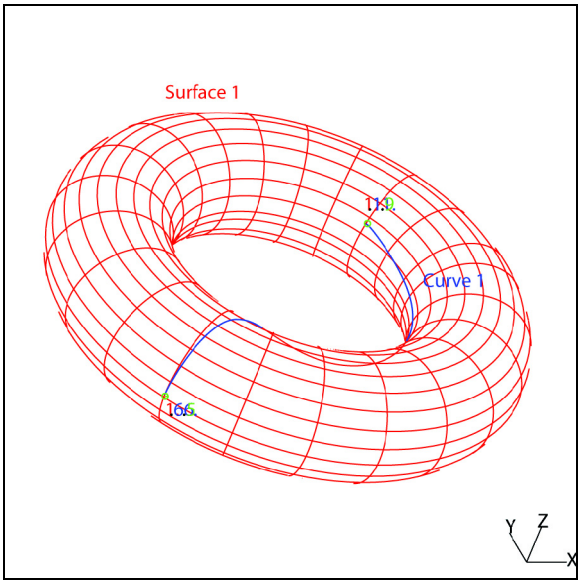

Figure 53 Contour Curve on Torus Surface

Example

You can also connect points from 2 different contours (66 and 11 in the i-index, 5 and 9 in the j-index).

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A curve is formed by seeking a short path (not necessarily the shortest path) between the first and last point. In a periodic surface, such as a torus, it will choose the shortest periodic path, but the smoothing algorithm may not converge to the to exact shortest path.

```
sd 1 ts 0 0 0 0 0 1 3 0 1           c torus surface definition
curd 1
contour 1.66.5 1.11.9;   c curve 1 defined by first

                         c point 1.66.5 and last point 1.11.9
```

csp3 create/append a 3D cubic spline curve

csp3 *option x1 y1 z1 ... xn yn zn* ; *trans* ;

where

option can be

loop close curve into a smooth loop

First Endpoint Natural Derivative

00 natural derivative for the Last Endpoint
01 *last_dx last_dy last_dz* - derivatives at the Last Point
02 *#_curve_2 derivative_mag_2* - match First Endpoint of Another Curve
03 *#_curve_2 derivative_mag_2* - match Last Endpoint of Another Curve

First Endpoint Specify Derivative

10 *first_dx first_dy first_dz*
11 *first_dx first_dy first_dz last_dx last_dy last_dz*
12 *first_dx first_dy first_dz #_curve_2 derivative_mag_2*
13 *first_dx first_dy first_dz #_curve_2 derivative_mag_2*

First Endpoint Match First Endpoint of Another Curve

20 *#_curve_1 derivative_mag_1*
21 *#_curve_1 derivative_mag_1 last_dx last_dy last_dz*
22 *#_curve_1 derivative_mag_1 #_curve_2 derivative_mag_2*
23 *#_curve_1 derivative_mag_1 #_curve_2 derivative_mag_2*

First Endpoint Match Last Endpoint of Another Curve

30 *#_curve_1 derivative_mag_1*
31 *#_curve_1 derivative_mag_1 last_dx last_dy last_dz*
32 *#_curve_1 derivative_mag_1 #_curve_2 derivative_mag_2*
33 *#_curve_1 derivative_mag_1 #_curve_2 derivative_mag_2*

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x1 y1 z1 ... xn yn zn - triplets of coordinates of Control Points

Theory

A cubic spline curve is composed from segments (**Figure 54**). Segment *i* is described by parametric algebraic equations:
where *t* is an independent variable, and *a,b,c,d,e,f,g,h,j,k,l,m* are unknown constants. There are 12

$$x(t)=a_i t^3+b_i t^2+c_i t+d_i$$
$$y(t)=e_i t^3+f_i t^2+g_i t+h_i$$
$$z(t)=j_i t^3+k_i t^2+l_i t+m_i$$

unknown constants for segment *i*. Suppose, there are *n* segments in the cubic spline. This produces *12n* unknowns. There are *6n* constraints since the endpoints of each section pass through the control points. There are *6n-6* additional constraints imposed so that the curve has equal left and right first derivatives at the interior control points.

You can choose the derivatives at the end points to impose the remaining 6 constraints.
You have 2 possible choices:

- **first derivatives** at the Endpoints
- **natural derivatives** at the Endpoints (second derivatives = 0)

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

Example

You can insert control point in between all ready defined control points from **Figure 54**. The result is shown in **Figure 55**.

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```
curd 1
csp3 00 c natural derivative
0 1 0
1 0 0
2 1.3 0
3 0 0
3 0.8 0  c inserted point
4 1 0;;;
```

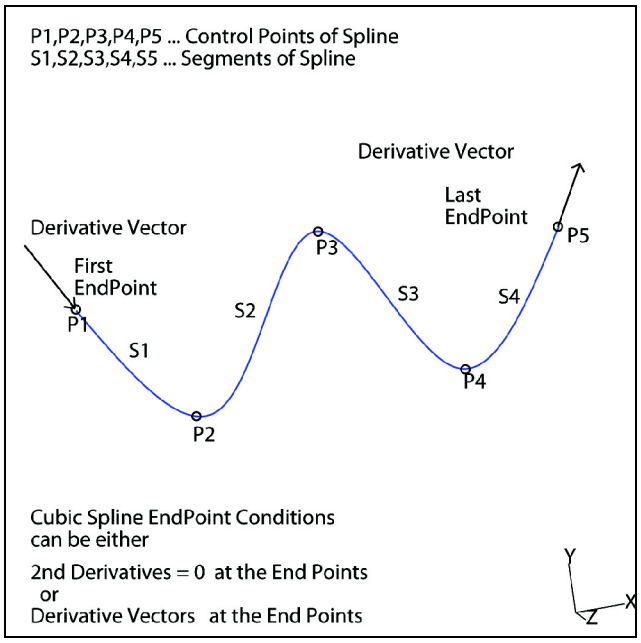

Figure 54 Spline Curve

Example

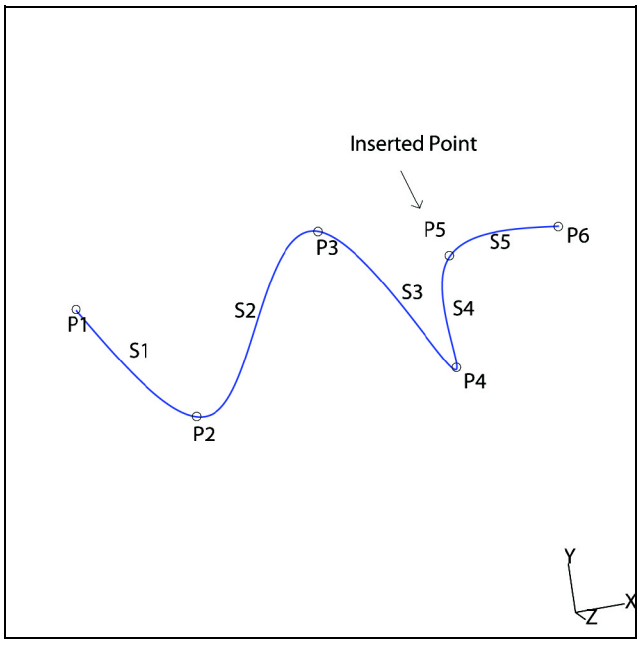

Figure 55 Inserting a Cubic Spline Control Point

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You can define a derivative condition for the previously defined control points in **Figure 54**. The result is shown in **Figure 56**.

```
curd 1
csp3 10 1 3 0 c derivatives
0 1 0
1 0. 0
2 1.3 0
3 0 0
4 1 0 ;;;
```

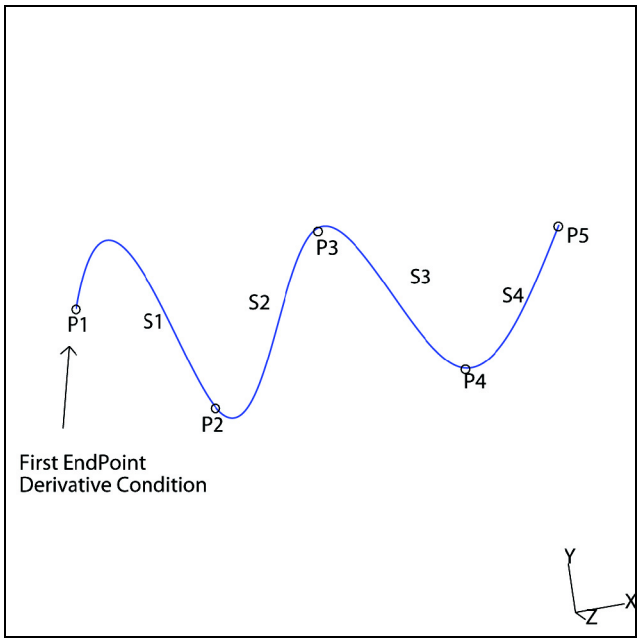

Figure 56 Endpoint Derivative Condition

bsp3 create/append a B-Spline curve

bsp3 $knot_1 \ldots knot_n$; $x_1\ y_1\ z_1 \ldots x_k\ y_k\ z_k$; *trans* ;

where

$knot_1 \ldots knot_n$ are the parametric coordinates of *n* knots

Remarks

The multiple knots have to be specified for the starting and ending point of the curve. The number of multiple knots for the starting and ending point determines the order of the curve (e.g. 2 - linear, 3 - quadratic, 4 - cubic, etc.). The multiplicity has to be the same for the starting and ending point.

$x_1\ y_1\ z_1 \ldots x_k\ y_k\ z_k$ are triplets of coordinates of control points.

where $k = n$ - *multiplicity*

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

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Example

```
curd 2 bsp3
     0 0 0 1 2 3 3 3; c knots
     0 0 0           c control
     1 1 1           c points
     2 1 1
     2 0 1
     3 2 1; ;

c bspline curve definition
c 4 knots
c 0 0 0 multiple knot
c 3 3 3 multiple knot
c 5 control points
c P1...P5
```

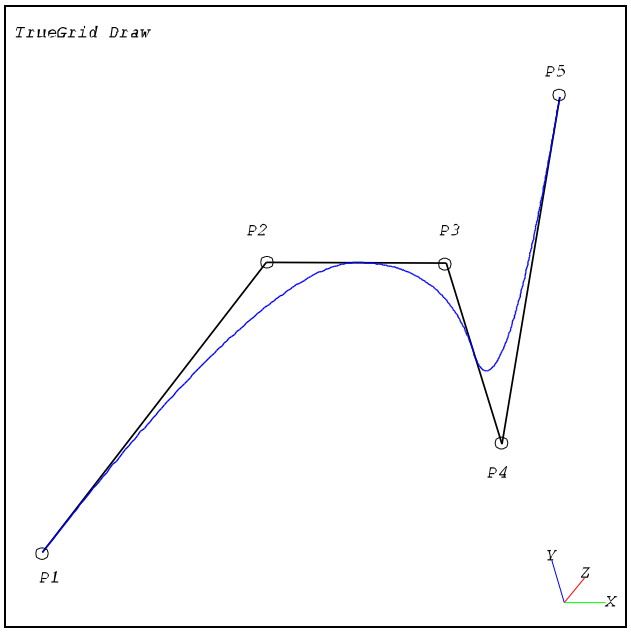

Figure 57 B-spline curve

nrb3 create/append a NURBS curve

nrb3 $knot_1 \ldots knot_n$; $weight_1 \ldots weight_k$; $x_1\ y_1\ z_1 \ldots x_k\ y_k\ z_k$; *trans* ;

where

$knot_1 \ldots knot_n$	are parametric coordinates of n knots
$weight_1 \ldots weight_k$	are weights for each control point
$x_1\ y_1\ z_1 \ldots x_k\ y_k\ z_k$	are triplets of coordinates of control points.

where $k = n$ - *multiplicity*.

Remarks

The **multiple knots** have to be specified for the starting and ending point of the curve. The number of multiple knots for the starting and ending point determines the order of the curve (e.g. 2 - linear, 3 - quadratic, 4 - cubic, etc.). The multiplicity has to be the same for the starting and ending point.

This command creates/appends a NURBS curve formed from an explicit list of knots, weights and control points.

You can shape the curve by changing the weights (**Figure 58**).

You can specify discontinuity at the control point by multiple internal knots (**Figure 59**).

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A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

Example

```
curd 2 nrb3
 0 0 0 1 2 4 5 5 5;   c knots
.1 1 1 5 1 1;         c weights
     0 1 0           c control
     1 0 0           c points
     2 1 0
     3 4 0
     4 2 0
     5 3 0;;
c curve definition - 2
c nurbs
c 0 0 0 multiple knot
c 5 5 5 multiple knot
c 6 control points
c P1 ...P6
c weight = 5 for P4 (x,y,z)

curd 3 nrb3
 0 0 0 1 2 4 5 5 5;   c knots
.1 1 1 .1 1 1;        c weights
     0 1 0           c control
     1 0 0           c points
     2 1 0
     3 4 0
     4 2 0
     5 3 0;;
c curve definition - 3
c nurbs
c 0 0 0 multiple knot
c 5 5 5 multiple knot
c 6 control points
c P1 ... P6
c weight = .1 for P4
c weight = 5 for P4
```

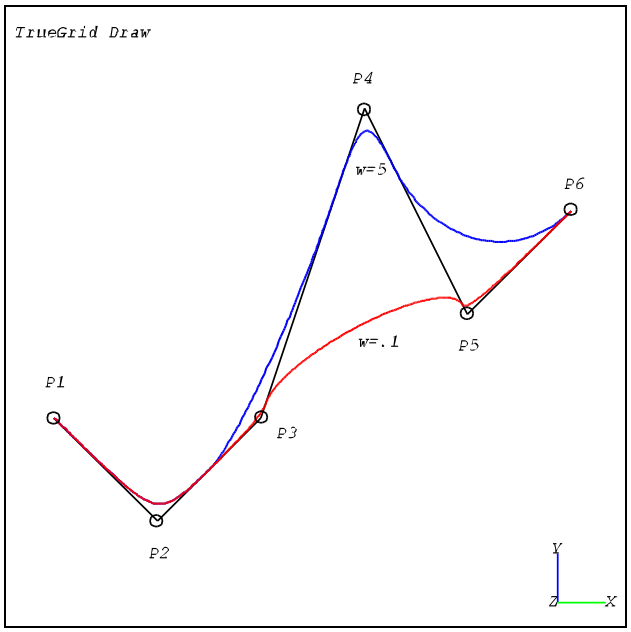

Figure 58 2 NURBS curves with various weights

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Example

```
curd 4 nrb3
0 0 0 1 1 2 4 5 5 5; c knots
.1 1 1 1 10 1 1;  c weights
     0 1 0   c control points
     0 .5 0
     1 0 0
     2 1 0
     3 4 0
     4 2 0
     5 3 0;;

c NURBS curve with
c discontinuity - multiple
c internal knots 1 1
c for P3
```

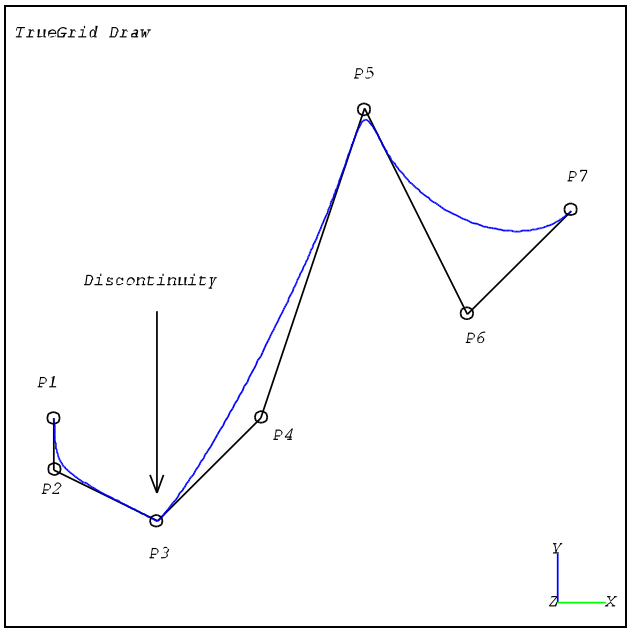

Figure 59 NURBS curve with discontinuity of the slope in P3

ld2d3d create/append a 2D curve converted to 3D

ld2d3d *2D_curve_# system coordinate start end transform ;*

where

system can be

rt for Cartesian coordinate interpolation,
cy for cylindrical coordinate interpolation, or
sp for spherical coordinate interpolation

coordinate can be

x for the first coordinate interpolated
y for the second coordinate interpolated
z for the third coordinate interpolated

start is the first value of the specified coordinate,
end is the last values of the specified coordinate, and
transform is an optional transformation that may be applied.

Remarks

The interpolation formula is: $x = (1-\alpha)X_0 + \alpha X_1$

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where x is the interpolated coordinate,

α is a parametric coordinate, it goes from 0 to 1, and

X_0, X_1 are the coordinates of the starting and ending points.

The parametric coordinate is segmented, with the segmentation based on the arc length of the curve. See the examples on the following page.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

Example

```
ld 1 lep 2 1 0 0 -100 100 0 ;
c 2d curve definition
c elliptic arc

curd 1 ld2d3d 1 rt x 2 2 ;
c 3d curve - 1 definition
c elliptic arc
c in plane

curd 2 ld2d3d 1 rt x 2 -2 ;
c 3d curve - 2 definition
c elliptic arc
c with nonuniformly
c interpolated coordinate x
```

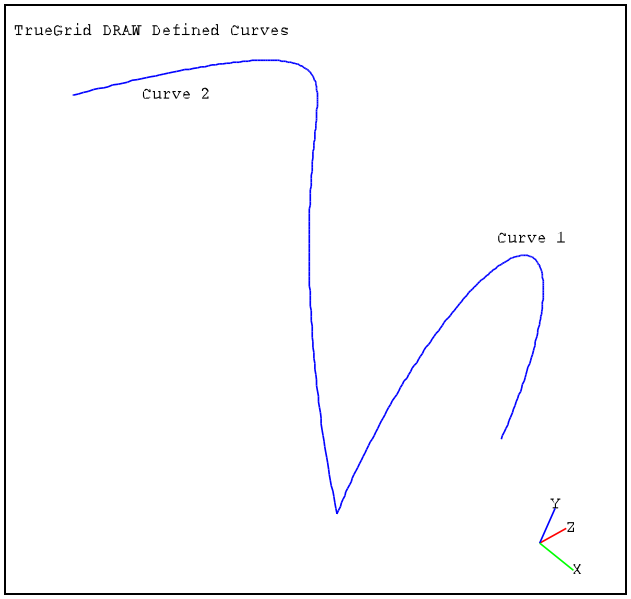

Figure 60 curve 1 and curve 2 by ld2d3d

intcur create/append a 3D curve by interpolation

intcur *3D_curve_1 3D_curve_2 interpolation_parameter transform ;*

where

3D_curve_1	is the ID number of a 3D curve,
3D_curve_2	is the ID number of the second 3D curve,
interpolation_parameter	is the distance from curve_1 (towards the second), and
transform	is a final transformation, which may be applied.

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Remarks

The **intcur** command is especially useful for defining the inside curves for a butterfly mesh which must be applied along curved geometry.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

Example

Curve **4** is created by interpolation between curves **2** and **3** with interpolation parameter .3 (**Figure 61**). Simplified command file:

```
curd 4 intcur 2 3 .3  ;
c definition of curve 4 by
c interpolation
c between curves 2 and 3
```

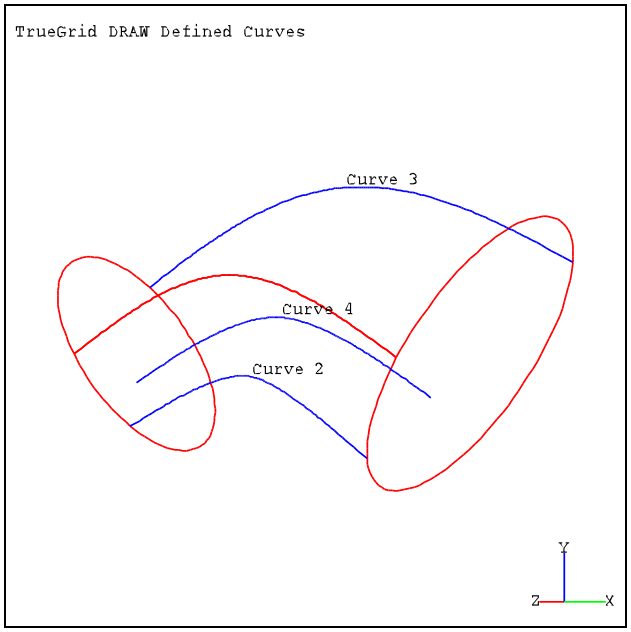

Figure 61 curve 4 by intcur

lp3pt create/append a 3D curve by pairs of defined points

lp3pt *point_id_1 ... point_id_n ; trans ;*

Remarks

The **lp3pt** command connects points on surface/curve by line segments. In the case of infinite surfaces (cylinder, plane, cone...), the surface must first be displayed. You can label these points by choosing ***Labels*** and ***Surf Point***. If you do not see a label where you expect one, then zoom in a bit further. If you click on a label you can then enter that label (or point ID) using the **F8** function key.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

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Example

```
sd 1 function -90 90 -90 90
   cos(u); cos(v);cos(u)  ;;
curd 1
lp3pt 1.50.3 1.40.6 1.30.50
     1.24.54 1.2.1  ;  ;
     c curve 1 is created
     c from segments
     c connecting points
     c 1.2.1  1.24.54  1.30.50
     c 1.40.6 1.50.3
```

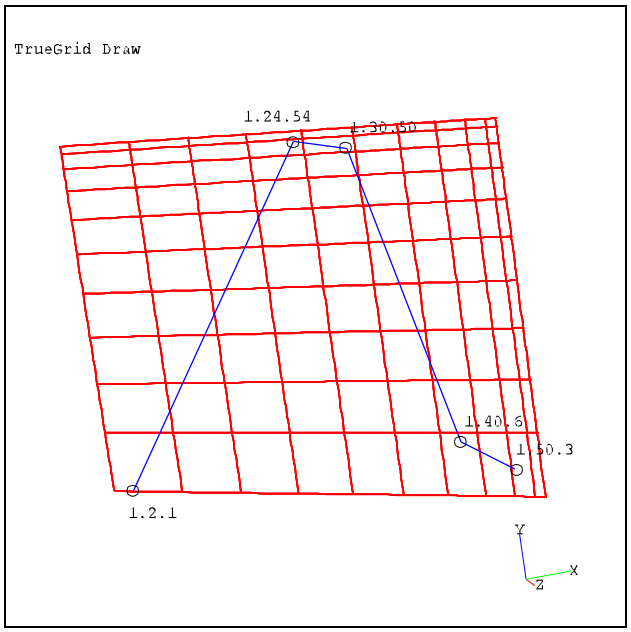

Figure 62 curve by lp3pt

3dfunc create/append a parameterized function curve

3dfunc *min_u max_u x_expres ; y-expres ; z-expres ; trans ;*

where

min_u	is the minimum value of the parameter, u,
max_u	is the maximum value of the parameter,
x_expres	is the function of u which defines the x-coordinates,
y-expres	is the function of u which defines the y-coordinates,
z-expres	is the function of u which defines the z-coordinates, and
trans	is an optional final transformation.

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Remarks

This command enables you to define a parameterized 3D curve over a specified domain.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

Example

```
curd 1 3dfunc -90 90 cos(u);
sin(u) ; sin(u)*cos(u) ; ;
c u goes from -90 to 90
c x = cos(u)
c y = sin(u)
c z = sin(u)*cos(u)
```

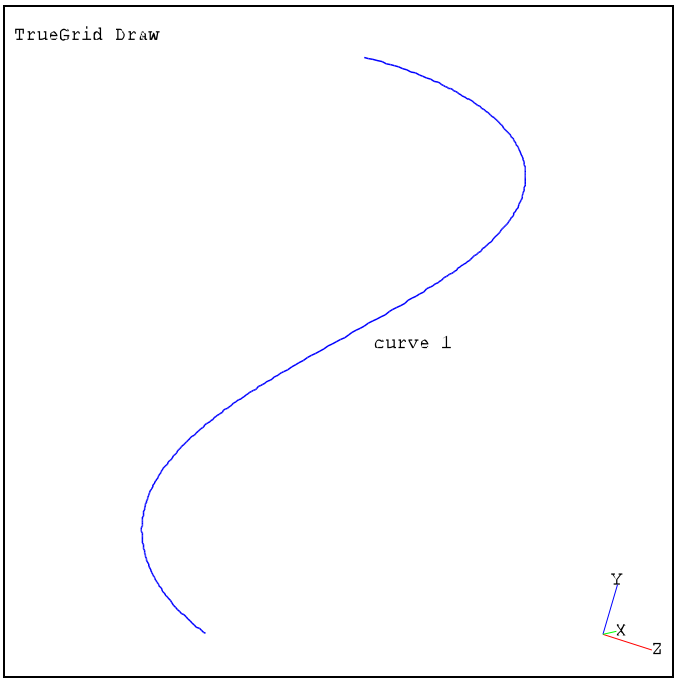

Figure 63 curve by 3dfunc

projcur create/append curve projected onto a surface

projcur *3D_curve_# surface_# trans ;*

where

3D_curve_# is the ID number of a previously-defined 3D curve,
surface_# is the ID number of a previously-defined surface, and
trans is an optional transformation.

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Remarks

The **projcur** command creates a curve by projecting an existing curve onto a surface. Projection is sensitive to surface curvature and/or the relative position of the curve to the surface. Also, see the following **pscur** command, which is similar to this command except that it also smooths the projected curve.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

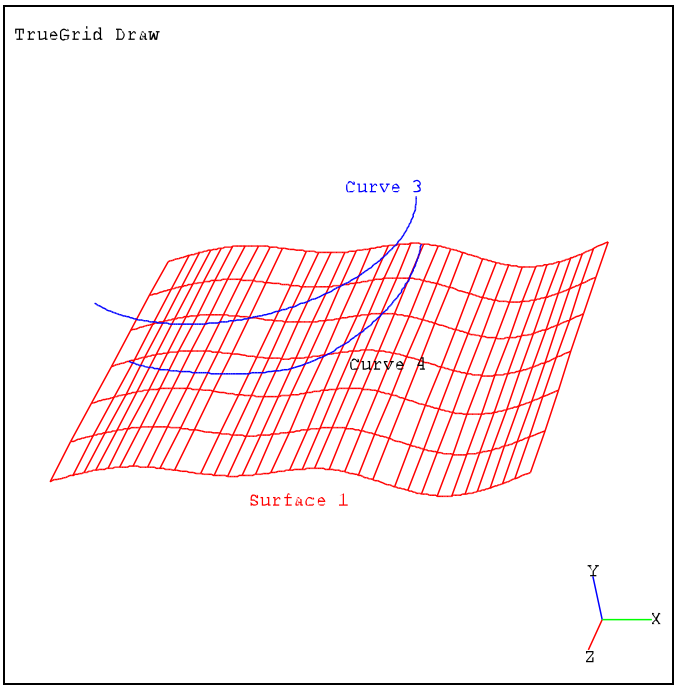

Figure 64 curve by **projcur**

Example

```
curd 4 projcur 3 1 ;
c curve 4 by projection of curve 3 onto surface 1
```

pscur create/append curve projected onto a surface and smoothed

pscur *3D_curve_# surface_# #_iterations tol trans ;*

where

3D_curve_#	is the ID number of an existing 3D curve,
surface_#	is the ID number of an existing surface,
#_iterations	is the number of iterations used,
tol	is the tolerance which limits the smoothing iterations, and
trans	is an optional final transformation.

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Remarks

This command projects the specified 3D curve onto the specified surface and then iteratively smooths the final curve. This iterative method continues until the differences between successive iterations are less than *tol*.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

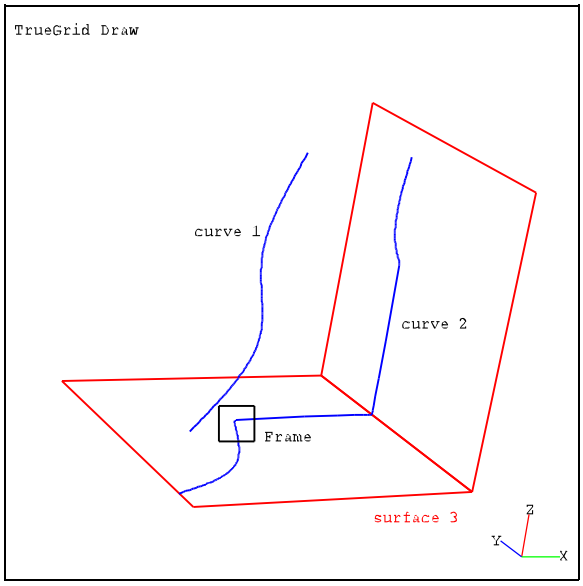

Figure 65 curve2 by pscur

Example

The **pscur** command is demonstrated by projection and smoothing of the spline curve onto discontinuous planar surface (**Figure 65**). **Figure 66** is created from **Figure 65** by magnification of the Frame. It shows the difference between the **pscur** and **projcur** command when applied to the same curve. **pscur** creates the smoother curve.

```
curd 2 pscur 1 3 100 .01  ;

     c curve 2 - by projection
     c and smoothing of
     c curve 1 onto surface 3
     c with maximum
     c 100 iterations
     c and with minimum
     c tolerance .1
```

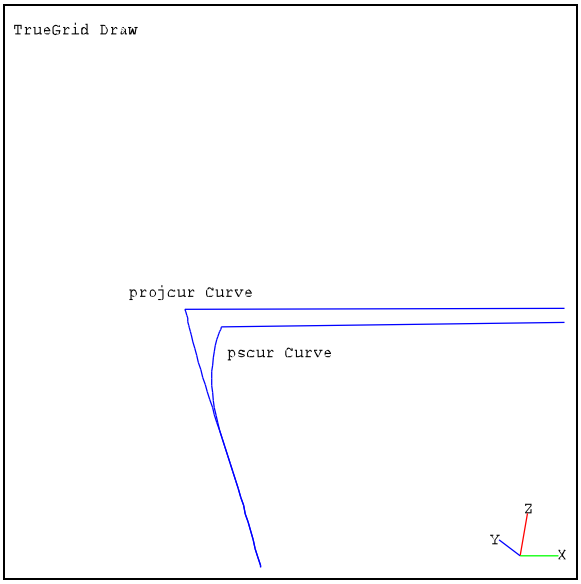

Figure 66 pscur vs projcur curve

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arc3 create/append arc of a circle

arc3 *option system point system point system point*

where

option can be

seqnc to specify the arc starting at the first point, passing through the second, and ending at the third point.

cmplt to specify the complementary arc of the circle specified in seqnc, or

whole to specify the entire circle, and

system and *point* can be

rt *x y z* for cartesian coordinates,

sp *rho theta phi* for spherical coordinates,

cy *rho theta z* for cylindrical coordinates,

pt *surface.i.j* for a point off of a surface, or

pt *curve.i* for a point off of a 3D curve.

Remarks

This command appends a 3D arc of a circle passing through 3 points. There are three options. An arc can be created passing from the first to the second to the third point - **seqnc** option (**Figure 68**). The second option - **cmplt** is the complement of the first option above (**Figure 68**). The third option - **whole** creates the entire circle (**Figure 69**).

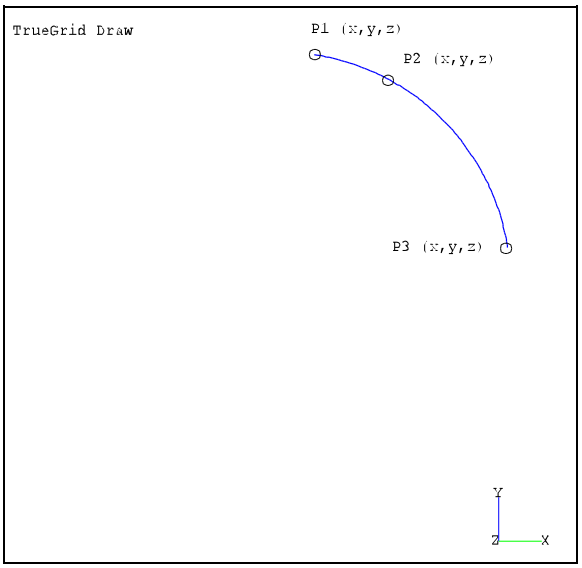

Figure 67 curve by arc3 seqnc

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Examples

```
curd 1 arc3  seqnc  rt 0 1 0  rt
.5 .8 0  rt 1 0 0 ;
c  circular arc - seqnc
c  given by points
c  P1(0,1,0)
c  P2(.5,.8,0)
c  P3(1,0,0)
```

```
curd 1 arc3
cmplt rt 0 1 0
rt .5 .8 0  rt 1 0 0 ;
c  circular arc - cmplt
c  given by points
c  P1(0,1,0)
c  P2(.5,.8,0)
c  P3(1,0,0) curd 1 arc3
 seqnc rt 0 1 0
 rt .5 .8 0  rt 1 0 0 ;
 c  circle - whole
 c  given by points
 c  P1(0,1,0)
 c  P2(.5,.8,0)
 c  P3(1,0,0)
```

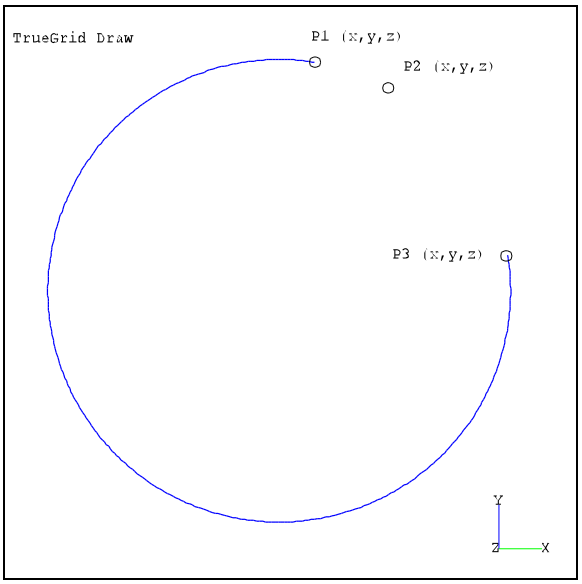

Figure 68 curve by arc3 cmplt

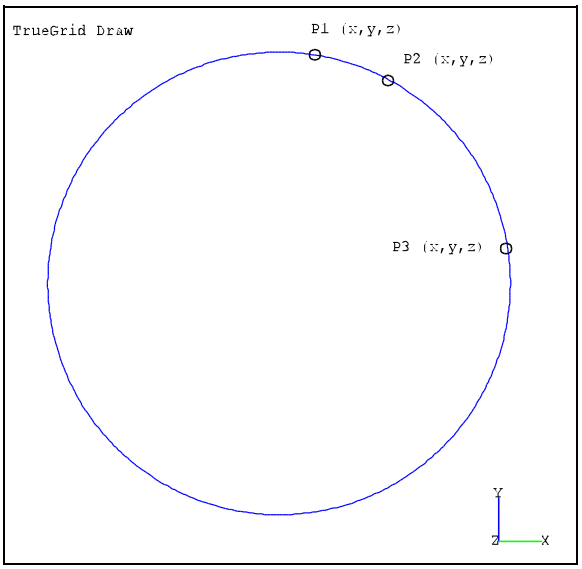

Figure 69 curve by arc3 whole

cpcd create/append a copy of a previously defined 3D curve.

cpcd *curve_# trans* ;

where

curve_# is the ID number of a previously-defined 3D curve, and

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trans is an optional final transformation.

Remarks

This option allows you to copy a previously-defined 3D curve and transform it.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

Example

```
c definition of spline curve
curd 1 csp3 loop
     0  -1   0
     3  -1.5 0
     5  -1   0
     5   0   0
     5.1 1   0
     3   .5  0
     0   1   0; ;
curd 2 cpcd 1 mz 5 csca .5 ;
c definition of curve 2
c  by copy from curve 1
c and moved in the z-direction
c and scaled to the half size
```

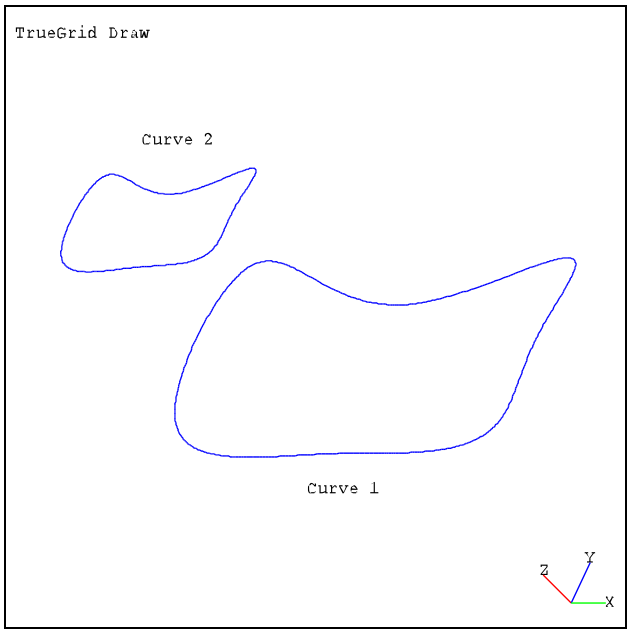

Figure 70 curve 2 by **cpcd** from curve 1

cpcds create/append a copy of previously defined 3D curves

cpcds *list_curve_#* *trans* ;

where

list_curve_# is an ordered list of 3D curve ID numbers, and

trans is an optional final transformation.

Remarks

The *list_curve_#* is ordered (**Figure 71**).

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

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Example

```
curd 5 cpcds 1 2 3 4;
rzx        c symetry - zx plane
my -1      c translation - y
mz .5      c translation - z
rz 10 ;    c rotation ar.- z

c definition of curve 5
c by copy of curves 1,2,3,4,5
```

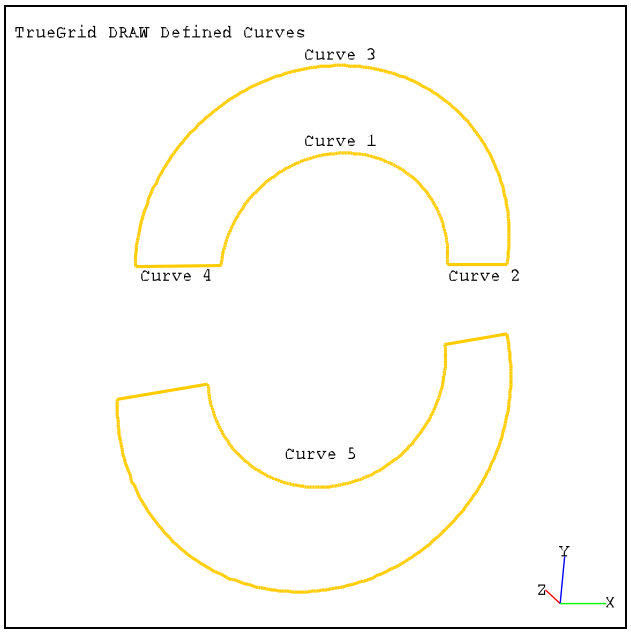

Figure 71 curve 5 by cpds

twsurf create/append the curve at the intersection of two surfaces

twsurf *surface_1 surface_2 x1 y1 z1 ... xn yn zn ; trans ;*

where

surface_1	is the ID number of the first surface,
surface_2	is the ID number of the second surface,
x1	is the approximate x-coordinate of the curve's beginning,
y1	is the approximate y-coordinate of the curve's beginning,
z1	is the approximate z-coordinate of the curve's beginning,
xn	is the approximate x-coordinate of the curve's end,
yn	is the approximate y-coordinate of the curve's end,
zn	is the approximate z-coordinate of the curve's end, and
trans	is an optional final transformation.

Remarks

This command creates/appends the curve at the intersection of two surfaces. Initial points are required to determine the endpoints and some intermediate points for surfaces with large curvature (greater than 120 degrees). These points need only be approximate.

A transformation is formed by a sequence of simple transformations such as rotate and translate. For a complete list of these simple transformations, see the **lct** command.

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Example

A curve is created at the intersection of a torus and a cylinder. The initial points are picked by ***Projection*** on Surface 1 (**Figure 72**). The points form an oriented curve, which is used as a first approximation of intersection. The curve is closed by coincident starting and ending initial points (**Figure 73**). The command file follows:

```
sd 1 ts 0 0 0 0 0 1 1 0 .5
     c definition of torus
     c - surface 1

sd 2 cy .9 .9 0 0 .0 1 .4
     c definition of
     c cylinder - surface 2

curd 1
twsurf 1 2
 5.47e-01 1.06e+00  4.60e-01
 4.99e-01 8.72e-01  4.99e-01
 7.70e-01 5.31e-01  4.95e-01
 1.20e+00 6.59e-01  3.29e-01
 1.26e+00 7.52e-01 -1.74e-01
 9.79e-01 4.96e-01 -4.90e-01
 5.77e-01 6.65e-01 -4.85e-01
 5.48e-01 1.09e+00 -4.47e-01
 7.76e-01 1.28e+00  4.41e-02
 5.47e-01 1.06e+00 4.60e-01;;;

     c definition of curve 1
     c at the intersection of
     c surface 1 and 2 with 9
     c initial points
     c (coincident starting
     c and ending initial
     c point)
```

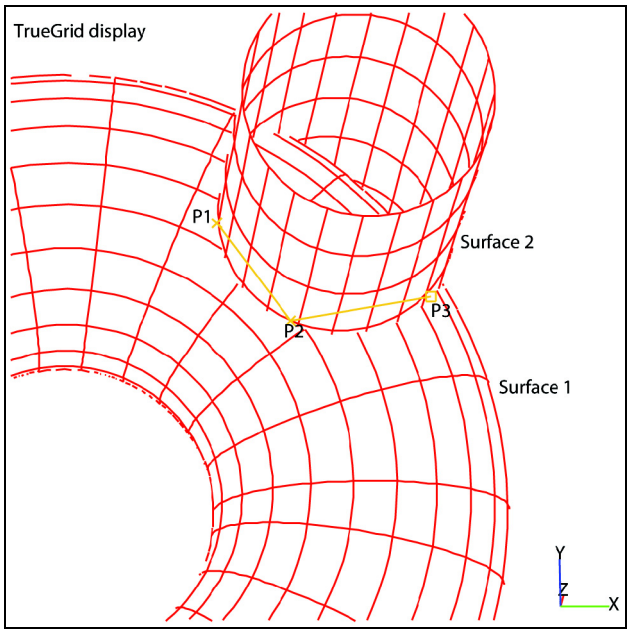

Figure 72 First 3 initial points of Curve 1 picked by Projection on Surface 1

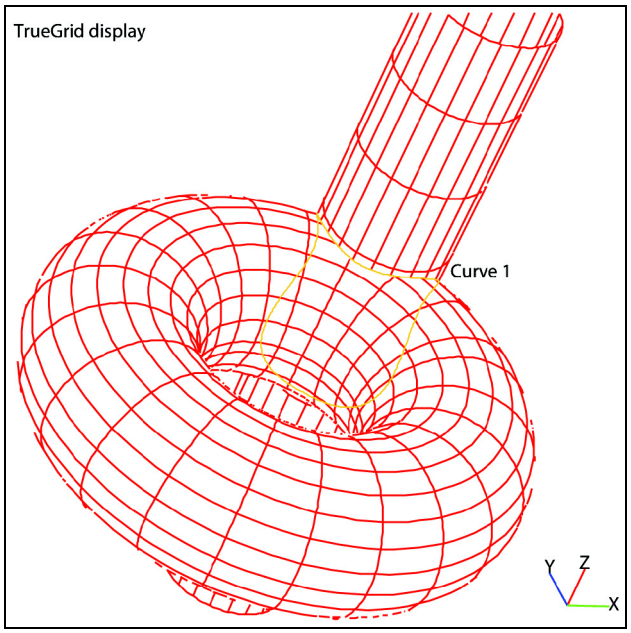

Figure 73 Curve by **twsurf**

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5. Display 3D Curve

dcd display a 3D curve

dcd *3d_curve_#*

dcds display a set of 3D curves from a list

dcds *3d_curve_list* ;

dacd display all 3D curves

dacd (no arguments)

acd add a 3D curve to the picture

acd *3d_curve_#*

acds add a list of 3D curves to the picture

acds *3d_curve_list* ;

rcd remove a 3D curve from the picture

rcd *3d_curve_#*

rcds remove a list of 3D curves from the picture

rcds *3d_curve_list* ;

racd remove all 3D curves from the picture

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racd (no arguments)

lacd list all of the active 3D curves in the picture

lacd (no arguments)

6. Print 3D Curves

cdinfo print information about the 3D curves

cdinfo <*no arguments*>

Remark

Prints the number of points, number segments, and the length of each 3D curve.

Example

```
igescd 42 45 3;      c evaluate IGES curves 42-45
labels sdedge        c show the edge labels to the surfaces
curd 1 sdedge 2.2 sdedge 2.4
rmseg
sdedge 2.1 sdedge 2.4
cdinfo c write information about the 3D curves
```

The **cdinfo** command writes the following table:

```
3D CURVE INFORMATION
 curve      1 has     71 points,    3 segments and length 2.10000
 curve      3 has    257 points,    1 segments and length 3.14159
 curve      4 has      2 points,    1 segments and length 0.10000
 curve      5 has      2 points,    1 segments and length 0.20000
 curve      6 has    188 points and      1 segments
```

7. Delete 3D Curve

rmseg remove last segment from 3D curve

rmseg <*no arguments*>

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Remark

Remove the last segment appended to a 3D curve by **curd**.

delcd delete a 3D curve from the True*Grid*® data base

delcd *3d_curve_#*

Remarks

The deletion of a 3D curve cannot be undone. If you wish to remove a 3D curve from the picture, use **rcd**.

delcds delete a list of 3D curves from the True*Grid*® data base

delcds *3d_curve_list* ;

Remarks

The deletion of a 3D curve cannot be undone. If you wish to remove a list of 3D curves from the picture, use **rcds**.

8. Surfaces

The following is a full description of all available methods for creating a surface. Commands that require a surface, such as the **sf** and **sfi** commands refer to a surface defined using the **sd**, **vpsd**, **iges**, **igessd**, **igespd**, or **nurbs** commands. Of this group of surface definition commands, only the **sd** command can specify the shape of the surface. The other surface definition commands import surfaces from a file and use the shape of the surface described in the file. Surfaces defined with the **sd** command have the advantage of, in many cases, of being simple, and in all cases can be formed using parameters and expressions so that you can have any degree of parameterization you desire. The disadvantage is that the parametric form of the surfaces can be less intuitive and require a more analytic approach to designing the desired shape.

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When a surface is defined, it must be tessellated to be drawn. This tessellation is a set of connected points forming linear polygons as an approximation to the surface. It is these polygons that appear in the drawing. In many cases, they look like a block mesh. These tessellation points are assigned labels of the form s.i.j where s is the surface number, i is the first index of the point in the grid of tessellation, and j is the second index. In a more complex surface with trimming curves or in a polygon surface, there are usually triangles. The same notation is used but with a different meaning. The first index is the polygon number. The second index is the sequence number of the point used to define the polygon.

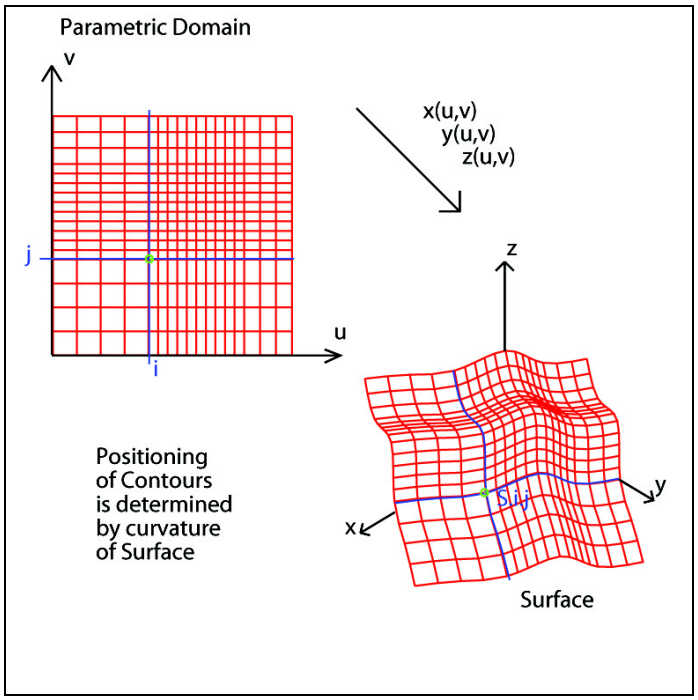

Figure 74 Tessellation

In some surface types (**csps**, **bsps**, **nrbs**, and **hermite**), you can specify boundary conditions (i.e. the derivatives along the edges). The easiest thing to do is to select the natural boundary condition. This means that the 2nd derivatives are set to zero. If you choose to specify the derivatives along an edge of the surface, you must specify the derivative for each control point along that edge. For example, suppose you choose to build a cubic spline surface with 6 control points in each row (u-direction) and 3 control points in each column (v-direction) of the table of control points. The edges are numbered. The edge corresponding to the minimum u is edge 1, maximum u is 2, minimum v is 3, and maximum v is 4. If you chose to specify derivatives along edge 1 or 2, you will need 3 derivatives. If you chose to specify derivatives along edge 3 or 4, you must specify 6 derivatives. Each derivative is a vector

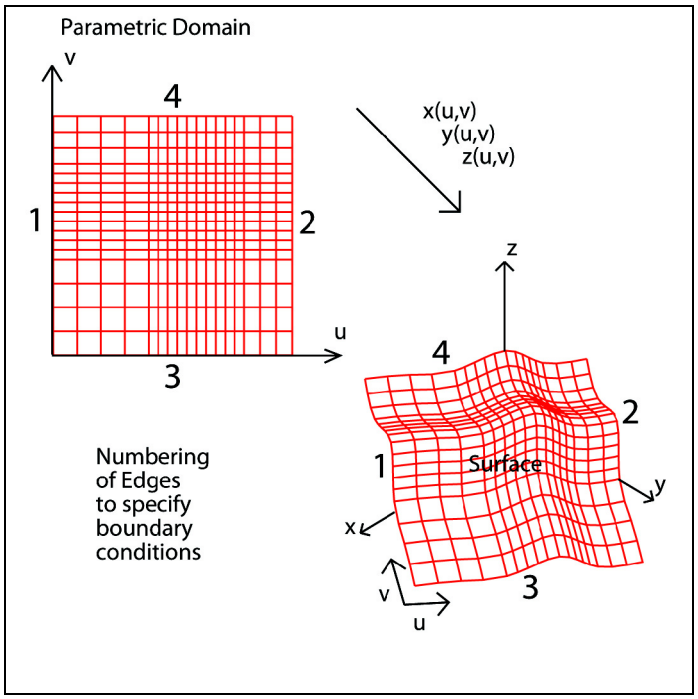

Figure 75 Numbered edges

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with three components forming the direction vector for the derivative. A flag formed by 4 digits are used to flag the type of boundary condition for each of the corresponding edges. 0 means a natural derivative along the edge and 1 means that the derivatives are being specified. All derivatives follow this 4 digit flag, in order. In some cases, you can specify a loop by setting either the first 2 digits in the flag or the last 2 digits to 2.

In the case of infinite surfaces (**plan**, **xyplan**, **yzplan**, **zxplan**, **pl3**, **pl3o**, **iplan**, **cy**, **cyr2**, **cyr3**, **xcy**, **ycy**, **zcy**, **cp**, **cn2p**, **pr**), the surface tessellation depend on the size of the bounding box for the graphics. The size of the bounding box is determined from the size of the finite surfaces in the model. In an infinite surface, the points used to tessellate the surface are recalculated after any change of the bounding box. The bounding box is recalculated with the **restore** command. The bounding box can be displayed by the **grid** command.

If an infinite surface is not drawn, which is usually the case when a command file is run in batch mode, then the points of tessellation will not be calculated. When you select points or edges of a surface symbolically (by label), be sure that surface has been drawn. When these points are selected by label, the coordinates are substituted in the save file, not its label.

accuracy set accuracy of surface projections

accuracy *acc*

where *acc* is the scale factor relative to the default accuracy of the projection

Remarks

The default accuracy parameter is 1.0 which produces approximately 3.5 digits of accuracy. The accuracy parameter must be positive. The projection to a single surface is more accurate when this parameter is increased. When an edge is projected to two surfaces or a vertex projected to three surfaces, then increasing this parameter will cause the intersection algorithm to be more accurate. There is a natural limit to this accuracy. On a 32 bit version, the maximum precision is approximately 7 digits. Therefore the accuracy parameter cannot exceed 500. On a 64 bit version, the accuracy parameter cannot exceed 5,000,000,000, which will try to get 14 digits 0f accuracy.

This accuracy is ideal. The iterative Newton algorithm has two termination criterium. The first is a maximum number of iterations. The second is a relative tolerance or change in coordinates between iterations. If either is reached, the algorithm terminates. When the accuracy parameter is increased, the maximum number of iterations is increased and the tolerance is decreased. It is possible the projection will reach maximum accuracy without increasing the accuracy parameter. It is also possible that the maximum accuracy is not reached after the maximum number of iterations. However, in this latter case, it is more likely that the change in coordinates from one iteration to the next falls below

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the tolerance before the maximum number of iterations is reached. This limitation can be blamed primarily on the truncation errors of digital floating point arithmetic.

There is also a limitation due to the geometry. It is worth mentioning that some surfaces are always at maximum accuracy, such as a sphere. A polygon surface is naturally limited to th accuracy due to the resolution of the polygons. Finally, it should be noted that if a portion of the mesh is projected to multiple surfaces and if these surfaces to not intersect, then it is impossible for the accuracy to be reached.

chkfolds check for folds in surfaces

chkfolds *option*

where the *option* can be

on	activate the fold checking projection algorithm
off	deactivate the fold checking projection algorithm (default)

Remarks

There is a problem in accuracy when projecting to a surface that has folds in it. To avoid additional costly calculations for all surface projections, this problem is corrected by issuing a flag only when a surface is folded and the projection is not done correctly. This flag is set using the **chkfolds** command. This is a very rare condition. It effects all projections within a part.

cosurf create a composite surface from displayed surfaces

cosurf *surface_# trans;*

Remarks

Use the Pick panel and the Display panel in the Environment window to select surfaces for display. Alternatively, use the surface display functions to choose surfaces for display. Be sure to choose a new surface number for this composite. It is not allowed to use a number for this new surface if the same number is in the surface Display list. The automatic parameter %nextsrf can be used to get the next available surface identification number.

delsd delete a surface

delsd *surface_#*

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where *surface_#* is the assigned surface identification number

Remarks

When a surface is deleted, it cannot be undone. Do not delete a surface that is being used by a part. Wait until the part is complete before deleting the surface. Actually, this is a useless command since one can avoid a surface by simply not showing it.

delsds delete a set of surfaces

delsds surface_list ;
where a *surface_#* is the assigned surface identification number

Remarks

When a surface is deleted, it cannot be undone. Do not delete a surface that is being used by a part. Wait until the part is complete before deleting the surface. Actually, this is a useless command since one can avoid a surface by simply not showing it.

fetol feature extraction from polygon surfaces

fetol *angle*
where *angle* is the maximum angle for a feature

Remarks

Interior edges are formed within a polygon surface when the angle between two facets or polygons fall below the specified maximum angle. The polygon surfaces affected by this command come from the **vpsd** and stlsd commands and the **sd** with the **mesh**, **face**, **faceset**, **poly**, **stl**, or **bstl** options. These interior edges are useful when attaching edges of the mesh to important features of the polygon surface.

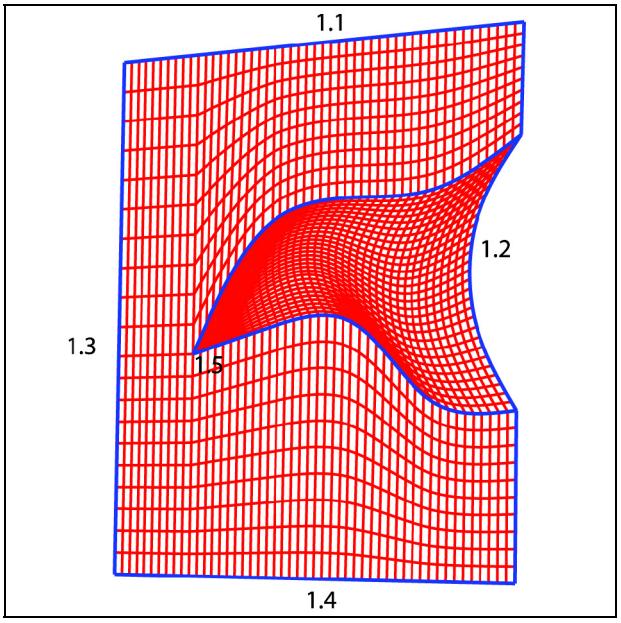

Figure 76 Interior feature edge 1.5

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The **fetol** command must precede the creation of the surface.

The default is no interior edges. To change **fetol** back to the default, use and angle greater than 180 degrees.

Example

```
fetol 163
vpsd 1 bed.cor bed.elm;
```

getol relative tolerance to tessellate curves and surfaces

getol *tolerance*
where *tolerance* is the relative tolerance

Remarks

The default relative tolerance for geometry is 1.0. This geometry tolerance is a factor used in the tessellation process. The actual tolerance is based on the total size of the geometry entity. Areas of greater curvature result in more points/polygons in the tessellation. When **getol** is increased, the number of points/polygons is decreased. Decrease **getol** to increase the number of points/polygons. This will also affect the amount of memory consumed for graphics and in some cases affect the accuracy of the projection when the **accuracy** parameter is set below 2.

lcsd list all composite surfaces with a specified surface

lcsd *surface_#*
where *surface_#* is a valid surface number

Remarks

This function helps identify a specific composite surface when there are many.

mvpn modify a surface node of a polygon surface

mvpn *point_id coor_id change(s)*
where
point_id identify the surface node which has the form *s.p.n*
where

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s	surface number
p	i-index or polygon number of the surface
n	j-index or node number of the polygon
coor_id	identify the coordinates to be changed
can be	
x	change the x-coordinate
y	change the y-coordinate
z	change the z-coordinate
xy	change the xy-coordinates
yz	change the yz-coordinates
xz	change the xz-coordinates
xyz	change the xyz-coordinates
change(s)	list of coordinate changes corresponding to *coor_id*.

Remarks

This command is used to move polygon surface nodes of types **faceset**, **poly**, **intp**, **stl**, and **bstl**. Polygon surfaces can be crude and error prone approximations to the actual surfaces, and as such, may produce polygon surfaces that are absurd to mesh. This feature can help you smooth the polygon surface to make the problem of meshing practical to solve. If you wish, you can save the modified polygon surface by using the **wrsd** command so that the next time you use this polygon surface, you do not have to repeat these polygon surface modifications.

Use the labels command to label the surface nodes on a polygon surface to aid in selecting the *point_id*.

npll set the quality of the projection to surfaces

npll *size*

where *size* is a relative curvature in a fold in the surface

Remarks

This command is almost never needed. In fact, there has been only one known case where this was needed. It is used in the special case that the surface being projected onto is extremely curved and almost folds onto itself. If the mesh does not project properly, this command will increase the quality of the projection search and perhaps fixes the problem. An alternative solution is to break the surface into smaller surfaces or use 3D curves with additional partitions in the mesh to constrain the mesh to the desired region of the surface. The default for this parameter is 30. It cannot be larger than 2000. If this number is increased, the length of time needed to project onto a surface will be increased. The cost for the projections should increase by something less than linear with respect to the number

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assigned with the command **npll**. If you are having difficulties, also try the **getol** command. This command is not found in the Graphical User Interface because it should only be used by experts who know to use it very sparingly.

project project a point to surface(s)

project *x y z list_of_surface_numbers* ;

where

x y z	coordinates of the initial point
list_of_surface_numbers	list of surface numbers

Remarks

Project uses the given 3D point as the initial point for the projection algorithm. This point is projected to the closest point in the intersection of the given surfaces.

This function returns the coordinates of the point of projection and the normal vector to each of the surfaces at that point of projection. These values can be used in expressions by referencing the following automatic parameters:

xprj - x-coordinate due to the **project** command
yprj - y-coordinate due to the **project** command
zprj - z-coordinate due to the **project** command
xnrm - x-component of the normal to the 1st surface from the **project** command
ynrm - y-component of the normal to the 1st surface from the **project** command
znrm - z-component of the normal to the 1st surface from the **project** command

This function is of little use in directly generating a mesh. Use the **sf** or the **sfi** command or click on the ***Project*** button to constrain (or project) a face of a mesh to a surface. The **project** command is intended to be used to demonstrate the projection method without the need to build a part. It can also be used to form parameters in creating geometry. For example, you can use the normal to define the tangent of a spline curve so that the curve is orthogonal to the surface. You can parameterize a normal offset from a surface, again by using the normal of the projection. You may find other creative uses for this function.

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For example, if one surface is given, then the result of the function is the coordinates of the point on that surface which is closest to the original point. If two surfaces are specified, then the result will be the point of projection along the intersecting curve of the two surfaces. If three surfaces are specified, then the result will be the point of intersection of the three surfaces.

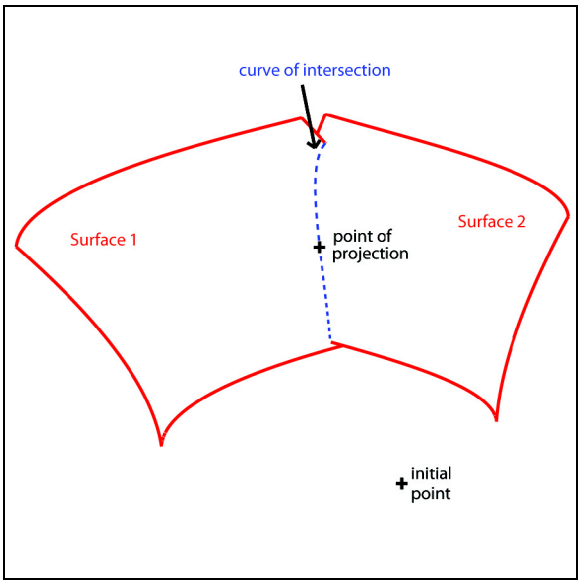

Figure 77 Project To Two Surfaces

Example

```
project 1.2 2.3 3.4 17 151;
```

This projects the point (1.2,2.3,3.4) to the intersection of surfaces 17 and 151. The results of this command are printed to the text window and the tsave file.

```
(x,y,z)=  -7.07107E-01  -3.33071E-08   7.07107E-01
 Table Of Distances To Surfaces
 distance to surface 17 is   0.00000E+00
 normal=-7.07107E-01 -3.33071E-08  7.07107E-01
 distance to surface 151 is   2.86341E-08
 normal= 5.77350E-01  5.77350E-01  5.77350E-01
```

pvpn place a surface node of a polygon surface

pvpn *point_id coor_id coordinate(s)*

where

point_id	identify the surface node which has the form *s.p.n*
where	
s	surface number
p	i-index or polygon number of the surface
n	j-index or node number of the polygon
coor_id	identify the coordinates to be changed
can be	
x	change the x-coordinate
y	change the y-coordinate
z	change the z-coordinate
xy	change the xy-coordinates

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yz	change the yz-coordinates
xz	change the xz-coordinates
xyz	change the xyz-coordinates
coordinate(s)	list of coordinates corresponding to *coor_id*.

Remarks

This command is used to move polygon surface nodes of types **faceset**, **poly**, **intp**, **stl**, and **bstl**. Polygon surfaces can be crude and error prone approximations to the actual surfaces, and as such, may produce polygon surfaces that are absurd to mesh. This feature can help you smooth the polygon surface to make the problem of meshing practical to solve. If you wish, you can save the modified polygon surface by using the **wrsd** command so that the next time you use this polygon surface, you do not have to repeat these polygon surface modifications.

Use the labels command to label the surface nodes on a polygon surface to aid in selecting the *point_id*.

Alternatively, you can use the graphical user interface to move a surface node of a polygon surface. There are two basic ways to do this. In both cases, choose the surface node of the polygon surface to be moved by selecting the ***Poly Surface*** button in the ***Move Pts.*** panel of the environment window. Then move the mouse close to the surface node that you wish to move and click on the **F5** function key.

If you choose a surface node using the ***Pick*** panel, you must then click on the ***Attach*** button to cause the surface node to move. This will produce a **pvpn** command that will be written to the text window and the tsave file.

If you choose the ***Move Pts.*** panel, you can then move the surface node with a click and drag of the left mouse button. When you release the mouse button, a **pvpn** command is generated.

sd surface definition

sd *[name] surface_# surface_type surface_parameters*

where

name	optional name for the surface
surface_#	positive integer to identify the surface

surface_type and *surface_parameter* can be any of:

Derivatives of other surfaces (see also **trsd**):

sds $sd_1\ sd_2\ ...\ sd_n\ ;$	combine several numbered surfaces into one
intp $sd_1\ sd_2\ fraction$	interpolate a surface between two surfaces

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Planes:

plan $x_0\ y_0\ z_0\ x_n\ y_n\ z_n$	infinite planar surface by a normal vector
xyplan *trans ;*	transform an infinite x-y plane
yzplan *trans ;*	transform an infinite y-z plane
zxplan *trans ;*	transform an infinite z-x plane
pl2 *system point system point*	plane specified by two points
pl3 *system point system point system point*	plane specified by three points
pl3o *system point system point system point offset*	plane specified by 3 points and an offset
iplan *a b c d*	plane defined by an implicit function

Spheres:

sp $x_0\ y_0\ z_0$ *radius*	sphere

Cylinders:

cy $x_0\ y_0\ z_0\ x_n\ y_n\ z_n$ *radius*	infinite cylindrical surface
xcy *radius trans ;*	transform an infinite x-axis cylinder
ycy *radius trans ;*	transform an infinite y-axis cylinder
zcy *radius trans ;*	transform an infinite z-axis cylinder
cy3 $x_1\ y_1\ z_1\ x_2\ y_2\ z_2\ x_3\ y_3\ z_3$	cylinder by 3 points
cyr2 $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$ *radius*	cylinder by 2 points and radius
cyr3 $x_1\ y_1\ z_1\ x_2\ y_2\ z_2\ x_3\ y_3\ z_3$ *radius*	cylinder by 3 points and radius

Conics:

cn2p $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r_1\ t_1\ r_2\ t_2$	infinite conical surface by two points
cone $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r\ \theta$	infinite conical surface by angle and radius
er $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r_1\ r_2$	ellipse revolved about its axis
pr $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r_1\ t_1\ r_2\ t_2\ r_3\ t_3$	infinite parabola revolved about an axis

Torus:

ts $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r_1\ t\ r_2$	torus

2D Curves, surfaces formed from:

crx *ln*	planar 2D curve rotated about the x-axis
cry *ln*	planar 2D curve rotated about the y-axis
crz *ln*	planar 2D curve rotated about the z-axis
cr $x_0\ y_0\ z_0\ x_n\ y_n\ z_n$ *ln*	planar 2D curve rotated about an arbitrary axis

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cp *ln trans ;* planar 2D curve extruded infinitely

rule2d y_1 ln_1 y_2 ln_2 *trans ;* ruled surface between two planar curves

swept ln_0 *direction* ln_1 α_1 ... ln_n α_n *; trans ;*
surface with planar cross-sections along a 2D curve

3D curves, surfaces formed from:

rule3d *3D-curve*$_1$ *3D_curve*$_2$ *trans ;*
ruled surface between two 3D curves

crule3d *3D-curve*$_1$ *3D_curve*$_2$ *trans ;*
cylindrical surface between two 3D curves

pipe *3d_curve radius*$_1$ *arc_length*$_1$ *radius*$_2$ *arc_length*$_2$ *... ; trans ;*
sweep a pipe shape along an arbitrary 3D curve

blend3 *3d_curve*$_1$ *3d_curve*$_2$ *3d_curve*$_3$ *trans ;*
blend three bounding 3D curves to form a patch

blend4 *3d_curve*$_1$ *3d_curve*$_2$ *3d_curve*$_3$ *3d_curve*$_4$ *trans ;*
blend four bounding 3D curves to form a patch

r3dc x_0 y_0 z_0 x_n y_n z_n *3D_curve start_angle end_angle trans ;*
3D curve rotated about an arbitrary axis

IGES (see also **iges**, **igesfile**, **igessd**, **igespd**, and **nurbsd**):

nurbs *nurbs* IGES NURBS (Non-Uniform Rational B-Spline)

igess *surface trans ;* IGES parametric surface

igesp *plane trans ;* IGES plane,

Algebraic surfaces:

function u_{min} u_{max} v_{min} v_{max} *x-expression ; y-expression ; z-expression ; trans ;*
surface by three algebraic expressions

csps *#_columns #_rows flag conditions* x_{11} y_{11} z_{11} x_{12} y_{12} z_{12} *... trans ;*
cubic spline surface

bsps *i-degree j-degree iknot*$_1$ *... iknot*$_{k1}$ *; jknot*$_1$ *.... jknot*$_{k2}$ *; xcontrol*$_1$ *ycontrol*$_1$ *zcontrol*$_1$ *... xcontrol*$_n$ *ycontrol*$_n$ *zcontrol*$_n$ *trans ;*
B-spline surface

nrbs *i-degree j-degree iknot*$_1$ *... iknot*$_{k1}$ *; jknot*$_1$ *.... jknot*$_{k2}$ *; weight*$_1$ *... weight*$_n$ *; xcontrol*$_1$ *ycontrol*$_1$ *zcontrol*$_1$ *... xcontrol*$_n$ *ycontrol*$_n$ *zcontrol*$_n$ *trans ;*
NURBS surface

hermite *#_columns #_rows flag conditions* x_{11} y_{11} z_{11} x_{12} y_{12} z_{12} *... trans ;*
2nd order, once differentiable, spline surface

Polygonal surfaces (see also **vpsd**, **wrsd**, **mvpn**, **pvpn**, and **fetol**):

mesh *m n* x_{11} y_{11} z_{11} x_{21} y_{21} z_{21} ... x_{mn} y_{mn} z_{mn} *trans ;* tabular data

face *region* face of the present part (Part phase only)

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faceset *set_name* face set
poly *polygon_set trans ;* polygon set
stl *file trans ;* read the standard ASCII STL file
bstl *file trans ;* read the standard binary STL file

trans is a product from left to right of the following operators

mx *x_offset*
my *y_offset*
mz *z_offset*
v *x_offset y_offset z_offset*
rx *theta*
ry *theta*
rz *theta*
raxis *angle x0 y0 z0 xn yn zn*
rxy
ryz
rzx
tf *origin x-axis y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* (Cartesian coordinates)
cy *rho theta z* (cylindrical coordinates)
sp *rho theta phi* (spherical coordinates)
pt *c.i* (label of a labeled point from a 3D curve)
pt *s.i.j* (label of a labeled point from a surface)

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* (Cartesian coordinates)
cy *rho theta z* (cylindrical coordinates)
sp *rho theta phi* (spherical coordinates)
pt *c.i* (label of a labeled point from a 3D curve)
pt *s.i.j* (label of a labeled point from a surface)

inv to invert the present transformation
csca scale all coordinates
xsca scale the x-coordinates
ysca scale the y-coordinates
zsca scale the z-coordinates

Remarks

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See the following section called Surface Dictionary for a full description of each surface type.

The surface name is optional. It still needs to be numbered. The automatic parameter %nextsrf can be used to get the next available surface identification number. All surfaces with the same name are automatically combine into a composite surface and is known by that name.

Names must have some alphabetic characters in them and no spaces. Other special characters like : and ; should also be avoided, although the implications of using none alphanumeric characters has not been explored. This feature is case insensitive. Only the first 8 characters of a name are used.

A surface can only have one name, so it can be conveniently included into only one composite surface by name. To include it into additional composite surfaces, use the surface's number in the **sd** command with the **sds** option.

All surfaces are defined in the Cartesian coordinate system regardless of the coordinate system used to define the mesh (**block** or **cylinder**).

Sd defines a surface to be used by other commands such as **sf**, **project**, **ms**, and **ssf**.

You can look at pictures of the surfaces you define, with such commands as **dsd**. In fact, surfaces defined by the **sd** command are the only surfaces which you can see displayed this way. Besides **dsd**, the relevant commands are **asd**, **rsd**, **dsds**, **asds, rsds, dasd**, and **rasd**. When a surface is defined, the surface is evaluated at many points so that the surface can be rendered graphically. Areas of the surface which have greater curvature will be evaluated at more points. This is done once when the surface is defined. Some surfaces, such as a plane, cylinder, rotated parabola, cone, or extruded 2D curve, are infinite and can not be completely rendered. They are cut off in order to fit into the picture with the rest of the surfaces. When such a surface is added to the picture, it may extend a little further than desired. A restore of the picture (***Rest***) will re-evaluate the infinite surfaces so that they are extended uniformly.

You can import surfaces from a CAD/CAM system. These surfaces are assigned surface numbers using the **iges**, **igessd**, **nurbsd**, and **igespd** commands or with the **igess**, **nurbs**, or **igesp** options of the **sd** command. Surfaces can also be selected by their level number inherited from the IGES interface for viewing using the **lv**, **alv**, **rlv**, and **dlvs** commands. Associativity groups may also be used through definitions found in the IGES file to view these surfaces using the **dgrp**, **agrp**, **rgrp**, and **dgrps** commands

Defined surfaces can also be used to construct 3D curves by extracting an edge or contour of a surface using the **sdedge** and **contour** options of the **curd** command, respectively.

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smgap small surface gap tolerance

smgap *tolerance*
where *tolerance* is n absolute (positive) distance

Remarks

The **smgap** command sets a tolerance to small gaps in surfaces. If a surface gap is found to be smaller than the value of **smgap**, the gap is removed. This command is almost very needed because the default works in about %99.9999 of the time. This may be useful only in the case when a mapped surface has an (almost) degenerate edge. If this command is used before the command that creates the surface is issued, then the degenerate edge will not appear in the surface.

trsd transform a surface definition

trsd *surface_id trans* ;
where
trans can be any of:

mx *x_offset*	to translate along the x-axis
my *y_offset*	to translate along the y-axis
mz *z_offset*	to translate along the z-axis
v *x_offset y_offset z_offset*	to translate in a given vector direction
rx *theta*	to rotate about the x-axis
ry *theta*	to rotate about the y-axis
rz *theta*	to rotate about the z-axis
raxis *angle x0 y0 z0 xn yn zn*	to rotate about a general axis
rxy	to reflect about the xy-plane
ryz	to reflect about the yz-plane
rzx	to reflect about the zx-plane
tf *origin x-axis y-axis*	to transform to a frame of reference
ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*	to transform from one frame of reference to another
inv	to invert the transformation.

Each argument in the **tf** and **ftf** transformations consists of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinated
cy *rho theta z*	cylindrical coordinates

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sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from surface

Remarks

The *surface_id* can be its number or name.

This command applies a transformation to the given surface. The new surface assumes the surface *surface_#id* and the pre-transformed surface is lost. This is significant in the case of a combined surface (sds) because every surface in the list is replaced by its transformed version.

The transform is order dependent. Note that in the example (**Figure 78**), applying the shift before the rotation is different from applying the rotation before the shift.

Example

```
c Set up first surface
sd   3 function
     0 90 0 360
     cos(u)*cos(v);
     cos(u)*sin(v);
     3 *sin(u);;
sd   4 function
    -90 0 0 360
     cos(u)*cos(v);
     cos(u)*sin(v);
     .5*sin(u);;
sd   1 sds   3   4;
c Set up second surface
c (This copy is needed
c  because each surface
c  in the surface list
c  is transformed)
sd 30 function
     0 90 0 360
     cos(u)*cos(v);
     cos(u)*sin(v);
     3 *sin(u);;
sd 40 function
    -90 0 0 360
     cos(u)*cos(v);
     cos(u)*sin(v);.
```

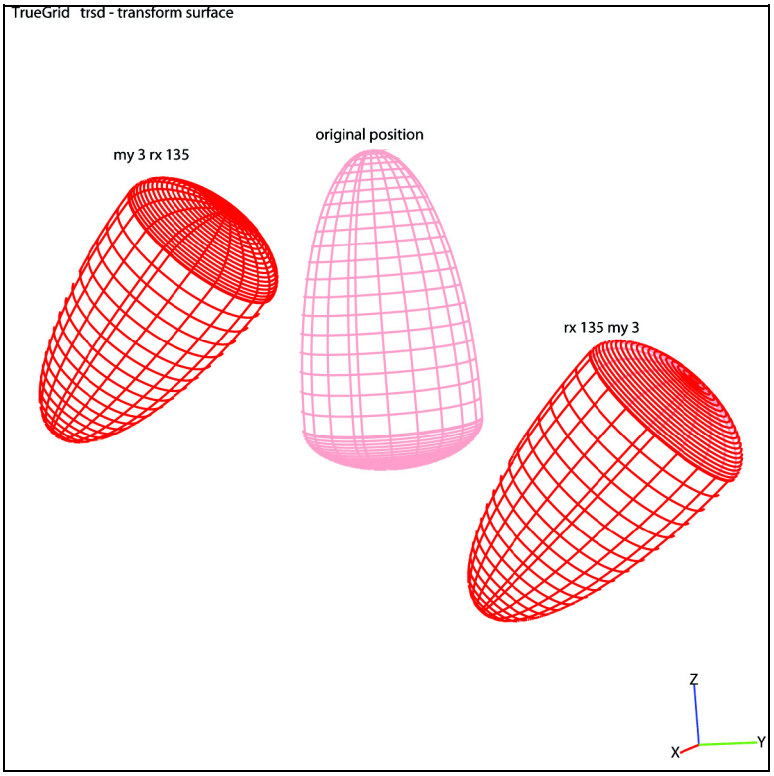

Figure 78: trsd - transforming a surface is order dependent

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```
    5*sin(u);;
sd 10 sds 30 40;
c Transform first surface by moving then rotating
trsd 1 my 3 rx 135;
c Transform second surface by rotating then moving
trsd 10 rx 135 my 3;
```

sdinfo list surfaces

sdinfo *(no arguments)*

Use this command to see a list of all the numbered surfaces. For each surface number, **True*Grid*®** gives a terse description of the surface.

vd define a volume

vd *number type parameters*

where

number is a unique volume number with which you can refer to the volume later, and

type and *parameters* can be any of the following:

sp *x y z radius* for a sphere with center *x, y, z* and radius *radius*.

csp *x y z radius* for the complement of a sphere with center *x, y, z* and radius *radius*.

el *x y z x_radius y_radius z_radius trans* ; for an ellipsoid aligned with the global coordinate axis and then transformed.

where (*x,y,z*) is the center of the ellipsoid
where *x_radius* is the radius in the x-direction
where *y_radius* is the radius in the y-direction
where *z_radius* is the radius in the z-direction

cel *x y z x_radius y_radius z_radius trans* ; for the compliment of an ellipsoid aligned with the global coordinate axis and then transformed.

where (*x,y,z*) is the center of the ellipsoid
where *x_radius* is the radius in the x-direction
where *y_radius* is the radius in the y-direction
where *z_radius* is the radius in the z-direction

cy *x y z xn yn zn radius* for an infinite cylinder of radius *radius* whose axis passes through *x, y ,z* in the direction of (*xn*,*yn*,*zn*).

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ccy *x y z xn yn zn radius* for the complement of an infinite cylinder of radius *radius* whose axis passes through *x, y, z* in the direction of (*xn*,*yn*,*zn*).

cr *x y z xn yn zn curve* for a rotated 2D curve. The curve number is *curve*, and the axis of rotation passes through *x, y, z* in direction (*xn*,*yn*,*zn*).

ccr *x y z xn yn zn curve* for the complement of a rotated 2D curve. The curve number is *curve*, and the axis of rotation passes through *x, y, z* in direction (*xn*,*yn*,*zn*).

cyf *x y z xn yn zn radius z_min z_max* for a finite cylinder of radius *radius* whose axis passes through *x, y ,z* in the direction of (*xn*,*yn*,*zn*). The values *z_min* and *z_max* are the distances along this axis to the start and end of the finite cylinder.

ccyf *x y z xn yn zn radius z_min z_max* for the complement of a finite cylinder of radius *radius* whose axis passes through *x, y ,z* in the direction of (*xn*,*yn*,*zn*). The values *z_min* and *z_max* are the distances along this axis to the start and end of the finite cylinder.

sd *surface distance* for a surface with thickness with thickness of the value of *distance*. *surface* is the surface number.

csd *surface distance* for the complement of a surface with thickness with thickness of the value of *distance*. *surface* is the surface number.

nsd *surface distance* for a surface with thickness of the value of *distance*. *surface* is the surface number. The thickness is only in the normal direction.

cnsd *surface distance* for the compliment of a surface with thickness of the value of *distance*. *surface* is the surface number. The thickness is only in the normal direction.

box *xm ym zm xx yx zx option* ; for a box with corners at (*xm,ym,zm*) and (*xx,yx,zx*)

where *option* is a LS-DYNA BOX feature and can be

adaptive *material level* for LS-DYNA BOX_ADAPTIVE

coarsen *inout_flag* for LS-DYNA BOX_COARSEN

cbox *xm ym zm xx yx zx option* ; for the complement of a box with corners at (*xm,ym,zm*) and (*xx,yx,zx*),

sphbox *xmin ymin zmin xmax ymax zmax option* ; for an SPH box in LS-DYNA

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with corners at (*xm,ym,zm*) and (*xx,yx,zx*)

where *option* can be

motion *xn yn yz load_curve_# flag*

where *flag* can be

0 for velocity

1 for displacement

Remarks

There are several uses of a volume. The **mtv** command uses a volume to modify the materials of elements within the volume. The Ls-dyna output uses the **box** and **sphbox** volume to write *DEFINE_BOX cards.

Examples

```
block 1 31;1 31;1 31;-1 1 -1 1 -1 1
vd 1 sp 0 0 0 1
mtv 1 1 1 2 2 2 1 4 2 ;
merge
dm 1
dm 2
```

In this example, elements with at least 4 nodes within a sphere are assigned material 2.

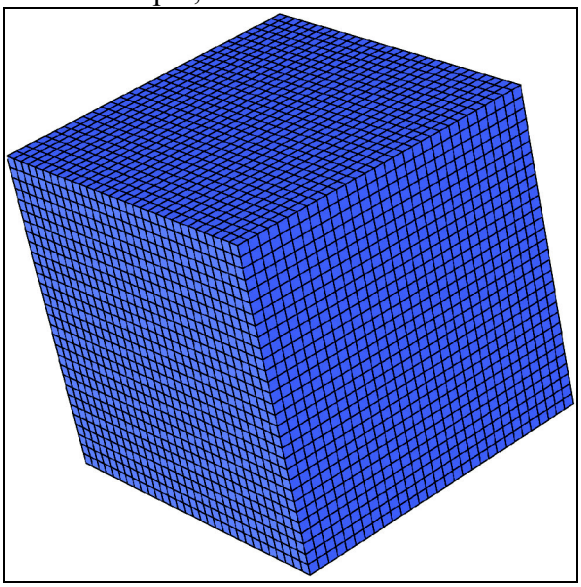

Figure 79 Material 1

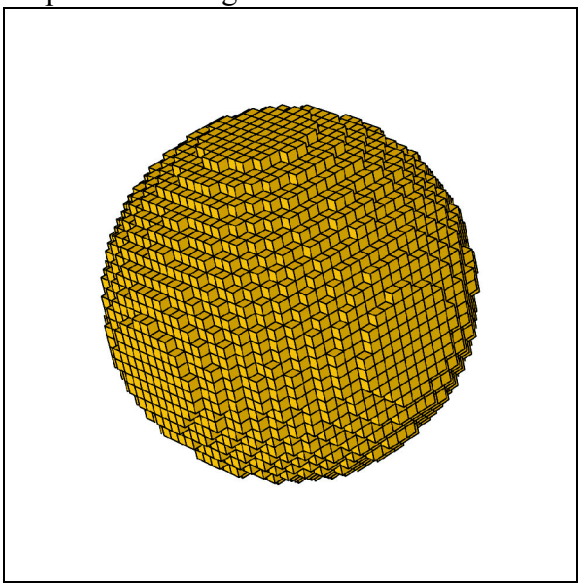

Figure 80 Material 2

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9. Surface Dictionary

The following is a list of surface types and parameters available, as options, under the **sd** command.

Many of the following surfaces refer to an axis to form a local coordinate system. The parameters (x_0 y_0 z_0) give a point through which the axis passes. The parameters (x_n y_n z_n) give the direction of the axis; that is, it is in the same direction as the line from the origin to the point (x_n y_n z_n). The distance to this point must be positive. Often they also refer to *r* and *t*; *r* is a distance from the axis and *t* is the distance along the axis from the point (x_0 y_0 z_0). Thus (*r*,*t*) form a local coordinate system in which we may define points, curves, etc..

Some of the following surfaces are based on a planar 2D curve defined with the **ld** command. They are assigned a number which is referred to by *ln*. The 2D curve must be defined first before it is referenced by the **sd** command.

Some of the following surfaces are based on a 3D curve defined with the **curd** command. They are assigned a number which is referred to by *3D_curve*. The 3D curve must be defined first before it is referenced by the **sd** command.

In some surface definitions, you make the surface in a convenient coordinate system and then immediately rotate or translate it to where you need it. This is indicated by the *trans ;* option in the surface description below. If you don't want any transformations, then you should simply omit any transformations and end the command with a semicolon.

blend3 blend three bounding 3D curves to form a patch

blend3 *3d_curve*$_1$ *3d_curve*$_2$ *3d_curve*$_3$ *trans ;*

where

3d_curve$_i$	3D curve number
trans	sequence of simple operators to transform the coordinates

Remarks

The three 3D curves used to form the boundaries of a surface patch are assumed to intersect. Care must be taken to insure that the 3D curves almost intersect. The algorithm can deal with gaps between the 3D curves, but the result can be unsatisfactory if the gaps are very large. The curves do not need to end at the desired corner points of the surface patch.

It is also assumed that the three curves surround only one 3-sided region. If, due to curvature, there is an ambiguity as to which enclosed region will be surfaced, the results may be unsatisfactory. In such

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a case, we recommend that you break up the 3D curves into smaller 3D curves so that the region to become a surfaced is obvious. The three curves do not need to be specified in any particular order.

The surface that is fit to the three bounding curves is a result of a transfinite interpolation, similar to the transfinite interpolation associated with the **tfi** command in the part phase.

This surface is similar to the **blend4** option.

Example

```
curd 1 csp3 00  -5.4e-01 3.6e-01
7.5e-01 6.1e-02  4.9e-01 8.6e-01
5.5e-01  3.2e-01 7.6e-01 ;;;
curd  2   csp3   00      -4.3e-01
4.8e-01     7.5e-01     -3.6e-01
-1.6e-02     9.2e-01     -4.1e-01
-5.9e-01     6.8e-01     -1.1e-01
-8.1e-01 5.7e-01 ;;;
curd 3 csp3 00 -2.7e-01 -7.9e-01
5.4e-01 2.3e-01 -6.2e-01 7.4e-01
4.0e-01  1.8e-01 8.9e-01 2.2e-02
7.8e-01 6.1e-01 ;;;
sd 1 blend3 2 1 3 ;
```

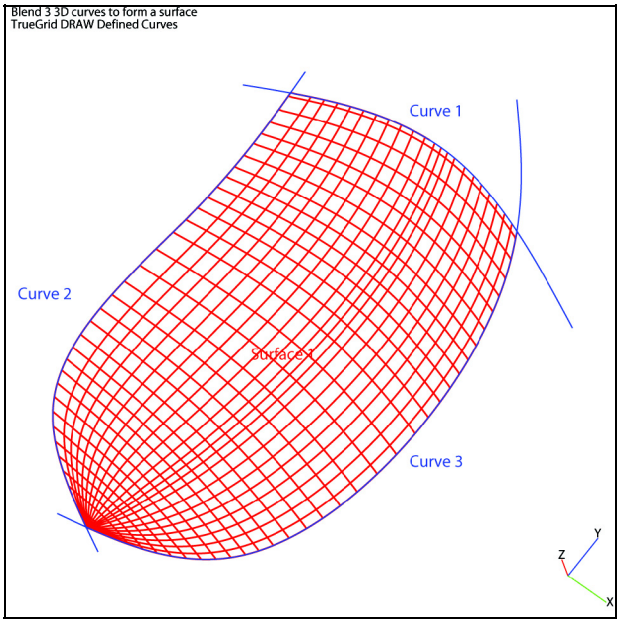

Figure 81 blended surface

blend4 blend four bounding 3D curves to form a patch

blend4 *3d_curve*$_1$ *3d_curve*$_2$ *3d_curve*$_3$ *3d_curve*$_4$ *trans* ;

where

3d_curve$_i$	3D curve number
trans	sequence of simple operators to transform the coordinates

Remarks

The four 3D curves used to form the boundaries of a surface patch are assumed to intersect. Care must be taken to insure that the 3D curves almost intersect. The algorithm can deal with gaps between the 3D curves, but the result can be unsatisfactory if the gaps are very large. The curves do not need to

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end at the desired corner points of the surface patch.

It is also assumed that the four curves surround only one 4-sided region. If, due to curvature, there is an ambiguity as to which enclosed region will be surfaced, the results may be unsatisfactory. In such a case, we recommend that you break up the 3D curves into smaller 3D curves so that the region to be surfaced is obvious. The four curves do not need to be specified in any particular order.

The surface that is fit to the four bounding curves is a result of a transfinite interpolation, similar to the transfinite interpolation associated with the **tfi** command in the part phase.

Selecting 3D curves to form a four sided surface patch is an art. Ideally, opposite curves should be parallel. Avoid sharp bends in the curves. A kink in a boundary curve usually indicates the need for several surface patches using pieces of the original 3D curves. Severe concave boundary curves can also be a problem. If this occurs, create several more 3D curves to bisect the region and build several surface patches instead of one.

This surface is similar to the **blend3** option.

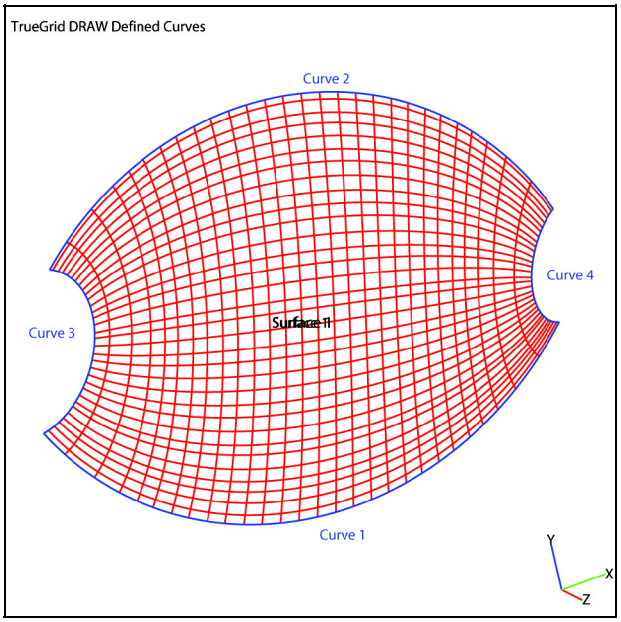

Figure 82 blended surface

Example

```
curd 1 arc3 seqnc   rt -6.7e+00 -5.2e-01  6.1e-01
                    rt  3.6e-01 -5.0e+00  3.1e-01
                    rt  6.4e+00 -2.5e+00  2.6e-01 ;
curd 2 arc3 seqnc   rt -5.7e+00  3.1e+00  8.2e-01
                    rt  4.0e-01  5.1e+00  5.9e-01
                    rt  6.6e+00  4.0e-01  7.5e-01 ;
curd 3 arc3 seqnc   rt -5.7e+00  3.1e+00  8.2e-01
                    rt -4.8e+00  1.3e+00  1.4e-01
                    rt -6.7e+00 -5.2e-01  6.1e-01 ;
curd 4 arc3 seqnc   rt  6.6e+00  4.0e-01  7.5e-01
                    rt  5.6e+00 -9.0e-01  9.7e-01
                    rt  6.4e+00 -2.5e+00  2.6e-01 ;
sd 1 blend4 1 2 3 4 ;
```

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bsps B-spline

bsps *i-degree j-degree*
iknot$_1$ ... *iknot*$_{k1}$; *jknot*$_1$ .... *jknot*$_{k2}$;
xcontrol$_{1,1}$ *ycontrol*$_{1,1}$ *zcontrol*$_{1,1}$ ... *xcontrol*$_{1,n2}$ *ycontrol*$_{1,n2}$ *zcontrol*$_{1,n2}$
. . .
xcontrol$_{n1,1}$ *ycontrol*$_{n1,1}$ *zcontrol*$_{n1,1}$ ... *xcontrol*$_{n1,n2}$ *ycontrol*$_{n1,n2}$ *zcontrol*$_{n1,n2}$;
trans ;
where

i-degree j-degree	degrees of surface polynomial in the i and j direction.
i*knot*$_1$ ... *iknot*$_{k1}$	parametric coordinates of *k1* knots (0..) in the i direction.
j*knot*$_1$ ... *jknot*$_{k2}$	parametric coordinates of *k2* knots (0..) in the j direction.
xcontrol$_{1,1}$ *ycontrol*$_{1,1}$ *zcontrol*$_{1,1}$ ... *xcontrol*$_{1,n2}$ *ycontrol*$_{1,n2}$ *zcontrol*$_{1,n2}$ *xcontrol*$_{n1,1}$ *ycontrol*$_{n1,1}$ *zcontrol*$_{n1,1}$ ... *xcontrol*$_{n1,n2}$ *ycontrol*$_{n1,n2}$ *zcontrol*$_{n1,n2}$;	triplets of coordinates of control points

where
$n_1 = k_1$ - *multiplicity*$_1$
$n_2 = k_2$ - *multiplicity*$_2$

trans	sequence of simple operators to transform the coordinates

Remarks

This type of surface is also known as a Basis Spline because it is defined as a linear combination of polynomial basis functions. The theory of B-Splines can be found in many text books. The control points are the coefficients forming the linear combinations of the basis functions and can form a crude caricature of the surface.

The multiple knots have to be specified for the starting and ending point of the curve. The number of multiple knots is called multiplicity. The multiplicity for the starting and ending point determines the order of the curve. (2 - linear, 3 - quadratic, 4 - cubic ...).

multiplicity = degree + 1

The multiplicity has to be the same for the starting and ending points.

B-spline surfaces are also available through the IGES interface.

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Example

The B-Spline surface, shown in **Figure 83**, is defined by the same control points as the NURBS Surface (**Figure 114**).

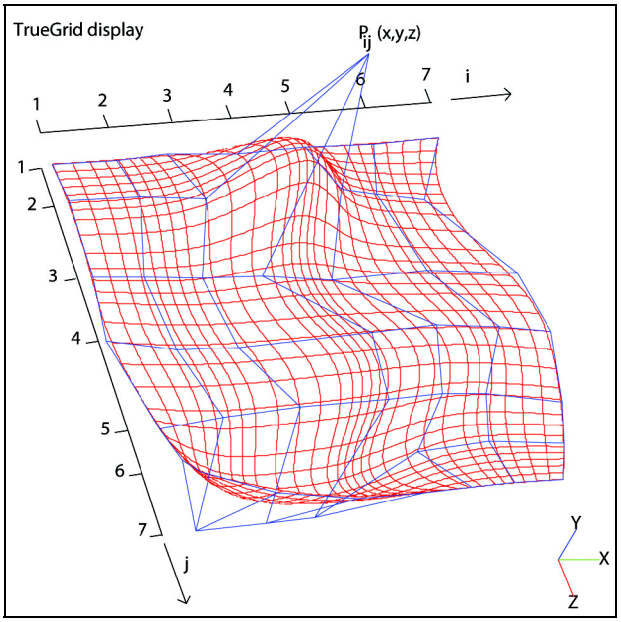

Figure 83 B-Spline Surface

```
sd 1 bsps 3 3
0 0 0 0 1 2 3 4 4 4 4;
0 0 0 0 1 2 3 4 4 4 4;
-2.17  4.90e-03 -1.80
-1.45  1.59e-02 -1.85
-8.54e-01       1.96e-02   -1.95
-4.35e-02  2.05e-02 -1.98
 8.01e-01  4.66e-03 -2.12
 1.65e+00 -5.72e-03 -2.19
 2.49e+00 -1.37e-03 -2.31
-2.14e+00  1.05e-02 -1.32
-1.35e+00  2.16e-02 -1.37
-6.19e-01       1.52e-02   -1.38
4.05e-01  2.00e+00 -1.50e+00
 1.01e+00  1.33e-04 -1.55e+00
1.41e+00  3.86e-03 -1.49e+00
 1.94e+00  2.21e-02 -1.40e+00
-2.17e+00 -2.57e-02 -3.81e-01 -1.58e+00 -6.88e-02 -4.22e-01
-9.27e-01 -1.13e-01 -4.49e-01 -2.38e-01 -1.24e-01 -4.82e-01
 5.19e-01 -9.57e-02 -4.24e-01  1.59e+00  1.64e-02 -4.53e-01
 2.60e+00  9.17e-02 -3.58e-01
-2.27e+00 -4.51e-02  3.97e-01 -1.48e+00 -1.52e-01 4.15e-01
-7.71e-01 -1.99e-01  3.57e-01  7.17e-01 -8.84e-02 2.68e-01
 1.43e+00  4.52e-03  2.52e-01  2.11e+00  6.75e-02 3.25e-01
 2.68e+00  9.55e-02  4.07e-01
-2.09e+00  3.17e-03  1.48e+00 -1.28e+00 -1.31e-01 1.26e+00
-3.79e-01 -1.57e-01  1.13e+00  5.72e-01 -6.70e-02 1.12e+00
 1.15e+00 -3.48e-03  1.11e+00  1.63e+00  3.45e-02 1.15e+00
 2.45e+00  5.85e-02  1.29e+00
-2.08e+00  1.66e-01  1.98e+00 -1.45e+00 -1.00e-00 1.93e+00
-6.71e-01 -1.00e-00  1.87e+00 -1.10e-01 -1.00e-00 1.83e+00
 5.99e-01 -2.52e-02  1.80e+00  1.42e+00  2.08e-02 1.73e+00
 2.28e+00  2.87e-02  1.78e+00
-2.03e+00  3.00e-01  2.21e+00 -1.14e+00  4.58e-02  2.29e+00
-2.59e-01 -2.78e-02  2.33e+00  4.70e-01 -1.12e-02 2.29e+00
 1.02e+00  4.84e-03  2.26e+00  1.52e+00  7.42e-03 2.221e+00
 2.10e+00  4.38e-04  2.19e+00 ;;
```

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bstl read the standard binary STL file

bstl *file trans* ;
where
file path and file name (case sensitive in UNIX/LINUX)
trans sequence of simple operators to transform the coordinates

Remarks

This command permits you to import triangular surfaces via StereoLithography (STL) files.

Use the **fetol** command before this command to extract interior features as edges.

The maximum number of polygons is 8,388,607. When a polygon has more than 4 nodes, it is broken into triangles and each triangle counts as a polygon.

Example

```
fetol 110
sd 1 bstl  prt_k1.bstl ;
```

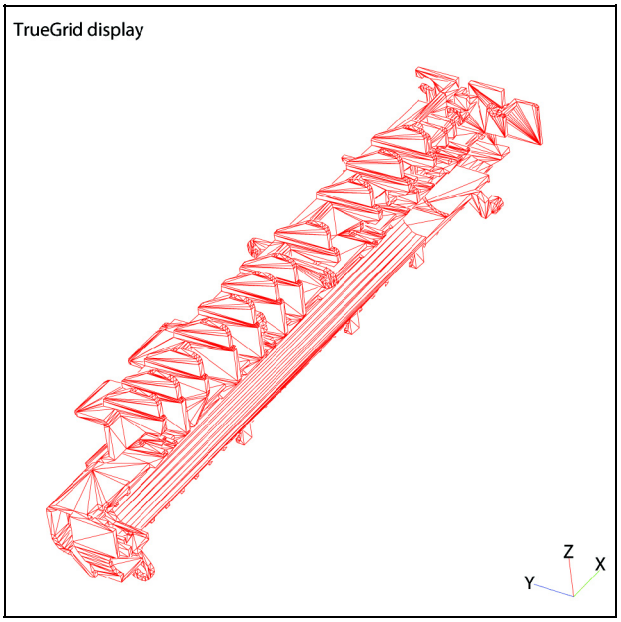

Figure 84 Surface imported from binary stl

cn2p infinite cone, defined by two points

cn2p $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r_1\ t_1\ r_2\ t_2$
where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of axis direction vector
y_n	second component of axis direction vector
z_n	third component of axis direction vector
r_1	radius at the cross section along axis at t_1
t_1	position along axis with radius r_1
r_2	radius at the cross section along axis at t_2
t_2	position along axis with radius r_2

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Remarks

(x_0,y_0,z_0) and (x_n,y_n,z_n) define the axis of symmetry. You complete the cone's definition by specifying two points (r_1,t_1) and (r_2,t_2) in any planar cross section containing the axis of symmetry. The t coordinate is measured along the axis from the point (x_0,y_0,z_0) in the direction of (x_n,y_n,z_n). The r coordinate is measured orthogonally from the axis of symmetry. For any plane containing the axis of symmetry, the line segment formed by the two points in the plane will be contained on the cone. The coordinate pairs (r_1,t_1) and (r_2,t_2) must satisfy the following conditions:

$$|r_1-r_2|\times 10^{-18} < |t_1-t_2|$$

and

$$|t_1-t_2|\times 10^{-18} < |r_1-r_2|$$

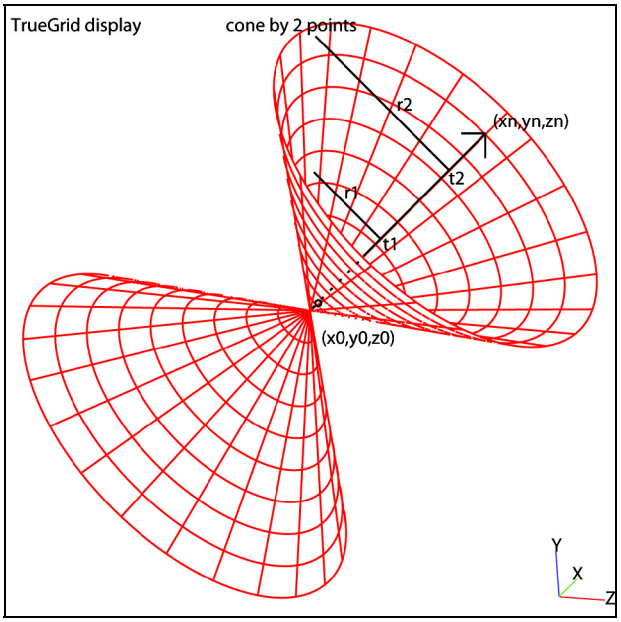

Figure 85 Cone by an axis and two points

A point projected to this surface should not start out on the axis of symmetry. See the example in **Figure 85**. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

Example

```
sd 2 cn2p 0 0 0 -1 1 1 1 1 2 2
```

cone infinite cone, defined by radius and angle

cone $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r\ \theta$

where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of axis direction vector
y_n	second component of axis direction vector
z_n	third component of axis direction vector
r	radius of the cross section along axis at the point $(x_0\ y_0\ z_0)$
θ	angle of the cone

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Remarks

A cone defined by radius and angle. The two points (x_0,y_0,z_0) and (x_n,y_n,z_n) define the axis of symmetry. The plane orthogonal to the axis at the point (x_0,y_0,z_0), slices through the cone to form a circle of radius r. r must be non-negative. Note that if $r=0$, then the point (x_0,y_0,z_0) is at the apex of the cone. The cone forms an angle θ relative to its axis of symmetry. This angle must be between -90 and 90, excluding -90, 0, and 90. A point projected to this surface can not be on the axis of symmetry. This is an infinite surface. The graphics will only show a portion of the surface. The portion shown in the graphics changes as the objects in the picture change.

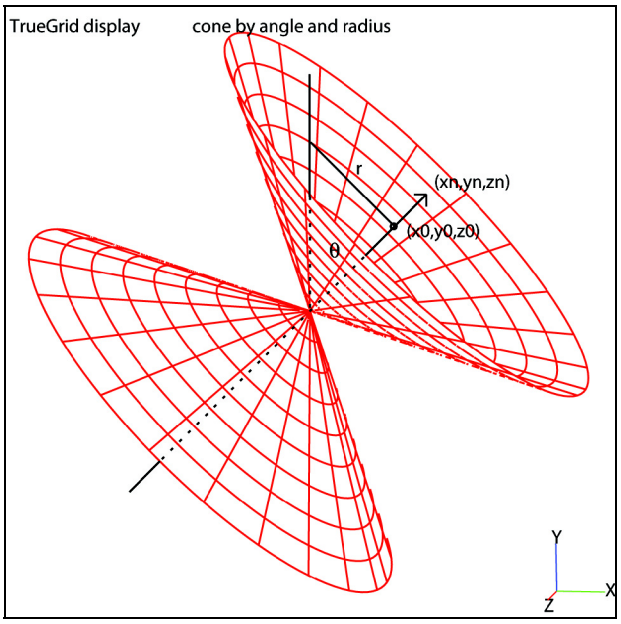

Figure 86 Cone specified by radius and angle

Example

```
sd 7 cone 0 1 2 1 1 0 1 45
```

cp infinite generalized cylinder (extruded or lofted curve)

cp *ln trans* ;

where

ln	2D curve number
trans	transform operators

Remarks

This option enables you to define a surface by sweeping a planar 2D curve through the third dimension, then transforming the result. The curve *ln* lies in the xz-plane and then is extended infinitely in the y direction to form a generalized cylinder.

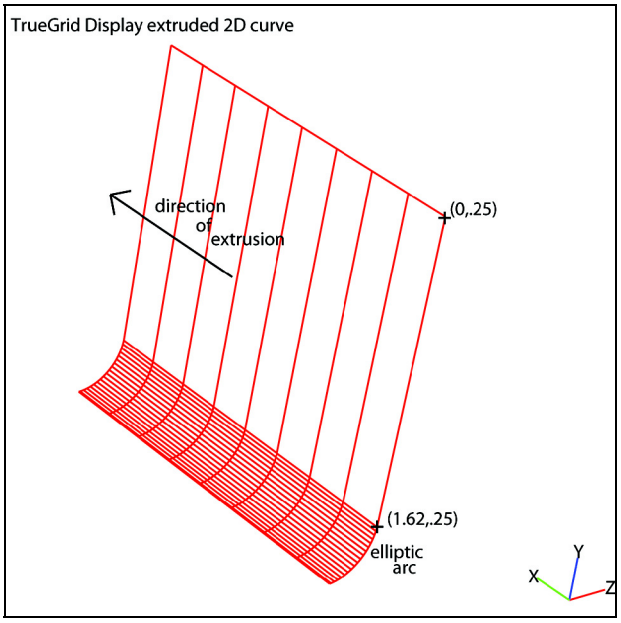

Figure 87 2D curve extruded in the x-direction

Example

```
ld 4 lp2 0 .25 1.62 .25;lep .25
.25 1.62 0 90 0 0 ;
sd 3 cp 4 rz -90;
```

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cr 2D curve revolved about an axis

cr x_0 y_0 z_0 x_n y_n z_n ln

where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of axis direction vector
y_n	second component of axis direction vector
z_n	third component of axis direction vector
ln	2D curve number

Remarks

This command forms a surface by rotating the 2D curve ln about an axis given by (x_0,y_0,z_0) and (x_n,y_n,z_n). As an example, the following commands were used to generate the picture below. The 2D curve must be defined before it is referenced by this command. The numbers in the picture correspond to the points in the 2D curve definition. A point projected to this surface should be away from the axis of symmetry.

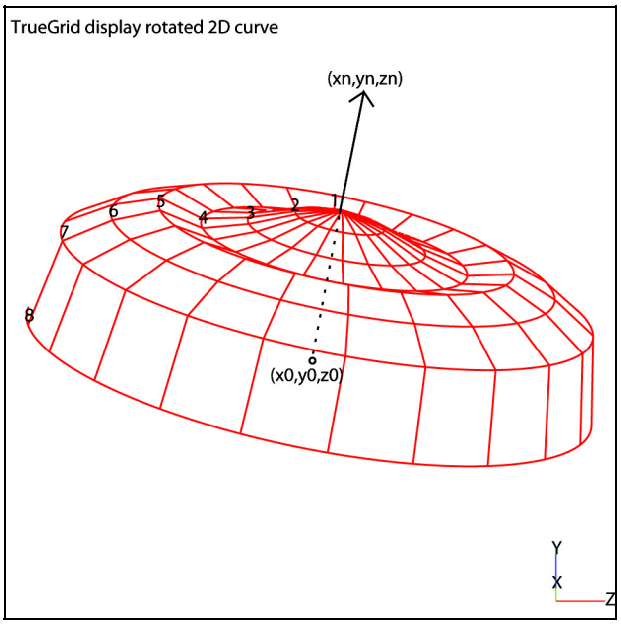

Figure 88 2D curve rotated about an axis

Example

```
ld 1 lp2 0 .2266 .0625 .2109 .125
.1875 .1875 .1719 .25 .1875 .3125
.164 .375 .125 .4 0;
sd 1 cr 0 4 1 0 1 .2 1
```

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crule3d cylindrical surface between two 3D curves

crule3d *3D-curve$_1$ 3D-curve$_2$ trans ;*

where

3D-curve$_1$	first boundary 3D curve
3D-curve$_2$	second boundary 3D curve
trans	sequence of simple operators to transform the coordinates

Remarks

This command forms a surface using linear interpolation in cylindrical coordinates between two 3D curves. At least one of the two curves must have a positive arc length. Both curves must be already defined. 3D curves are defined with the **curd** command or read from an **IGES** file. The direction of the curves is critical to the interpolation of the surface. Check the order of the points in each of the curves by displaying the points with the **label** command. In order to reverse the order of a curve, specify the curve definition number with a negative number. The surface is constructed by pairing each point on one curve with a point on the other curve. These two points are connected by a line segment in cylindrical coordinates. Points along the two curves are paired by relative arc length.

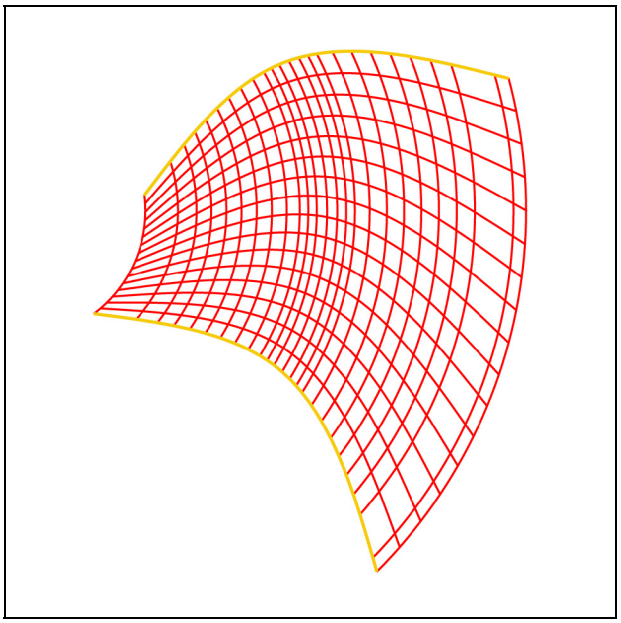

Figure 89 Cylindrical ruled surface

Example

```
curd 1 csp3 00 .3 -.4 0 1 -.6 0 1.4 -1.4 0;;
curd 2 cpcd 1 rz 60;
sd 1 crule3d 1 2 ; ;
```

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crx rotate 2D curve about x-axis

crx *ln*

where *ln* is the 2D curve number

Remarks

The 2D curve *ln* must be defined before it is referenced by this command using the **ld** command. A point projected to this surface should be away from the axis of symmetry. The following commands were used to generate the surface in the picture below. The first two points of the 2D curve are labeled in the picture. A circular arc is appended to the line segment between the two points.

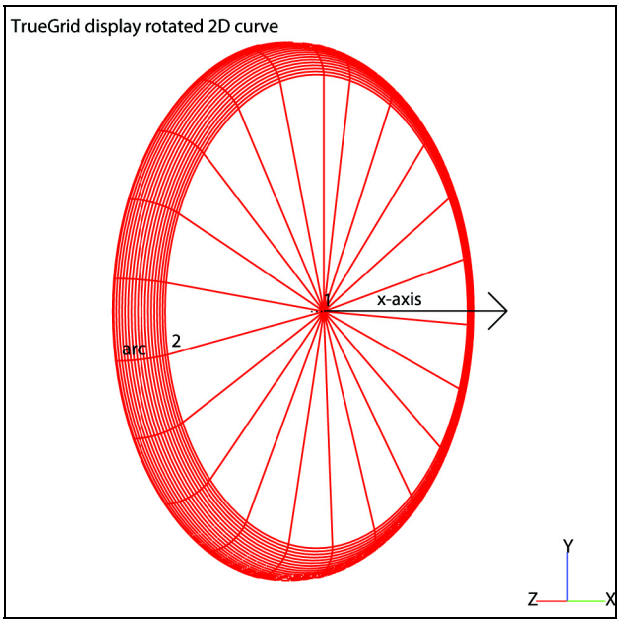

Figure 90 2D curve rotated about the x-axis

Example

```
ld 4 lp2 0 .25 1.62 .25; lep .25
.25 1.62 0 90 0 0 ;
sd 3 crx 4
```

cry rotate 2D curve about y-axis

cry *ln*

where *ln* is the 2D curve number

Remarks

This command defines a surface by rotating a 2D curve about the y-axis. The 2D curve *ln* must be defined before it is referenced by this command using the **ld** command. A point projected to this surface should be away from the axis of symmetry. The following commands were used to create the surface in the picture below. The four points in the **lp2** command are annotated in the picture (**Figure 91**).

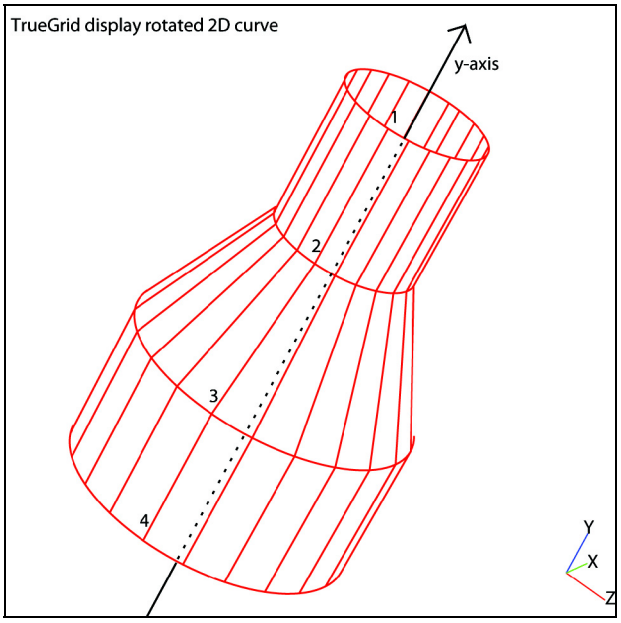

Figure 91 2D curve rotated about the y-axis

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Example

```
ld 1 lp2 1 1 1 2 .5 3 .5 4;
sd 1 cry 1
```

crz rotate 2D curve about z-axis

crz *ln*

where *ln* is the 2D curve number

Remarks

This command defines a surface by rotating a 2D curve about the z-axis. The 2D curve *ln* must be defined before it is referenced by this command using the **ld** command. A point projected to this surface should be away from the axis of symmetry. The following commands created the surface in the picture below. The two points in the **lp2** command are annotated.

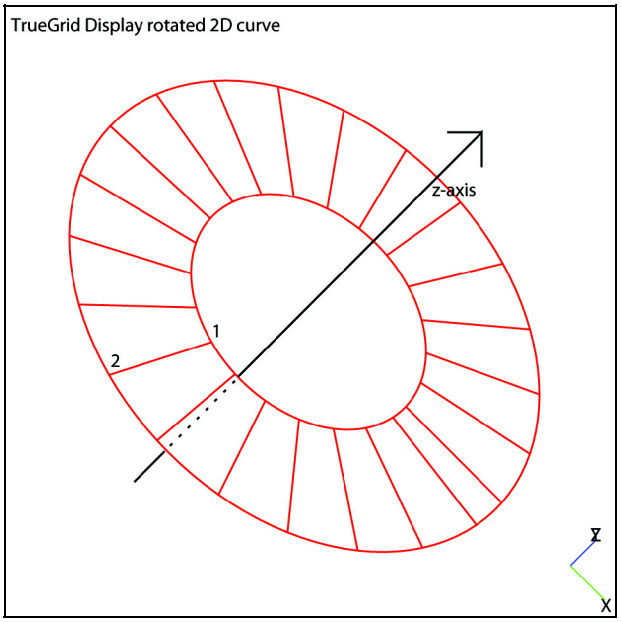

Figure 92 2D curve rotated about the z-axis

Example

```
ld 1 lp2 1 1 2 1;
sd 1 crz 1
```

csps fit a cubic spline surface through 3D data

csps *#_columns #_rows flag conditions* $x_{11}\ y_{11}\ z_{11}\ x_{12}\ y_{12}\ z_{12}$ ... *trans ;*

where

#_columns	number of control points in each column of the table of control points
#_rows	number of control points in each row of the table of control points
flag	4 digits formed using **0** for natural boundary condition, **1** for

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	specified boundary derivatives, one digit for each of the four edges
conditions	specifies the derivatives along the edges, in order, when a flag of 1 is used in the corresponding digit in the *flag*
where each *condition* is formed by derivatives	
dx_{ij}, dy_{ij}, dz_{ij}	derivative triplet, one for each control point along the edge
x_{ij} y_{ij} z_{ij}	list of coordinates for the control points on the surface
trans	sequence of simple operators to transform the coordinates

Remarks

There are four binary digits forming the flag. Each digit corresponds to an edge of the surface. The first edge is formed by the first column. The second edge is formed by the last column. Edge 3 is the first row. Edge 4 is the last row.

The theory of cubic splines allows for a derivative to be specified at each end of a spline curve. In a cubic spline surface, derivatives can be specified at both ends of each row and column. Each derivative is specified as a vector with three components, *dx*, *dy*, and *dz*. The magnitude of the derivatives can have a significant affect on the shape of the surface. Experimentation may be required. If derivatives are not specified, then the natural derivative is used. This means that the second derivative is set to zero which uses the remaining degrees of freedom with no first derivatives needed.

In a laborious way, one can form a cubic spline surface from a set of 3D cubic spline curves. Using a text editor on the tsave file that recorded the definition of the 3D curves, one can cut and paste to form the table of control points for the cubic spline surface. This requires that all of the 3D curves have the same number of control points and oriented in the same general direction. The curve control points have to placed into the table in the proper order. Derivatives at the end control points (if any) become the derivatives along an edge. The ordering is critical. Care is also needed in selecting corresponding interior control points in each 3D curve that represent corresponding features in the shape of the surface.

The transformation *trans* can include the normal offset. The normal offset must be the first operator in the list of operators that form the transformation. This normal offset is intended for small offsets only, otherwise the resulting surface will be self-intersecting.

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Examples

In this first example, one of the edges is degenerate and all of the edge derivatives are specified. In this case, the end derivatives of the degenerate edge are zero.

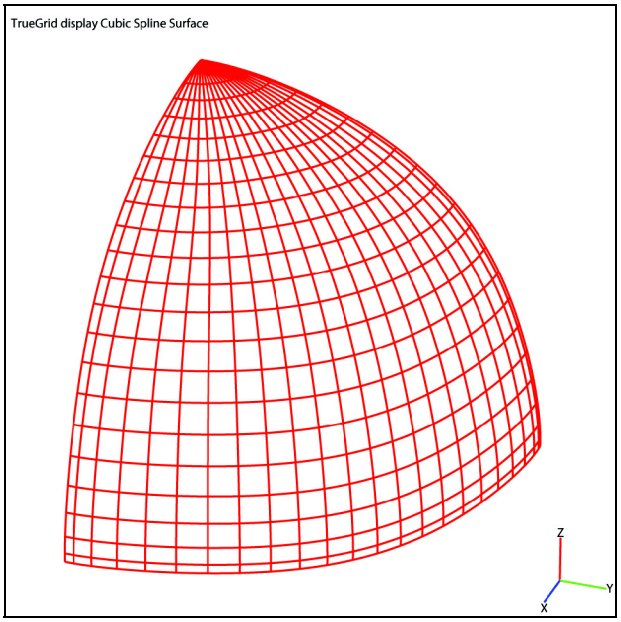

Figure 93 Spline surface with degenerate edge

```
c  The magnitude of the derivatives are parametric
para am1 1 am2 1 am3 1 am4 1.3;
c  All derivatives are specified in this example
sd 1 csps 4 4 1111
c  x,y,z components of the derivatives on edge 1
0 %am1 0 0 %am1 0 0 %am1 0 0 0 0
c  x,y,z components of the derivatives on edge 2
[-%am2] 0 0 [-%am2] 0 0 [-%am2] 0 0 0 0 0
c  x,y,z components of the derivatives on edge 3
0 0 %am3 0 0 %am3 0 0 %am3 0 0 %am3
c  x,y,z components of the derivatives on edge 4
[-%am4] 0 0 [-%am4*cos(30)] [-%am4*sin(30)] 0
[-%am4*cos(60)] [-%am4*sin(60)] 0 0 [-%am4] 0
c  x,y,z control points for row 1
1 0 0 [cos(30)] [sin(30)] 0 [cos(60)] [sin(60)] 0 0 1 0
c  x,y,z control points for row 2
[sqrt(3)/2] 0 .5 [cos(30)*sqrt(3)/2] [sin(30)*sqrt(3)/2] .5
[cos(60)*sqrt(3)/2] [sin(60)*sqrt(3)/2] .5 0 [sqrt(3)/2] .5
c  x,y,z control points for row 3
.5 0 [sqrt(3)/2] [cos(30)/2] [sin(30)/2] [sqrt(3)/2]
[cos(60)/2] [sin(60)/2] [sqrt(3)/2] 0 .5 [sqrt(3)/2]
```

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```
c  x,y,z control points for row 4
0 0 1 0 0 1 0 0 1 0 0 1 ;;
```

```
sd 1 csps 5 5 0000
c number of control points
c in the i and j-direction
c(rows and columns)
c Natural Boundary Conditions
1 1 1 2 1 1 2.5 1 1 3 1 1 4 1 1
1 2 1 1.79 1.79 2 2.5 1.5 2
3.2 1.79 2 4 2 1 1 2.5 1
1.5 2.5 2 2.5 2.5 1 3.5 2.5 2
4 2.5 1 1 3 1 1.79 3.2 2
2.5 3.5 2 3.2 3.2 2 4 3 1 1 4 1
2 4 1 2.5 4 1 3 4 1 4 4 1;;
```

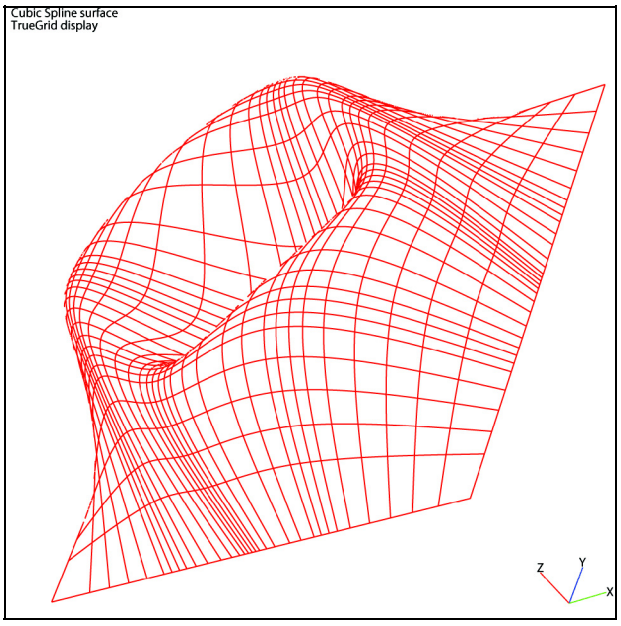

Figure 94 cubic spline surface

The last example is used with a normal of 1 applied to it to form the next surface. Care is needed in using the normal offset. The normal distance must not exceed the inverse of the curvature or the resulting surface will be self-intersecting.

```
sd 1 csps 5 5 0000
1 1 1 2 1 1 2.5 1 1 3 1 1 4 1 1
1 2 1 1.79 1.79 2 2.5 1.5 2
3.2 1.79 2 4 2 1 1 2.5 1
1.5 2.5 2 2.5 2.5 1 3.5 2.5 2
4 2.5 1 1 3 1 1.79 3.2 2
2.5 3.5 2 3.2 3.2 2 4 3 1 1 4 1
2 4 1 2.5 4 1 3 4 1 4 4 1 normal
1;;
```

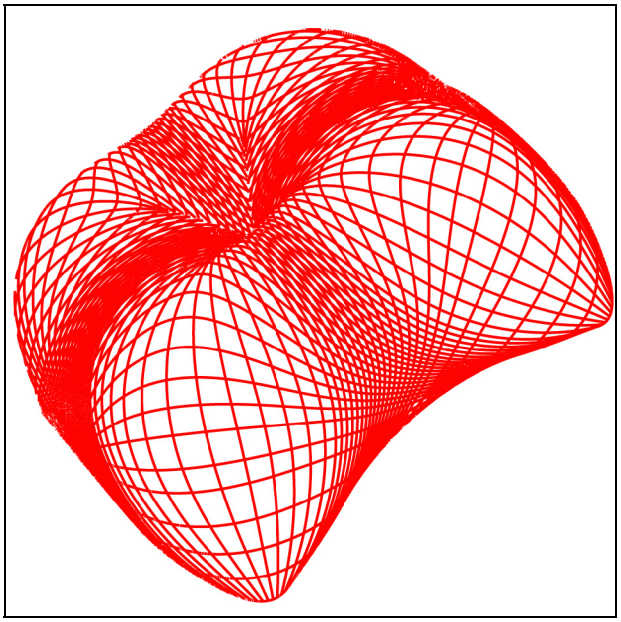

Figure 95 Cubic spline surface w/ derivatives

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In this next example, two surfaces are constructed from 3D cubic spline curves. All of the control points are from the unit sphere. The boundary control points are the same for both surfaces. Only the interior control point differs. This pair of surfaces is intended to demonstrate the need to select points in the cross section curves so that they correspond to the same features. But this example also demonstrates the absurdity of trying to approximate a sphere with a cubic spline surface with only 9 control points. It also shows that the boundary edges do not follow the 3D cubic spline curves that pass through the same control points.

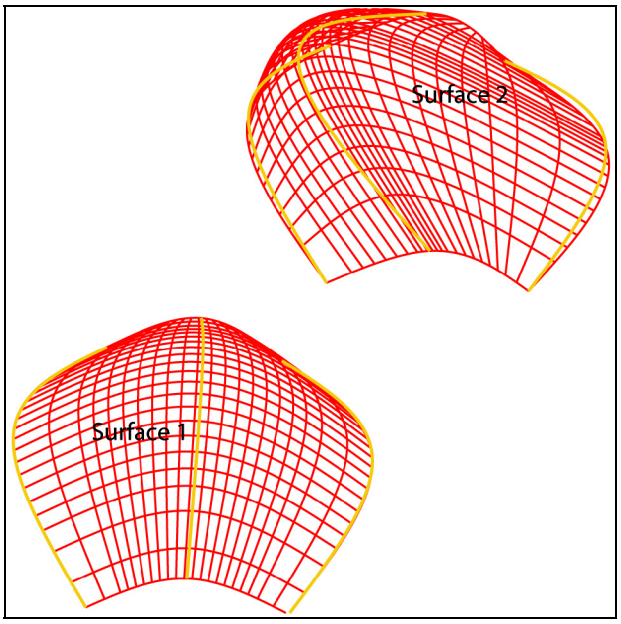

Figure 96 Cubic spline surface from curves

```
curd 1 csp3 00
  -5.2497667e-01  -2.9622331e-07
8.5111660e-01
  -1.0000000e+00  -5.6425995e-07
2.2724271e-07
 -5.2497709e-01 -2.9622353e-07 -8.5111636e-01 ;;;
curd 2 csp3 00
 -2.6248786e-01 -4.5464340e-01 8.5111660e-01
 -4.9999908e-01 -8.6602592e-01 2.2724271e-07
 -2.6248807e-01 -4.5464376e-01 -8.5111636e-01 ;;;
curd 3 csp3 00
 2.4999917e-01 -4.3301326e-01 8.6602539e-01
 4.9999827e-01 -8.6602640e-01 2.2724271e-07
 2.8678745e-01 -4.9673271e-01 -8.1915170e-01 ;;;
sd 1 csps 3 3 0000
 -5.2497667e-01 -2.9622331e-07 8.5111660e-01
 -1.0000000e+00 -5.6425995e-07 2.2724271e-07
 -5.2497709e-01 -2.9622353e-07 -8.5111636e-01
 -2.6248786e-01 -4.5464340e-01 8.5111660e-01
 -4.9999908e-01 -8.6602592e-01 2.2724271e-07
 -2.6248807e-01 -4.5464376e-01 -8.5111636e-01
 2.4999917e-01 -4.3301326e-01 8.6602539e-01
 4.9999827e-01 -8.6602640e-01 2.2724271e-07
 2.8678745e-01 -4.9673271e-01 -8.1915170e-01 ;;;
curd 4 csp3 00
 -5.2497667e-01 -2.9622331e-07 8.5111660e-01
 -1.0000000e+00 -5.6425995e-07 2.2724271e-07
 -5.2497709e-01 -2.9622353e-07 -8.5111636e-01 ;;;
```

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```
curd 5 csp3 00
 -2.6248786e-01 -4.5464340e-01 8.5111660e-01
 -7.5146174e-01 -2.2497362e-01 6.2023556e-01
 -2.6248807e-01 -4.5464376e-01 -8.5111636e-01 ;;;
curd 6 csp3 00
 2.4999917e-01 -4.3301326e-01 8.6602539e-01
 4.9999827e-01 -8.6602640e-01 2.2724271e-07
 2.8678745e-01 -4.9673271e-01 -8.1915170e-01 ;;;
sd 2 csps 3 3 0000
 -5.2497667e-01 -2.9622331e-07 8.5111660e-01
 -1.0000000e+00 -5.6425995e-07 2.2724271e-07
 -5.2497709e-01 -2.9622353e-07 -8.5111636e-01
 -2.6248786e-01 -4.5464340e-01 8.5111660e-01
 -7.5146174e-01 -2.2497362e-01 6.2023556e-01
 -2.6248807e-01 -4.5464376e-01 -8.5111636e-01
 2.4999917e-01 -4.3301326e-01 8.6602539e-01
 4.9999827e-01 -8.6602640e-01 2.2724271e-07
 2.8678745e-01 -4.9673271e-01 -8.1915170e-01 ;;;
```

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cy infinite cylinder

cy $x_0\ y_0\ z_0\ x_n\ y_n\ z_n$ *radius*

where

$x_0\ y_0\ z_0$	point on the cylinder's axis
$x_n\ y_n\ z_n$	direction vector which defines the cylinder's axis
radius	radius of the cylinder

Remarks

This command defines a cylinder with radius, *radius*, and an axis given by (x_0,y_0,z_0) and (x_n,y_n,z_n), as shown in the picture below. Specifically, (x_0,y_0,z_0) is any point on the cylinder's axis and the cylinder's axis is parallel to the vector which extends from the origin to the point (x_n,y_n,z_n). The radius of the cylinder must be positive. When projecting a point onto this surface, be sure that it is initialized somewhere away from the axis of symmetry to make it clear which direction to project. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

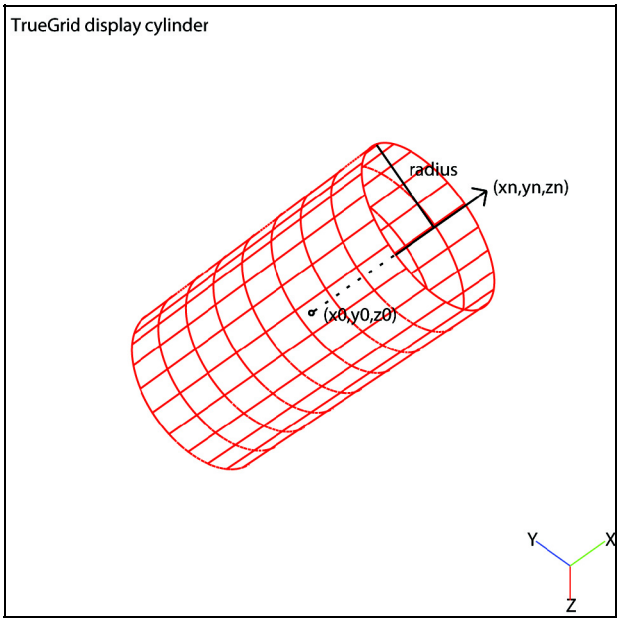

Figure 97 Cylinder by an axis and a radius

Example

```
sd 5 cy 0 0 0 1 0 0 12.5
```

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cy3 cylinder from 3 points

cy3 $x_1\ y_1\ z_1\ x_2\ y_2\ z_2\ x_3\ y_3\ z_3$

where

$(x_1\ y_1\ z_1)$	global coordinates of the first point
$(x_2\ y_2\ z_2)$	global coordinates of the second point
$(x_3\ y_3\ z_3)$	global coordinates of the third point

Remarks

The three points are assumed to form a cross section of the cylinder, perpendicular to the axis of symmetry.

Example

```
curd 1 arc3 whole rt .1 .2 .3
                  rt -.2 .3 1
                  rt -1 -2 -1 ;
sd 1 cy3
 -.9190374 -.05553633 1.7810655
    -.23211999     -1.1207378
-1.0815152
 -2.073 -2.403686 .35669363
```

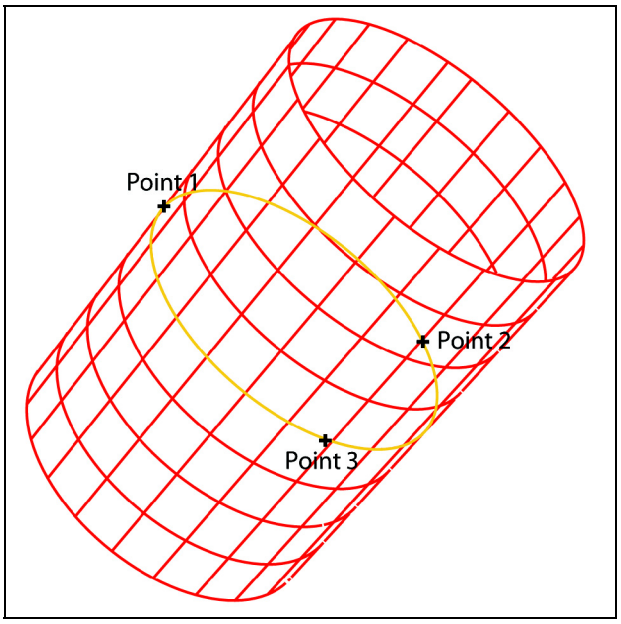

Figure 98 Cylinder by 3 points

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cyr2 cylinder from 2 points and radius

cyr2 $x_1\ y_1\ z_1\ x_2\ y_2\ z_2\ radius$

where

$(x_1\ y_1\ z_1)$	global coordinates of the first point
$(x_2\ y_2\ z_2)$	global coordinates of the second point
radius	radius of the cylinder

Remarks

The two points are used to establish the axis of rotation. They are selected in global Cartesian coordinates.

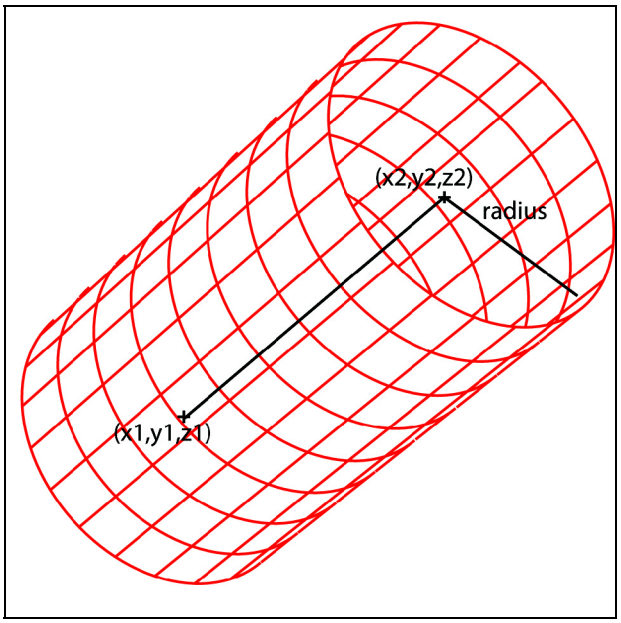

Figure 99 2 points form the axis

Example

```
sd 1 cyr2 1 2 3 1.5 4 5 1 ;
```

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cyr3 cylinder from 3 points and radius

cyr3 $x_1\ y_1\ z_1\ x_2\ y_2\ z_2\ x_3\ y_3\ z_3$ *radius*

where

$(x_1\ y_1\ z_1)$	global coordinates of the first point
$(x_2\ y_2\ z_2)$	global coordinates of the second point
$(x_3\ y_3\ z_3)$	global coordinates of the third point
radius	radius of the cylinder

Remarks

Three points are selected on a planar cross section, perpendicular to the axis, are used to establish the axis. The radius of the cylinder that passes through these three points is ignored and uses the one specified. This makes it easy to create a plug to a hole with a gap between the plug and the hole. The points are selected in global Cartesian coordinates.

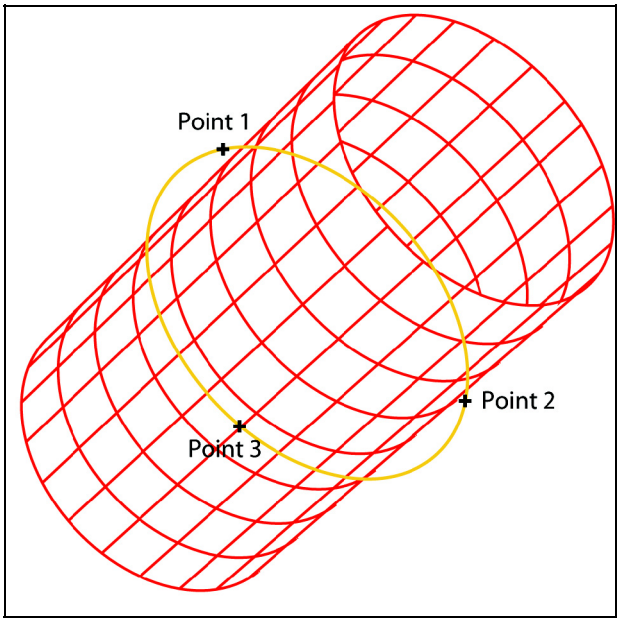

Figure 100 3 points from a circle

Example

```
curd 1 arc3 whole rt .1 .2 .3
                  rt -.2 .3 1
                  rt -1 -2 -1 ;
circent
 -.9190374 -.05553633 1.7810655
    -.23211999    -1.1207378
-1.0815152
 -2.073 -2.403686 .35669363
c  center =  -1.01674E+00  -1.07451E+00   4.24898E-01
c  radius =    1.6991307
c  normal =  -7.11559E-01   5.85396E-01  -3.88582E-01
sd 1 cyr3
 -.9190374 -.05553633 1.7810655
 -.23211999 -1.1207378 -1.0815152
 -2.073 -2.403686 .35669363
 [1.6991307-.25]
```

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er ellipsoid (ellipse revolved about an axis)

er x_0 y_0 z_0 x_n y_n z_n r_1 r_2

where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of axis direction vector
y_n	second component of axis direction vector
z_n	third component of axis direction vector
r_1	radius in the axial direction
r_2	radius orthogonal to the axis

Remarks

An ellipse is rotated about an axis given by (x_0,y_0,z_0) and (x_n,y_n,z_n). To specify the shape of the ellipse, you give the lengths of its two axes: r_1 is the length of the axis perpendicular to the axis of rotation, and r_2 is the length of the axis along the axis of rotation. Both lengths must be positive. A point projected onto this surface from within should be away from the axis of symmetry. See **Figure 101**.

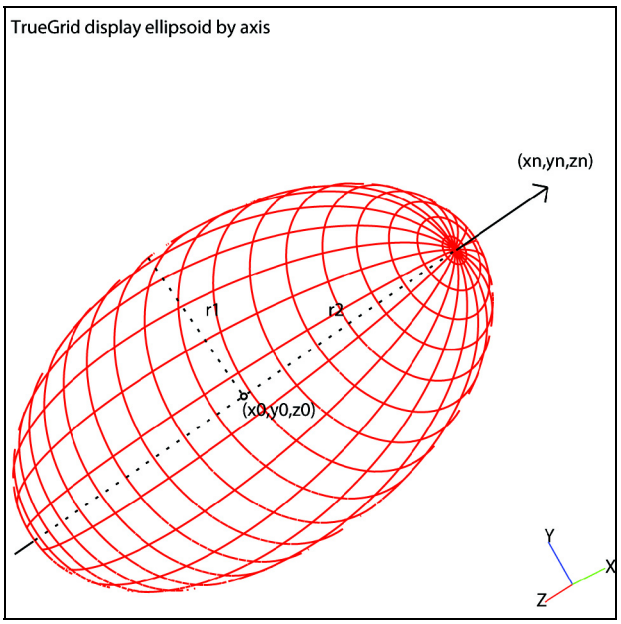

Figure 101 ellipsoid by axis and two dimensions

Example

```
sd 6 er 0 0 0 1 0 0 1 2
```

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face face of the present part (Part Phase only)

face *region*

Remarks

The region should be a face of the current part; the surface is defined to match the face as shown in **Figure 102**. This surface becomes available after the **endpart** command finishes the definition of the part.

Use the **fetol** command before this command to extract interior features as edges.

Example

```
sd 1 face 1 2 1 1 5 6
```

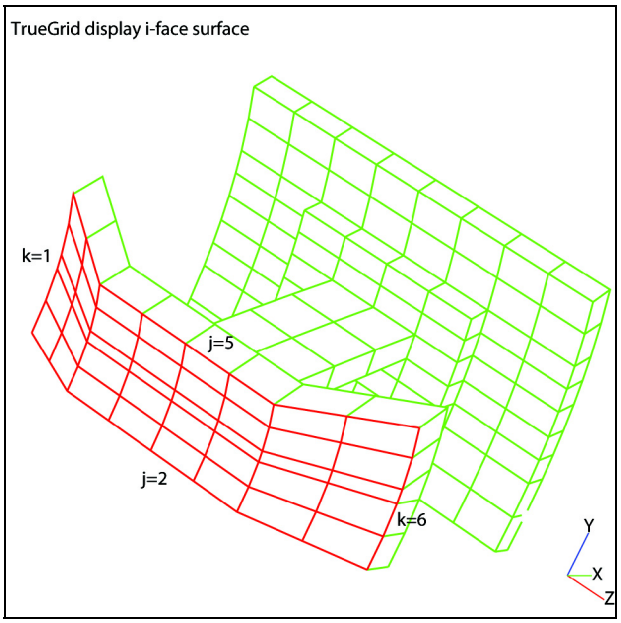

Figure 102 Face of a Part as a Surface

faceset convert a face set into a surface

faceset *set_name*

where *set_name* is the name of a face set

Remarks

This option is only available in the merge phase. A face set can be created using the set operations in the part or merge phase. The part phase method is parametric and the merge phase is very general. The command to use is **fset**. Alternatively, select the ***Pick*** button in the merge phase environment window followed by the ***Sets*** and ***Faces*** buttons. Use the lasso to select the faces. These methods can be combine to form the desired face set. The set is then easily converted to a surface with this command.

When 8-noded faces are used, they are converted to four 4-noded faces using the centroid as a 9th node.

Use one of the merging commands such as **stp** before using this command. This causes near duplicate nodes to be removed and the surface will have no extra or strange interior edges.

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Use the **fetol** command before this command to extract interior features as edges.

This surface is ideal when a hex mesh is to be attached to an existing model. For example, a NASTRAN model can be read in and the exterior of that mesh can be converted to a surface using this method. This is required when using the **blude** part command.

Example

A NASTRAN mesh of a simple ship is import using the **readmesh** command and one of the sides of the ship (wet surface only) is selected using the interactive set creation features. This set is then converted to a surface using:

```
sd faceset starboard

c starboard is the face set highlighted in red.
c These faces where selected using the
c lasso after selecting pick, sets, and faces.
```

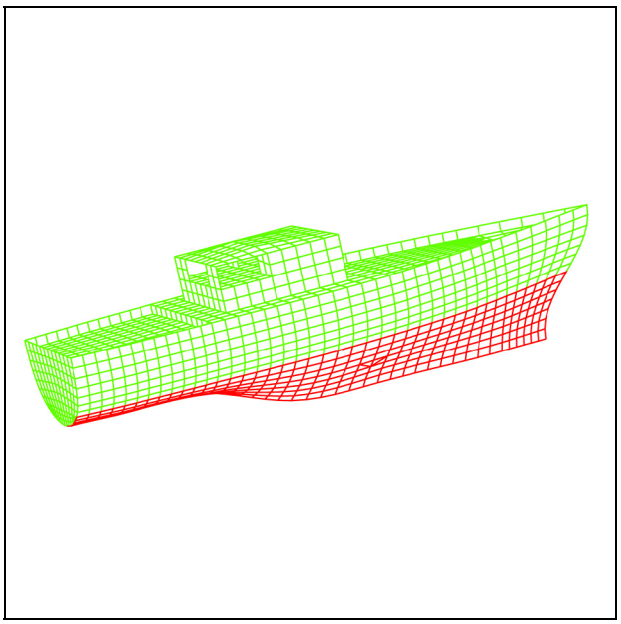

Figure 103 Faceset surface

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function surface by three algebraic expressions

function u_{min} u_{max} v_{min} v_{max} *x-expression ; y-expression ; z-expression ; trans ;*

where

u_{min}	minimum value of the first independent variable
u_{max}	maximum value of the first independent variable
v_{min}	minimum value of the second independent variable
v_{max}	maximum value of the second independent variable
x-expression	algebraic expression for first 3D coordinate
y-expression	algebraic expression for second 3D coordinate
z-expression	algebraic expression for third 3D coordinate
trans	sequence of simple operators to transform the coordinates

Remarks

Function defines a surface with algebraic forms referencing independent variables u and v in the rectangle:

$$u_{min} < \mathbf{u} < u_{max}, \qquad v_{min} < \mathbf{v} < v_{max}.$$

Three Fortran-like expressions map the rectangle to the coordinates x, y, and z in the global coordinate system.

Example

```
sd 2 function -1 1 -1 1
u;v;u*u-v*v;;
```

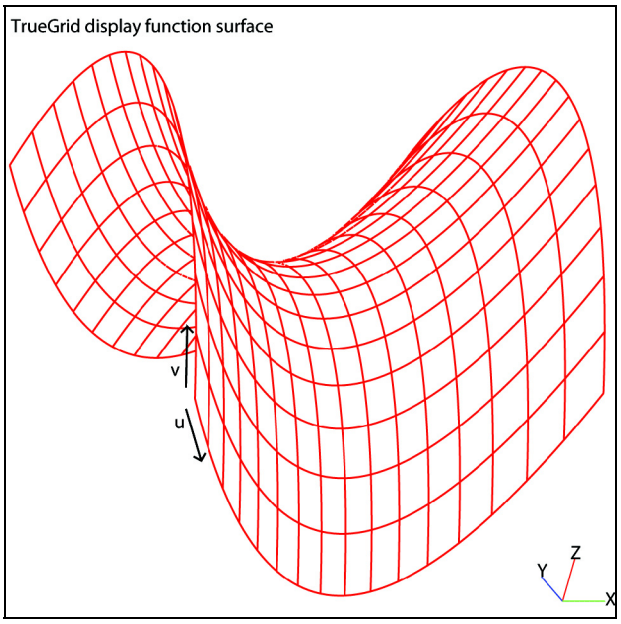

Figure 104 Saddle surface using functions

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hermite precision 2nd order spline surface

hermite *#_columns #_rows flag conditions* $x_{11}\ y_{11}\ z_{11}\ x_{12}\ y_{12}\ z_{12}$ ... [normal δ] *trans ;*

where

#_columns	number of control points in each column of the table of control points
#_rows	number of control points in each row of the table of control points
flag	4 digits formed using **0** for natural boundary condition, **1** for specified boundary derivatives, **2** for a loop, one digit for each of the four edges

where *conditions* specifies the list of derivatives along the edges, in order, when a flag of 1 is used in the corresponding digit in the *flag*

where each *condition* is formed by derivatives

dx_{ij}, dy_{ij}, dz_{ij}	derivative triplet, one for each control point along the edge
$x_{ij}\ y_{ij}\ z_{ij}$	list of coordinates for the control points on the surface
normal δ	optional argument for a normal offset
trans	sequence of simple operators to transform the coordinates

Remarks

This is designed to place a smooth surface through a very large number of regularly spaced points. It has 2 derivatives at each point. The resulting surface is very accurate in that it passes approximately through all of the points within a very small tolerance. This surface is different from a cubic spline surface in that it is not 2 times differentiable everywhere. This 2nd derivative, which is typically found in a cubic spline surface, has been sacrificed to get a surface that is very fast and very accurate. It can easily handle millions of control points and has at least 1 derivative everywhere. The form of this data for this surface is identical to a cubic spline surface, so, if after trying a cubic spline and not getting the desired results, you can easily try this type of spline surface.

A normal offset can be applied to this surface by issuing the normal option followed by the offset as the first operator in the transformation *trans*.

Use the **wiges** command to write surfaces of **hermite** type to an IGES file as a parametric spline surface.

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Examples

In the examples below, all have the same set of control points. Only the edge derivatives are different.

In this first example, the natural derivatives are used.

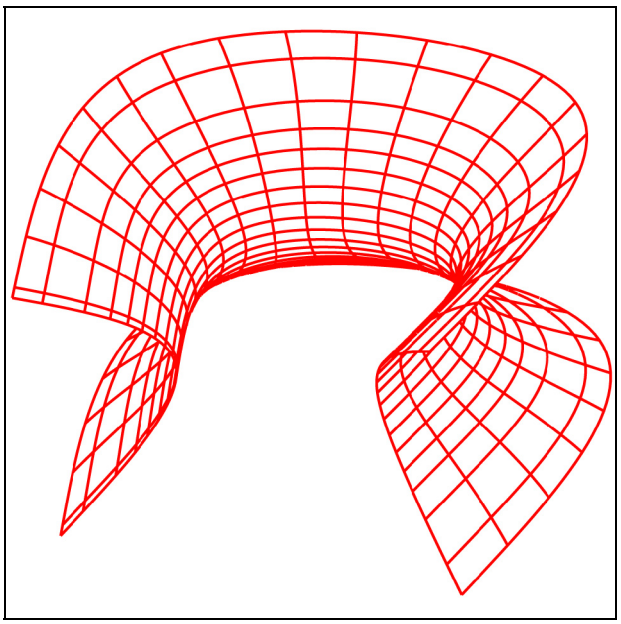

Figure 105 Natural derivatives

```
sd 1 hermite 4 3 0000
 2 0 0 0 2 0 -2 0 0 0 -2 0
 1 0 1 0 1 1 -1 0 1 0 -1 1
 2 0 2 0 2 2 -2 0 2 0 -2 2 ; ;
```

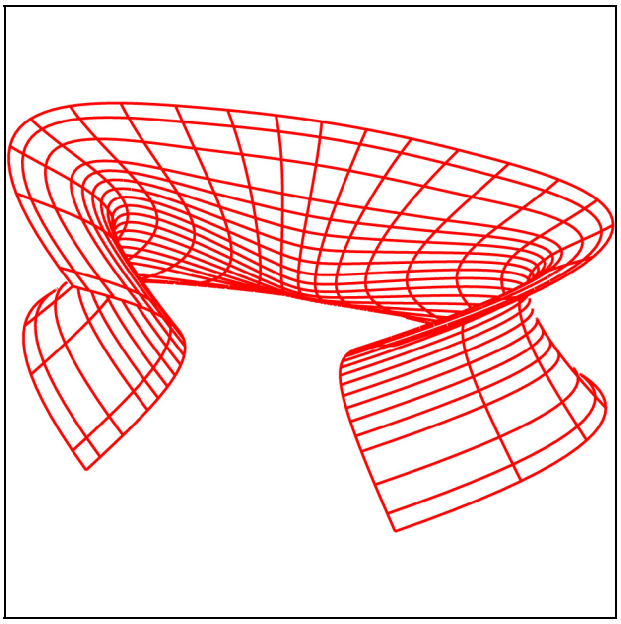

Figure 106 Edge derivatives

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In the next example, derivatives have been selected. The magnitude of the derivatives can be changed to vary the effect of the derivatives.

```
para d 10;
sd 1 hermite 4 3 1100
 0 %d 0 0 %d 0 0 %d 0
 %d 0 0 %d 0 0 %d 0 0
 2 0 0 0 2 0 -2 0 0 0 -2 0
 1 0 1 0 1 1 -1 0 1 0 -1 1
 2 0 2 0 2 2 -2 0 2 0 -2 2 ; ;
```

Two edges can be forced to match, including their derivatives.

```
sd 1 hermite 4 3 2200
 2 0 0 0 2 0 -2 0 0 0 -2 0
 1 0 1 0 1 1 -1 0 1 0 -1 1
 2 0 2 0 2 2 -2 0 2 0 -2 2 ; ;
```

Both pairs if edges can be matched to form a variation of a torus.

```
sd 1 hermite 4 3 2222
 2 0 0 0 2 0 -2 0 0 0 -2 0
 1 0 1 0 1 1 -1 0 1 0 -1 1
 2 0 2 0 2 2 -2 0 2 0 -2 2 ; ;
```

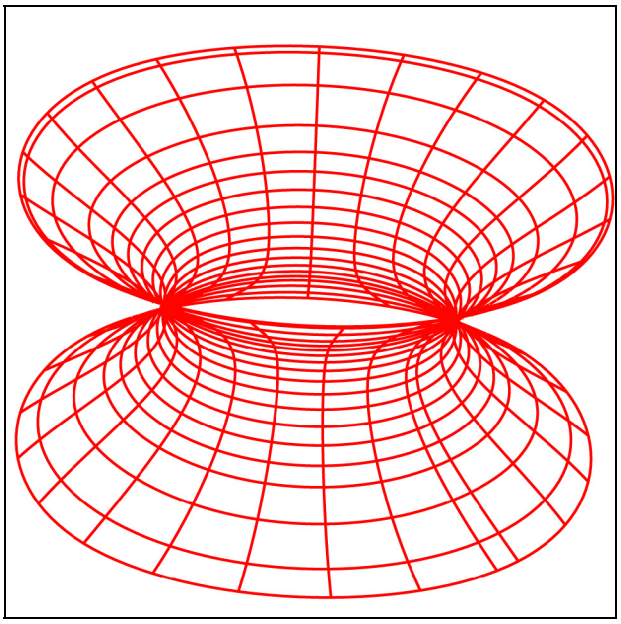

Figure 107 Periodic in 1 direction

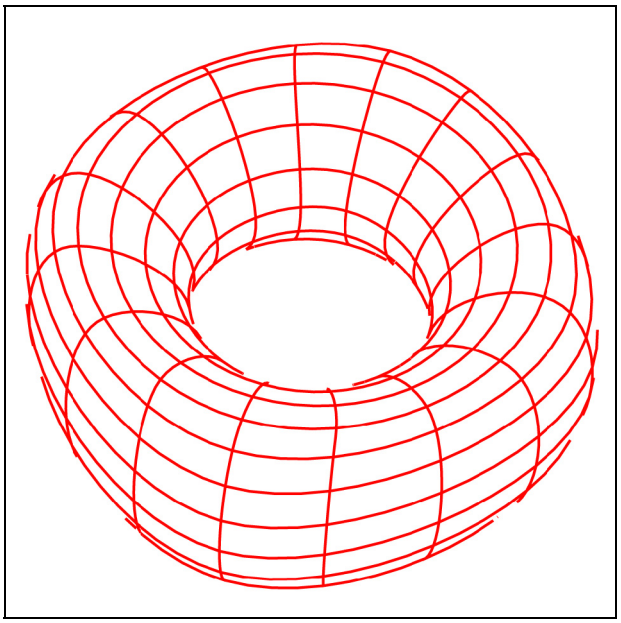

Figure 108 Periodic in both directions

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igess import a parametric IGES surface

igess *seq_# trans* ;

where

seq_#	surface's sequence number in the IGES file
trans	list of transformations to be applied to it

Remarks

You make these surfaces with a CAD system, and write them to a file in IGES format. Then you can read the entity directory from the IGES file using the **igesfile** command (or less appropriately the **iges** command). The **igesfile** command assigns a sequence number to each of the parametric surfaces it encounters in the IGES file. The argument *seq_#* in this option is the parametric surface sequence number.

IGES surfaces have sequence numbers based on whether they are a NURBS, plane, or parametric type. This command imports one of the parametric type.

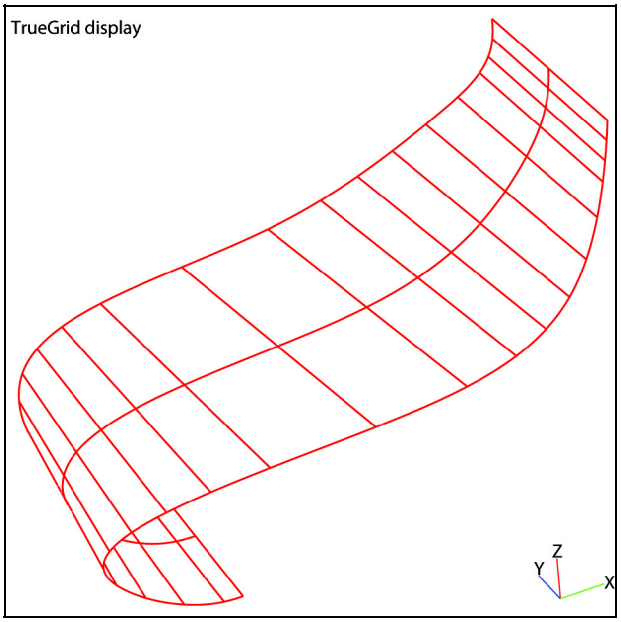

Figure 109 Imported parametric IGES Surface

When this IGES surface is assigned a surface definition number using this option, the surface data are read and the surface is rendered. The time it takes to render is a function of the curvature of the surface and of the trimming curves, because the surface must be evaluated more frequently in the areas of higher curvature.

When surfaces are trimmed, in which case all surfaces are usually trimmed, the surface will be counted twice, once as an untrimmed surface and once as a trimmed surface. All of the untrimmed surfaces are given sequence numbers first. Then the trimmed versions of the same surfaces are assigned sequence numbers in the same order as the untrimmed surfaces.

Alternatively, you can import all of the surfaces from an IGES file using the **iges** command. The **iges** command is much easier to use but you have to render all of the surfaces. You can use the **igessd** command to import a sequence of parametric surfaces instead of issuing the **sd** command many times.

There may be times when you use the **iges** command to import all of the surfaces and then want to import a single surface with the **sd** command. You may wish to do this because you want the

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untrimmed version of the surface or you may want another copy of the surface with a transformation applied. The **igess** option of the **sd** command can also be applied to the parametric surfaces assigned sequence numbers from the **iges** command (i.e. there is no need to issue the **igesfile** command if you have already issued the **iges** command).

Example

```
igesfile surfaces.igs
```

The text window and the tsave file will show the sequence numbers for the surfaces found in the IGES file.

```
c  IGES file contained   114 surfaces from     1 to   114
c  IGES file contained    24 NURBS surfaces from     1 to    24
c  IGES file contained     4 planes from     1 to     4
```

The sequence numbers reported are then used to select the surface that is desired.

```
sd 1 igess 101;
```

If you read another IGES file, the sequence numbers are added to the existing sequence numbers.

```
igesfile more_surfaces.igs
```

The report to the text window and the tsave is:

```
c  IGES file contained    11 surfaces from   115 to   125
c  IGES file contained     4 NURBS surfaces from    25 to    28
c  IGES file contained     4 planes from     5 to     8
```

```
sd 2 igess 125 rx 30 rxy ;
```

In this last example, the surface is transformed by rotating about the x-axis by 30 degrees and reflected through the xy-coordinate plane before it is rendered.

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igesp import an IGES plane

igesp *seq_# trans* ;

where

seq_#	plane's sequence number in the IGES file
trans	list of transformations to be applied to it

Remarks

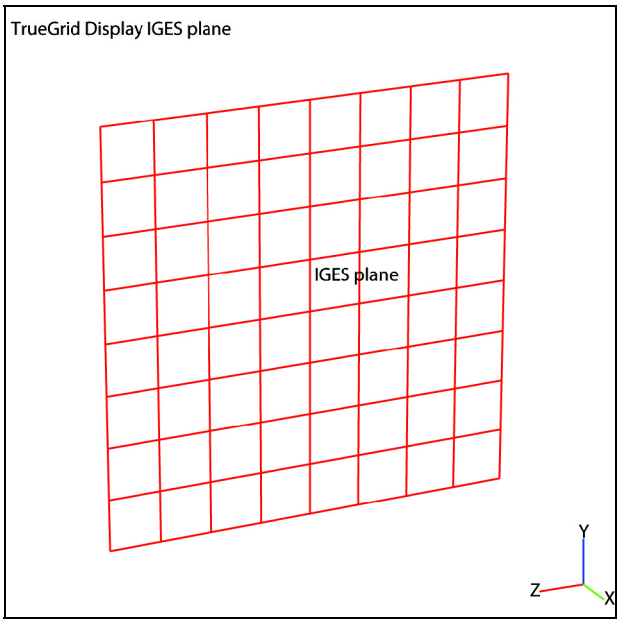

Figure 110 Imported IGES Plane

You make these surfaces with a CAD system, and write them to a file in IGES format. Then you can read the entity directory from the IGES file using the **igesfile** command (or less appropriately the **iges** command). The **igesfile** command assigns a sequence number to each of the plane surfaces it encounters in the IGES file. The argument *seq_#* in this option is the plane sequence number.

IGES surfaces have sequence numbers based on whether they are a NURBS, plane, or other type. This command imports one of the planar type.

When a planar surface is assigned a surface definition number using this option, the surface parameters are read and the surface is rendered. The time it takes to render is a function of the curvature of the trimming curves, because the surface must be evaluated more frequently in the high curvature areas.

When surfaces are trimmed, in which case all surfaces are usually trimmed, the surface will be counted twice, once as an untrimmed surface and once as a trimmed surface. All of the untrimmed surfaces are given sequence numbers first. Then the trimmed versions of the same surfaces are assigned sequence numbers in the same order as the untrimmed surfaces.

Alternatively, you can import all of the surfaces from an IGES file using the **iges** command. The **iges** command is much easier to use but you have to render all of the surfaces. You can use the **igespd** command to import a sequence of planar surfaces instead of issuing the **sd** command many times.

There may be times when you use the **iges** command to import all of the surfaces and then want to import a single surface with the **sd** command. You may wish to do this because you want the

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untrimmed version of the surface or you may want another copy of the surface with a transformation applied. The **igesp** option of the **sd** command can also be applied to the planar surfaces assigned sequence numbers from the **iges** command (i.e. there is no need to issue the **igesfile** command if you have already issued the **iges** command).

Example

```
igesfile surfaces.igs
```

The text window and the tsave file will show the sequence numbers for the surfaces found in the IGES file.

```
c  IGES file contained   114 surfaces from     1 to   114
c  IGES file contained    24 NURBS surfaces from     1 to    24
c  IGES file contained     4 planes from     1 to     4
```

The sequence numbers reported are then used to select the surface that is desired.

```
sd 1 igesp 1;
```

If you read another IGES file, the sequence numbers are added to the existing sequence numbers.

```
igesfile more_surfaces.igs
```

The report to the text window and the tsave is:

```
c  IGES file contained    11 surfaces from   115 to   125
c  IGES file contained     4 NURBS surfaces from    25 to    28
c  IGES file contained     4 planes from     5 to     8

sd 2 igesp 6 rx 10 ry 10 ;
```

In this last example, the surface is transformed by rotating about the x-axis by 10 degrees and rotating about the y axis 10 degrees before it is rendered.

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intp interpolate a surface between two surfaces

intp sd_1 sd_2 *fraction*

Remarks

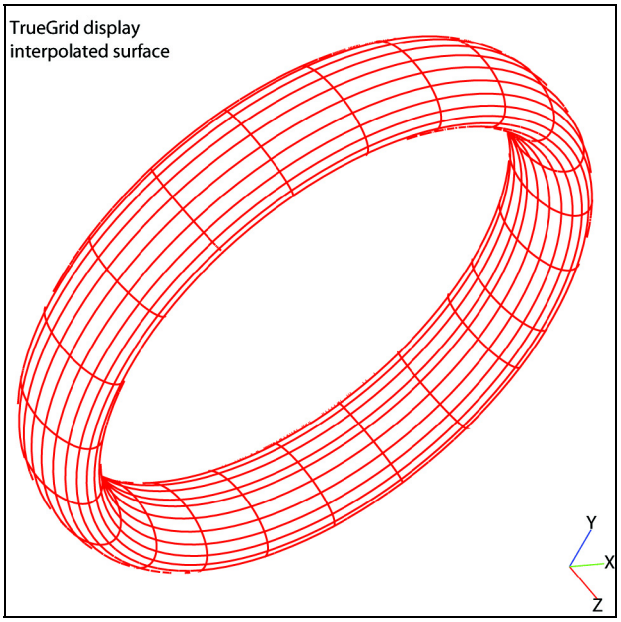

Figure 111 Surface Interpolated Between Two Other Surfaces

The two surfaces sd_1 and sd_2 must be defined before they are referenced by this command. The first surface is offset toward the second surface by projecting the points from the first surface onto the second surface. The offset distance is specified by a fraction of the distance from the first surface to the second surface. The first surface must be one of the following types: a rotated 2D curve (**cr**, **crx**, **cry**, or **crz**), a sphere (**sp**), a rotated ellipse (**er**), a structured mesh (**mesh**), a linear interpolation between two curves (**rule2d**), an interpolation between 2 3D curves (**rule3d**), a torus (**tr**), 2D cross sections placed along a 2D curve (**swept**), another **intp**, an IGES surfaces other than a plane, or a composite surface formed from the surface types above. Care is needed when choosing the second surface or unexpected results may occur. For example if the second surface has a sharp extrusion, many of the points from the first surface will project to that point. The **intp** surface is the only surface type valid for the variable shell surface command **ssf**.

Example

```
sd 1 ts 0 0 0 0 0 1 5 0 1
sd 2 cy 0 0 0 0 0 1 5
sd 3 intp 1 2 .5
```

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iplan infinite plane defined by an implicit function

iplan *a b c d*

where the arguments are coefficients in the equation below.

Remarks

This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture changes. The plane is defined by the

$$a\,x + b\,y + c\,z = d$$

set of points that satisfy
where the coordinate triplet *(x,y,z)* is any point on the plane and where at least one of the variables *a*, *b*, or *c* is different from zero as shown in **Figure 112**.

```
sd 3 iplan 1 -1 1 1
```

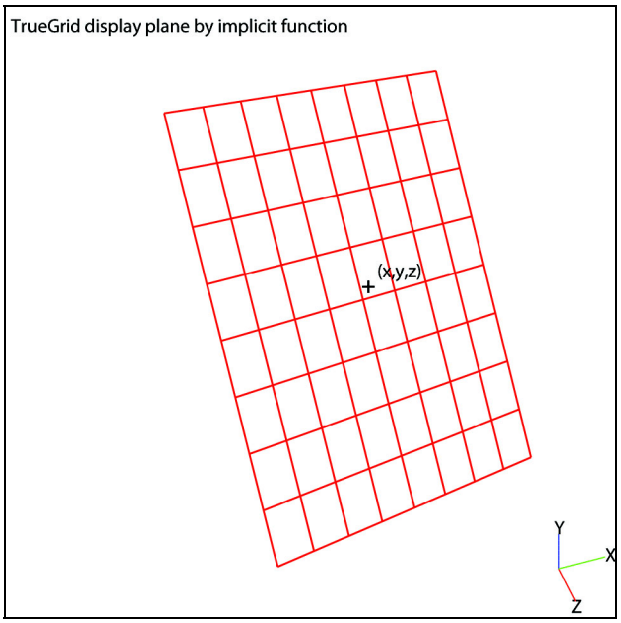

Figure 112 Plane specified by an implicit function

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mesh surface by tabular points

mesh *m n* x_{11} y_{11} z_{11} x_{21} y_{21} z_{21} ... x_{mn} y_{mn} z_{mn} *trans ;*

where

m	number of rows in the table or 2D array
n	number of columns in the table or 2D array
(x_{ij}, y_{ij}, z_{ij})	3D coordinates at position (i,j) in the table or 2D array
trans	standard coordinate transformation (sequence of operators)

Remarks

Mesh defines a surface by specifying a table or 2D array of points to be part of the surface. The rest of the surface is implicitly defined by bilinear interpolation between those points. This type of surface provides a general way to use experimental or computational data from any source.

The first two integers, *m* and *n*, give the number of points in the rows and columns, respectively. These parameters must be greater than 1. The remaining arguments give the positions of each mesh point in physical space. These points should be ordered the way Fortran would order a 2D array with *m* rows and *n* columns, i.e. with:

(i,j) = (1,1), (2,1), ... , (*m*,1), (1,2), ... , (*m*,2), ... , (1,*n*) , ... , (*m*,*n*).

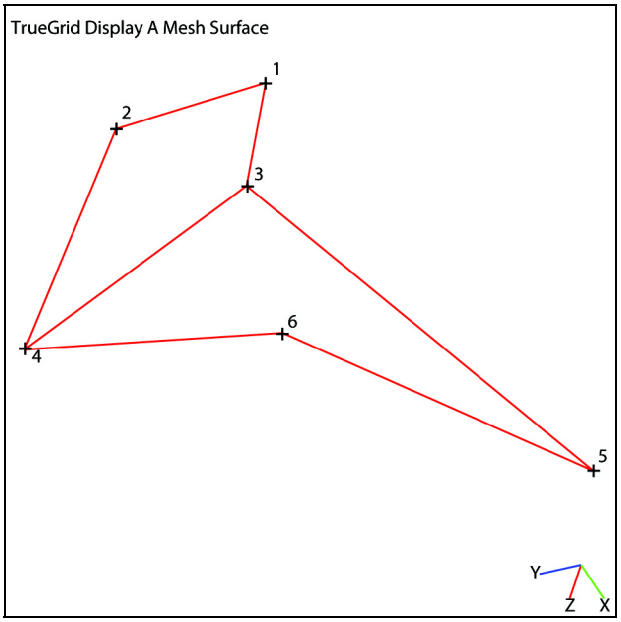

Figure 113 2 by 3 Mesh Surface

Example

```
sd 1 mesh 2 3
1.125 0.999 1.269
1.566 1.875 0.869
2.092 1.753 0.644
2.087 2.377 1.170
3.053 0.936 1.119
2.606 1.891 0.712 ;
```

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nrbs NURBS

nrbs *i-degree j-degree*
iknot$_1$ ... *iknot*$_{k1}$; *jknot*$_1$ .... *jknot*$_{k2}$;
weight$_{1,1}$ *weight*$_{1,1}$ *weight*$_{1,1}$ ... *weight*$_{1,n2}$ *weight*$_{1,n2}$ *weight*$_{1,n2}$
...
weight$_{n1,1}$ *weight*$_{n1,1}$ *weight*$_{n1,1}$ ... *weight*$_{n1,n2}$ *weight*$_{n1,n2}$ *weight*$_{n1,n2}$;
xcontrol$_{1,1}$ *ycontrol*$_{1,1}$ *zcontrol*$_{1,1}$ ... *xcontrol*$_{1,n2}$ *ycontrol*$_{1,n2}$ *zcontrol*$_{1,n2}$
...
xcontrol$_{n1,1}$ *ycontrol*$_{n1,1}$ *zcontrol*$_{n1,1}$ ... *xcontrol*$_{n1,n2}$ *ycontrol*$_{n1,n2}$ *zcontrol*$_{n1,n2}$;
trans ;
where

i-degree j-degree	degrees of the surface polynomial in the i and j direction.
i*knot*$_1$ ... *iknot*$_{k1}$	parametric coordinates of *k1* knots (0..) in the i direction.
j*knot*$_1$ ... *jknot*$_{k2}$	parametric coordinates of *k2* knots (0..) in the j direction.
weight$_{1,1}$ *weight*$_{1,1}$ *weight*$_{1,1}$ ... *weight*$_{1,n2}$ *weight*$_{1,n2}$ *weight*$_{1,n2}$ *weight*$_{n1,1}$ *weight*$_{n1,1}$ *weight*$_{n1,1}$ ... *weight*$_{n1,n2}$ *weight*$_{n1,n2}$ *weight*$_{n1,n2}$;	weights for each control point
xcontrol$_{1,1}$ *ycontrol*$_{1,1}$ *zcontrol*$_{1,1}$ ... *xcontrol*$_{1,n2}$ *ycontrol*$_{1,n2}$ *zcontrol*$_{1,n2}$ *xcontrol*$_{n1,1}$ *ycontrol*$_{n1,1}$ *zcontrol*$_{n1,1}$ ... *xcontrol*$_{n1,n2}$ *ycontrol*$_{n1,n2}$ *zcontrol*$_{n1,n2}$;	triplets of coordinates of control points

where
$n_1 = k_1$ - *multiplicity*$_1$
$n_2 = k_2$ - *multiplicity*$_2$

trans	sequence of simple operators to transform the coordinates

Remarks

First, a few notes on the parameters. The multiple knots have to be specified for the starting and ending point of the curve. The number of multiple knots is called multiplicity. The multiplicity for the starting and ending point determines the order of the curve. (2 - linear, 3 - quadratic, 4 - cubic ...).

multiplicity = degree + 1

The multiplicity has to be the same for the starting and ending points.

A normal offset can be applied to this surface by issuing the normal option followed by the offset as the first operator in the transformation *trans*.

NURBS surfaces are also available through the IGES interface.

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Example

The NURBS surface (**Figure 114**) is defined by the same control points as the B-Spline Surface (**Figure 83**). The weights are specified for each control point. Every weight is 1.0 except 4 control points with weights 10.0 (**Figure 114**). The NURBS surface is passing closer to the control points with w=10.

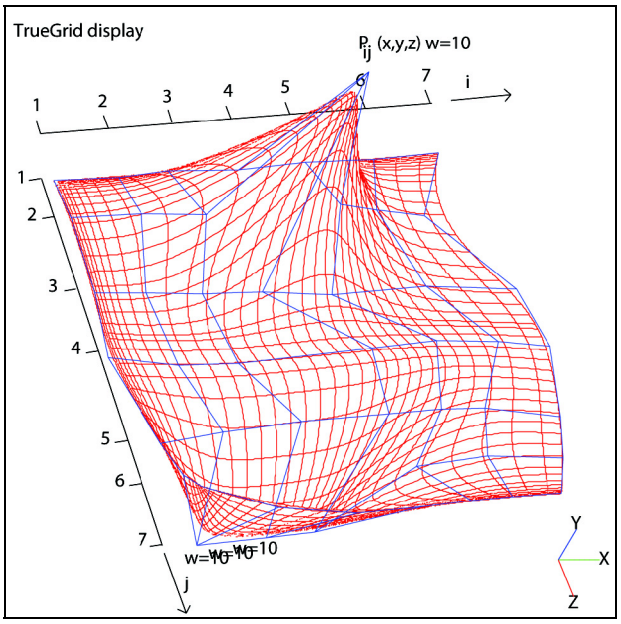

Figure 114 NURBS Surface

```
sd 1 nrbs 3 3
0 0 0 0 1 2 3 4 4 4 4;
0 0 0 0 1 2 3 4 4 4 4;
1 1 1 1 1 1 1
1 1 1 10 1 1 1
1 1 1 1 1 1 1
1 1 1 1 1 1 1
1 1 1 1 1 1 1
1 10 10 10 1 1 1
1 1 1 1 1 1 1;
...                  c x,y,z coordinates of control points
...                  c organized into 7 rows with 7 columns
...                  c the same as for BSPS surface
```

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nurbs import a NURBS surface

nurbs *seq_# trans* ;

where

seq_#	NURBS's surface sequence number in the IGES file
trans	list of transformations to be applied to it

Remarks

You make these surfaces with a CAD system, and write them to a file in IGES format. Then you can read the entity directory from the IGES file using the **igesfile** command (or less appropriately the **iges** command). The **igesfile** command assigns a sequence number to each of the NURBS surfaces it encounters in the IGES file. The argument *seq_#* in this option is the NURBS sequence number.

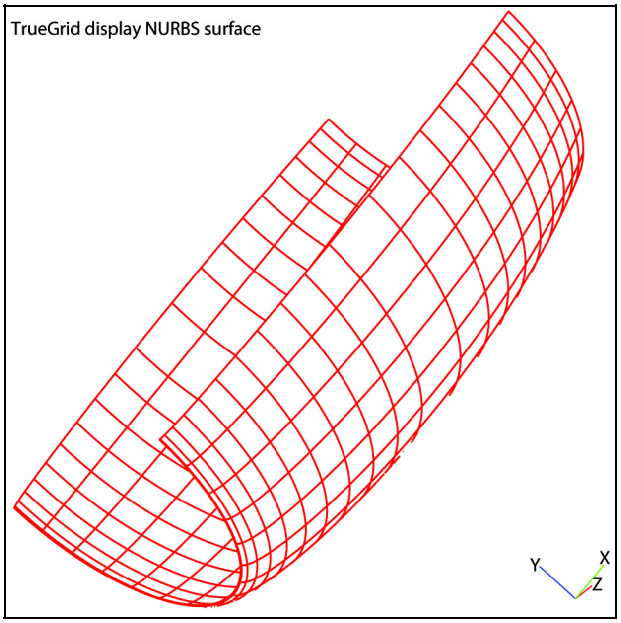

Figure 115 NURBS Surface from an IGES File

IGES surfaces have sequence numbers based on whether they are a NURBS, plane, or parametric type. This option imports one of the NURBS type.

When a NURBS surface is assigned a surface definition number using this option, the surface parameters are read and the surface is rendered. The time it takes to render is a function of the curvature of the surface and the curvature of the trimming curves, because the surface must be evaluated more frequently in the high curvature areas.

When surfaces are trimmed, in which case all surfaces are usually trimmed, the surface will be counted twice, once as an untrimmed surface and once as a trimmed surface. All of the untrimmed surfaces are given sequence numbers first. Then the trimmed versions of the same surfaces are assigned sequence numbers in the same order as the untrimmed surfaces.

Alternatively, you can import all of the surfaces from an IGES file using the **iges** command. The **iges** command is much easier to use but you have to render all of the surfaces. You can use the **nurbsd** command to import a sequence of NURBS surfaces instead of issuing the **sd** command many times.

There may be times when you use the **iges** command to import all of the surfaces and then want to

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import a single surface with the **sd** command. You may wish to do this because you want the untrimmed version of the surface or you may want another copy of the surface with a transformation applied. The **nurbs** option of the **sd** command can also be applied to the NURBS surfaces assigned sequence numbers from the **iges** command (i.e. there is no need to issue the **igesfile** command if you have already issued the **iges** command).

Example

```
igesfile surfaces.igs
```

The text window and the tsave file will show the sequence numbers for the surfaces found in the IGES file.

```
c  IGES file contained   114 surfaces from     1 to   114
c  IGES file contained    24 NURBS surfaces from     1 to    24
c  IGES file contained     4 planes from     1 to     4
```

The sequence numbers reported are then used to select the surface that is desired.

```
sd 1 nurbs 16;
```

If you read another IGES file, the sequence numbers are added to the existing sequence numbers.

```
igesfile more_surfaces.igs
```

The report to the text window and the tsave is:

```
c  IGES file contained    11 surfaces from   115 to   125
c  IGES file contained     4 NURBS surfaces from    25 to    28
c  IGES file contained     4 planes from     5 to     8

sd 2 nurbs 27 rz 90 mx .1 ;
```

In this last example, the surface is transformed by rotating about the z-axis by 90 degrees and translating in the x direction .1 before it is rendered.

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pipe sweep a pipe shape along an arbitrary 3D curve

pipe *3d_curve* $radius_1$ arc_length_1 $radius_2$ arc_length_2 ... ; *trans* ;

where

3d_curve	3D curve that is used to sweep along
$radius_i$ arc_length_i	radius and arc length pairs specified along the 3D curve
trans	optional transformation at the end

Remarks

This command allows you to define a pipe with varying cross-section. The arc length varies from a value of 0 at one end of the 3D curve to a value of 1 at the other end. In other words, the arc length is the fractional value of the entire 3D-curve length at the point where the circular cross-sectional radius is defined. You may specify as many cross-sections as you like.

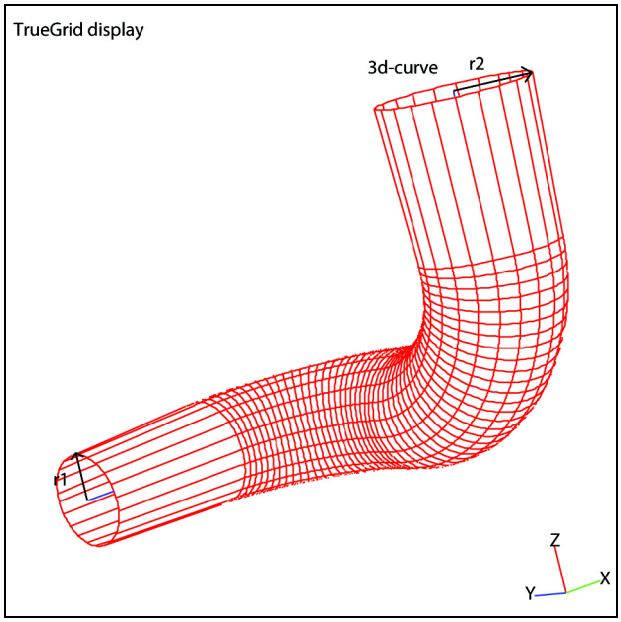

Figure 116 pipe surface

Example

```
curd 1 lp3 -1 1 1 0 1 1;;
3dfunc 0 90 cos(u;sin(u);1;;
3 d f u n c      1 8 0      2 7 0
1;sin(u);cos(u)+2;;
lp3 1 -1 3; ;
sd 1 pipe 1 .3 0 .5 1;;
```

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pl2 plane specified by two points

pl2 *system point system point*

where

system		one of the three coordinate systems:
	rt	for Cartesian
	sp	for spherical
	cy	for cylindrical
point		set of 3 coordinate in the appropriate coordinate system

Remarks

This option defines the plane using 2 points. The first point is on the plane, the second point in not on the plane and on the line normal to the plane through the first point.

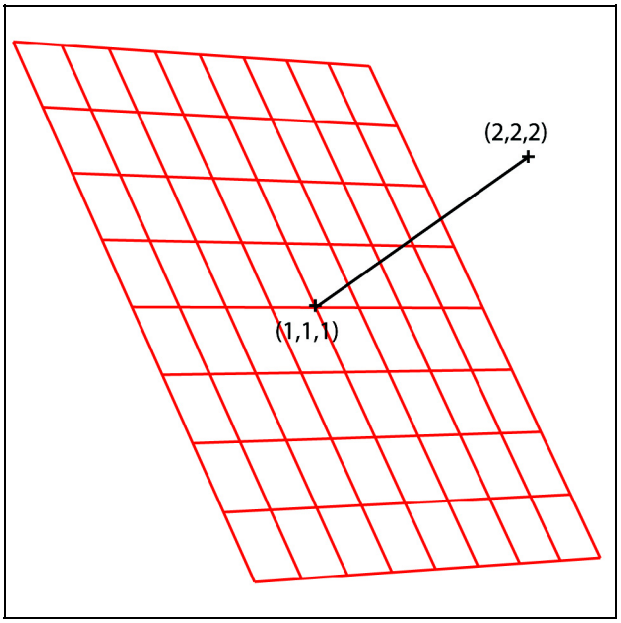

Figure 117 Plane by 2 points

Example

```
sd 1 pl2 rt 1 1 1 rt 2 2 2 ;
```

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pl3 plane specified by three points

pl3 *system point system point system point*

where

system		one of the three coordinate systems:
	rt	for Cartesian
	sp	for spherical
	cy	for cylindrical
point		set of 3 coordinate in the appropriate coordinate system

Remarks

In Cartesian coordinates, the point is designated by its x, y, and z coordinate values. In cylindrical coordinates, the radius is followed by the angle and z-coordinates. In spherical coordinates, the radius is first, followed by the polar and azimuthal angles. All angles are in degrees. Remember, the three points must not be co-linear. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

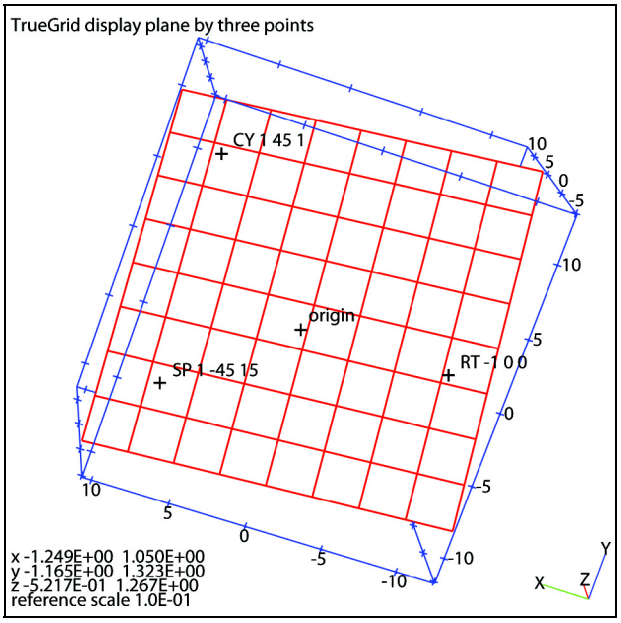

Figure 118 Plane specified by 3 points

Example

```
sd 1 pl3 cy 1 45 1 sp 1 -45 15
rt -1 0 0
```

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pl3o plane specified by 3 points and an offset

pl3o *system point system point system point offset*

where

system	one of the three coordinate systems:
rt	for Cartesian
sp	for spherical
cy	for cylindrical
point	set of 3 coordinate in the appropriate coordinate system
offset	normal offset distance

Remarks

This command defines a plane specified by three points and a positive normal offset. The direction of the normal is defined by the right hand rule through the 3 points.

The point is designated by its x, y, and z coordinate values. In cylindrical coordinates, the radius is followed by the angle and z-coordinates. In spherical coordinates, the radius is first, followed by the polar and azimuthal angles. All angles are in degrees.

This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

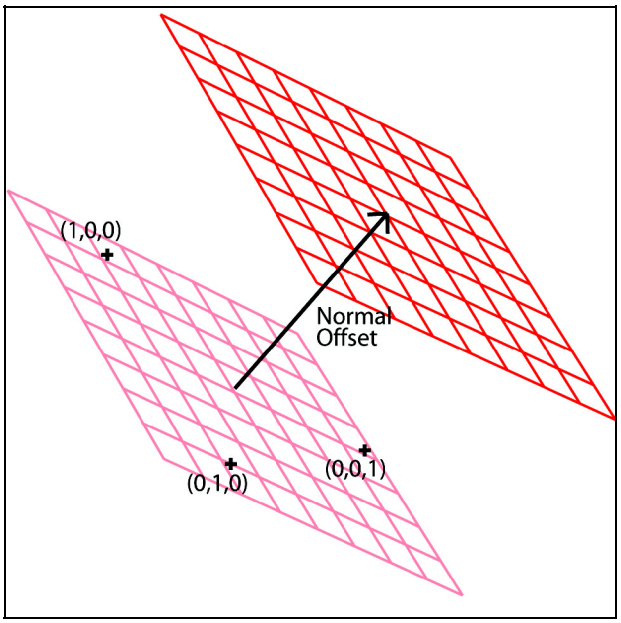

Figure 119 Normal offset plane

Example

```
sd 1 pl3o rt 1 0 0 rt 0 1 0 rt 0 0 1 1 ;
```

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plan infinite plane

plan $x_0\ y_0\ z_0\ x_n\ y_n\ z_n$

where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of the normal vector
y_n	second component of the normal vector
z_n	third component of the normal vector

Remarks

A plane containing the point (x_0,y_0,z_0) and normal to the vector (x_n,y_n,z_n) as shown in **Figure 120**.

```
sd 1 plan 0 0 0 1 0 0
```

This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

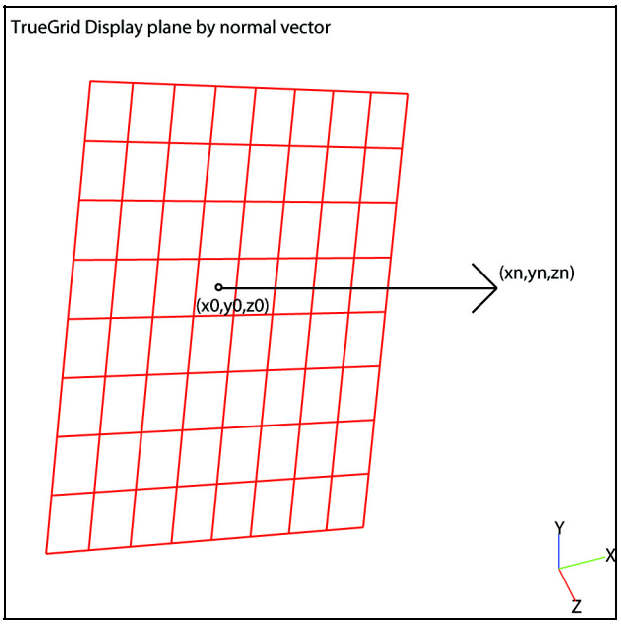

Figure 120 Plane by a point and a normal vector

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poly convert a polygon set into a surface

poly *polygon_set trans ;*

where

polygon_set	named set of polygons
trans	sequence of simple operators to transform the coordinates

Remarks

You can create a polygon surface using the **pset** command and the interactive feature to select or modify the polygons in a set with the ***Sets*** window. The ***Sets*** window can be activated in the merge phase from the ***Pick*** panel in the Environment Window. These features, alongwith the **wrsd** command, can be used to sort out complex polygon surfaces and split them into multiple surfaces or remove features. This can also be used to create a normal offset by using the normal option as one of the transformation primitives.

Use the **fetol** command before this command to extract interior features as edges.

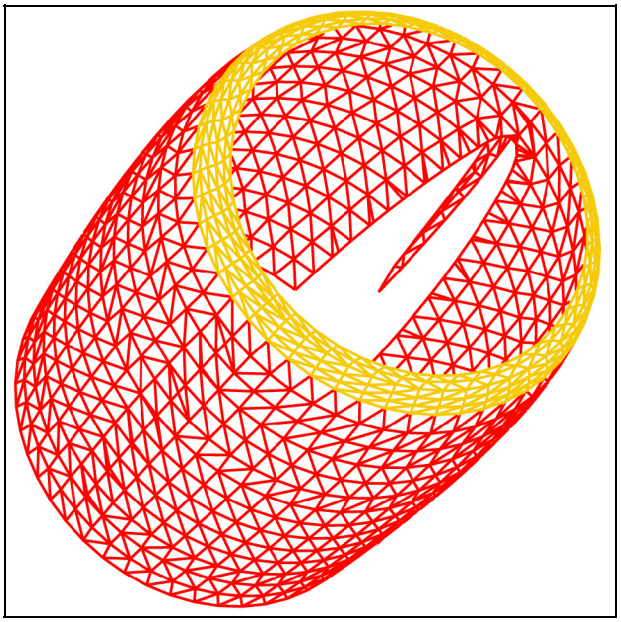

Figure 121 Surface from a polygon set

Example

It is assumed in this example that a polygon set was formed using the Sets window in the Pick panel of the Environment window. This polygon set was named p1. Then the command to create a surface form this set is:

```
sd 2 poly p1;
```

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pr paraboloid (parabola revolved about an axis)

pr $x_0\ y_0\ z_0\ x_n\ y_n\ z_n\ r_1\ t_1\ r_2\ t_2\ r_3\ t_3$

where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of axis direction vector
y_n	second component of axis direction vector
z_n	third component of axis direction vector
r_1	radius at the cross section along axis at t_1
t_1	position along axis with radius r_1
r_2	radius at the cross section along axis at t_2
t_2	position along axis with radius r_2
r_3	radius at the cross section along axis at t_3
t_3	position along axis with radius r_3

Remarks

A parabola is rotated about an axis given by (x_0,y_0,z_0) and (x_n,y_n,z_n). You define the parabola by specifying three points that it passes through in any planar cross section containing the axis of symmetry: (r_1,t_1), (r_2,t_2), and (r_3,t_3). The t coordinate is measured along the axis, starting at the point (x_0,y_0,z_0). The r coordinate is measured perpendicular to the axis. The three points can be in any order. These points must uniquely specify a parabola. This means that the coordinates must satisfy:

$$|t_1^2\times(t_3-t_2)+t_2^2\times(t_1-t_3)+t_3^2\times(t_2-t_1)| > 10^{-18}$$

A point projected to this surface should not be on the axis of symmetry. See the example in **Figure 122**. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

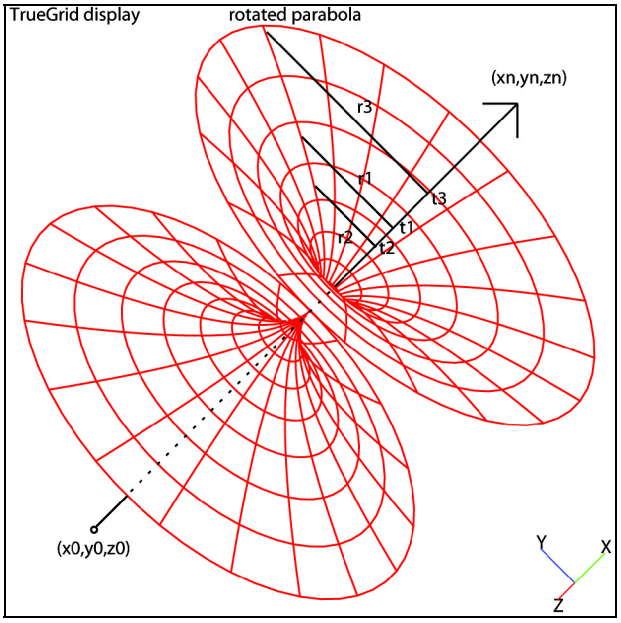

Figure 122 Parabola specified by three points

Example

```
sd 4 pr 0 0 0 1 0 0 .75 8 1.2675 8.3 3 9
```

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r3dc 3D curve revolved about an axis

r3dc $x_0\ y_0\ z_0\ x_n\ y_n\ z_n$ *3D_curve begin_angle end_angle tran ;*

where

(x_0,y_0,z_0)	point on the axis of rotation
(x_n,y_n,z_n)	vector parallel to the axis of rotation
3D_curve	3D curve being rotated (generatrix)
begin_angle	beginning angle for the rotational sweep
end_angle	ending angle for the rotational sweep
trans	sequence of simple operators to transform the coordinates

Remarks

This command forms the surface of rotation determined by the generatrix, *3D_curve*, the axis of rotation (defined by the point (x_0,y_0,z_0) and the direction vector (x_n,y_n,z_n)) and the starting and ending angles (*begin_angle* and *end_angle*, respectively).

The surface is created by rotating the generatrix around the axis of rotation from the beginning angle to the ending angle. For any point on the generatrix, the rotational arc created lies on the circle centered along the axis rotation, orthogonal to (x_n,y_n,z_n) and through that point on the generatrix. The portion of the circle retained is measured counter-clockwise starting *begin_angle* degrees from the generatrix point to *end_angle*.

The best results occur when the generatrix is coplanar with the axis, and the worst results occur along segment tangent to the generated arc.

To improve results add a point to the generatrix near to the tangent points. The geometry tolerance (set with **getol**) may need to be reduced so that the line thinning algorithm does not remove new points.

Example

```
curd 1 lp3 6.01 0 3.35  6.1 0 3.5 6
0 3.5 6 0 .5  10 0 0  11 0 2.5 11.1
0 2.5  11.2 0 3 11.1 0 3 12.1 0 5.5
12.05 0 5.5 11.995  0 5.25;;
sd 1 r3dc 0 0 0 0 0 1 1 0 60;
curd 2 cpcd 1 my -10 mz .1;
```

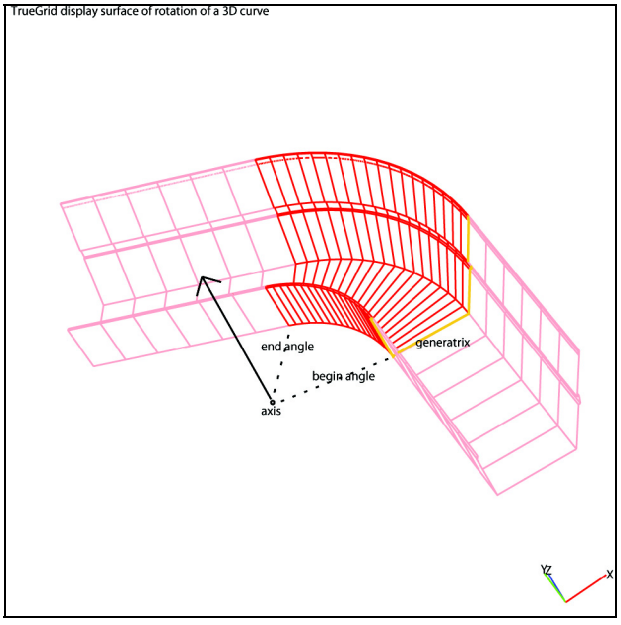

Figure 123 Surface of revolution of a 3D curve

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```
sd 2 rule3d 1 2;;
curd 3 cpcd 1 rz 60;
curd 4 cpcd 3 mx -8.66 my 5 mz -.1;
sd 3 rule3d 3 4;;
```

rule2d ruled surface between two 2D curves

rule2d y_1 ln_1 y_2 ln_2 *trans* ;

where

y_1	y-coordinate of the plane for the first 2D curve
ln_1	2D curve ID which will form one edge of the surface
y_2	y-coordinate of the plane for the second 2D curve
ln_2	2D curve ID which will form the opposite edge
trans	sequence of simple operators to transform the coordinates

Remarks

This surface is formed by a linear interpolation between two 2D curves. Prior to using this command, you must have defined the 2D curves numbered ln_1 and ln_2, using commands like **ld**. These 2D curves would originally lie in the plane y=0 (the xz-plane), but **rule2d** will translate them in the y-direction to the y=*y1* and y=*y2* planes, respectively. The interpolation amounts to constructing a straight line from each point of one translated curve to the corresponding point of the other translated curve. The order of points along the two curves (orientation) must be the same. To reverse the order of the points in one of the curves, specify that 2D curve definition number as negative. The points that correspond to each other are those which have the same relative arc length position on their respective curves. At least one of the two 2D curves must have positive arc length.

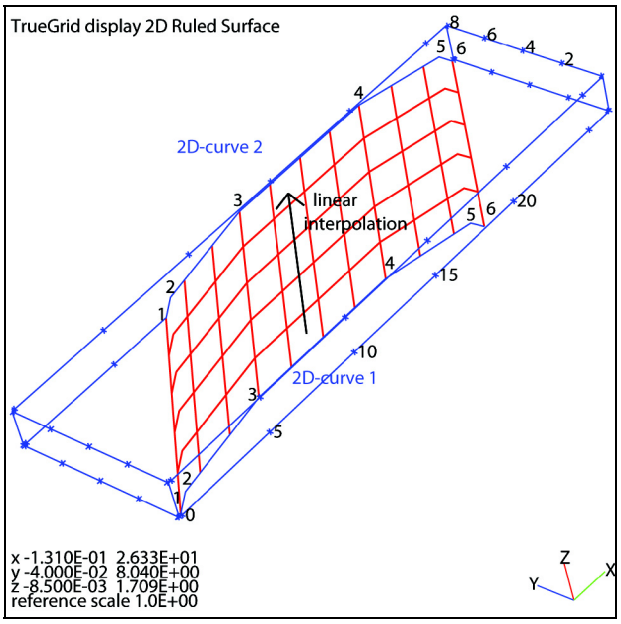

Figure 124 Ruled Surface From Two 2D Curves

Example

```
ld 1 lp 0 0  .5 .7  5 1.7 12.5 1.7 17.5 .7 18.2 0;
```

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```
ld 2 lp 8 0 8.5 .7 13 1.7 20.5 1.7 25.5 .7 26.2 0;
sd 1 rule2d 0 1 8 2;
```

rule3d ruled surface between two 3D curves

rule3d *3D-curve$_1$ 3D-curve$_2$ trans ;*

where

3D-curve$_1$	first boundary 3D curve
3D-curve$_2$	second boundary 3D curve
trans	sequence of simple operators to transform the coordinates

Remarks

A surface is formed by a linear interpolation in Cartesian coordinates between two 3D curves. At least one of the two curves must have a positive arc length. Both curves must be already defined. 3D curves are defined with the **curd** command or read from an **IGES** file. The direction of the curves is critical to the interpolation of the surface. Check the order of the points in each of the curves by displaying the points with the **label** command. In order to reverse the order of a curve, specify the curve definition number with a negative number. The surface is constructed by pairing each point on one curve with a point on the other curve. These two points are connected by a line segment. Points along the two curves are paired by relative arc length. The following is an example of three 3D curves, which are annotated below, and the two surfaces interpolated between pairs of curves.

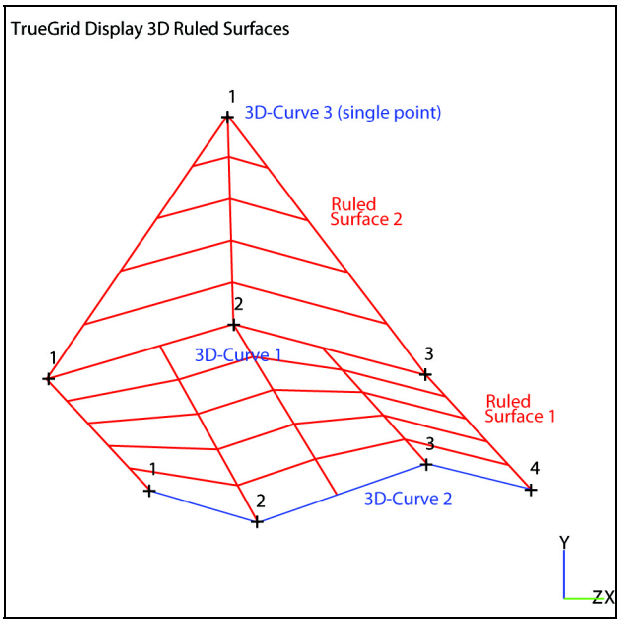

Figure 125 3D Ruled Surfaces From 3D Curves

Example

```
curd 1 lp3 1 2 1 3 2.5 1.1 5 2 1.5;;
curd 2 lp3 1.5 1 2 2 .75 3 3.5 1.25 3.5 4.5 1 3.75;;
curd 3 lp3 3 4.5 1;;
sd 1 rule3d 1 2;
sd 2 rule3d 1 3;
```

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sds union of surfaces

sds s_1 s_2 ... s_n ;

where s_i is the number of a surface previously defined

Remarks

The surfaces that form this union do not need to meet perfectly. There can be large gaps and overlappings.

This is the most important surface type when dealing with CAD models. Typically, the CAD model is broken into many surfaces for construction purposes. No consideration is made in the construction of the surfaces for mesh generation. In most cases, it is more convenient, if not essential, to form a mesh across many of these surfaces. Use this option to form a composite surface and it will be as if it were one surface. The projection of a face of the mesh works the same for both single and composite surfaces.

Sometimes, this surface type can be used to group surfaces when there are many surfaces to sort (i.e. it is very useful when working with IGES input files).

When a node is projected to a multiple surface, it is projected to each of the component surfaces. The point of projection which is the shortest distance among all of the component surface projections is defined as the projection point for the multiple surface. This makes it possible for the surfaces to overlap or not meet perfectly. A closest point will always be found.

A multiple surface list can include surfaces defined with the **sds** option.

An effective way to prepare a list of surfaces is to first display all surface definitions (**dasd**). Then, in the Environment Window, use the ***Display List***, ***Surface*** button in the Apply Action To: menu, and the ***Remove*** button to remove those surfaces which you do not want in your grouped surface. When you only have those surfaces displayed which you wish to group together, use the **lasd** command to list all of the displayed surfaces. Type

sd *some_number* sds

in the text window and cut and paste the listed surfaces, followed by a semi-colon. (Where *some_number* is the definition number of the surface that you are creating.)

Care is needed when using this option to define several multiple surfaces, each containing some of the same component surfaces. The intersection of the two multiple surfaces includes the surfaces that they both have in common. This would be an extreme case of two surfaces being tangent. This can be an

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effective technique when you want an edge of the mesh to lay along the middle of a surface common to both composites.

Example

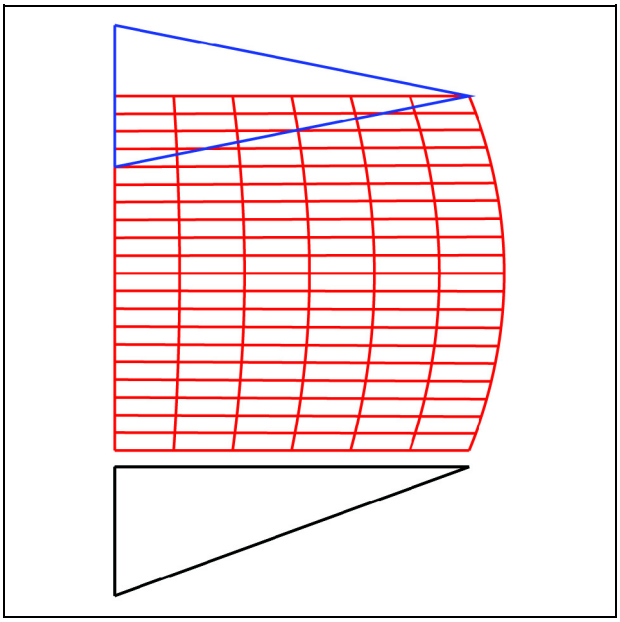

Figure 126 Union of 3 surfaces with errors

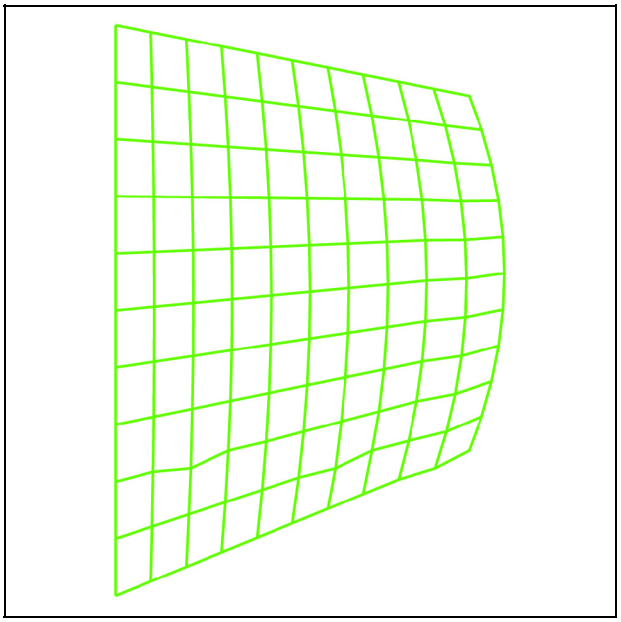

Figure 127 Mesh projected to union of surfaces

Typically, a CAD model will have inaccuracies in the geometry such as gaps between surfaces or surfaces that overlap. This example is an exaggeration of this common problem to demonstrate the robust nature of the **sds** option and the projection method. These two features go hand in hand. In this example, three surfaces form a composite surface. There is a gap between two of the surfaces. There is an overlap between two surfaces. When a face of the mesh is projected to this union of surfaces, the only effect these inaccuracies have on the mesh is that some of the interior nodes are perturbed because they cannot project into the gap. This is usually not a problem because the gaps are usually not that large.

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In this next example, two surfaces form a composite surface. The **sds** option is a natural way to combine tangent surfaces so that you can avoid intersecting tangent surfaces.

```
sd 1 function 0 90 0 360
cos(u)*cos(v);
cos(u)*sin(v);
3*sin(u);;
sd 2 function -90 0 0 360
cos(u)*cos(v);
cos(u)*sin(v);
.5*sin(u);;
sd 3 sds 1 2;
```

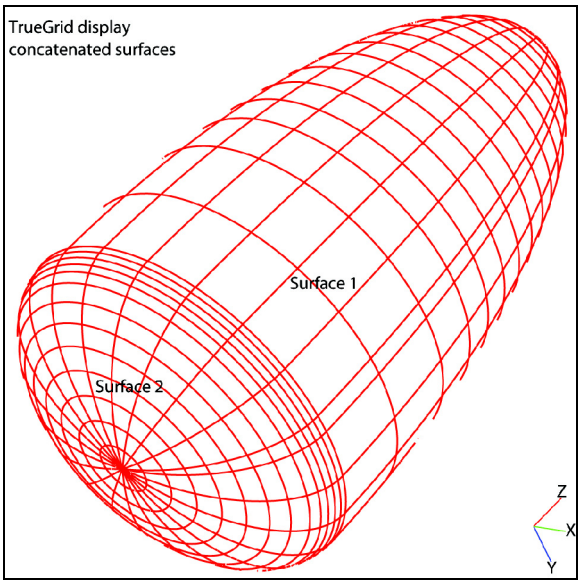

Figure 128 Two surfaces concatenated

sp sphere

sp x_0 y_0 z_0 *radius*

where

x_0	first coordinate of the center of the sphere
y_0	second coordinate of the center of the sphere
z_0	third coordinate of the center of the sphere
radius	radius of the sphere

Remarks

This command defines a sphere with center (x_0,y_0,z_0) and radius *radius*, as shown in the following picture. The radius must be positive. When projecting a point onto the sphere, be sure that the point is away from the center of sphere. Projections are made along rays emanating from the center of the sphere.

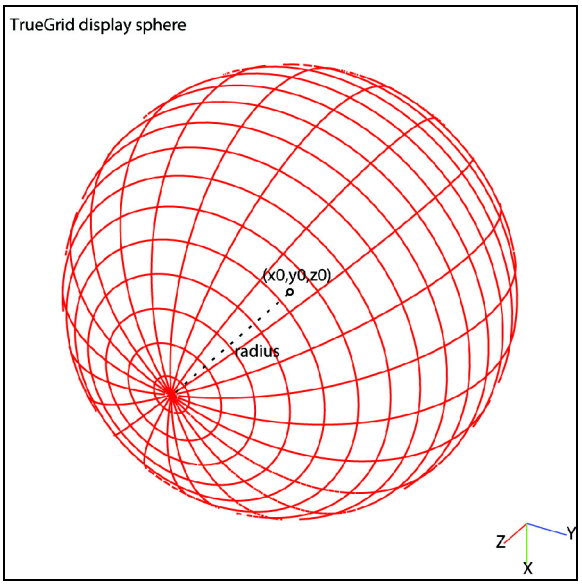

Figure 129 Sphere by center and radius

Example

```
sd 10 sp 1 0 -1 9
```

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stl read the standard ASCII STL file

stl *file trans* ;

where

file	path and file name (case sensitive in UNIX/LINUX)
trans	sequence of simple operators to transform the coordinates

Remarks

This command permits you to import triangular surfaces via StereoLithography (STL) files. Both the ASCII and binary versions of the file are supported.

Use the **fetol** command before this command to extract interior features as edges.

The maximum number of polygons is 8,388,607. When a polygon has more than 4 nodes, it is broken into triangles and each triangle counts as a polygon.

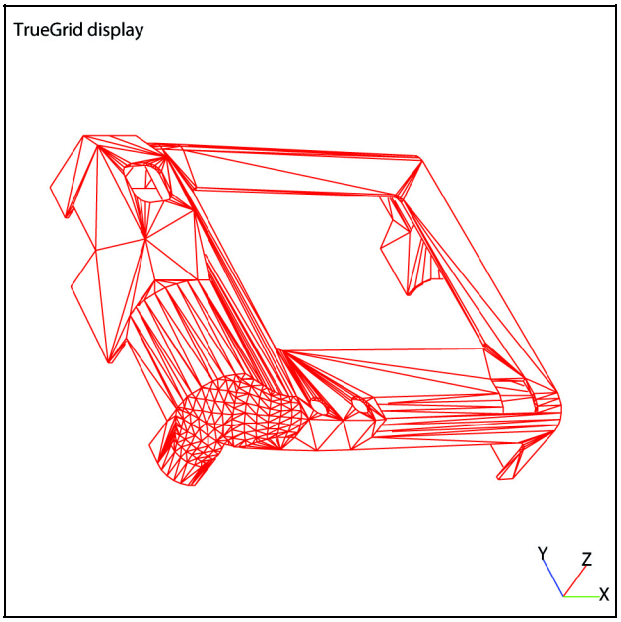

Figure 130 Surface imported from stl file

Example

```
fetol 110
sd 1 stl lcab2b.stl ;
```

swept sweep 2D curves along a 2D curve

swept ln_0 *direction* $ln_1\ \alpha_1 \ldots ln_n\ \alpha_n$ *; trans ;*

where

ln_0	2D curve number
direction	either **r** for the right side or **l** for the left side
$ln_i\ \alpha_i$	pairs of 2D curve numbers and relative arc lengths
trans	sequence of simple operators to transform the coordinates

Remarks

This is a surface which exactly matches various 2D curves at an arbitrary number of cross-sections.

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The 2D curve ln_0 is used as the base of the surface. This 2D curve is on the xz-plane. Positions are measured along this curve by relative arc length. The first point on the base curve has a relative arc length of zero. The last point on the base curve has a relative arc length of 1.0. A cross section of the surface can be specified at any point along the base curve by specifying the relative arc length and the 2D curve which forms the cross section at that point. The 2D curves ln_1, ... ln_n form the cross sections. The parameters α_1 ...α_n are the relative arc lengths of the cross sections, respectively. These arc length parameters must all be between 0 and 1, inclusive. The plane containing the 2D cross section curve will be placed normal to the curve such that the origin in the 2D cross section plane matches the selected point on the base curve. If the *direction* is **r** for right, then the right side of the base curve is the positive direction in the cross section. If *direction* is specified as **l** for left, then the left side of the base curve is the positive direction for the planar cross section. The 2D cross section curves must be specified in order along the base curve. The surface between the cross sections is constructed by linearly interpolating the specified cross sections. The ordering of the points in the 2D cross section curves must match. The arc length of all of the 2D curves must be positive. Care must be taken when using this surface because it is easy to specify cross sections that cause the surface to overlap itself when the base curve has a lot of curvature. All 2D curves must be defined before they are referenced by this command. A transformation can be applied to this surface.

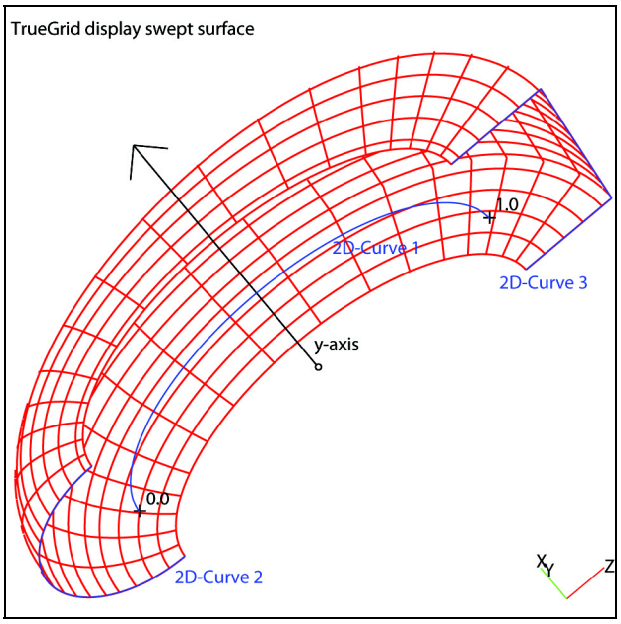

Figure 131 Swept surface with 2 cross sections

Example

In this example the base curve is half of an ellipse. The first cross section is a half of a circle. The last cross section is a half of a cube. The surface transitions smoothly from the smooth arc to 90 degree corners.

```
ld 1 lep 1 2 1 2 -90 90 0;
ld 2 lep 1 1 0 0 -90 90 0;
ld 3 lp2 0 -1 1 -1 1 1 0 1;;
sd 1 swept 1 r 2 0 3 1;;
```

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ts torus

ts x_0 y_0 z_0 x_n y_n z_n r_1 t r_2

where

x_0	first coordinate of a point on the axis of rotation
y_0	second coordinate of a point on the axis of rotation
z_0	third coordinate of a point on the axis of rotation
x_n	first component of axis direction vector
y_n	second component of axis direction vector
z_n	third component of axis direction vector
r_1	larger radius
t	distance from the pont (x_0 y_0 z_0) to the center of the torus
r_2	cross section radius

Remarks

(x_0,y_0,z_0) and (x_n,y_n,z_n) give the axis of symmetry, which passes through the center of the hole in the middle. t is the distance (along the axis) from (x_0,y_0,z_0) to the center point of the torus. The torus has major radius r_1 and minor radius r_2, and both must be positive. A point projected onto this surface should be away from the axis of symmetry and away from the circle at the center of the interior of the torus.

Example

```
sd 4 ts 0 0 0 1 0 1 4 6 1
```

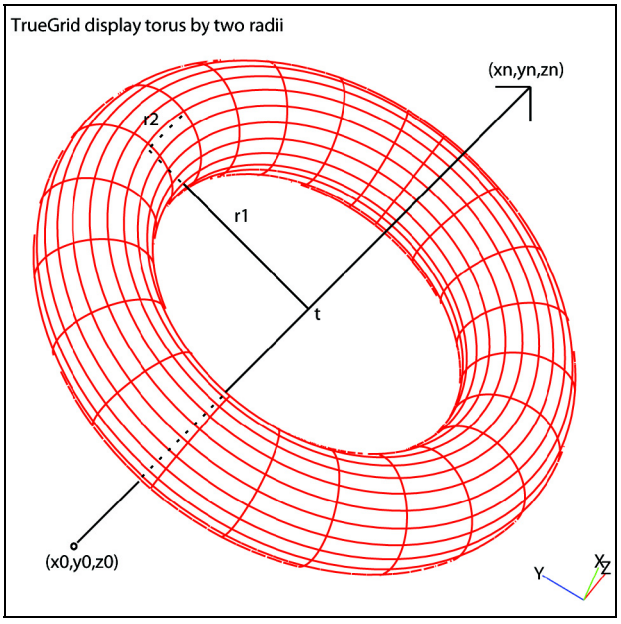

Figure 132 Torus specified by two radii

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xcy to transform an infinite x-axis cylinder

xcy *radius trans* ;

where

radius	is the radius of the cylinder, and
trans	sequence of simple operators to transform the coordinates

Remarks

An infinite cylinder starts out with its axis aligned with the x-axis. The radius must be positive. The transformations make it convenient to start with this cylinder and then translate and rotate it to the required position. Translations are applied to the point which started at the origin. For the syntax of the transformations, see the sections on replication and transformation of parts. When projecting a point onto this surface, be sure that it is initialized somewhere away from the axis of symmetry in order to make it clear which direction to project. See **Figure 133**. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

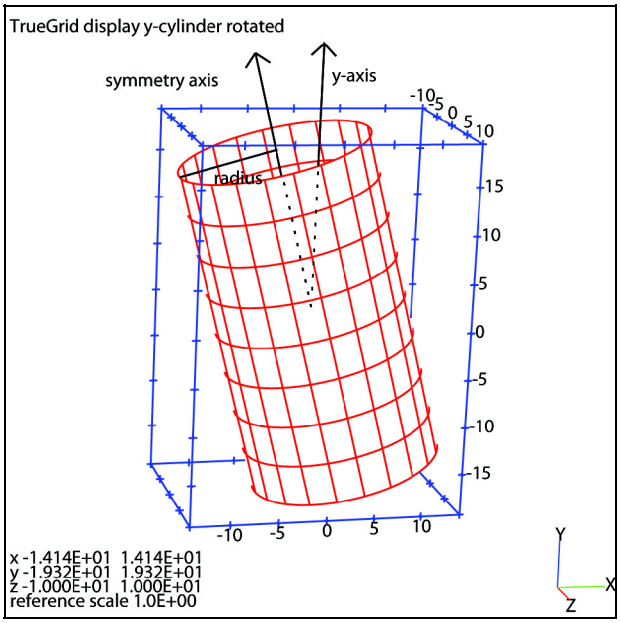

Figure 133 An x-cylinder translated and rotated

Example

```
sd 1 xcy 1 my 1 ry 15 ;
```

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xyplan transform an infinite xy-plane

xyplan *trans* ;

where *trans* is a sequence of simple operators to transform the plane.

Remarks

This is the infinite plane defined by $z = 0$. The transform makes it convenient to start with this plane and then translate or rotate it to the desired position. For the syntax of the transformations, see the sections on replication and transformation of parts. The following example produced the picture below. Translations are applied to the plane at the point which originated at *(0,0,0)*.

This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

Example

```
sd 1 xyplan rx -15;
```

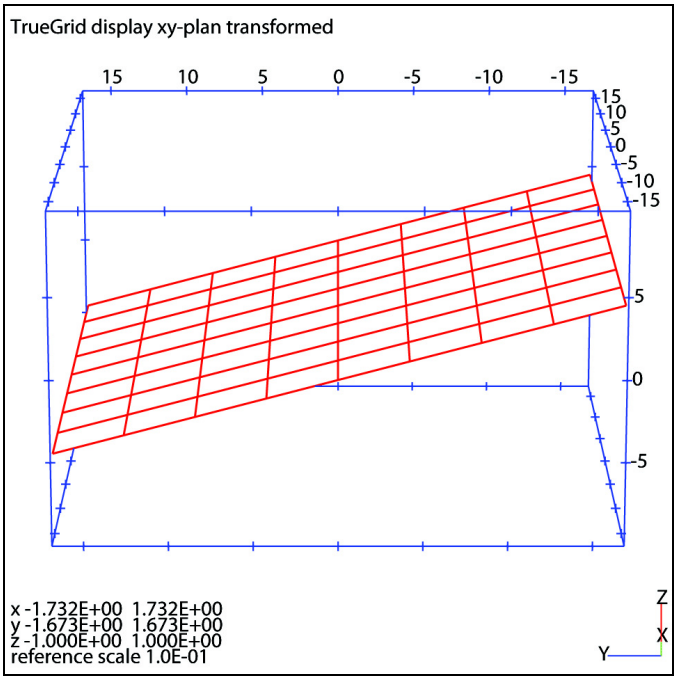

Figure 134 xy-plane rotated

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ycy to transform an infinite y-axis cylinder

ycy *radius trans* ;

where

radius	radius of the cylinder
trans	sequence of simple operators to transform the coordinates

Remarks

An infinite cylinder starts out with an axis the same as the y-axis. The radius must be positive. The transformations make it convenient to start with this cylinder and then translate and rotate it to the required position. Translations are applied to the point which started at the origin. For the syntax of the transformations, see the sections on replication and transformation of parts. When projecting a point onto this surface, be sure that it is initialized somewhere away from the axis of symmetry to make it clear which direction to project. The following example produced the cylinder below. See **Figure 135**. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

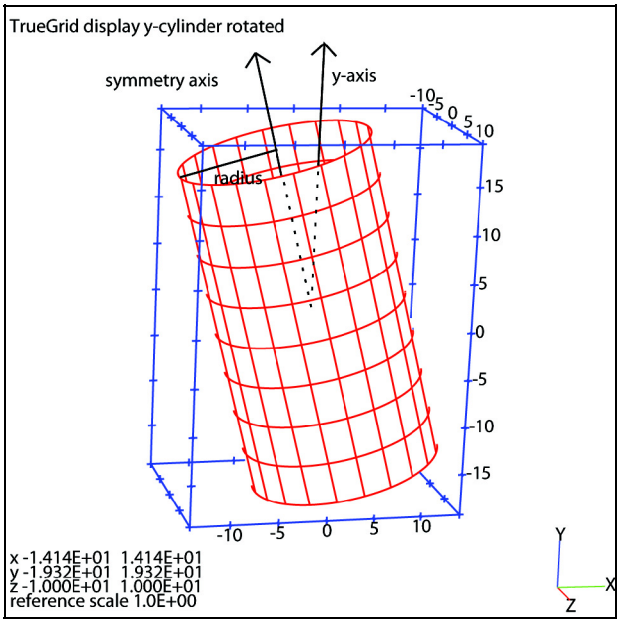

Figure 135 y-cylinder rotated about the z axis

Example

```
sd 1 ycy 10 rz 15;
```

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yzplan to transform an infinite yz-plane

yzplan *trans* ;

where *trans* is a sequence of simple operators to transform the plane.

Remarks

This is the infinite plane defined by $x = 0$. The transform makes it convenient to start with this plane and then translate or rotate it to the desired position. For the syntax of the transformations, see the sections on replication and transformation of parts. The following example produced the picture below. Translations are applied to the plane at the point which originated at *(0,0,0)*.

This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

Example

```
sd 1        yzplan mx 1 rz 45;
```

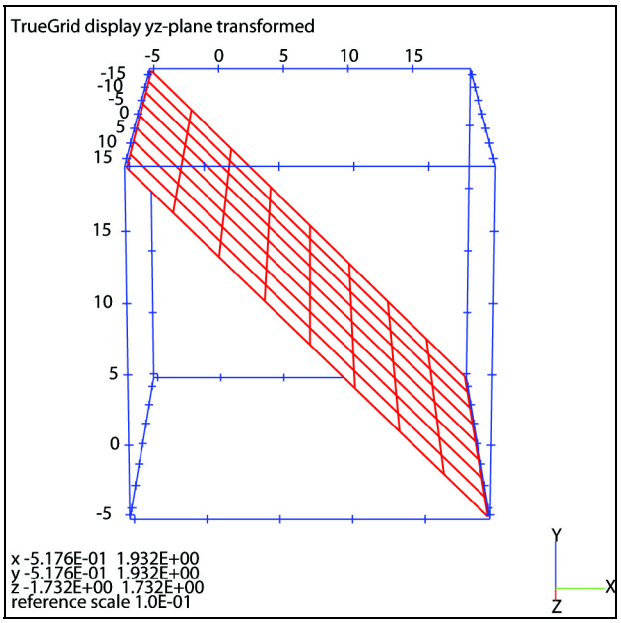

Figure 136 yz-plane moved and rotated

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zcy to transform an infinite z-axis cylinder

zcy *radius trans* ;

where

radius	radius of the cylinder
trans	sequence of simple operators to transform the coordinates

Remarks

An infinite cylinder starts out with an axis the same as the z-axis. The radius must be positive. The transformations make it convenient to start with this cylinder and then translate and rotate it to the required position. Translations are applied to the point which started at the origin. For the syntax of the transformations, see the sections on replication and transformation of parts. When projecting a point onto this surface, be sure that it is initialized somewhere away from the axis of symmetry to make it clear which direction to project. The following example produced the picture below. See **Figure 137**. This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

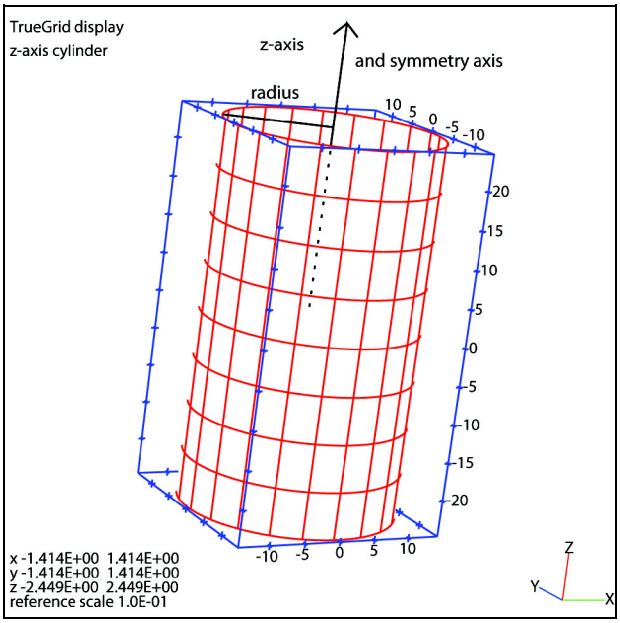

Figure 137 z-cylinder on the z-axis

Example

```
sd 1 zcy [sqrt(2)];
```

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zxplan to transform an infinite zx-plane

zxplan *trans* ;

where *trans* is a sequence of simple operators to transform the plane.

Remarks

This is the infinite plane defined by $y = 0$. The transform makes it convenient to start with this plane and then translate or rotate it to the desired position. For the syntax of the transformations, see the sections on replication and transformation of parts. The following example produced the picture below. Translations are applied to the plane at the point which originates at *(0,0,0)*.

This is an infinite surface. The graphics will only show a portion of the surface. That portion shown in the graphics changes as the objects in the picture change.

Example

```
sd 1 xzplan rx -15 mz -1;
```

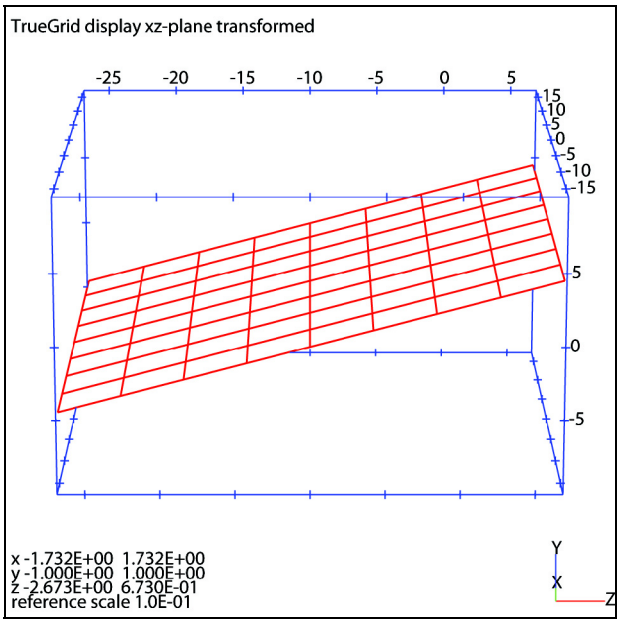

Figure 138 xz-plane rotated and moved

10. Surface Display Commands

These commands let you control which surfaces are displayed in the picture.

When you issue a command to display a surface, you will not necessarily see it. This is because it may lie off-screen. If you issue a surface display command and cannot see the surface, try moving the picture around or zooming in and out. When all else fails, issue a **restore** command (hit the ***Rest*** button).

Many of these commands can be executed using the GUI. In the Environment Window, select the ***Display List*** button. Then choose the ***Surface*** button under the **Apply Action To:** label. The **Action:** buttons will then enable you to reproduce many of the following commands. Some of the display commands in this section are essential. For example, the **asd** command has not been replaced by a button in the GUI. The display commands for many objects, such as 3D curves, parts, and materials, are similar.

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Table I

DISPLAY COMMANDS

Menu	Surface	3D Curve	Parts	Material	Interface	Cad	Cad
Type	Surfaces (sd)	3Dcurves (cd)	Parts (p)	Material (m)	Boundary (bb)	Groups (grp)	Levels (lv)
Display 1 (d∗)	dsd #	dcd #	dp #	dm #	dbb #	dgrp #	dlv #
Add 1 (a∗)	asd #	acd #	ap #	am #	abb #	agrp #	alv #
Remove 1 (r∗)	rsd #	rcd #	rp #	rm #	rbb #	rgrp #	rlv #
Display Many (d∗s)	dsds list;	dcds list ;	dps list ;	dms list ;	dbbs list ;	dgrps list;	dlvs list ;
Add Many (a∗s)	asds list ;	acds list ;	aps list ;	ams list ;	abbs list ;	---	---
Remove Many (r∗s)	rsds list ;	rcds list ;	rps list ;	rms list ;	rbbs list ;	---	---

A sequence in a list can be abbreviated by inserting a colon between the first and last numbers. Surfaces can be named. If a surface has been named, then the name can be used instead of its number.

asd add a surface to the picture

asd *surface_#*

Remarks

Surface *surface_#* is added to the display.

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asds add surfaces to the picture

asds *surface_numbers* ;

Remarks

The surfaces in the list are added to the displayed.

ansd add neighboring surfaces to the picture

ansd *d θ repeats*

where

d	distance tolerance
θ	tangent plane angle tolerance
repeats	number of times to repeat the process of adding surfaces

Remarks

The purpose of this command is to display the surfaces near to the ones already displayed.
All surfaces are searched other than infinite surfaces without edges, such as a cylinder or paraboloid.

All edges of a surface of a candidate surface are checked for points within a distance *d* of an existing displayed surface. For such nearby points it compare's the candidate surface's tangent plane with the existing surface's tangent plane (at the closest point). If the two tangent planes intersect at an angle less than *θ*, then the two surfaces are near each other.

This is repeated so that the next iteration uses the surfaces just added to the display list to find the next set of near surfaces. If any iteration fails to add any surfaces to the display list, the process terminates.

This command is useful in creating composite surfaces. This is the first step in a four step process.

1. select a set of seed surfaces for display
2. use the **ansd** command to find all neighboring surfaces which are nearly tangent
3. use the **lasd** to list the surface numbers
4. cut and paste this list of surfaces into the **sds** option of the **sd** command

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Examples

This figure shows how the angles between surfaces are measured. The normals to surfaces 1 and 4 differ by 9 degrees. Surface 7 is rejected because the angle is too great.

```
ansd 1 10 1
```

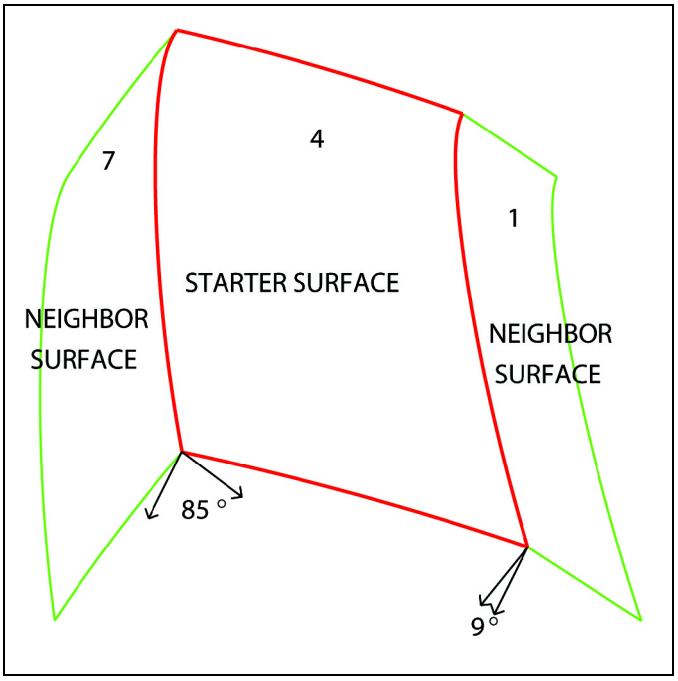

Figure 139 Surface 1 selected adjacent to 4

If you provide a repeat count greater than one, then the search will be repeated. The picture will grow outwards each iteration from the original group of surfaces.

In this example, only surface 1080 was displayed. Then the command

```
ansd .1 3 100
```

was issued to select this set of nearly tangent surfaces.

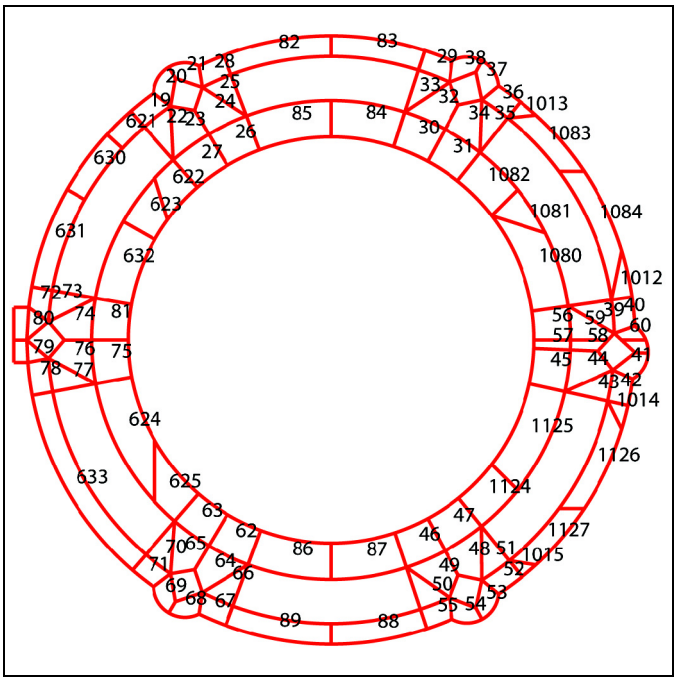

Figure 140 Surface 1080 seeded ANSD

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dasd display all surfaces in the picture

dasd *(no arguments)*

dsd display a surface in the picture

dsd *surface_#*

Remarks

Displays only one surface – removing all others, if necessary.

dsds display several surfaces in the picture

dsds s_1 s_2 ... s_n ;
where s_i is a numbered surface defined by an **sd** command

Remarks

Displays only the surfaces listed.

lasd list the surfaces in the picture

lasd *(no arguments)*

Remarks

This command is useful in creating composite surfaces. This is the second step in a three step process.

1. select a set of seed surfaces for display
2. use the **lasd** to list the surface numbers
3. cut and paste this list of surfaces into the sds option of the **sd** command

rasd remove all surfaces from the picture

rasd *(no arguments)*

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rsd remove a surface from the picture

rsd *surface_#*

11. Importing Geometry from CAD Programs

IGES is a standard format for CAD geometry and can be used to pass geometry from a CAD program to **True*Grid*®**. It is advantageous to use IGES geometry because most CAD systems are better than **True*Grid*®** at creating complex geometry. Also, you may already have an IGES model and there is no need to reproduce the effort in creating the geometry in **True*Grid*®**. IGES geometry has the disadvantage that it may be overly complex and reducing the geometry to that which is needed for meshing is more difficult than building the geometry within **True*Grid*®**. This point is mitigated since you have the option, in **True*Grid*®**, to ignore certain features by not projecting a portion of the mesh to certain surfaces. Also, you may not have easy access to a CAD model or you may wish not to learn how to use an available CAD system. Also significant to experienced **True*Grid*®** users is that CAD geometry is not easily made parametric while every aspect of geometry built in **True*Grid*®** can be parametric.

The projection method applied to IGES geometry is superior to other mesh generation techniques in two significant ways. IGES geometry is usually broken into many small components that do not meet perfectly. These small gaps between the geometric components are ignored by the projection method. Also, many meshing techniques may fail or do poorly on some geometry, requiring that the CAD system be reused to break the geometry into smaller pieces with the hope that the smaller pieces will be easily meshed. If successful, these other meshing techniques then must merge the individual parts together. In contrast, the projection method encourages building composite surfaces, done very simply using the **sds** option of the **sd** command for surfaces and the **coedge** command for edges of surfaces, to simplify the geometry and the block structure of the mesh. The mesh generation methods used in **True*Grid*®** have never been shown to fail on any geometry.

There are numerous options when creating an IGES file in a CAD system since there is more than one way to define geometry. In our experience, most CAD systems default to creating an IGES file consisting of solids. This is not acceptable to **True*Grid*®**. **True*Grid*®**'s methods were developed for trimmed surfaces and curves, not solids. Also, in our experience, most CAD operators who create the IGES files are not aware of the options for writing an IGES file. In most cases you need to insist that the IGES file be produced using trimmed surfaces, not solids.

You can import this IGES geometry created with a CAD program using the **iges** command. This interface is a little different from the interfaces of most CAD systems in that the **True*Grid*®** internal

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data base is complex and may require significant computations. For this reason, you can save this data base, once generated, in a binary IGES file. In the first instance, you use the **iges** command to import the geometry followed by the **saveiges** command to save the binary IGES data base. Subsequently, you use the **useiges** command to import the binary IGES data base prior to using the same **iges** command that you used in the first instance. If you take these extra steps, you may save yourself considerable time.

First usage:

```
iges file.igs 1 1;
saveiges file.bin
```

All subsequent usage:

```
useiges file.bin
iges file.igs 1 1;
```

Although IGES is an international standard, there is no enforcement of the standard and no quality assurance. CAD systems, on occasions, produce faulty IGES files because of bugs in the CAD system, deviations from the IGES standard, and mathematical abuses. Our policy is to develop algorithms that work around these problems produced by the CAD systems. Since the CAD systems are constantly evolving, **True*Grid*®** must be constantly evolving as well. So we ask that if you find a problem when importing an IGES file, please contact XYZ Scientific applications, Inc.. The best way to help is to send the IGES file. This may be difficult because the IGES file may be too large or it may be proprietary or classified. Alternatively, you can use the WINDOWS® IGES Inspection Utility (**igesfind.exe**) or the UNIX®/LINUX® **igesfind** program supplied by XYZ Scientific Applications, Inc. to extract a specific surface from an IGES file. Then send just that one surface to XYZ Scientific Applications, Inc. by email to info@truegrid.com. For example, suppose that surface 123 of an IGES file called widget.igs causes a problem in **True*Grid*®**. With the IGES Inspection Utility on WINDOWS®, select that numbered surface and click on *save as* to create a new IGES file with just the surface numbered 123. On a UNIX®/LINUX® system type

igesfind widget.igs -surface 123 > s123.igs

to produce an IGES file s123.igs with only the surface numbered 123.

Geometric entities in an IGES file can depend on other entities, forming a hierarchy. When the geometric entities are numbered, room is left in the sequence for those entities which form the basis for others. If you use the **iges** command, then entities which form the basis for other entities are not rendered. For example, the curves used to trim a surface are not evaluated since the entity will be embedded in the trimmed surface as one of its edges. You can use one of the other IGES evaluation

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commands discussed below if you wish to use these basis entities. There are two important entities which fall in this category. They are composite curves (entity type number 102) and trimmed surfaces (entity type number 144). You will notice, when importing an IGES file using the **iges** command, that the underlying composite and component curves as well as the untrimmed surface are all assigned numbers, but they are not rendered. There is no loss of data, since all of the information about these component entities are built into the trimmed version of the surface.

The following is a list of the supported IGES entity types. The bounded surface entity (143) and its associated bounded entity (type 141), which are used to form solids, are not supported.

3D curves	line	entity type 108
	circular arc	entity type 100
	conic arc	entity type 104
	parametric spline	entity type 112
	NURBS curve	entity type 126
	composite	entity type 102
	copious data	entity type 106
Surfaces	plane including bounded plane	entity type 108
	ruled surface	entity type 118
	surface of revolution	entity type 120
	tabulated cylinder	entity type 122
	parametric spline	entity type 114
	NURBS surface	entity type 128
	trimmed surface	entity types 144 & 142
	offset surface	entity type 140
Others	transformation	entity type 124
	level	N/A
	associativity	entity type 302
	subfigure	entity type 308

NURBS (Non-Uniform Rational B-Spline) surfaces are a type of surface that has become, in most CAD systems, the preferred type of surface because of their versatility. Planar surfaces can also be found in an IGES file. There are other types of surfaces such as ruled, parametric spline, revolved, and tabulated. These three basic types of surfaces, NURBS, planes, and others, are each assigned an IGES sequence number, according to their type, so that they can be referenced for rendering purposes. These IGES sequence numbers are different from the number you assign a surface. For example, the first NURBS IGES surface may be assigned to be surface number 2 in **True*Grid*®** because you may already have a surface number 1. The untrimmed NURBS surfaces are assigned sequence numbers first, followed by the trimmed versions of these surfaces. As is noted above, the untrimmed version of a NURBS surface will not be automatically rendered if there is also a trimmed version of that surface. In the same manner, planes and other surface types are sequenced and rendered. Then IGES

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curves are numbered and rendered in the order that they are found in the IGES file. Although an IGES curve may be 2D or 3D, they are all treated as 3D curves, with 2D curves being assigned a third constant coordinate. Since most 2D curves are usually used to define a trimmed surface, they are not automatically rendered.

When there are multiple IGES files being used, the IGES sequence numbering of NURBS, planes, other surfaces, and curves start where they left off on from the previous IGES surface.

You may have difficulty reading an IGES file created on a different machine. This is because WINTEL systems use a different end-of-line and carriage return from UNIX/LINUX systems. One way to solve this problem is to use a text editor to remove the extra character at the end of the line and rewrite it. There may also be utilities on your system to do this conversion. If you use ftp to transfer the file, be sure to leave it in ASCII or character mode (not binary).

You can display rendered IGES curves with commands such as **dcd**, **dacd**, **acd**, **acds**, **rcd**, **rdds**, and **dcds**. You can display rendered IGES surfaces with commands such as **dsd**, **dasd**, **asd**, **asds**, **rsd**, **rsds**, **dsds**. You can place the edge of the mesh along a rendered IGES curve with the **curs** command. You can place a face of the mesh on a rendered IGES surface with the **sf** command.

The default accuracy in rendering the geometry is approximately 3.5 digits of accuracy. This rendering tolerance is relative to the size of the geometric entity being rendered. The accuracy of the rendering affects the time it takes to do the rendering and the amount of memory needed to store the rendering data. This also affects the size of an IGES binary file written using the **saveiges** command. It also affects the accuracy of the default projection to surfaces. Both the accuracy of the rendering and the accuracy of the projection can be controlled, independently, using the **getol** and **accuracy** commands, respectively. In particular, you may wish to decrease the accuracy of the rendering to speed the rendering and decrease the memory requirements while increasing the projection accuracy. This is a classic example of the tradeoff of memory and computations.

You can also read in geometry which has been stored in a simple polygon data format using the **vpsd** command, the **stlsd** command, or the **stl** or **bstl** options of the **sd** command. These types of surfaces are not affected by the **getol** command because the polygons dictate the accuracy.

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iges render geometry in an IGES file

iges *file surface_# curve_# transformations ;*

where

file IGES file name including the path relative to the working directory

surface_# surface number to be assigned to the first surface in the file

curve_# 3D curve number to be assigned to the first 3D curve in the file

transformations list of coordinate transformations to be applied to all geometry

where a transform can be

mx *x_offset* for translation in the x direction

my *y_offset* for translation in the y direction

mz *z_offset* for translation in the z direction

v *x_offset y_offset z_offset* for a general translation

scv *xn yn zn offset* to translate a given distance along a vector

dv x_1 y_1 z_1 x_2 y_2 z_2 to translate by a vector given by two points

rx θ for rotation about the x axis

ry θ for rotation about the y axis

rz θ for rotation about the z axis

raxis *angle x0 y0 z0 xn yn zn* axis of rotation

rxy for reflection about the x-y plane

ryx for reflection about the y-x plane

rzx for reflection about the z-x plane

tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* Cartesian coordinates

cy *rho theta z* cylindrical coordinates

sp *rho theta phi* spherical coordinates

pt *c.i* label of a labeled point from a 3D curve

pt *s.i.j* label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* Cartesian coordinates

cy *rho theta z* cylindrical coordinates

sp *rho theta phi* spherical coordinates

pt *c.i* label of a labeled point from a 3D curve

pt *s.i.j* label of a labeled point from a surface

inv invert the transformation preceding this option

csca *scale_factor* to scale all coordinates

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xsca *scale_factor* to scale the x coordinate
ysca *scale_factor* to scale the y coordinate
zsca *scale_factor* to scale the z coordinate

Remarks

It is inefficient to start with a high number for the curves or surfaces.

The number of NURBS, planes, other surface, and curves found in an IGES file is reported in the text window and the tsave file.

This command is the simplest way to import IGES geometry. You may use the **igesfile** to be selective on which entities to render. Although this may be faster, because you only render the geometry you want, you must have already decided what curves and surfaces you want. This can only be decided by using the **iges** command first so that you can see the surfaces and curves to make your choices.

To save the rendering calculations for future re-use, use the **saveiges** command. The next time you want the surfaces and 3D curves from this IGES file, issue the **useiges** command before issuing the **iges** or other IGES rendering command.

In the WINDOWS operating system, when you click on the iges command in the menus, you will get a browser that will aid you in selecting the IGES file.

igesfile open an IGES file

igesfile *filename*

Remarks

When you enter an **igesfile** command, the IGES file is opened. The geometry in the IGES file is reported, but not rendered. To render curves, use the **curd** command using the **igc** option, or with the **igescd** command. To render a NURBS surface, use the **sd** command using the **nurbs** option, or with the **nurbsd** command. For a plane, evaluate it with the **sd** command using the **igesp** option, or with the **igespd** command. For any other type of IGES surface, evaluate it with the **sd** command using the **igess** option, or with the **igessd** command.

This command should only be used by an experienced user. It requires prior knowledge of the geometry in the specified file. This command has the advantage, over the iges command, of selecting out a subset of the geometric entities for rendering. It is advantageous only when the subset is relatively small.

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The number of NURBS, planes, other surface, and curves found in an IGES file is reported in the text window and the tsave file.

igescd render a sequence of IGES curves

igescd *curve$_1$ curve$_n$ first_curve transformations ;*

where

curve$_1$ is the sequence number in the IGES file of the first curve to be evaluated

curve$_n$ is the sequence number in the IGES file of the last curve to be evaluated

first_curve is the **True*Grid*®** curve number to be assigned to the first IGES curve

transformations is a list of any of the following coordinate transformations:

mx *x* for x translation

my *y* for y translation

mz *z* for z translation

v *x y z* for a general translation

scv *xn yn zn offset* to translate a given distance along a vector

dv $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$ to translate by a vector given by two points

rx θ for rotation about the x-axis

ry θ for rotation about the y-axis

rz θ for rotation about the z-axis

raxis *angle x0 y0 z0 xn yn zn* axis of rotation

rxy for reflection about the x-y plane

ryz for reflection about the y-z plane

rzx for reflection about the z-x plane

tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve

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pt *s.i.j*	label of a labeled point from a surface
inv	invert the present transformation
csca *scale_factor*	to scale all coordinates
xsca *scale_factor*	to scale the x coordinate
ysca *scale_factor*	to scale the y coordinate
zsca *scale_factor*	to scale the z coordinate

Remarks

First specify the IGES file with the **igesfile** or the **iges** command.

It is inefficient to start with a high number for the curves.

A unique **True*Grid*®** 3D curve number will be assign to each rendered IGES curve, starting with the specified number. All specified curves will be rendered, even those used in the construction or rendering of more complex geometric entities.

igeslbls use IGES surface labels to name surfaces

igeslbls *flag*
where *flag* can be **on** or **off**

Remarks

The default is **off**. Sometimes, an IGES file includes labels for the surfaces. When several surfaces are found to have the same label, they are treated as a composite surface. The result would be the same if the **sd** command were issued using the **sds** option to combine the several surfaces into one. Surface numbers are assigned for internal bookkeeping. It will grab the next available surface number to do this. You will be notified for each composite surface. Names must have some alphabetic characters in them and no spaces. Other special characters like : and ; should also be avoided, although the implications of using none alphanumeric characters has not been explored. This feature is case insensitive. Only the first 8 characters of a name are used.

One can add surfaces to a composite by simply naming it in the **sd** command when the surface is defined.

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igespd render a sequence of IGES planes

igespd *plane$_1$ plane$_n$ first_surface transformations*

where

plane$_1$ sequence number in the IGES file of the first plane to be evaluated

plane$_n$ sequence number of the last plane to be evaluated

first_surface **True*Grid*®** surface number to be assigned to the first plane

transformations is a list of any of the following coordinate transformations:

mx *x* for x translation

my *y* for y translation

mz *z* for z translation

v *x y z* for a general translation

scv *xn yn zn offset* to translate a given distance along a vector

dv $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$ to translate by a vector given by two points

rx θ for rotation about the x-axis

ry θ for rotation about the y-axis

rz θ for rotation about the z-axis

raxis *angle x0 y0 z0 xn yn zn* axis of rotation

rxy for reflection about the x-y plane

ryz for reflection about the y-z plane

rzx for reflection about the z-x plane

tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface

inv invert the present transformation

csca *scale_factor* to scale all coordinates

xsca *scale_factor* to scale the x coordinate

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ysca *scale_factor* to scale the y coordinate
zsca *scale_factor* to scale the z coordinate

Remarks

First specify the IGES file with the **igesfile** or **iges** command.

It is inefficient to start with a high number for the surfaces.

A unique **True*Grid*®** surface number will be assign to each rendered IGES plane, beginning with the specified number.Everyl requested plane will be rendered, even if it is untrimmed and forms the basis for a trimmed surface.

igessd render a sequence of IGES surfaces

igessd *surface$_1$ surface$_n$ first_surface transformations*

where

surface$_1$ is the first IGES surface number to be selected
surface$_n$ is the last IGES surface number to be selected
first_surface is the first **True*Grid*®** surface definition number
transformations is a list of any of the following coordinate transformations:

mx *x* for x translation
my *y* for y translation
mz *z* for z translation
v *x y z* for a general translation
scv *xn yn zn offset* to translate a given distance along a vector
dv *x_1 y_1 z_1 x_2 y_2 z_2* to translate by a vector given by two points
rx θ for rotation about the x-axis
ry θ for rotation about the y-axis
rz θ for rotation about the z-axis
raxis *angle x0 y0 z0 xn yn zn* axis of rotation
rxy for reflection about the x-y plane
ryz for reflection about the y-z plane
rzx for reflection about the z-x plane
tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates

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sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface

inv invert the present transformation
csca *scale_factor* to scale all coordinates
xsca *scale_factor* to scale the x coordinate
ysca *scale_factor* to scale the y coordinate
zsca *scale_factor* to scale the z coordinate

Remarks

First specify the IGES file with the **igesfile** or **iges** command.

It is inefficient to start with a high number for the surfaces..

A unique **True*Grid*®** surface number will be assign to each rendered IGES surface, beginning with the specified number. Each specified surface will be rendered, even if it is untrimmed and forms the basis for a trimmed surface.

To evaluate NURBS surfaces, use the **nurbsd** command. In order to evaluate planes, use the **igespd** command. This command is for all other IGES surfaces that **True*Grid*®** supports.

nurbsd render a sequence of IGES NURBS surfaces

nurbsd $surface_1$ $surface_n$ *first_surface transformations ;*
where
$surface_1$ is the IGES file sequence number of the first NURBS surface used
$surface_n$ is the IGES file sequence number of the last NURBS surface used
first_surface is the first **True*Grid*®** surface definition number
transformations is an optional list of transformations to be applied to the surfaces:
mx *x* for x translation

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my *y* for y translation
mz *z* for z translation
v *x y z* for a general translation
scv *xn yn zn offset* to translate a given distance along a vector
dv x_1 y_1 z_1 x_2 y_2 z_2 to translate by a vector given by two points
rx θ for rotation about the x-axis
ry θ for rotation about the y-axis
rz θ for rotation about the z-axis
raxis *angle x0 y0 z0 xn yn zn* axis of rotation
rxy for reflection about the x-y plane
ryz for reflection about the y-z plane
rzx for reflection about the z-x plane
tf *origin x-axis y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface

inv invert the present transformation
csca *scale_factor* to scale all coordinates
xsca *scale_factor* to scale the x coordinate
ysca *scale_factor* to scale the y coordinate
zsca *scale_factor* to scale the z coordinate

Remarks

This command applies to the IGES file you selected with the **igesfile** or **iges** command.

It is inefficient to start with a high number for the surfaces.

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A unique **True*Grid*®** surface will be assigned to each rendered IGES NURBS surface, starting with the specified number. Each specified surface will be rendered, even if it is untrimmed and forms the basis for a trimmed surface.

In order to evaluate planes, use the **igespd** command. Use the **igessd** command to render other IGES surface types.

saveiges save IGES binary data

saveiges *file*

where *file* is the file name to contain the binary IGES surface and 3D curve data

Remarks

If your IGES file has a large number of surfaces and 3D curves that you need to reuse, it may be time-consuming for **True*Grid*®** to render them each time you rerun your batch file. With this command you can save the evaluations so that in the future you can quickly access the surfaces and curves. This command is used after the IGES geometry has been rendered. All IGES geometric entities (surfaces and 3D curves) that have been rendered will be stored in the binary data file. Each surface or 3D curve from an IGES file is assigned a key in this binary data file. When this binary data is retrieved using the **useiges** prior to either the **iges** or **igesfile** command, the surfaces and curves in the binary data file will be matched to the corresponding geometric entities from the appropriate IGES files. The name of the IGES file is used to form this key for each surface and curve and it is required that the IGES file name not be changed so that the correct correspondence between IGES entities and the binary data can be established.

The binary data file that is formed by this command contains the detailed information resulting from the rendering process. The **getol** command controls this rendering process. When a binary file is used (see the **useiges** command), the IGES surfaces and curves inherit the results of the rendering process at the time the binary file was written. If you want to change the rendering, such as changing the **getol** parameter, you cannot use the previously saved IGES binary file.

trimming controls the trimmed surface algorithm

trimming *option*

where the *option* can be

on	this activates the trimming
off	this deactivates the trimming

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Remarks

This feature is not usually needed. It is used for backward compatibility and for debugging. The default is **on**.

ltrim set the surface trimming work space

ltrim *size*

where *size* is the number of words to use in the trimming algorithm

Remarks

This command is not usually needed. The default is 2000000 words. Only use this command to increase the work space if the code issues a warning that it had difficulties trimming a surface and perhaps increasing the work space may help.

```
warning - not enough work space to trim surface.
          try increasing the work space from 2000000
          by using the ltrim command.
```

This can occur if there are some small details in the trimming curves or if there are many interior trimming curves. **True*Grid*®** tries to trim the surface by first using the domain curves. Sometimes these trimming curves are not supplied in the IGES file. If they are not supplied or if they fail, then **True*Grid*®** attempts to trim the surface with the model space curves.

You can also try reducing the resolution of the trimming using the **getol** command.

Often, when this problem occurs, the IGES file includes a poorly defined trimmed surface. For instance, the surface component trimming curves do not meet at their respective end points or the trimming curve has a loop causing the curve to intersect itself. These problems violate the IGES standard and sometimes occur in IGES files. Most of these problems with trimming curves are automatically fixed by **True*Grid*®** and you are never warned about it.

The best way to determine if the trimming curves are bad is to use the **igescd** to render the trimming curves. Then inspect the curves for these flaws. If you have a large IGES file, it will be easier if you first extract the offending surface using the IGESFIND utility that is supplied with each **True*Grid*®** distribution.

If the trimming curves are bad, perhaps you can modify the IGES file using the CAD system which created the trimmed surfaces originally. If it appears that there is no problem with the trimming

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curves, and you have tried increasing the size of the work space with no results, please contact XYZ.

useiges use saved binary IGES data

useiges *file*

where *file* is the name of the binary IGES file made by the **saveiges** command.

Remarks

If your IGES file has a large number of surfaces and 3D curves that you need to use, it may be time-consuming for **True*Grid*®** to evaluate them all. You can save the evaluations with **saveiges** and then quickly read them back in with this command.

This command is used before the IGES file is opened using either **iges** or **igesfile**. During a previous run with **True*Grid*®**, any IGES geometric entities (surfaces and curves) that were rendered (were visible) and were stored in the binary data file will be retrieved. Each surface or curve from an IGES file was assigned a key in this binary data file. When this binary data is retrieved using the **useiges** followed by either the **iges** or **igesfile** command, the surfaces and curves in the binary data file are matched to the corresponding geometric entities from the appropriate IGES files. The name of the IGES file is used to form this key for each surface and curve and it is required that the IGES file name not be changed so that the correct correspondence between the IGES entities and the binary data can be established. Up to 100 binary files can be read in with this command.

stlsd import StereoLithography surfaces

stlsd *first_surface stl_file transformations ;*

where

first_surface	is the **True*Grid*®** surface number to be assigned to the first surface read from the file,
stl_file	is the StereoLithography file,
transformations	is a list of coordinate transformations to be applied to all the curves and surfaces. Choose them from among:

mx *x_offset*	for translation in the x direction
my *y_offset*	for translation in the y direction
mz *z_offset*	for translation in the z direction
v *x_offset y_offset z_offset*	for a general translation along a vector
scv *xn yn zn offset*	to translate a given distance along a vector
dv $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$	to translate by a vector given by two points

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rx θ for rotation about the x axis
ry θ for rotation about the y axis
rz θ for rotation about the z axis
raxis *angle x0 y0 z0 xn yn zn* axis of rotation
rxy for reflection about the x-y plane
ryx for reflection about the y-x plane
rzx for reflection about the z-x plane
tf *origin x-axis y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface
ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type and coordinate data:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface
inv invert the present transformation
csca *scale_factor* to scale all coordinates
xsca *scale_factor* to scale the x coordinate
ysca *scale_factor* to scale the y coordinate
zsca *scale_factor* to scale the z coordinate

Remarks

This command is ideal for importing geometry from a biological source. Sometimes these polygon surfaces are not ideal. To split a polygon surface into several polygon surfaces, choose the Sets button in the Pick panel in the Environment window, form a Polygon set, and then use the **sd** command to convert the polygon set to a surface. There are three ways to move a surface polygon point. In the first method, move the mouse close to the point to be moved and click the **F5** function key. Then select a new geometric point, such as Pick>Z-Buffer in the Environment window and click on the Attach button in the Environment window. Alternatively, use the **mvpn** or **pvpn** command to move or change the polygon point. In the final method, after selecting a surface polygon point with the **F5** key, select the Poly Surface button on the left side of the Move Pts. Panel in the Environment window and use the options on the right side to interactively move the surface polygon point by clicking the left mouse button and dragging the mouse.

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If you make changes to a polygon surface, you may wish to save the new surface(s) instead of relying on the saved session file to be rerun every time you want to use the surface(s). You can save them as a ViewPoint surface using the **wrsd** command. Use the **vpsd** command to import the modified surface(s).

An edge of a polygon is considered a polygon boundary edge if that edge is not shared by another polygon. These polygon boundary edges are sewn together to form surface boundary edges. A corner is identified when two polygon boundary edges from the same polygon meet. Most surface edges start and end at a corner. If a surface boundary edge is closed (i.e. the start and end points are the same), it will contain no corners. It is also possible that a surface will have no boundary edges.

An edge of a polygon is considered a polygon interior feature if that edge is shared by another polygon and if the angle between the two polygons is less than the feature tolerance parameter (use the **fetol** command to set this parameter). These polygon feature edges are sewn together to form surface interior edges in the same way that surface boundary edges are formed.

Both types of surface edges can be labeled using the **la** command or by selecting labels from the Environment window. Points can be selected from these edges by using the Projection button in the Environment window. Edges of the mesh can be attached to one of these surface edges using the **edge** command or by using the Attach button in the Environment window. These edges can be combine into composite curves using the **se** option of the **curd** command or by using the interactive **coedge** button under the 3D Curves menu.

Surface points that are (nearly) coincident will produce artificial boundary edges. This can be avoided by using the command **smgap**. When **smgap** is used, (nearly) coincident points are merged.

vpsd import ViewPoint surfaces

vpsd *first_surface nodal_file connectivity_file transformations ;*

where

first_surface	is the **True*Grid*®** surface number to be assigned to the first surface read from the file,
nodal_file	is the ViewPoint nodal file,
connectivity_file	is the ViewPoint connectivity file, and
transformations	is a list of coordinate transformations to be applied to all the curves and surfaces. Choose them from among:

mx *x_offset*	for translation in the x direction
my *y_offset*	for translation in the y direction
mz *z_offset*	for translation in the z direction
v *x_offset y_offset z_offset*	for a general translation along a vector

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scv *xn yn zn offset* to translate a given distance along a vector
dv x_1 y_1 z_1 x_2 y_2 z_2 to translate by a vector given by two points
rx θ for rotation about the x axis
ry θ for rotation about the y axis
rz θ for rotation about the z axis
raxis *angle x0 y0 z0 xn yn zn* axis of rotation
rxy for reflection about the x-y plane
ryx for reflection about the y-x plane
rzx for reflection about the z-x plane
tf *origin x-axis y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface
ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type and coordinate data:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface
inv invert the present transformation
csca *scale_factor* to scale all coordinates
xsca *scale_factor* to scale the x coordinate
ysca *scale_factor* to scale the y coordinate
zsca *scale_factor* to scale the z coordinate

Remarks

If you need to write your own software to generate polygonal surfaces, you can write your output as ViewPoint files. These file formats are a good choice because of their simplicity. The first file contains the point data for the polygons. Each record is a point in three-dimensional space. There are 4 fields, each field separated by a comma. The first field is the node number, followed by the x, y, and z-coordinates. The second file contains the connectivity of the nodes. Each record defines a polygon. The fields are space delimited. The first field is the name of the set containing this polygon. This is followed by the ordered list of node numbers defining the polygon.

The polygons in these surfaces should be almost planar. When they are read in, they are reduced to

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triangles, unless they are quadrilaterals. The shape of the interior of a many sided polygon will be ambiguous, if the nodes are not co-planar.

Each unique name in the connectivity file produces a surface in **True*Grid*®**. The polygons do not have to be connected.

This command is ideal for importing geometry from a biological source. Sometimes these polygon surfaces are not ideal. To split a polygon surface into several polygon surfaces, choose the Sets button in the Pick panel in the Environment window, form a Polygon set, and then use the **sd** command to convert the polygon set to a surface. There are three ways to move a surface polygon point. In the first method, move the mouse close to the point to be moved and click the F5 function key. Then select a new geometric point, such as Pick>Z-Buffer in the Environment window and click on the Attach button in the Environment window. Alternatively, use the **mvpn** or **pvpn** command to move or change the polygon point. In the final method, after selecting a surface polygon point with the F5 key, select the Poly Surface button on the left side of the Move Pts. Panel in the Environment window and use the options on the right side to interactively move the surface polygon point by clicking the left mouse button and dragging the mouse.

If you make changes to a polygon surface, you may wish to save the new surface(s) instead of relying on the saved session file to be rerun every time you want to use the surface(s). You can save them as a ViewPoint surface using the **wrsd** command. Use the **vpsd** command to import the modified surface(s).

An edge of a polygon is considered a polygon boundary edge if that edge is not shared by another polygon. These polygon boundary edges are sewn together to form surface boundary edges. A corner is identified when two polygon boundary edges from the same polygon meet. Most surface edges start and end at a corner. If a surface boundary edge is closed (i.e. the start and end points are the same), it will contain no corners. It is also possible that a surface will have no boundary edges.

An edge of a polygon is considered a polygon boundary edge if that edge is not shared by another polygon. These polygon boundary edges are sewn together to form surface boundary edges. A corner is identified when two polygon boundary edges from the same polygon meet. Most surface edges start and end at a corner. If a surface boundary edge is closed (i.e. the start and end points are the same), it will contain no corners. It is also possible that a surface will have no boundary edges.

An edge of a polygon is considered a polygon interior feature if that edge is shared by another polygon and if the angle between the two polygons is less than the feature tolerance parameter (use the **fetol** command to set this parameter). These polygon feature edges are sewn together to form surface interior edges in the same way that surface boundary edges are formed.

Both types of surface edges can be labeled using the **la** command or by selecting labels from the

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Environment window. Points can be selected from these edges by using the Projection button in the Environment window. Edges of the mesh can be attached to one of these surface edges using the **edge** command or by using the Attach button in the Environment window. These edges can be combine into composite curves using the **se** option of the **curd** command or by using the interactive **coedge** button under the 3D Curves menu.

Surface points that are (nearly) coincident will produce artificial boundary edges. This can be avoided by using the command **smgap**. When **smgap** is used, (nearly) coincident points are merged.

Example

The following example creates 2 surfaces, **inner** and **outer**. The nodal and connectivity files are listed below. These files were read with the command

```
vpsd  1  vpsd.nodes vpsd.elem ;
```

See below for a picture of the resulting surfaces.

Nodal file: vpsd.nodes

```
1,0.87,0.50,3.00
2,0.00,1.00,3.00
3,-0.87,0.50,3.00
4,-0.87,-0.50,3.00
5,0.00,-1.00,3.00
6,0.87,-0.50,3.00
7,1.30,0.75,2.75
8,0.00,1.50,2.75
9,-1.30,0.75,2.75
10,-1.30,-0.75,2.75
11,0.00,-1.50,2.75
12,1.30,-0.75,2.75
13,1.97,-0.35,2.5
14,1.97,0.35,2.5
15,1.29,1.53,2.5
16,0.68,1.88,2.5
17,-0.68,1.88,2.5
18,-1.29,1.53,2.5
19,-1.97,0.35,2.5
20,-1.97,-0.35,2.5
21,-1.29,-1.53,2.5
22,-0.68,-1.88,2.5
```

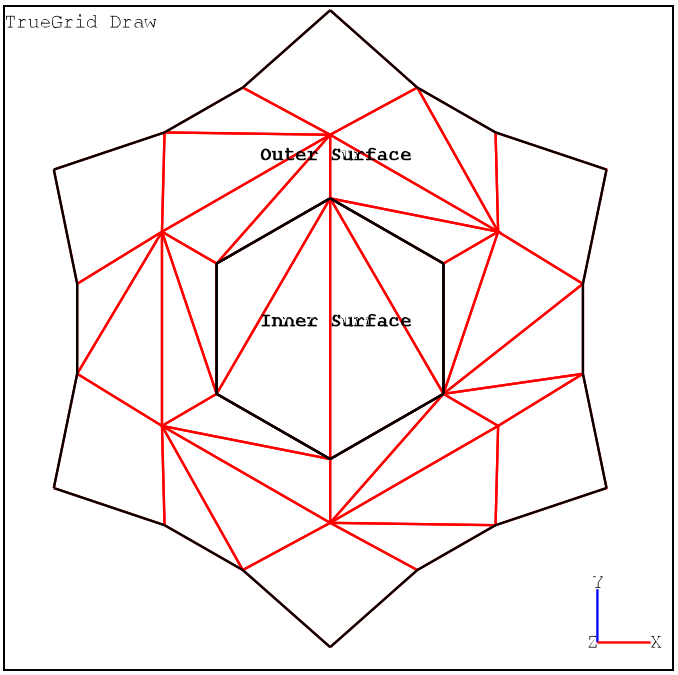

Figure 141 Polygon Surfaces

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```
     23,0.68,-1.88,2.5
     24,1.29,-1.53,2.5
     25,2.17,1.25,2.25
     26,0.00,2.50,2.25
     27,-2.17,1.25,2.25
     28,-2.17,-1.25,2.25
     29,0.00,-2.50,2.25
     30,2.17,-1.25,2.25
Connectivity file: vpsd.elem

     outer 1 6 12 13 14 7
     outer 1 7 15 16 8 2
     outer 2 8 17 18 9 3
     outer 3 9 19 20 10 4
     outer 4 10 21 22 11 5
     outer 5 11 23 24 12 6
     inner  1 2 3 4 5 6
     outer 7 14 25 15
     outer 8 16 26 17
     outer 9 18 27 19
     outer 10 20 28 21
     outer 11 22 29 23
     outer 12 24 30 13
```

wrsd write surface using the ViewPoint format

wrsd *node_file element_file surface_list* ;

where

node_file	file name to contain the nodal data for the surfaces
element_file	file name to contain the connectivity data for the surfaces
list	list of **True*Grid*®** surface numbers

Remarks

The files produced with this command can be read into **True*Grid*®** using the **vpsd** command.

12. Displaying Geometry from CAD Programs

IGES is the acronym for the Initial Graphics Exchange Specification which is an international standard interface for communicating geometry between Computer Aided Design (CAD) programs. When

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True*Grid*® reads such a file, it writes a short report of the geometry found in the file. Two types of objects that might be found in an IGES file are used to organize the geometry into sets. These objects must be created by the CAD operator before **True*Grid*®** can take advantage of them. The first is a level. Every geometric entity in an IGES file can belong to a level. **True*Grid*®** maintains the levels and you can select levels for graphics. The associativity group can be any set of geometric entities. A geometry entity can belong to many associativity groups. **True*Grid*®** maintains these groups and you can select groups for graphics. The following example of reading an IGES file demonstrates the identification of levels and groups.

```
iges maison.igs 1 1;

   The following IGES levels were found:
      1    2
   IGES file contained    34 NURBS surfaces from     1 to    34
   IGES file contained   192 curves from     1 to   192
   IGES file contained     2 groups from     1 to     2
```

When a sequence of levels or groups are needed, you only need to give the first and last numbers separated by a colon. When there are many surfaces in an IGES model, the first step in creating a mesh is to organize the geometry. Usually, there are many smaller surfaces that should be grouped to form larger composite surfaces because they touch each other tangentially. These larger surfaces become a more intuitive decomposition of the geometry. These larger composite surfaces also make it easier to design and build the block topology of the mesh because there is no longer the need to match the mesh boundaries to arbitrary component surface boundaries. If the CAD operator can use the associativity groups and levels in anticipation of building the mesh, the time spent using **True*Grid*®** to organize the geometry can be greatly reduced. If a set of surfaces are in the same level or group, to be formed into a composite surface, take the following steps:

1. Display only that level using the **dlv** command or that group using the **dgrp** command.
2. Use the **lasd** command to list all the surfaces in the picture.
3. Use the **sds** option of the **sd** command and cut and paste the list of surfaces into the command.

alv add a CAD level to the picture

alv *level*

where *level* is the number of the CAD level to be added.

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Remarks

This adds to the picture all defined surfaces resulting from CAD IGES surfaces and all 3D curves resulting from CAD IGES curves defined within the specified level.

dlv display a single CAD level in the picture

dlv *level*
where *level* is the number of the CAD level to be displayed

Remarks

This will display all defined surfaces resulting from CAD IGES surfaces and all 3D curves resulting from CAD IGES curves defined within the specified level. This command will remove all other surfaces and 3D curves from the picture.

dlvs display several CAD levels in the picture

dlvs $level_1$ $level_2$ ... $level_n$ **;**
where
$level_i$ is the number of one of the CAD levels to be displayed.

Remarks

This command will display all defined surfaces resulting from a CAD IGES surface and all 3D curves resulting from a CAD IGES curve defined within any of the specified levels. This will remove all other surfaces and 3D curves from the picture.

rlv remove a CAD level from the picture

rlv *level*
where *level* is the number of the CAD level to be removed.

Remarks

This command will stop displaying the defined surfaces resulting from any CAD IGES surfaces and all 3D curves resulting from any CAD IGES curves defined within the specified level.

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agrp add a CAD group to the picture

agrp *group*
where *group* is the number of the CAD group to be added.

Remarks

This command will add to the picture all defined surfaces resulting from CAD IGES surfaces and all 3D curves resulting from CAD IGES curves defined within the specified group.

dgrp display a single CAD group in the picture

dgrp *group*
where
group is the number of the CAD group to be displayed

Remarks

This command will display all defined surfaces resulting from CAD IGES surfaces and all 3D curves resulting from CAD IGES curves defined within the specified group. This command will remove all other surfaces and 3D curves from the picture.

dgrps display several CAD groups in the picture

dgrps $group_1\ group_2 \ldots group_n$ **;**
where $group_i$ is the number of one of the CAD groups to be displayed.

Remarks

This command will display all defined surfaces resulting from a CAD IGES surface and all 3D curves resulting from a CAD IGES curve defined within any of the specified groups. This command will remove all other surfaces and 3D curves from the picture.

rgrp remove a CAD group from the picture

rgrp *group*
where *group* is the number of the CAD group to be removed.

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Remarks

This command will stop displaying the defined surfaces resulting from any CAD IGES surfaces and all 3D curves resulting from any CAD IGES curves defined within the specified group.

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II. Assembly Commands - Merge Phase

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1. Merging Parts

In the Merge Phase, nodes that are close to one another are merged into a single node. The merging commands allow you to define how close is close. All tolerances are in absolute distances. There are commands for specifying tolerances for the general merging of all nodes over all parts or just nodes on the exterior faces of the mesh. There are commands for specifying the tolerances for the special merging of nodes between parts or within a part. These special tolerances override the general ones. If no tolerance commands are specified, then no merging is done. However, the Merge Phase must be entered in order to build the node map which is used to generate the output.

Invocation of a tolerance command (**t**, **tp**, **st**, **stp**) within the Merge Phase causes an immediate merging of nodes. These commands can also be invoked within any phase; when the Merge Phase is entered, those tolerance commands are immediately executed or re-executed as the case may be. The merge process is always performed on the nodes in their original (prior to any merging) state. Merging is not cumulative. If you leave the Merge Phase and reenter it, all merging is recalculated with what ever new parts that have been added. This lets you interactively experiment with merging and tolerances. Setting a tolerance to a negative value is an easy way to restore the nodes to their original states. Graphical displays of the mesh in the Merge Phase always reflect the results of any merging.

Nodes are merged depending on the distance between them. If a node lies within a tolerance distance of more than one other node, then it is merged with the closest one. When merging several nodes into one node, the first-defined node survives. This can be overridden by the **bptol** command. Nodes within a joint and across the two sides of a sliding surface are not merged. When the first merging of nodes occurs, a sliding interface table is calculated which is used in the merging process. This table is written to the screen and to the save file and is intended as diagnostics. The following is a sample of that table:

SLIDING INTERFACE SUMMARY

Surf	S-node	S-lseg	S-qseg	M-node	M-lseg	M-qseg
1	105	84	0	468	418	0
2	232	0	52	468	418	0
3	221	0	0	390	304	0
4	221	0	0	390	304	0
5	158	0	0	120	88	0
6	158	30	30	120	88	0
7	204	102	0	204	102	0
8	232	0	52	90	52	0
9	101	18	18	161	132	0
10	101	18	18	161	132	0
11	548	120	120	3216	0	1056
12	133	84	0	161	132	0

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```
 13      133       84        0      161      132        0
 14      308      240        0     3216        0     1056
```

This table is organized by the sliding interface number on the right. Columns 2, 3, and 4 are datum pertaining to the slave side on the interface; columns 5, 6, and 7 to the master side. Columns 2 and 5 (`S-node` and `M-node`) are node counts. Columns 3 and 6 (`S-lseg` and `M-lseg`) are linear face counts, and columns 4 and 7 (`S-qseg` and `M-qseg`) are quadratic face counts.

A table of merged nodes is always written after the **tp** or **stp** commands are executed.

```
                MERGED NODES SUMMARY
  12 nodes merged between parts        1 and        2
  16 nodes merged between parts        2 and        2
  12 nodes merged between parts        1 and        3
  16 nodes merged between parts        3 and        3
 216 nodes merged between parts        4 and        4
  30 nodes merged between parts        7 and        7
  88 nodes merged between parts        8 and        8
 390 nodes were deleted by tolerancing
```

Up through 4000 parts can be merged under general tolerancing (i.e. no use of the **ptol** or **bptol** commands). 300 parts can be merged under special tolerancing (**ptol** and **bptol**).

The following is a common error to avoid. Suppose you create three parts that meet as shown in 142 and 143. Then define a sliding interface between parts 1 and 2 and also between parts 1 and 3. No nodes will be merged between parts 1 and 2 and between parts 1 and 3. However, nodes can be merged between parts 2 and 3. Sometimes you need to look closely in the graphics or carefully check the Merged Nodes Summary to detect this error. To fix this error, if indeed it is an error, use a dummy sliding interface between parts 2 and 3 to force no merging between those parts. Alternatively, use the **bptol** command with a negative number to avoid merging between those parts. You should also consider extending both interfaces 1 and 2 across to parts 3 and 2, respectively, because they may come in contact. This is an ambiguous situation since there are equally plausible situations where parts 2 and 3 should be merged together.

```
sid 1 sv;
sid 2 sv;
block 1 3;1 3;1 3;
     1 2 1 2 1 2
sii -2;;;1 s;
```

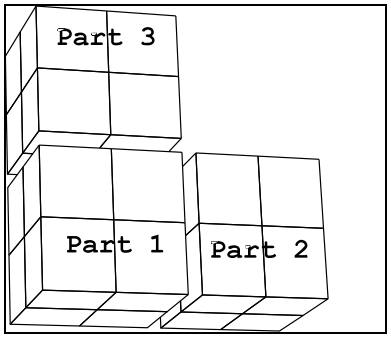

Figure 142 Before **stp**

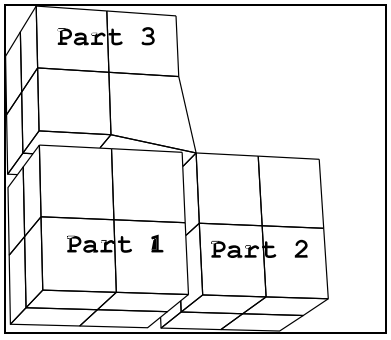

Figure 143 After **stp**

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```
sii ;-2;;2 s;
block 1 3;1 3;1 3;
     2.1 3 1 2 1 2
sii -1;;;1 m;
block 1 3;1 3;1 3;
     1 2 2.1 3 1 2
sii ;-1;;2 m;
merge
stp .2
```

For other commands affecting how nodes are merged, see the "Merging Parts" section in the chapter "Global Commands".

mnl write the merged nodes list

mnl *(no arguments)*

Remarks

The merge table is written out to the save file. This may be useful when you are using **True*Grid*®** to get intermediate results and need to process the mesh data further. The table consists of two columns. The first column is the node number before merging. The second column is the node number after merging. Two rows with the same second node number means the two corresponding nodes have been merged together

Example

The following butterfly block structure is a good example of merging within one part.

```
sd 1 cy 0 0 0 0 0 1 3
block 1 2 3 4;1 2 3 4;-1;-1 -1 1 1 -1 -1 1 1 0
dei 1 2 0 3 4;1 2 0 3 4; -1;
sfi -1 0 -4;;-1;sd 1
sfi ;-1 0 -4;-1;sd 1
merge
stp .001
mnl
```

This produces the merged nodes table below. Notice that the original node number 5 is merged to node 1. This causes the original node 6 to be renamed as node 5. The original node 9 is merged to node 2, causing the original node 10 to be renamed node 8. The original node 11 is merged to node 5, which

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was originally node 6. The original node 12 is merged to node 8, which was originally node 10.

```
c MERGED NODES TABLE
c          1          1
c          2          2
c          3          3
c          4          4
c          5          1
c          6          5
c          7          6
c          8          7
c          9          2
c         10          8
c         11          5
c         12          8
```

pn place a node at a new location

pn *node x y z*

where

node	node number
x y z	Cartesian coordinates of the new location

Remarks

This command is used automatically when the *Move Pts.* and *node* button are selected in the Merge Phase environment window.

The node numbers change when a different merge command is issued. If two nodes are merged into one and assigned a new number, it is this number that is used in this command. If, after this node is moved using the **pn** command, a new merge command is issued with a negative value so that the two nodes are no longer merged together, both nodes will have been moved to the same location so that they are coincident but separate nodes.

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2. Diagnostics Commands

ajnp find node near point

ajnp *x_coordinate y_coordinate z_coordinate*

Remarks

The automatic parameter, %node, is set to the node number of the node that is located nearest (Euclidean distance) the given point. This parameter is then available for future commands. This feature has been provided so that the node number can be used in batch mode.

sajnp silently find node near point

sajnp *x_coordinate y_coordinate z_coordinate*

Remarks

The automatic parameter, %node, is set to the node number of the node that is located nearest (Euclidean distance) the given point. This parameter is then available for future commands. This feature has been provided so that the node number can be used in batch mode.

ajel find element near point

ajel *element_type x_coordinate y_coordinate z_coordinate*

where the *element_type* can be:

1d	for linear beams
2d	for linear shells
3d	for linear bricks
1dq	for quadratic beams
2dq	for quadratic shells
3dq	for quadratic bricks

Remarks

The automatic parameter, %element, is set to the element number of the element that is located nearest (Euclidean distance) the given point. This parameter is then available for future commands. This feature has been provided so that the element number can be used in batch mode. The resulting element

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number is printed to the text window.

The measurement is made from the point at (*x_coordinate,y_coordinate,z_coordinate*) to the approximate center of mass of the element. Both linear and quadratic elements are measured using the same number of nodes to calculate the center of mass. For example, the 8 nodes of a linear hexa element are used to calculate the center of mass. If the same element is measured as a quadratic brick element (i.e. the quadratic command was issued before the part was created) then only the 8 linear nodes at the corners of the hexa element are used to calculate the center of mass. This way, the same numbered element will be chosen no matter if the element is linear of quadratic.

sajel silently find element near point

sajel *element_type x_coordinate y_coordinate z_coordinate*

where the *element_type* can be:

1d	for linear beams
2d	for linear shells
3d	for linear bricks
1dq	for quadratic beams
2dq	for quadratic shells
3dq	for quadratic bricks

Remarks

The automatic parameter, %element, is set to the element number of the element that is located nearest (Euclidean distance) the given point. This parameter is then available for future commands. This feature has been provided so that the element number can be used in batch mode. No message is sent to the test window when this command is executed.

The measurement is made from the point at (*x_coordinate,y_coordinate,z_coordinate*) to the approximate center of mass of the element. Both linear and quadratic elements are measured using the same number of nodes to calculate the center of mass. For example, the 8 nodes of a linear hexa element are used to calculate the center of mass. If the same element is measured as a quadratic brick element (i.e. the quadratic command was issued before the part was created) then only the 8 linear nodes at the corners of the hexa element are used to calculate the center of mass. This way, the same numbered element will be chosen no matter if the element is linear or quadratic.

cenref restore reference center for moment calculations

cenref *(no arguments)*

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Remarks

Restore the reference center for moment calculations to its default, the origin of the coordinate system.

Normally, the origin of the coordinate system is used as a reference center in moment calculations. You can change that with **reference** command. The purpose of this command is simply to restore the default afterwards.

centroid moments and inertia

centroid *(no arguments)*

Remarks

Print the moments and inertias of each part and material.

elm highlight elements within a *measure* interval

elm *minimum maximum*

where

minimum	is the lower property limit, and
maximum	is the upper limit on the values of the property.

Remarks

This command is designed to be used after you have determined the quality level of your mesh using the **measure** command. First, you measure some property (orthogonality, aspect ratio, Jacobian, etc.) using the **measure** command. Then select a range of values for this property. The elements in that range will be highlighted. For example:

```
measure orthogon
elm -90 -45
```

highlights elements which contain angles less than 45 degrees. Notice that the orthogonality option of the measure reports the deviation from 90 degrees – not the absolute angles.

If no element lies inside the selected range of the **elm** command the element(s) whose measure are closest to that range are highlighted. (e.g., if no element has a negative measure and the command **elm -2 -1** is issued, the element(s) with smallest measure will be highlighted.)

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If, after merging nodes, an element is collapsed to form an illegal element (e.g. a 7 noded brick), many of the options in the **measure** command (from the volume, jacobian, stiffn, and pointvol options only) will report elements with a measure of -1.E10. To view these illegal elements set the minimum value in the **elm** command to the illegal element value. The orthogon, warp, and triangle option use -91 to indicate illegal elements are present.. The avolume, smallest, and aspect options use -1 to flag illegal elements.

elmoff turn off highlighting from the elm command

elmoff *(no arguments)*

info mesh model summary

info *(no argument)*

Example

The **info** command for the following input produced the table that follows.

```
title Spring Example
spd 1 le 1.1 spd 2 le 2.2 spd 3 dhpt 3.3 4.4
block 1 3 5;1 3 5;1 3 5;1 3 5;1 3 5;1 3 5;
spdp 1 1 3 2 2 3 1 m 1 dx dy dz;
spring 2 v1 3 3 3 dx1 sddn 3 amp 2.12;
npm 2 0 0 0 1.2 inc 2 dx rx;lct 1 rz 180;lrep 0 1;
block 1 3 5;1 3 5;1 3 5;1 3 5;1 3 5;6 8 10;
spdp 1 1 1 2 2 1 1 s dx dy dz;spring 2 v2 3 3 1 dx2;
lct 1 rz 180;lrep 0 1;merge
spring 1 n1 125 n2 335 rx1 ry1 rz1 rx2 ry2 rz2 sddn 2 amp 1;;
info
```

```
           2 parts
         502 nodes after merging
         256 linear solids
          18 bulk springs/dampers
           3 numbered springs/dampers
           2 point masses
```

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mass mass table

mass *(no arguments)*

Print the masses, volumes and center of gravities of each part and material.

Remarks

The density of the material in this calculation is based on the material model. Be sure to define the material model before using this command. Densities are used only for materials defined for the following codes: ABAQUS, ALE3D, ANSYS, DYNA3D, ENIKE3D, ES3D, NASTRAN, LSDYNA, LSNIKE, MARC, NEUTRAL, NIKE3D, and TOPAZ3D.

measure choose a way to measure mesh quality at every element

measure *option*

where

option can be:

volume	to integrate the volume of each element
avolume	to integrate the absolute volume
jacobian	to compute the determinant of the Jacobian
orthogon	to measure deviations from orthogonality (90 degrees) in quadrilaterals faces of elements
smallest	for the smallest dimension of each element
pointvol	to calculate the volume with a one point integration formula
aspect	to calculate the aspect ratio for each element
triangle	to measure deviations from optimum (60 degrees) in triangular faces of elements
warp	to measure the angle between opposite corners of each element face
stiffn	stiffness or the condition number of the Jacobian
dvi	compare the volumes of neighboring elements in the i-direction - only for structured meshes
dvj	compare the volumes of neighboring elements in the j-direction - only for structured meshes
dvk	compare the volumes of neighboring elements in the k-direction - only for structured meshes
dli	compare the lengths of neighboring edges in the i-direction - only for structured meshes
dlj	compare the lengths of neighboring edges in the j-direction - only for

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	structured meshes
dlk	compare the lengths of neighboring edges in the k-direction - only for structured meshes
subvol	volumes of sections of elements
anjac	average determinant of the Jacobian for every node
annjac	average determinant of the relative Jacobian for every node
anstif	average stiffness of the Jacobian for every node
wannjac	average determinant of the relative Jacobian for every node

Remarks

A histogram is drawn to show the profile of the mesh according to the selected measure. The abscissa is the measure and the ordinate is the number of elements when there is one measurement per element or element segments when there are several measurements per element. The range of the measurement, the mean, and the standard deviation is written to the save file and it is displayed in the text window during an interactive session. These 4 values are assigned to the minmea, maxmea, mean, and standev automatic parameters, respectively.

This measure can detect invalid elements if a tolerance is specified using one of the merging commands (**t**, **tp**, **stp**). This tolerance is also used to identify the element geometry (hex, tet, wedge, pyramid, quad, or triangle). Many of the options depend on a shape function which is different for each type of element geometry. The order of the element (linear, quadratic, and triquadratic) also determines the shape function used in the measure. There are 14 different shape functions based on the element types generated by **True*Grid*®**. For more details on the mathematics, see "Application of An *a priori* Jacobian-based Error Estimate Metric to the Accuracy Assessment of 3D Finite Element Simulations" by Robert B. Rainsberger, Jeffrey T. Fong, and Pedro V. Marcal published in the "Proceedings of the ASME Pressure Vessels and Piping Conference", PVP2018-84784, July 15-20, 2018, Prague, Czech Republic. See the **vigq** command to change the number of Gaussian quadrature points from the default of 2.

The 4 node based options (**anjac**, **annjac**, **anstif**, and **wannjac**) calculate the Jacobian at the nodes instead of at interior points. Note that this implies that the **vigq** command has no effect on the results of these 4 options. All of the measurements for a single node are averaged. This is done for every node.

The **volume** option integrates the volume of a brick element using the appropriate shape function to estimate the volume. It is possible for the volume to be negative in some regions of the element; in that case the net volume will not be realistic. Shell elements are given thickness and the same method is then used to calculate the volume. If the shell element was not assigned a thickness, then the default of 1 is used. An illegal element is assigned a measure of -1.0e+10.

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The **avolume** option has the advantage that it is not affected by negative volumes since the absolute volume is integrated. Shells are given thickness as for the volume option above and then treated like a brick element. Illegal elements are assigned a measure of -1.

The **Jacobian** option measures the distortion of each element by sampling the Jacobian matrix of the appropriate shape function. In order to graph this data, the Jacobian matrix is reduced to a single number by first determining the eigenvalues of the matrix. The eigenvalue whose modulus is found between the other two is used to scale the Jacobian matrix. The matrix is divided by the cube of this modulus. This factors out the scale of the element. The determinant of the resulting matrix is graphed. Shell elements are given thickness to make this measurement. An illegal element is assigned a measure of -1.0e+10.

The **orthogonal** option measures all angles at all corners of a quad face of a brick element. It then graphs the deviation from 90 degrees. Each quad shell element is measured at each of the four corners. Illegal elements are flagged with -91.

The **smallest** option determines the smallest dimension of an element as the measurement. Illegal elements are flagged with -1.

The **pointvolume** option approximates the volume of an element using the Jacobian at a single point in the center of the element. The shell elements are given thickness and treated the same as bricks (see above). An illegal element is assigned a measure of -1.0e+10.

The **aspect** ratio is defined as the ratio of the largest diagonal to the smallest diagonal of an element. If the largest diagonal is zero, then the ratio is set to zero. If the smallest diagonal is zero, the ratio is set to a very large constant. Illegal elements are flagged as -1.

The **triangles** option measures the 3 angles of all triangles in an element. It then graphs the deviation from 60 degrees. Illegal elements are flagged as -91 degrees.

The **warp** option measures the angle between the normals at opposing nodes of each quad face. Illegal elements are flagged with -1.

The **dvi** option measures the ratio of volumes of consecutive elements along every thread of elements in the i-direction. This is only for structured meshes.

The **dvj** option measures the ratio of volumes of consecutive elements along every thread of elements in the j-direction. This is only for structured meshes.

The **dvk** option measures the ratio of volumes of consecutive elements along every thread of elements in the k-direction. This is only for structured meshes.

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The **dli** option measures the ratio of edge lengths of consecutive elements along every thread of elements in the i-direction. This is only for structured meshes.

The **dlj** option measures the ratio of edge lengths of consecutive elements along every thread of elements in the j-direction. This is only for structured meshes.

The **dlk** option measures the ratio of edge lengths of consecutive elements along every thread of elements in the k-direction. This is only for structured meshes.

The **stiffn** options measures the stiffness or condition number of the Jacobian using the appropriate shape function. Shell elements are given thickness to make this measurement. An illegal element is assigned a measure of -1.0e+10.

The **subvol** option displays the minimum and maximum element sample volumes used to calculate the total volume of an element. This makes it easy to identify those elements with regions of negative volume.

The **anjac** option means for each node, the determinant of the Jacobian is calculated for each element that contains this node, and the average is calculated.

The **annjac** option means for each node, the determinant of the relative Jacobian is calculated for each element that contains this node, and the average is calculated. The relative Jacobian is the Jacobian divided by the cube of the middle singular value of the Jacobian. In plane English, size of the element is factored out.

The **anstif** option means for each node, the stiffness of the Jacobian is calculated for each element that contains this node, and the average is calculated.

The **wannjac** option means for each node, the determinant of the relative Jacobian is calculated for each element that contains this node, and the weighted average is calculated. The relative Jacobian is the Jacobian divided by the cube of the middle singular value of the Jacobian. In plane English, size of the element is factored out. The weight in the averaging is the volume of the contributing brick element or the area of the contributing shell element.

With the **elm** and **elmoff** commands, you can see the locations in the mesh of the most interesting elements; e.g. you can use **measure** to measure volume and then **elm** to highlight the biggest elements.

nodcor retrieve coordinates of a node

nodcor *node_#*

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Remarks

This function prints the coordinates of a node. Note that node numbers may change as nodes are merged. The automatic parameters %xprj, %yprj, and %zprj will be assigned the coordinates of the node.

snodcor silently retrieve coordinates of a node

snodcor *node_#*

Remarks

This function prints the coordinates of a node. Note that node numbers may change as nodes are merged. The automatic parameters %xprj, %yprj, and %zprj will be assigned the coordinates of the node.

pmass active part mass

pmass *(no arguments)*

Remarks

The density of the material in this calculation is based on the material model. Be sure to define the material model before using this command. Densities are used only for materials defined for the following codes: ABAQUS, ALE3D, ANSYS, DYNA3D, ENIKE3D, ES3D, NASTRAN, LSDYNA, LSNIKE, MARC, NEUTRAL, NIKE3D, and TOPAZ3D.

reference reference point for moments and inertia

reference x_0 y_0 z_0

where

(x_0,y_0,z_0) is the reference point

Remarks

This command changes the center reference point used in moment calculations. If you like, you can later issue the **cenref** command to restore this reference point to its default value: the origin of the coordinate system.

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size dimensions of the mesh

size *(no arguments)*

smags detect detached small groups of elements

smags *element_set maximum*

Remarks

This command detects small detached groups of elements. These elements are put into a set. A cut off is specified and any detached group of elements with that or fewer elements will be placed into the garbage element set. This command is useful after reading a mesh with the **dynain** option of the **readmesh** command.

tmass total mass

tmass *(no arguments)*

Remarks

The density of the material in this calculation is based on the material model. Be sure to define the material model before using this command. Densities are used only for materials defined for the following codes: ABAQUS, ALE3D, ANSYS, DYNA3D, ENIKE3D, ES3D, NASTRAN, LSDYNA, LSNIKE, MARC, NEUTRAL, NIKE3D, and TOPAZ3D.

vigq quadrature rule for mass calculations

vigq *n*

where *n* is the number of Gaussian points used in the **measure** command.

Remarks

The default is 2.

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3. Graphics Commands

backplane toggles back plane removal

backplane on
or
backplane off

Remarks

This affects the efficiency of the hidden line display graphics and the number of polygons in the draw. Specifically, it instructs the drawing algorithm to ignore all of those polygons which are "facing" towards the "back" of the image, thus greatly reducing re-draw time. This command is especially useful when re-drawing large, complicated models or when looking at all the IGES surfaces.

dpic dumps all picture parameters

dpic *(no arguments)*

Remarks

This is used to save a view to be read back into **True*Grid*®** using **rpic**.

rpic reads the picture parameters

rpic *picture_parameters*

Remarks

The picture parameters that were dumped by **dpic**.

4. Tokens in the Picture

This section describes tokens in the picture which identify features in the model. The **condition** command helps display boundary conditions, constraints, and the like. The **labels** command displays indices, surface numbers, and so forth. The **mlabs** command combines the **condition** and **labels** command and allows for multiple options.

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Some of these labels and conditions may not be available in a filled (**tvv**) display. No tokens are available in OpenGL graphics (***Fast Gr.*** graphics).

mlabs multiple labels and conditions displayed

mlabs *options* ;

where an *option* can be any option available in the **labels** (**la**) or **condition** (**co**) commands.

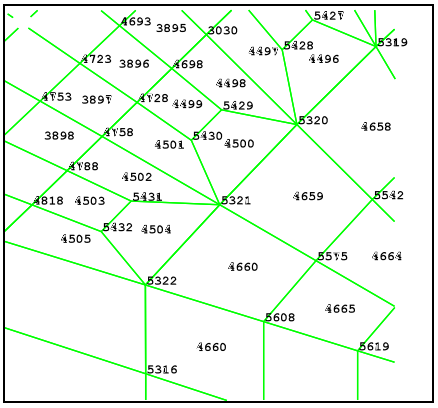

Figure 144 Multiple Labels

Remarks

Any combination of options listed under both the labels and condition commands can be used in this command. Warning messages are issued when doing facial conditions. The messages consist of the number of degenerate faces encountered.

Example

```
mlabs nodes 3d ;
```

condition specify type of condition/constraint to be displayed

condition *option*

or

co *option*

where *option* can be:

		Relevant commands:
dx	for nodes with fixed translation in x	**b, bi**
dy	for nodes with fixed translation in y	**b, bi**
dz	for nodes with fixed translation in z	**b, bi**
rx	for nodes with fixed rotation in x	**b, bi**
ry	for nodes with fixed rotation in y	**b, bi**
rz	for nodes with fixed rotation in z	**b, bi**
mom *load_curve direction*	for nodes with prescribed torque	**mom, momi**

where *direction*

x	around the x-axis
y	around the y-axis
z	around the z-axis

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fmom *load_curve* follower moments **fmom**

fc *load_curve* for point load vectors at nodes **ll, fc, fci**

fd *load_curve* for forced displacement vectors at nodes **fd, fdi, fdc, fdci, fds, fdsi**

pr *load_curve* for pressure surface amplitude vectors **pr, pri**

sy *symmetry_plane* for nodes on symmetry planes **plane**

where

symmetry_plane symmetry plane number

si *sliding_interface type* sliding interfaces or contact surfaces **sid, si, sii**

where

sliding_interface sliding interface number

type can be

m	master
s	slave
b	both

rb for boundary radiation orientation vectors on faces **rb, rbi**

re for radiation enclosure orientation vectors on faces **re, rei**

fl for boundary flux orientation vectors on faces **fl, fli**

cv for boundary convection orientation vectors on faces **cv, cvi, vcv, vcvi**

cvt convection thermal loads **cvt, cvti**

tm for initial temperature at nodes **tm, tmi, vtm, vtmi**

ft for prescribed temperature at nodes **ft, fti, vft, vfti**

fv *load_curve_number* for prescribed velocity at nodes **fv, fvi, fvc, fvci, fvs, fvsi, fvv, fvvi, fvvc, fvvci, fvvs, fvvsi**

sw *stone_wall_number* for nodes assigned to stone walls **plane, sw, swi**

nr for nonreflecting boundary faces **nr, nri**

jt *joint_number* for nodes in joints **jd, jt**

iss for interface save segments **iss, issi**

ve initial velocities **velocity, ve, vei**

efl for electric flux orientation vectors on faces **efl, efli**

vhg *load_curve* volumetric heat generation **vhg, vhgi, vvhg, vvhgi**

or sys to display element local coordinate system axis **or**

where

sys can be any one of the following:

r	to label the local element r-axis
s	to label the local element s-axis
t	to label the local element t-axis
rs	to label the local element r-axis and s-axis
tr	to label the local element r-axis and t-axis

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st to label the local element s-axis and t-axis

rst to label the local element r-axis, s-axis, and t-axis

syf *plane* for nodes assigned to a symmetry plane with failure **plane, syf, syfi**

where

plane number of the symmetry plane with failure

bv prescribed boundary velocities **bv, bvi**

ol outlets **ol, oli**

il inlets **il, ili**

npb nodal print blocks **npb**

epb element print block **epb**

sp *spring/damper_#* springs and dampers **sp, spdp**

where a spring or damper number can be

0 for all numbered springs

n for a numbered spring

-1 for all unnumbered springs

pm point masses **pm, npm**

interfac *slide_#* ABAQUS interface elements **sid, si, sii, bb**

thic shell thicknesses **th, thi, thic, ssf, ssfi**

mdep momentum deposition **mdep**

sfb *constraint* surface oriented constraints **sfb**

where

constraint can be:

dx constraint in the local x-direction

dy constraint in the local y-direction

dz constraint in the local z-direction

rx constrain rotation about the local x-axis

ry constrain rotation about the local y-axis

rz constrain rotation about the local z-axuis

tepro *load_curve* temperature profile **tepro**

sc smoothing constraints **sc**

acc *load_curve* for prescribed acceleration **acc, acci, accc, accci, accs, accsi, vacc, vacci, vaccs, vaccsi, vaccc, vaccci**

n for shell element outward normal vectors **n**

resn for radiation enclosure surface numbers **re, rei**

grtol *grid# patch#* glued interfaces for CFX **supblk, CFX4, stp, st, tp, p**

where

grid# is the assigned part number for a grid

patch# is the sequence number of the patch associated with that grid. The table of grid and patches is written to the screen each time

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	the model is merged with either **stp**, or **tp** and this option is invoked. This is used only for the CFX output option. Be sure to merge the model before using this option or writing the output.	
spw *spotweld_#*	numbered spotweld	**spw, spwd**
spwf *material_#*	Ls-dyna material 100 spotweld	**spwf**
ffc *load_curve*	follower point loads	**ffc**
bf *number*	bulk fluid	**bf, bfi, bfd**
frb *op load_curve*	prescribed nodal rotations	**frb**
	where *op* can be	
	v for velocity	
	a for acceleration	
	d for displacement (LS-DYNA only)	
	dofv nodal DOF velocity	
	dofa nodal DOF acceleration	
	dofd nodal DOF displacement (LS-DYNA only)	
detp	detonation points	**detp**
trp	tracer particles	**trp**
fbc *zone_num*	Fluent boundary conditions	
starbc *region_id*	Starcd boundary conditions	
kivabc *type*	Kiva boundary conditions	
cfxbc *label*	Cfx5 boundary conditions	
sph	smooth particle hydro elements	
dem	discrete element spheres	
mpce	multiple point constraint equations	
off	to turn off condition display	
size *scale*	to scale the size of the tokens displaying conditions	
angle θ	to change cone angle for tokens displaying conditions	

Remarks

The tokens used to represent conditions on the model may be different when viewing the same condition in the wire and the hide mode graphics. For example, when viewing a sliding interface, the wire graphics circles every node on the interface. The hide shows an arrow at the center of every face on the interface, pointing in the outward direction. Both can be useful information. Warning messages are issued when doing facial conditions. The messages consist of the number of degenerate faces encountered.

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Example

co fc display concentrated nodal loads

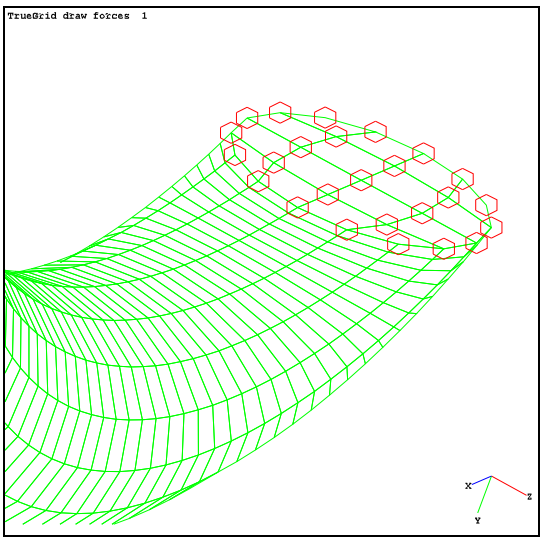

Figure 145 Wire frame version of CO FC

Example

```
cylinder 1 5;1 226;1 5;
    1.5 2.5 0 900 -.5 .5
sfi -1 -2;1 2;-1 -2;
     ts 0 0 0 0 0 1 2 0 1
z=z+0.05*j
fc 1 2 1 2 2 2 1
  1.2 .1 .2 .3
endpart
merge
co size 3.2
co fc 1
```

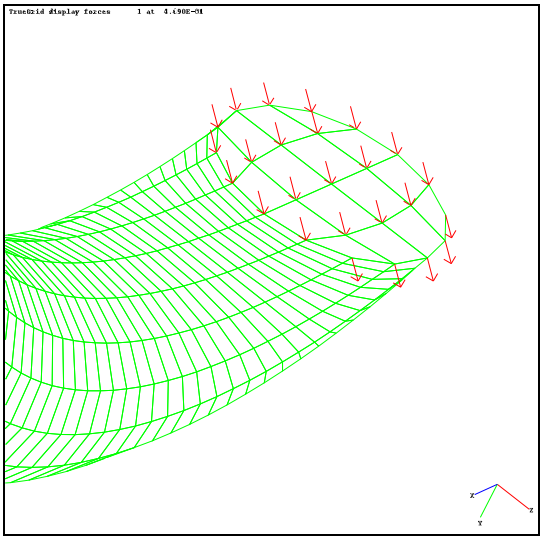

Figure 146 Hidden Line CO FC

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Remarks

The **co acc** (display accelerations), **co fmom** (display follower nodal moment), and **co fd** (display fixed displacements) options have a similar display.

co pr **display pressure**

Remarks

The **co bf** (display bulk fluid) is displayed in a similar fashion.

Example

```
gct 1 mx 4 my 4 rz 15;
lev 1 grep 0 ;levct 11 rz 30;
     repe 11;;
lev 2 grep 0 1;;
pslv 1
pslv 2
block 1 4;1 4;1 4;5 7 5 7 5 7
sfi -1 -2;-1 -2;-1 -2;sp 6 6 6 1
lcd 1 0 0 1 1;
pr 2 1 1 2 2 2 1 1
endpart
pplv
pplv
merge
co pr 1 co size 3.2
```

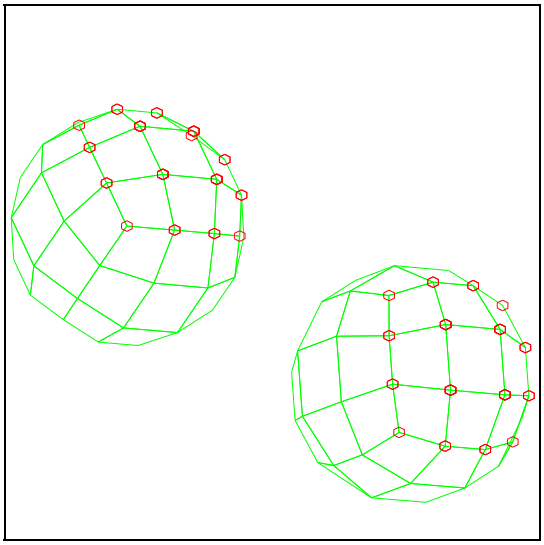

Figure 147 Wire frame version CO PR

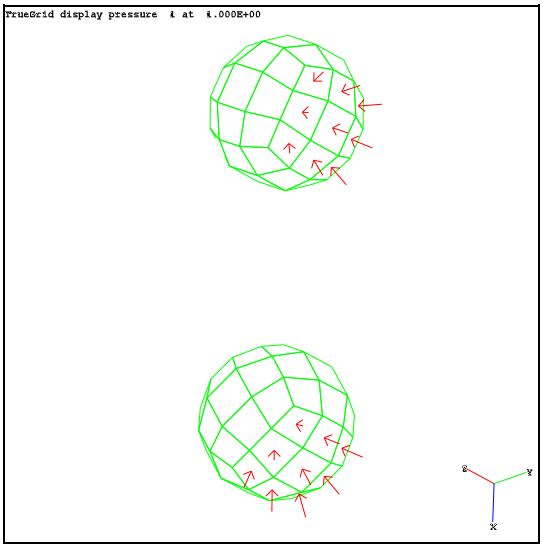

Figure 148 Hidden Line version CO PR

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co si **display sliding interfaces**

Examples

co rb **display radiation boundary condition**

Remarks

To assign this condition, see the **rb**, **rbi**, **vrb**, and **vrbi** commands in the part phase and the **rb** command in the merge phase.

Example

```
block 1 3 5 7 9;
      1 3 5 7 9;
      1 3 5 7 9;
      -1 -1 0 1 1;
      -1 -1 0 1 1;
      -1 -1 0 1 1;
dei 1 2 0 4 5; 1 2 0 4 5;;
dei 1 2 0 4 5;;1 2 0 4 5;
dei ;1 2 0 4 5;1 2 0 4 5;
sfi -1 -5;-1 -5;-1 -5;sp 0 0 0 5
sfi -2 -4;-2 -4;-2 -4;sp 0 0 0 3
de 2 2 2 4 4 4
de 1 0 0 3 0 0
orpt - 0 0 0
rbi 3 -5;-1 -5;-1 -5;1 1 2 1
merge
co rb co size 3
```

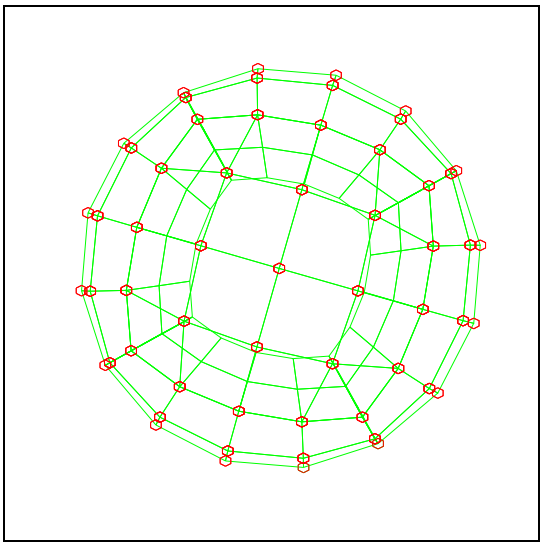

Figure 149 Wire frame version CO RB

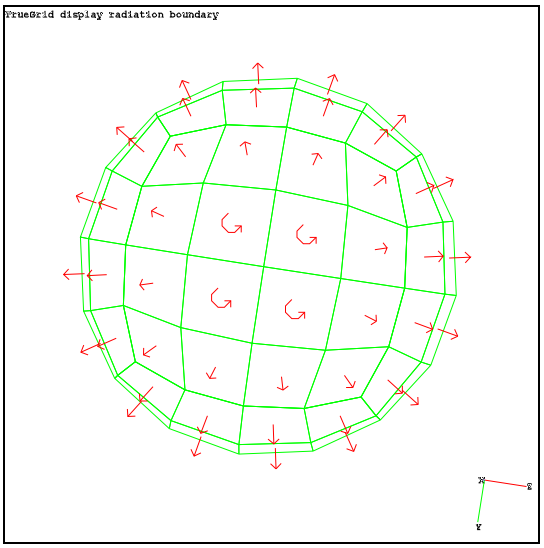

Figure 150 Hidden Line version CO RB

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Example

```
block 1 3 5 7 9;
      1 3 5 7 9;
      1 3 5 7 9;
      -1 -1 0 1 1;
      -1 -1 0 1 1;
      -1 -1 0 1 1;
dei 1 2 0 4 5;1 2 0 4 5;;
dei 1 2 0 4 5;;1 2 0 4 5;
dei ;1 2 0 4 5;1 2 0 4 5;
sfi -1 -5;-1 -5;-1 -5;sp 0 0 0 5
sfi -2 -4;-2 -4;-2 -4;sp 0 0 0 3
de 2 2 2 4 4 4
de 1 0 0 3 0 0
orpt + 0 0 0
rei 3 -4;-2 -4;-2 -4;0 2.2 no
merge
```

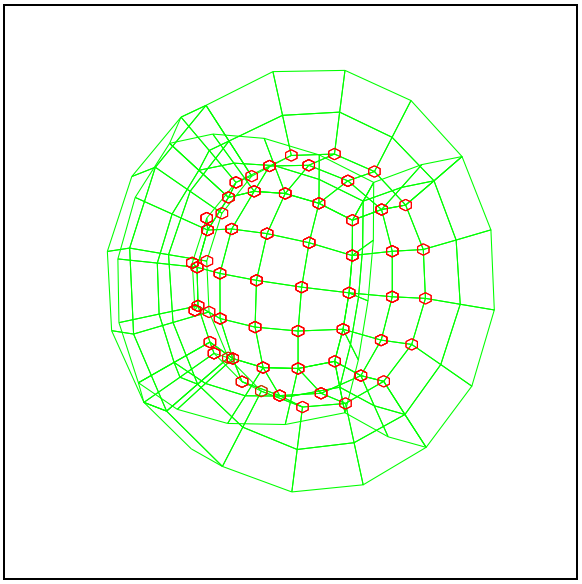

Figure 151 Wire frame version CO RE

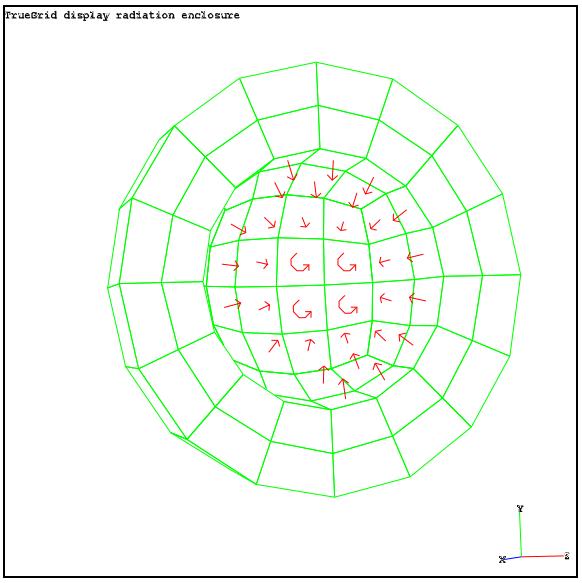

Figure 152 Hidden Line version CO RE

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co fl **display**

Example

```
block 1 13 17 31 35 47;
      1 5 9 13;1 25;
      1 13 17 31 35 47;
      1 5 9 13;1 25;
dei 2 5; 1 2 0 3 4;;
mate 1
mti 1 3 ; 1 2 ; ; 2
mti 1 3 ; 3 4 ; ; 3
mti 4 6 ; 1 2 ; ; 4
mti 4 6 ; 3 4 ; ; 5
mbi -2; -1 0 -4; 1 2;x -1
mbi -5; -1 0 -4; 1 2;x 1
bb 2 2 1 3 2 2 1;
bi ;; -1;dz 1;
fdi -6;1 2 0 3 4;1 2;1 1 1 0 0
fdi -1;1 2 0 3 4;1 2;1 1 -1 0 0
lcd 1 0 0 1 1;
orpt - 24 7 13
fli -1 2 0 5 -6;-1 -4;-1 -2;1 1
merge
co fl co size 2.6
```

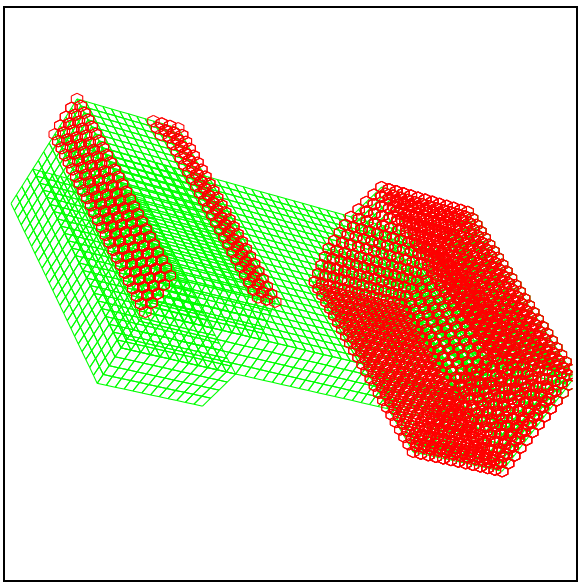

Figure 153 Wire frame version CO FL

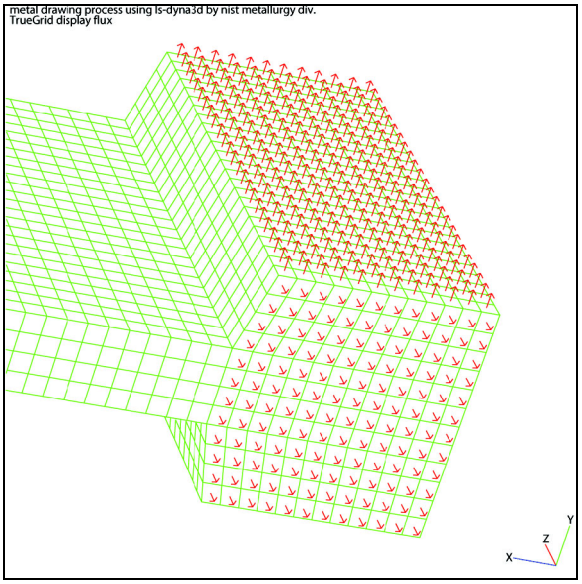

Figure 154 Hidden Line version CO FL

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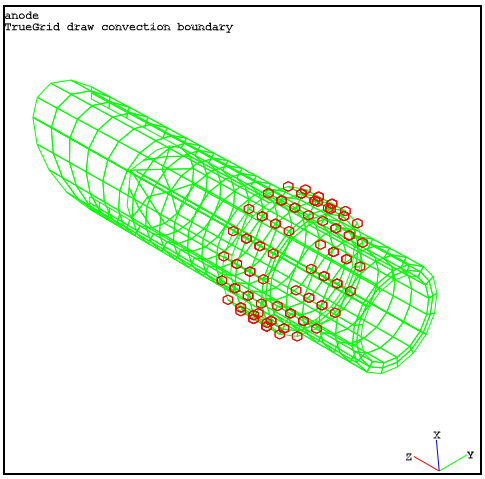

Figure 155 Wire version CO CV

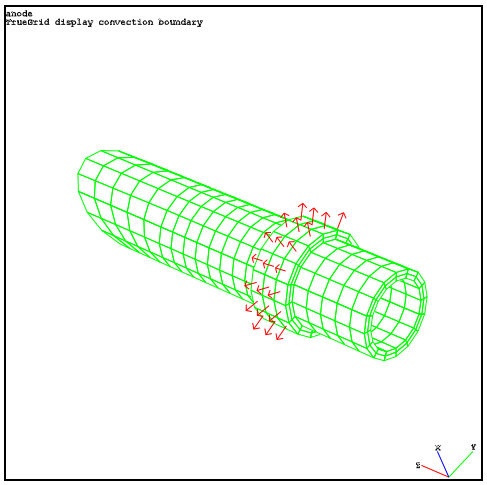

Figure 156 Hidden version CO CV

co cv **display convection conditions**

co cvt **display convection thermal loads**

Example

```
block -1 5 -9;-1 5 -9;-1 5 -9;-1 0 1;-1 0 1;-1 0 1;
sfi -1 -3 ;-1 -3 ;-1 -3 ; sp 0 0 0 5 de 1 0 0 2 0 0 orpt + 0 0 0 cvti 1
-3;-1 -3;-1 -3;.12 55 merge co cvt co size 2.4
```

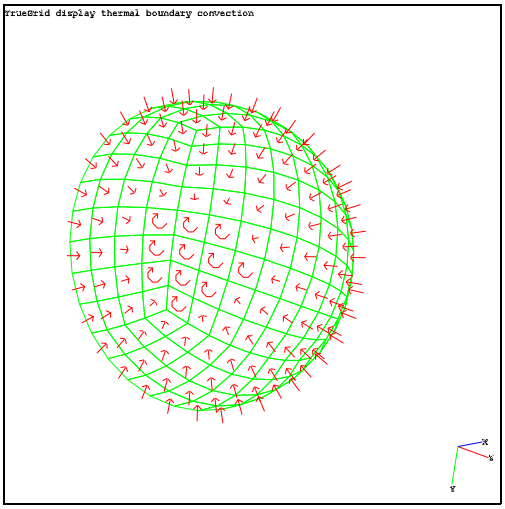

Figure 157 Hidden Line CO CVT

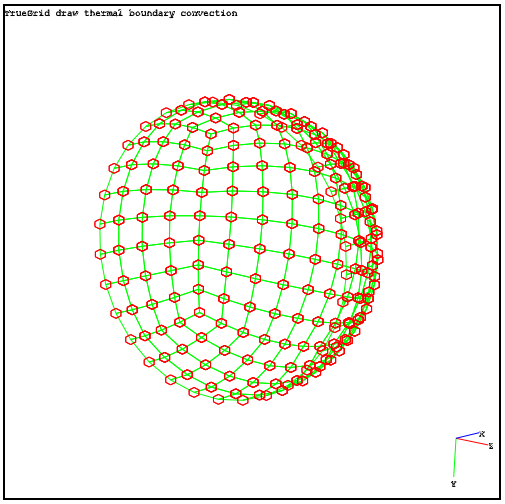

Figure 158 Wire frame CO CVT

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co tm **display initial temperature**

Example

```
sd 1 cy 0 0 0 0 0 1 5
sd 2 cy 0 0 0 0 0 1 15
sd 3 plan 0 0 0 0 1 0
sd 4 plan 0 0 0 1 0 0
sd 5 plan 0 0 0 0 0 1
sd 6 plan 0 0 10 0 0 1
block 1 11;
      1 11;
      1 3;
      0 10 0 10 0 10
tr 1 2 1 2 2 2 my -10
               rz 90
               my 5 ;
mb 1 1 1 2 1 2 x 5
sfi -1; ; ;sd 1
sfi -2; ; ;sd 2
sfi ;-1; ;sd 3
sfi ;-2; ;sd 4
sfi ; ;-1;sd 5
sfi ; ;-2;sd 6
tfi ; ;-1 0 -2;
tf 1 1 1 2 2 2
tm 2 1 1 2 2 2 100
endpart
merge
rx 20 ry 20
co tm
co size 2
```

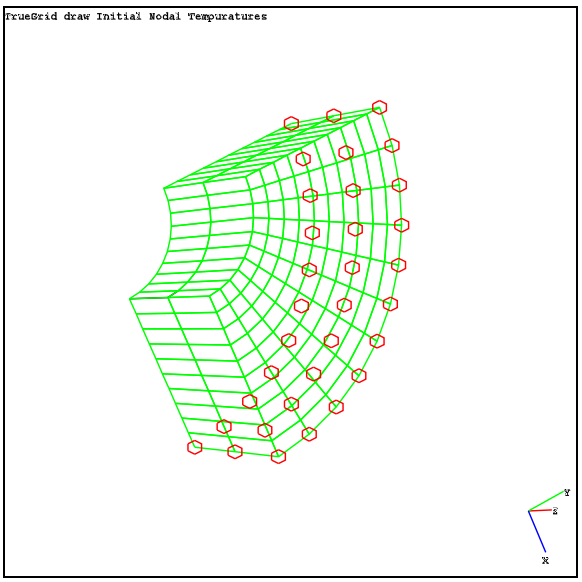

Figure 159 Wire frame version CO TM

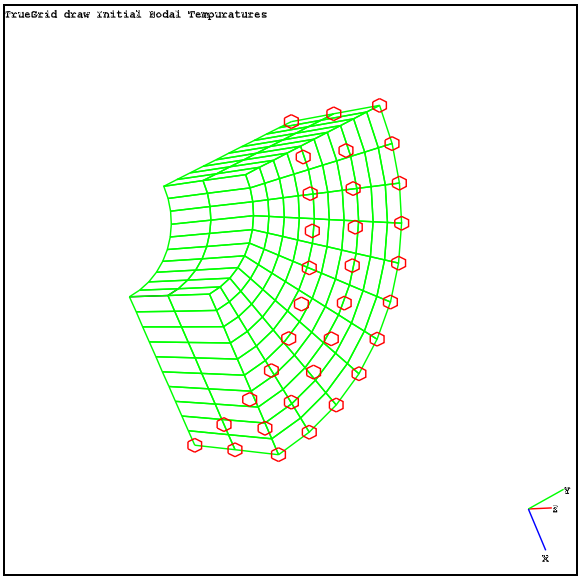

Figure 160 Hidden Line version CO TM

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co ft **display boundary temp**

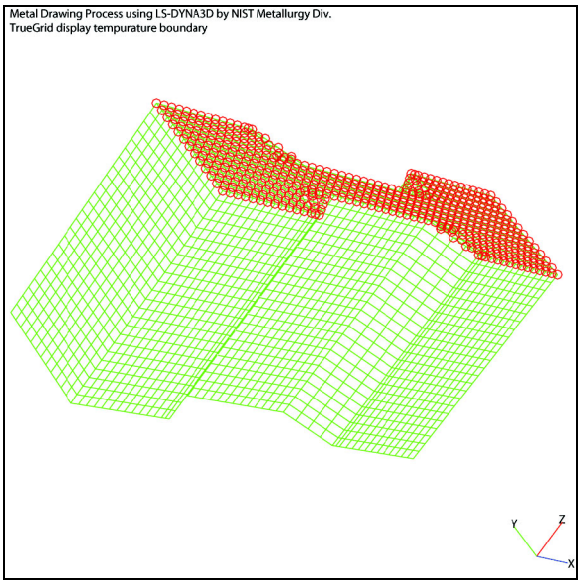

Figure 161 Hidden Line version CO FT

co fv **display nodal velocities**

Example

```
block 1 11; 1 11; 1 11; -1 1; -1 1;
-1 1;
sfi -1 -2; -1 -2; -1 -2; sp 0 0 0 1
lcd 1 0 0 .1 1 1 1;
fv 2 1 2 2 1 2 1 1 .57 -.57 .57 fv 2 2 2 2 2 2 1 1 .57 .57 .57
fv 1 2 2 1 2 2 1 1 -.57 .57 .57 fv 1 1 2 1 1 2 1 1 -.57 -.57 .57
merge co fv 1 co size 2
```

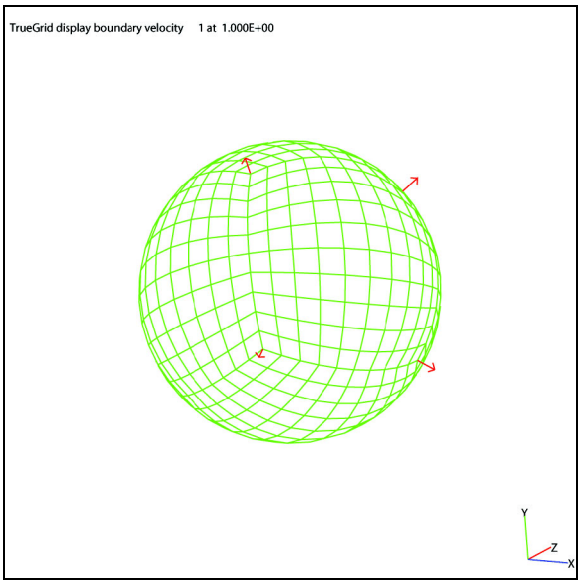

Figure 162 Hidden Line version CO FV

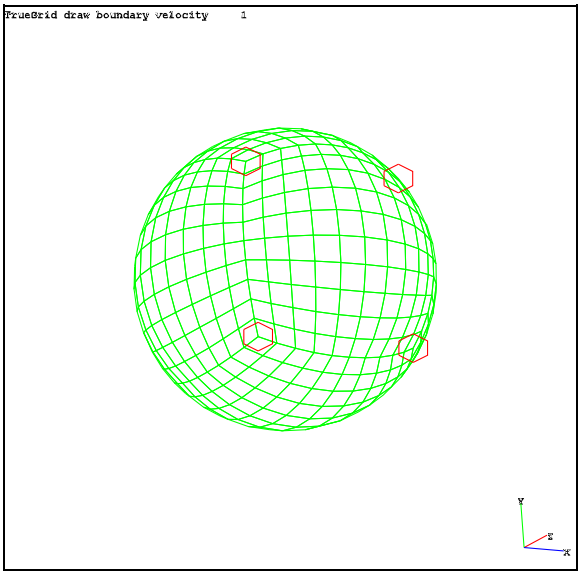

Figure 163 Wire frame version CO FV

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co sw **display stone wall**

Example

```
sd 1 cy 0. 0. 3.8 1 0 0 1.
sd 2 cy 0. 0. 3.8 1 0 0 1.6
sd 3 cy 0. 0. 5.8 0 1 0 .25
sd 4 cy 0. 0. 5.8 0 1 0 .68
sd 5 cy 0. 0. 3.8 0 1 0 .125
block 1 4 6 8 11 14 17 19;1 3 5 7 9;1
6 12 14 18 20 22 26 28 31 33 35
37;-2.5 -1.9 -1.7 -1.5 -1 -.6 -.2 0
-1.5 -1 -.7 -.3 0 0 .7 2.2 3.15 3.65
3.8 3.95 4.45 5.4 5.6 5.8 6 6.2
dei 1 5;;2 13;;
dei 5 6;;2 3 0 9 13;
dei 7 8;;5 7 0 10 12;
dei ;1 2;2 13;dei ;3 5;4 8;
dei ;4 5;7 13;dei 2 4;3 5;1 2;
sfi ;-3 5;-4 -8;sd 1
sfi ;-2 5;-3 -9;sd 2
sfi ;-2;9 13;plane 0 -1.1 0 0 1 0
sfi ;-4;8 13;plane 0 -.08 0 0 1 0
sfi ;-2;2 3;plane 0 -1. 0 0 1 0
sfi ;-3;8 13;plane 0 -.6 0 0 1 0
sfi ;-3;3 4 ;plane 0 -.6 0 0 1 0
sfi ;-4;3 4 ;plane 0 -.3 0 0 1 0
sfi -7 8; ; -10 -12; sd 3
sfi -6 8 ; ; 11 -13; sd 4
sfi -7 8 ; ; -5  -7; sd 5
sfi -6;;8 11;plan -.68 0 0 1 0 0
sfi -2 -4;-3 4;;cy -1.7 -.187 0 0 0 1
.125
sfi -2;4 5;;plan -1.825 0 0 1 0 0
sfi -4;4 5;;plan -1.575 0 0 1 0 0
relaxi 6 8;-2;9 13;50 .001 1
relaxi 6 8;-4;9 13;50 .001 1 relaxi -8;2 5;3 9;50 .001 1
relaxi 5 8;-2;3 9;50 .001 1 relaxi 5 8;-3;4 8;50 .001 1
relaxi 1 6;1 5;-1;50 .001 1 relaxi 1 6;1 5;-2;50 .001 1
relaxi -5;2 5;3 9;50 .001 1 lct 3 ryz;rxz;ryz rxz;
lrep 0 1 2 3;endpart merge stp .001 nset n1 = l 997 998 1650 1651 1668
1669 2986 2987 3004 3005 3657 3658;sw nset n1 1
```

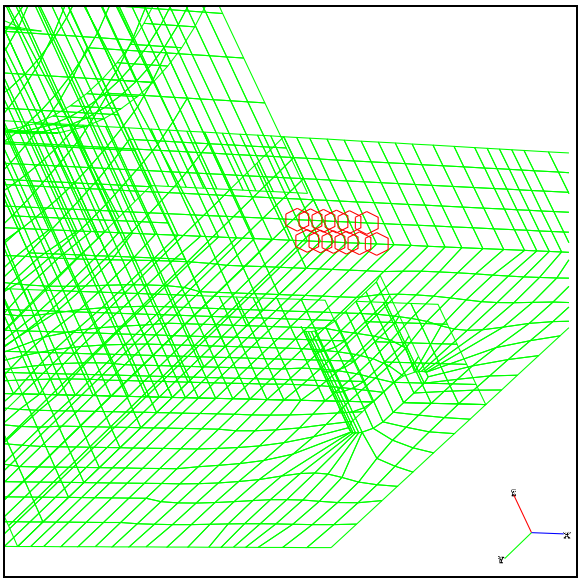

Figure 164 Wire frame version CO SW

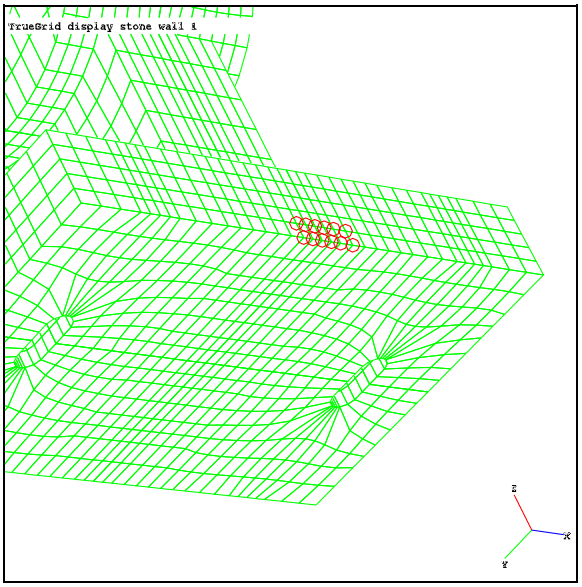

Figure 165 Hidden Line version CO SW

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co nr **display non-reflecting**

Remarks

The **co iss** command works the same way.

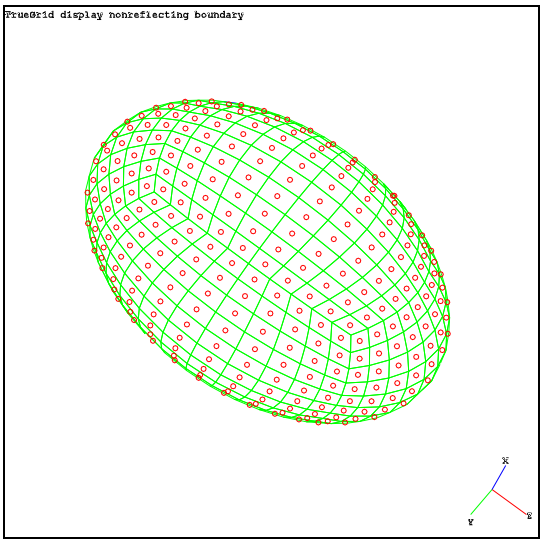

Figure 166 Hidden Line version CO NR

Example

```
block 1 11;1 11;1 11;
      -1 1;-1 1;-1 1;
sfi -1 -2;-1 -2;-1 -2;sp 0 0 0 1
z=1.5*z
nri -1 -2;-1 -2;-1 -2;merge
```

co ve **display initial velocities**

Example

```
block 1 3 5 7; 1 3 5 7;1 15;2 2 3 3;2 2
3 3;1 8;
sd 1 cyli 2.5 2.5 2.5 0 0 1 2 dei 1 2 0 3 4;1 2 0 3 4;;
sfi -1 -4; -1 -4;;sd 1 ve 1 1 2 4 4 2 .1 .1 1
```

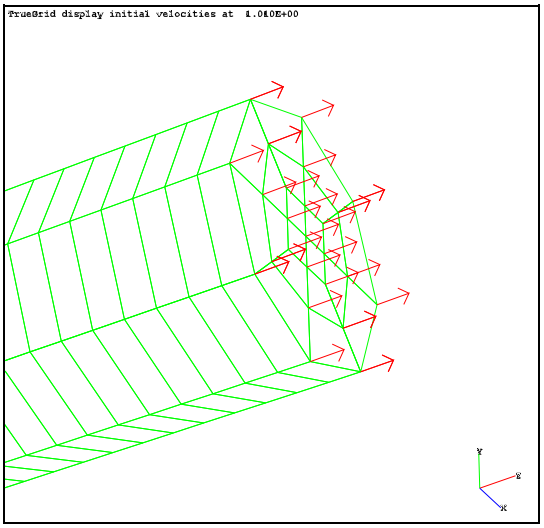

Figure 167 Hidden Line CO VE

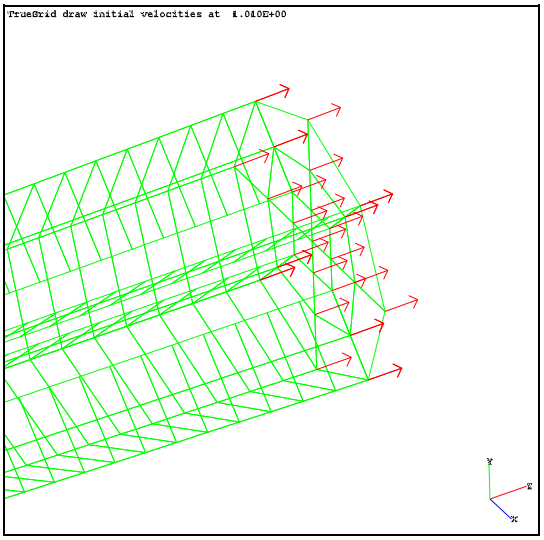

Figure 168 Wire frame CO VE

```
merge stp .0001 co ve co size 5
```

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co efl **display electric flux**

Example

```
ld 1 lp2 1 1;
  lfil 45 1 2 -45 .25
  lp2 1 2;
sd 1 crz 1
cylinder -1;
         1 4 7 10 91;
         1 23;
         1;
         0 12 24 36 360;
         1 2;
sfi -1;; 1 2;sd 1
efl 1 1 1 1 2 2 1.3
efl 1 2 1 1 3 2 2.3
efl 1 3 1 1 4 2 1.3
endpart
merge
rx -30 ry 30 zf 1.3 tvv
```

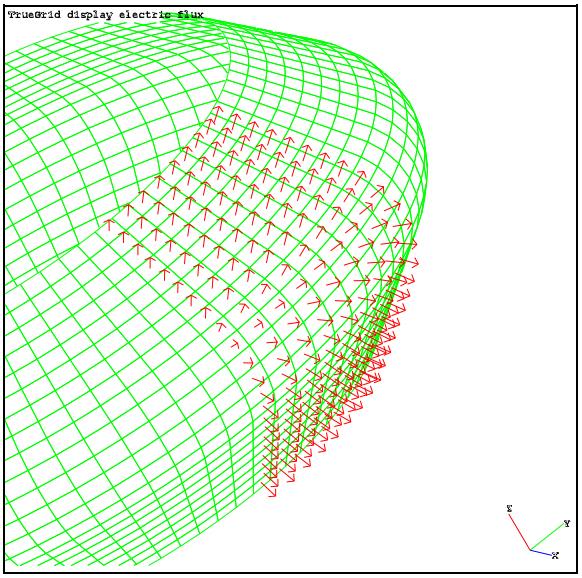

Figure 169 Hidden Line version CO EFL

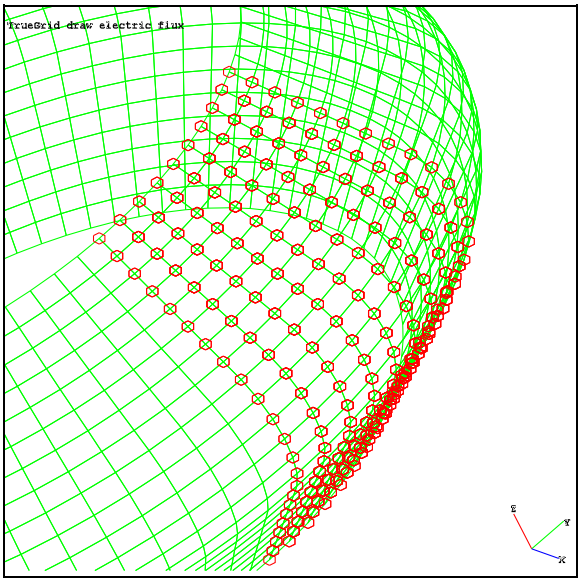

Figure 170 Wire frame version CO EFL

```
co efl co size 2.8
```

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co vhg **display volumetric heat generation**

Example

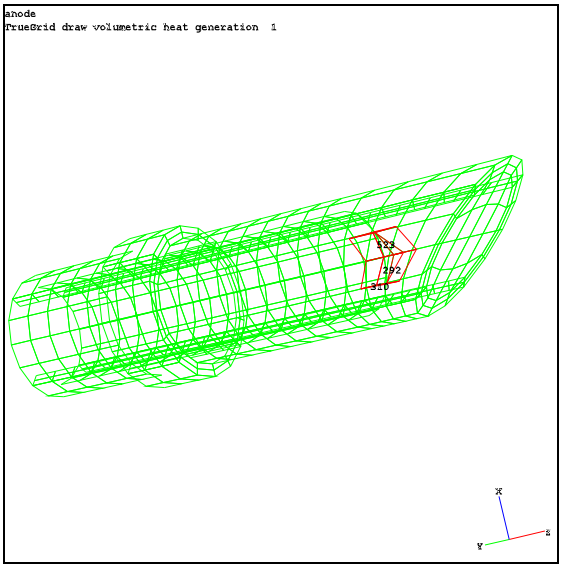

Figure 171 Wire frame version CO VHG

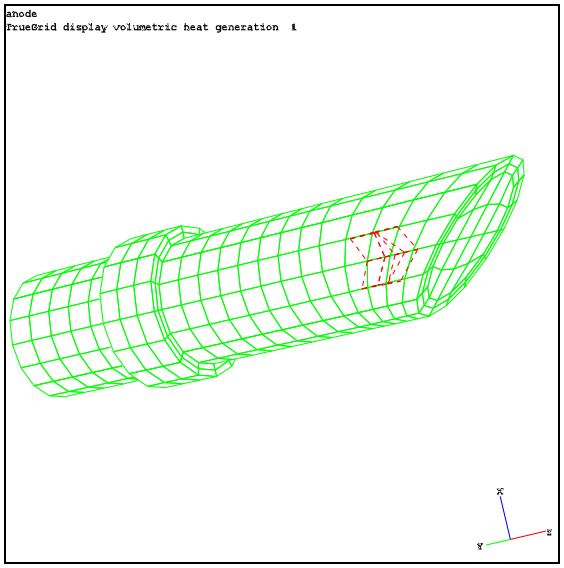

Figure 172 Hidden Line version CO VHG

This example is based on the anode.tg batch file in the EXAMPLES directory.

co sfb **display surface boundary conditions**

Remarks

The **co or** options works in a similar fashion.

Example

```
ld 1 csp2 00 1 -1 1.2 0 1 1; ;
sd 1 crz 1
cylinder -1;1 6;1 10;
         1 -30 30 -1 1
sfi -1;;;sd 1
sfb 1 1 1 1 2 2 surface j dy 1 ;
merge
```

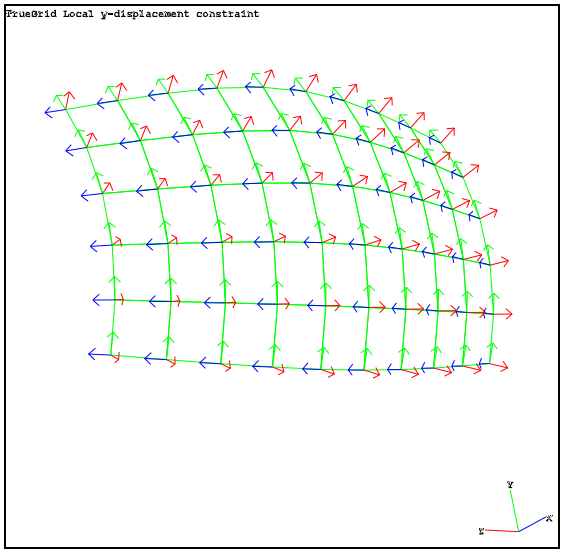

Figure 173 Wire frame version CO SFB

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co syf **display symmetry plane with failure**

Example

```
sd 1 sp 0 0 0 5
sd 2 sp 0 0 0 5.5
block 1 3 5 7 9 11 13;
      1 3 5 7 9 11 13;
      1 3 5 7 9 11 13;
  -2.5 -2.5 -2.5 0 2.5 2.5 2.5;
  -2.5 -2.5 -2.5 0 2.5 2.5 2.5;
  -2.5 -2.5 -2.5 0 2.5 2.5 2.5;
dei ;1 3 0 5 7; 1 3 0 5 7;
dei 1 3 0 5 7;; 1 3 0 5 7;
dei 1 3 0 5 7; 1 3 0 5 7;;
sfi -2 -6;-2 -6;-2 -6;sd 1
sfi -1 -7;-1 -7;-1 -7;sd 2
sfi -4;;;plan 0 0 0 1 0 0
de 1 0 0 4 0 0
plane 1 0 0 0 1 0 0 .0001 syf
syf 4 1 1 4 7 7 1 .9
merge
```

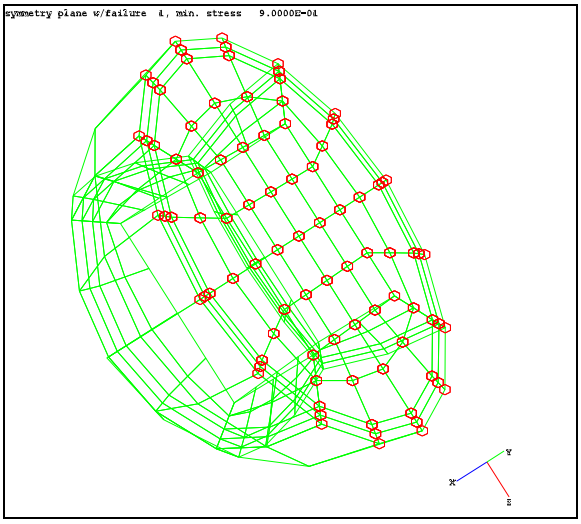

Figure 174 Wire frame version CO SYF

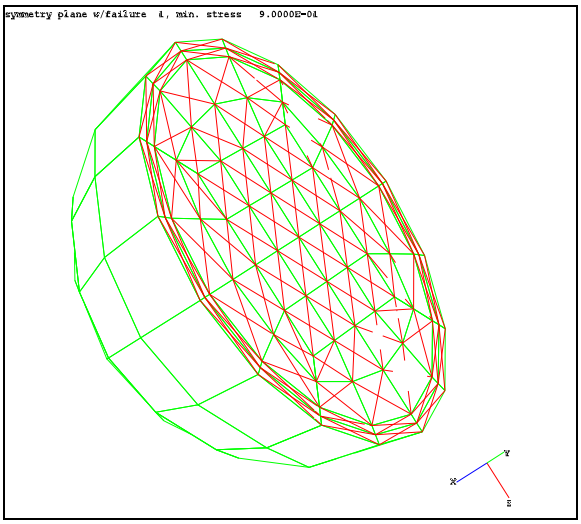

Figure 175 Hidden Line version CO SYF

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co bv **display boundary velocities**

Example

```
cyli 1 3 5 7 9;
     1 41;
     1 3 5 7 9;
     1 2 3 4 5;
     0 360;
     0 1 2 3 4;
bv 1 1 5 5 2 5 .3 0 .5
b 1 1 1 5 2 1 dx 1 dy 1 dz 1 ;
merge
co bv co size 2.9
```

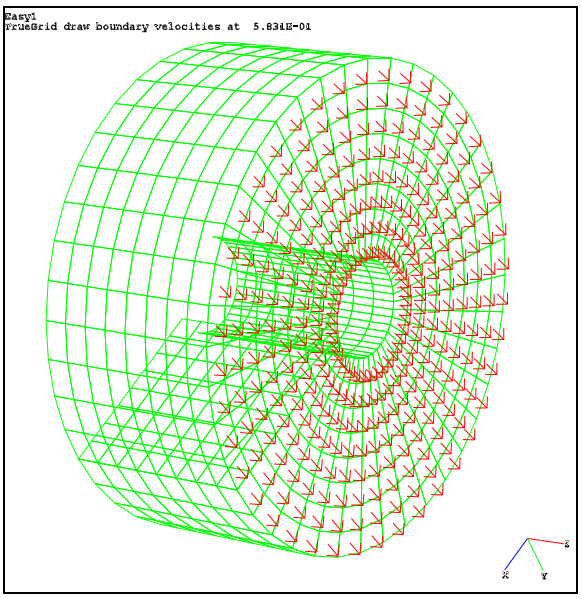

Figure 176 Wire frame version CO BV

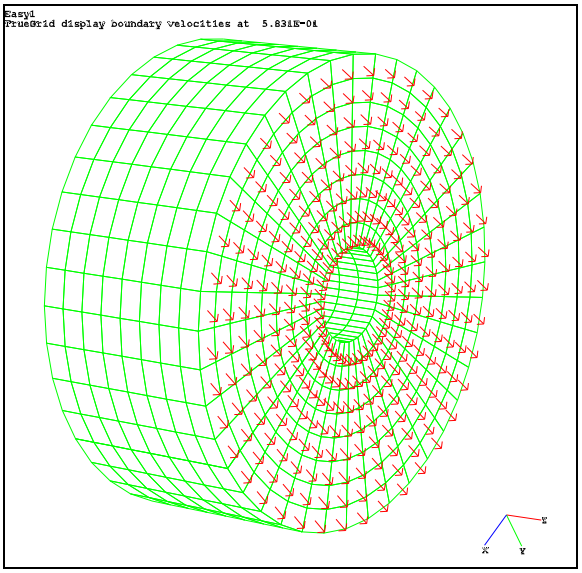

Figure 177 Hidden Line version CO BV

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co npb **display nodal print blocks**

Example

```
sd 1 sp 0 0 0 5
sd 2 plan 0 0 3 0 0 1
sd 3 plan 0 0 -3 0 0 1
block -1 -5 -9;
      -1 -5 -9;
       1 5 9;
       -1 0 1;
       -1 0 1;
       -1 0 1;
sfi -1 -3 ;-1 -3 ;1 3 ; sd 1
sfi ;; -3;sd 2
sfi ;; -1;sd 3
merge
nset nodes1 = 1 112 117 200 253 268;
npb nset nodes1
```

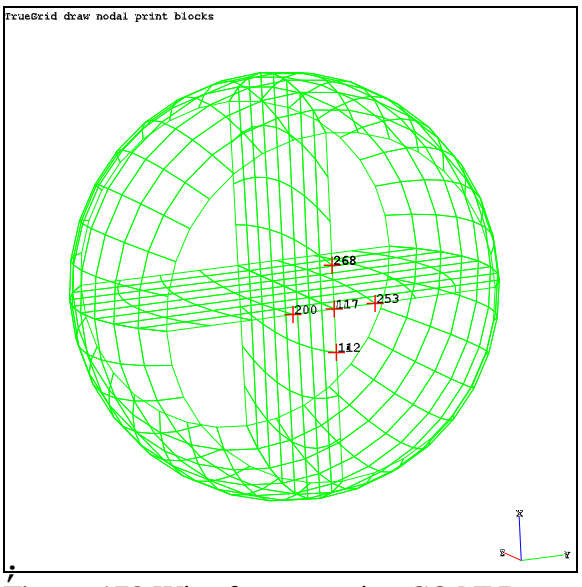

Figure 178 Wire frame version CO NPB

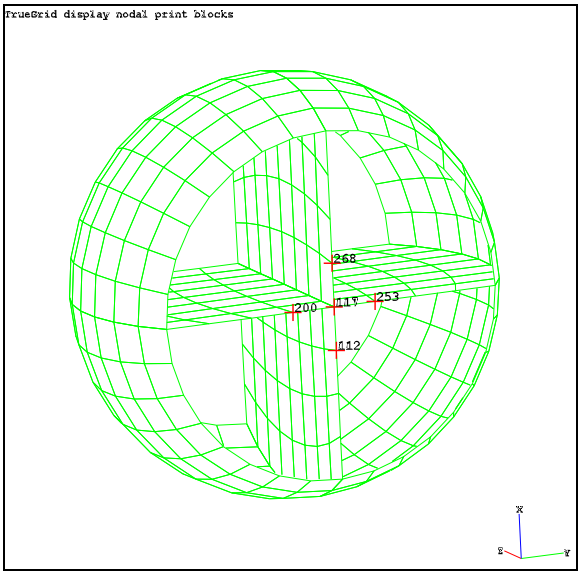

Figure 179 Hidden Line version CO NPB

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Example

```
sd 1 sp 0 0 0 5
sd 2 plan 0 0 3 0 0 1
sd 3 plan 0 0 -3 0 0 1
block -1 -5 -9;-1 -5 -9;1 5 9;
       -1 0 1;-1 0 1;-1 0 1;
sfi -1 -3 ;-1 -3 ;1 3 ; sd 1
sfi ;; -3;sd 2
sfi ;; -1;sd 3
merge
eset elements1 rpl ls
 4 20 65 132 177 289 324 353;
epb s elements1
```

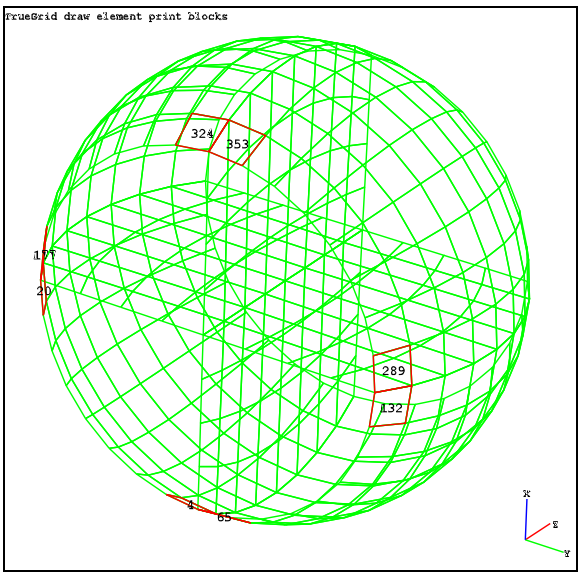

Figure 180 Wire frame version CO EPB

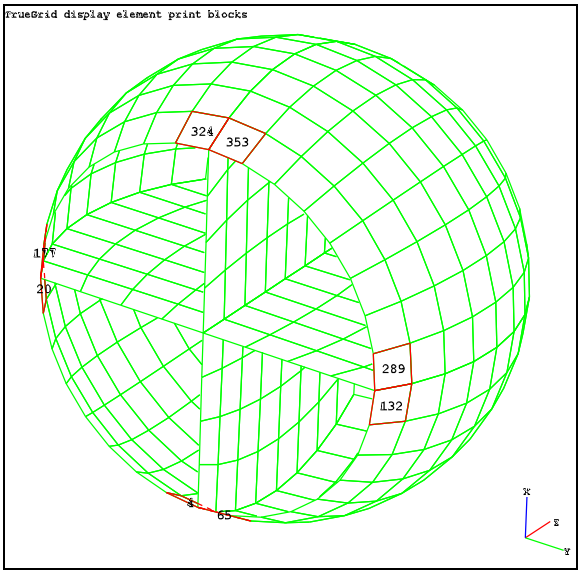

Figure 181 Hidden Line version CO EPB

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co sp **display springs**

Example

```
cylinder 1 3;1 226;1 3;
   1.5 2.5 0 900 -.5 .5
sfi -1 -2;1 2;-1 -2;ts 0 0 0 0 0 1 2
0 1
z=z+0.05*j
jbm 1 1 1 2 2 2 1 1 1 i 1 ;
mate 0
merge
spring 1 n1 1 n2 448 sddn 1 ; ;
```

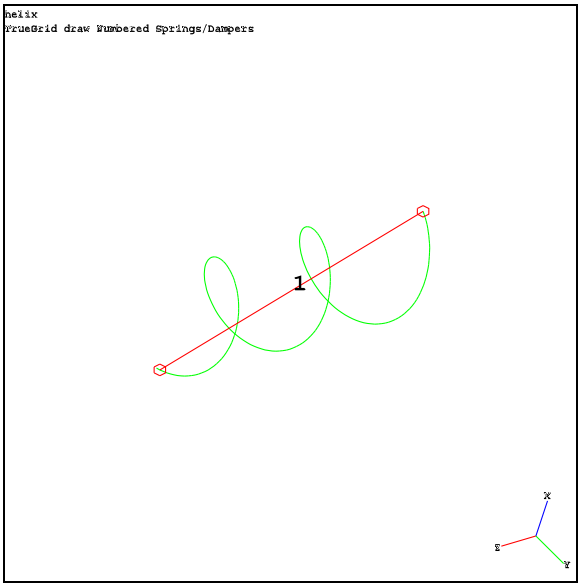

Figure 182 Wire frame version CO SP

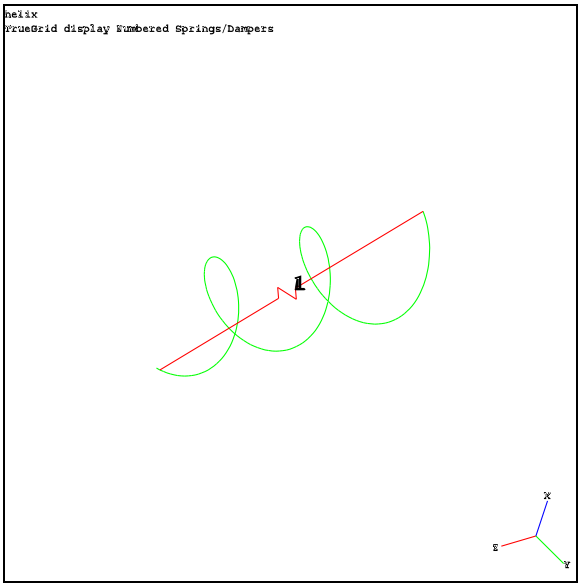

Figure 183 Hidden Line verison CO SP

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co pm **display point masses**

Example

```
sd 1 cy 0 0 0 0 0 1 1
sd 2 cy 0 0 0 0 1 0 1
sd 3 plan 0 0 0 0 1 0
sd 4 plan 0 0 0 0 1 1
sd 5 plan 0 0 0 0 1 -1
block 1 3 5 7;1 3 5 7;1 3;
-.3 -.3 .3 .3 -.3 -.3 .3 .3 1 2
dei 1 2 0 3 4; 1 2 0 3 4;;
sfi -1 -4; -1 -4;;sd 1
insprt 1 4 2 1
sfi ; -3;;sd 3
sfi -1 0 -4;; -1;sd 2
sfi 2 3; -1 0 -5; -1;sd 2
sfi ; 1 3; -1;sd 4
sfi ; 3 5; -1;sd 5
sd 6 plan .3 -.3 2 0 1 0
sd 7 plan -.3 .3 2 0 1 0
sd 8 plan .3 -.3 2 1 0 0
sd 9 plan -.3 .3 2 1 0 0
sfi -2; 2 4;;sd 9
sfi -3; 2 4;;sd 8
sfi 2 3; -2;;sd 6
sfi 2 3; -4;;sd 7
mseq i 4 6 4
mseq j 4 3 3 4
mseq k 6
lct 3 rx 90;rx 180;rx 270;
lrep 0 1 2 3 ;
pm 1 3 2 1 3 2 1 ;
endpart
merge
stp .001
pm 37 2 ;
npm 1 -2 0 0 2 ;
bm n1 37 n2 9262 ;
```

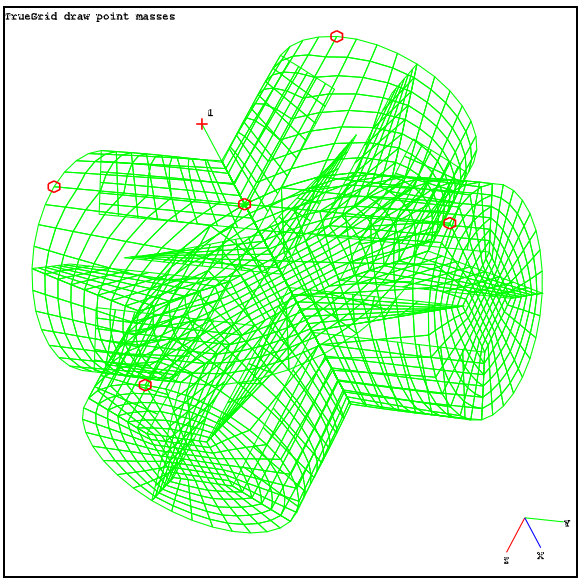

Figure 184 Wire frame version CO PM

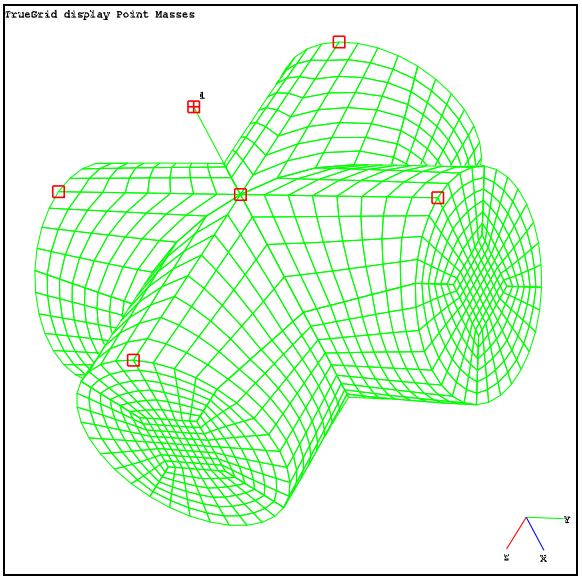

Figure 185 Hidden Line version CO PM

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Example

```
sd 1 function -1 1 -1 1 u+v ;
              u-v ;
               (u+v)*(u-v) ; ; ;
sid 1 inter;
block 1 11;1 11;1 3 4;
     -1 1 -1 1 -1 1 1
tr 1 1 1 2 2 3 rz 45;
patch 1 1 2 2 2 2 1
bb 1 1 2 2 2 2 1;
bb 1 1 1 2 2 1 1 mz .5;
bb 1 1 3 2 2 3 1 normal .3;
si 1 1 2 2 2 2 1 m ;
mti ;;2 3;2
block 1 11;1 11;1 3;-1 1 -1 1 1 5
tr 1 1 1 2 2 2 rz -45;
bb 1 1 1 2 2 1 1 normal .3;
bb 1 1 2 2 2 2 1 normal .9;
si 1 1 1 2 2 1 1 s ;
merge
```

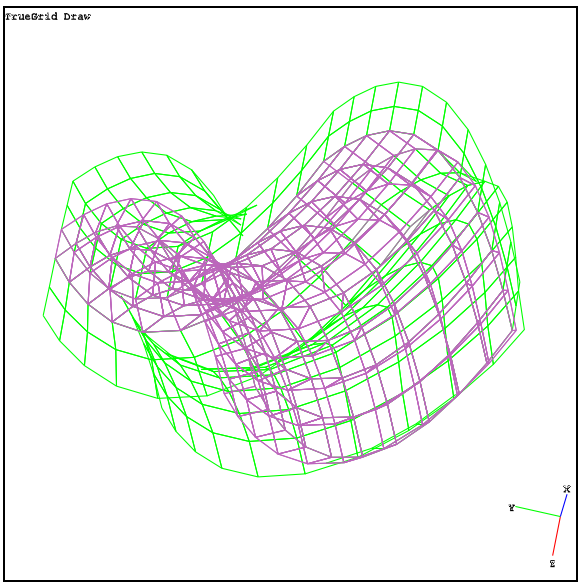

Figure 186 Wire frame version CO Interfac

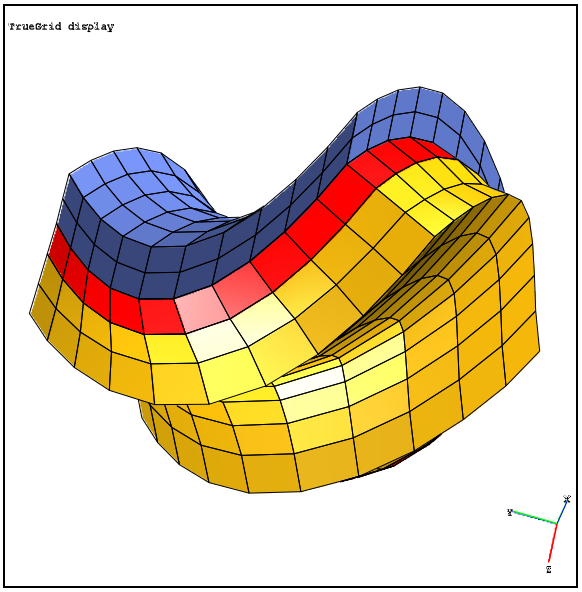

Figure 187 Fill verison CO Interfac

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co thic **display shell thicknesses**

Example

```
ld 1 lep 1 2 3 0 110 180 0 ;
sd 1 crz 1
sd 2 cy 0 0 0 0 0 1 1.9
sd 3 intp 1 2 .5 ;
block -1 -11;-1 -11;1 6;
          -2 2 -2 2 0 1
ssfi -1 -2; -1 -2;;3
edge 1 1 1 2 1 1 3.3
edge 1 2 1 2 2 1 3.3
edge 2 1 1 2 2 1 3.3
edge 2 1 1 2 2 1 3.3
edge 2 1 2 2 2 2 3.1
edge 1 1 2 1 2 2 3.1
edge 1 1 2 2 1 2 3.1
edge 1 2 2 2 2 2 3.1
merge
```

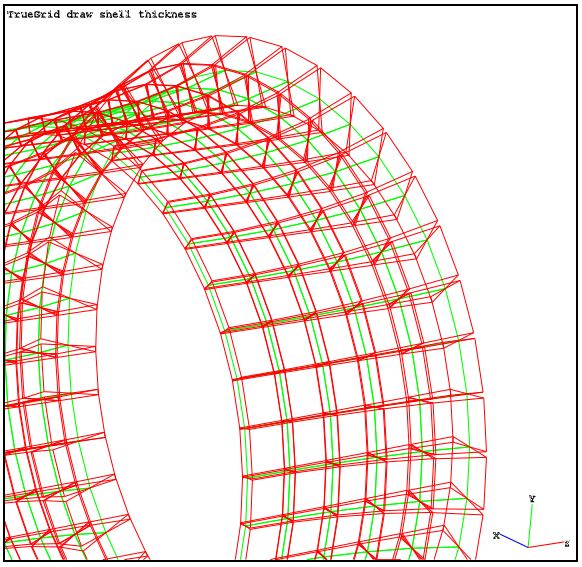

Figure 188 Wire frame version CO THIC

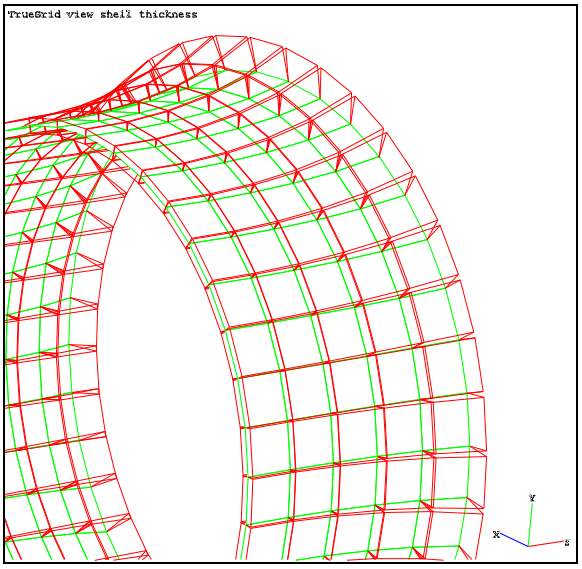

Figure 189 Hidden Line version CO THIC

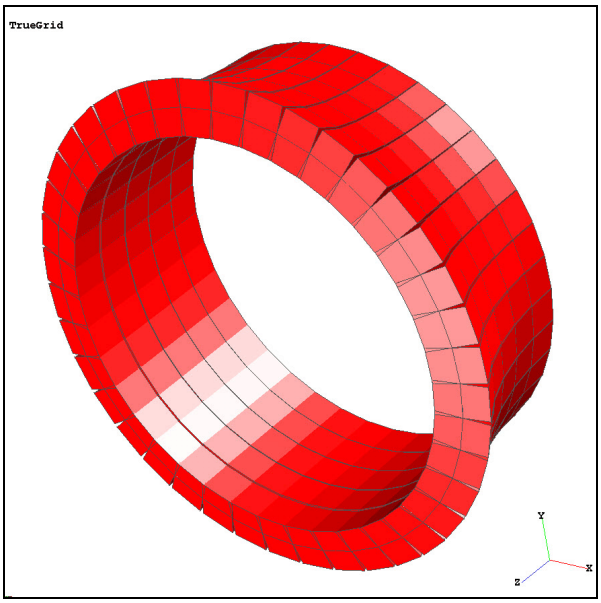

Figure 190 Fill version of CO THIC

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Example

```
sd 1 sp 0 0 0 1
sd 2 sp 0 0 0 2
sd 3 plan 0 0 0 -1 0 1
sd 4 plan 0 0 0 1 0 1
sd 5 plan 0 0 0 -1 1 0
sd 6 plan 0 0 0 1 1 0
block 1 5;1 5;1 5;
      1 2 -1 1 -1 1
sfi -1;;;sd 1
sfi -2;;;sd 2
sfi ;;-2;sd 3
sfi ;;-1;sd 4
sfi ;-2;;sd 5
sfi ;-1;;sd 6
lcd 1 0 0 1 1;
tepro   1   1   1   2   2   2   1
1.5*sqrt((x-2)^2+y*y+z*z)     ;
atan2(y,x) ;
merge
```

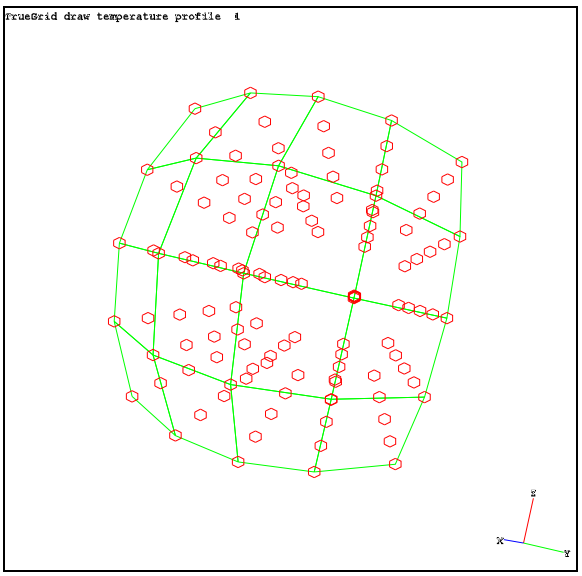

Figure 191 Wire frame verison CO TEPRO

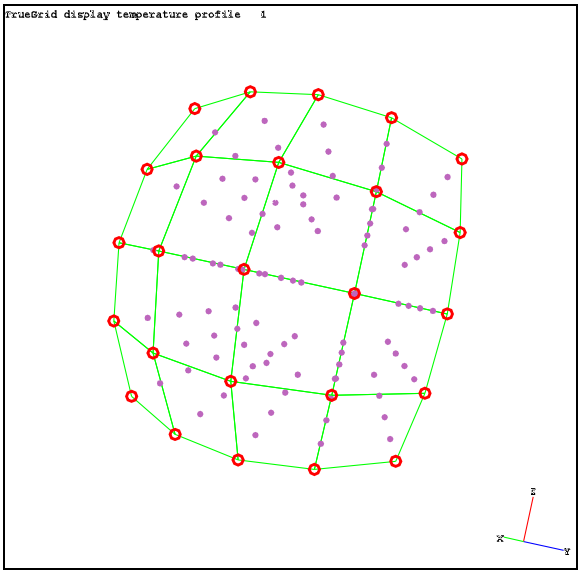

Figure 192 Hidden Line version CO TEPRO

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co sc **display smoothing constraints**

Example

```
sd 1 cy 0 0 0 0 0 1 2
block -1 2 5 -6;1 4;1 4;
      -1 -1 1 1 -1 1 -1 1
sfi -1 0 -2 0 -3 0 -4;;;sd 1
sc 2 1 1 3 2 2 i
merge
```

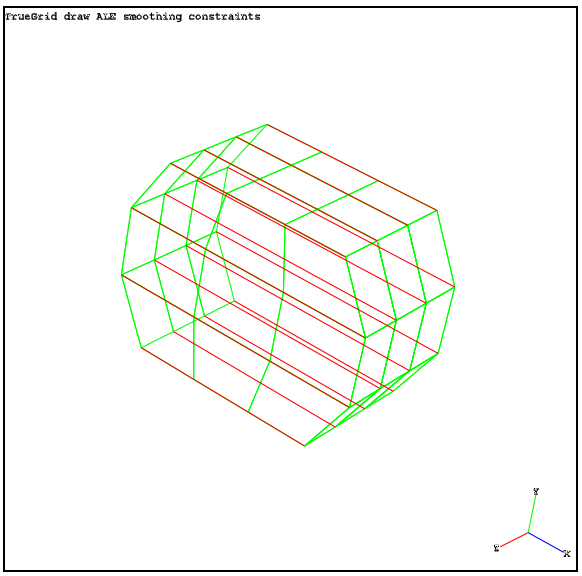

Figure 193 Wire frame version CO SC

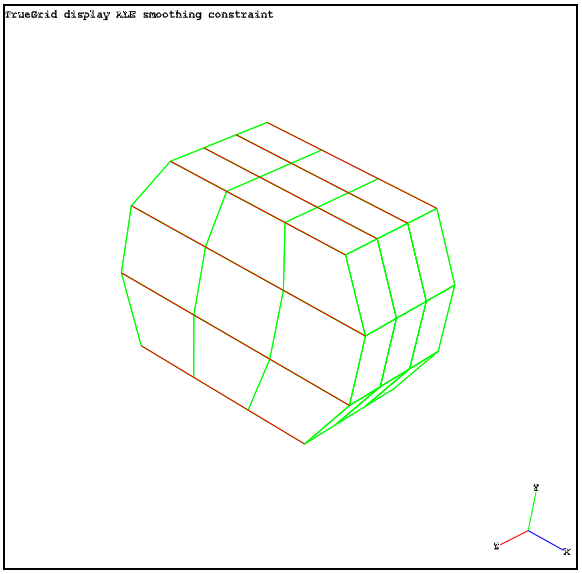

Figure 194 Hidden Line version CO SC

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co n **display shell normals**

Example

```
block -1 -6;-1 -6;-1 -6;
      -1 1 -1 1 -1 1
sd 1 sp 0 0 0 1
sfi -1 -2; -1 -2; -1 -2;sd 1
orpt - 0 0 0
n 1 1 1 1 2 2
n 2 1 1 2 2 2
n 1 1 1 2 1 2
n 1 2 1 2 2 2
n 1 1 1 2 2 1
n 1 1 2 2 2 2
merge
```

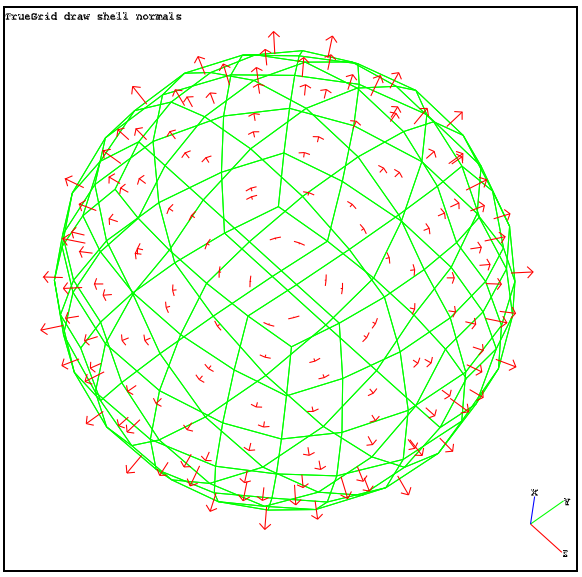

Figure 195 Wire frame version CO N

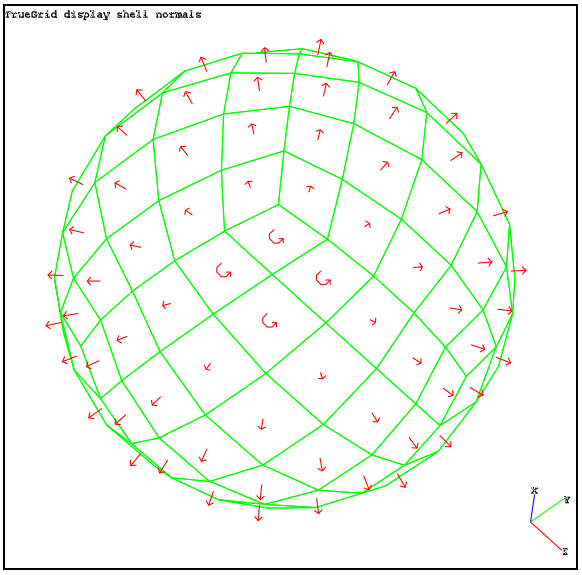

Figure 196 Hidden Line version CO N

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co resn **display radiation enclosures labeled**

Example

```
block 1 3 5 7 9;
      1 3 5 7 9;
      1 3 5 7 9;
      -1 -1 0 1 1;
      -1 -1 0 1 1;
      -1 -1 0 1 1;
dei 1 2 0 4 5; 1 2 0 4 5;;
dei 1 2 0 4 5;;1 2 0 4 5;
dei ;1 2 0 4 5;1 2 0 4 5;
sfi -1 -5;-1 -5;-1 -5;sp 0 0 0 5
sfi -2 -4;-2 -4;-2 -4;sp 0 0 0 3
de 2 2 2 4 4 4
de 1 0 0 3 0 0
orpt  + 0 0 0
rei 3 -4;-2 -4;-2 -4;0 2.2 no
merge
```

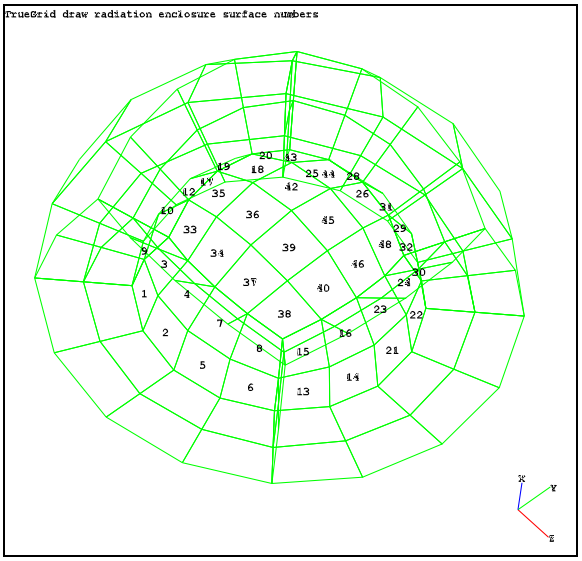

Figure 197 Wire frame version CO RESN

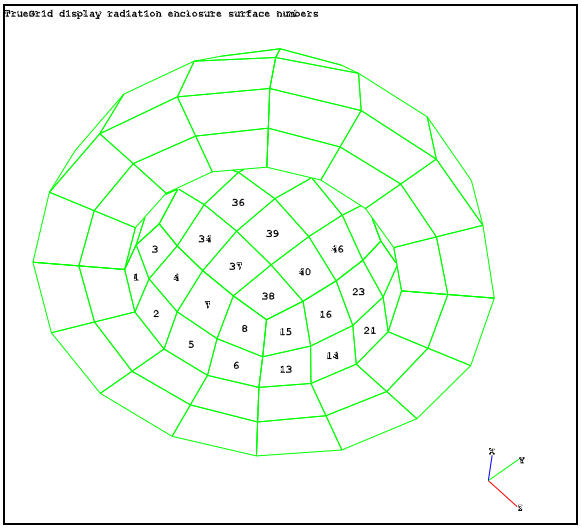

Figure 198 Hidden Line version CO RESN

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co grtol **display glued interface for CFX**

Remarks

See the **supblk** command.

Example

A butterfly mesh of the fluid in the pipe is created. The **supblk** command was used to define grid 1. The interface 2 for grid 1 was displayed by the **co grtol** command (**Figure 199**). The simplified command file follows:

... transformations, projections of mesh ...

```
supblk 1 1 2 3 4 3

merge
```

... display of part 1...

```
co grtol 1 2   c display interface 2 for grid 1
```

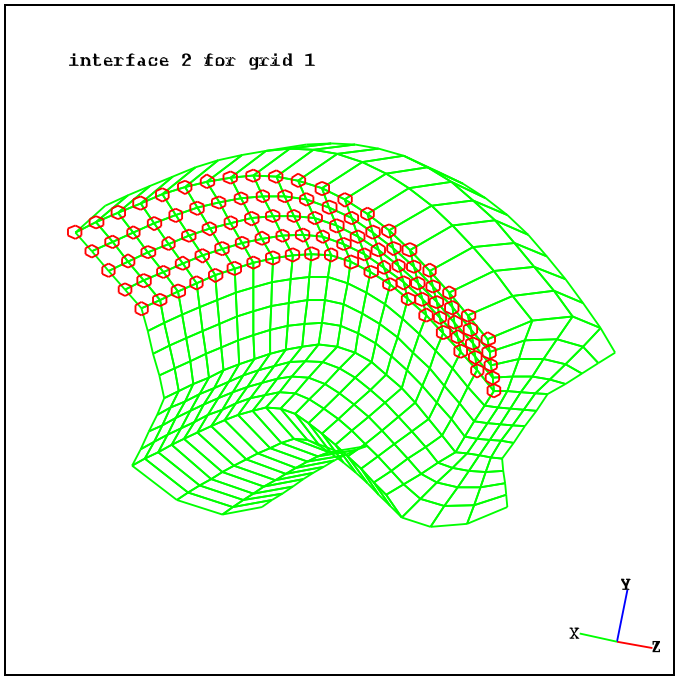

Figure 199 interface between grids

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Example

```
sd 1 plan 6 0 0 1 0 0
sd 2 plan 8 0 0 1 0 0
sd 3 plan 0 6 0 0 1 0
sd 4 plan 0 8 0 0 1 0
sd 5 cy 7 7 7 0 0 1 7
sd 6 sp 7 7 2 7
sd 7 sp 7 7 12 7
lcd 1 0 0 1 1;
block 1 6 7 9 10 15;
      1 6 7 9 10 15;
      1 6 10 14 18 23;
      5 5 6 8 9 9;
      5 5 6 8 9 9;
      -2 -2 2 12 16 16;
dei 1 2 0 5 6;1 2 0 5 6;;
dei 1 2 0 5 6;;1 2 0 5 6;
dei ;1 2 0 5 6;1 2 0 5 6;
dei 3 4;3 4;;
sfi -3;3 4;;sd 1
sfi -4;3 4;;sd 2
sfi 3 4;-3;;sd 3
sfi 3 4;-4;;sd 4
sfi -1 -6;-1 -6;3 4;sd 5
sfi -1 -6;-1 -6;-1 3;sd 6
sfi -1 -6;-1 -6;4 -6;sd 7
mom 3 3 1 4 4 1 1 1 x
endpart
merge
```

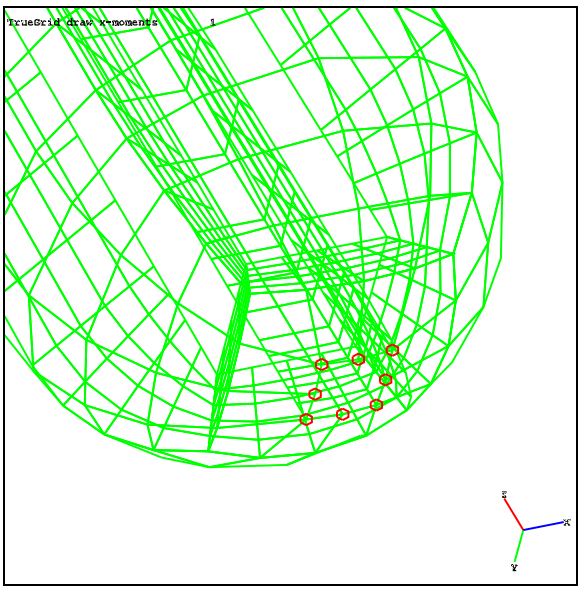

Figure 200 Wire drawing with moments

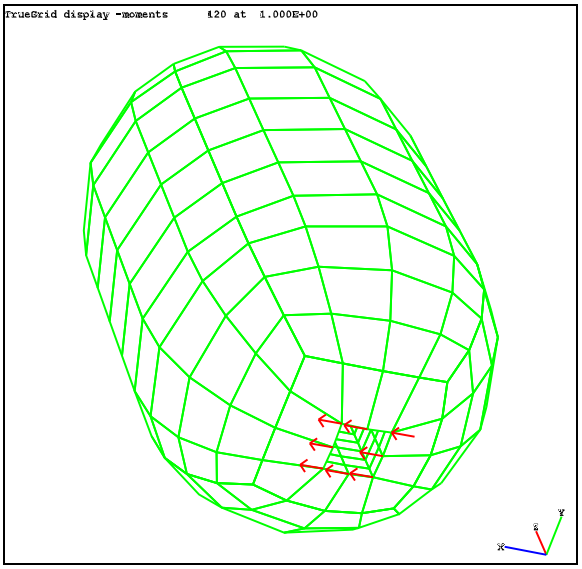

Figure 201 Hidden view with moments

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Example

This example takes two files. The second one is named row and is included in the first file.

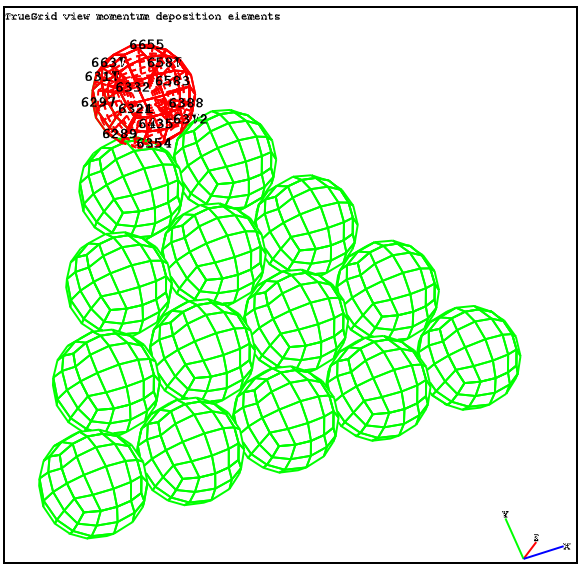

Figure 202 Parametric replication

```
sd 1 sp 0 0 0 3
para n 5
     x0 0
     y0 0;
include row
para n 4
     x0 3
     y0 [3*tan(60)];
include row
para n 3
     x0 [2*3]
     y0 [2*3*tan(60)];
include row
para n 2
     x0 [3*3]
     y0 [3*3*tan(60)];
include row
para n 1
     x0 [4*3]
     y0 [4*3*tan(60)];
include row
```

The second file, called row, contains the following commands:

```
block 1 5 9 13;1 5 9 13;1 5 9 13;
-1 -1 1 1 -1 -1 1 1 -1 -1 1 1
dei 1 2 0 3 4;1 2 0 3 4;;
dei 1 2 0 3 4;;1 2 0 3 4;
dei ;1 2 0 3 4;1 2 0 3 4;
sfi -1 -4;-1 -4;-1 -4;sd 1
if(%n.eq.1)then
 mdep 1 1 1 4 4 4 .2 2 0 .01
endif
gct 1 mx %x0 my %y0;grep 1;
lct %n mx 6;repe %n;lrep 1:%n;
```

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```
endpart
```

co sy **display symmetry plane conditions**

Example

```
accuracy 3

ld 1 lp2 0 0 1 1 2 4;;
     lp2 3 4 4 1 5 0;;
sd 1 crz 1
sd 2 plan 0 0 1 0 0 1
sd 3 plan 0 0 2.75 0 0 1
sd 4 cy 0 0 0 0 0 1 2.5
sd 5 plan 0 0 0 1 1 0
sd 6 plan 0 0 0 -1 1 0
sd 7 plan 0 0 0 0 0 1
block -1 -9 15 -21 -29;
      -1 -9 15 -21-29;
      1 5 -11;
      -5 -1 0 1 5
      -5 -1 0 1 5
      0 2 4
dei -2 -4;-2 -4;-1 -2;
dei -2 0 -4;1 2 0 4 5;;
dei 1 2 0 4 5;-2 0 -4;;
dei 2 4; 2 4; -3;
dei -1 2 0 4 -5;-1 2 0 4 -5;1 -3;
sfi -1 -2 -4 -5; -1 -2 -4 -5; 1 -3;sd 1
sfi ;; -2;sd 2
sfi -1 0 -5; 2 4; -3;sd 3 sfi 2 4; -1 0 -5; -3;sd 3
sfi -2 0 -4; 2 4; -3;sd 4 sfi 2 4; -2 0 -4; -3;sd 4
sfi -2; 4 5; 1 3;sd 5 sfi 1 2; -4; 1 3;sd 5
sfi -4; 1 2; 1 3;sd 5 sfi 4 5; -2; 1 3;sd 5
sfi -2; 1 2; 1 3;sd 6 sfi 1 2; -2; 1 3;sd 6
sfi -4; 4 5; 1 3;sd 6 sfi 4 5; -4; 1 3;sd 6
sfi ;; -1;sd 7 dei ; 1 3;;
plane 1 0 0 0 0 1 0 .001 symm ;
endpart
merge
grid on
```

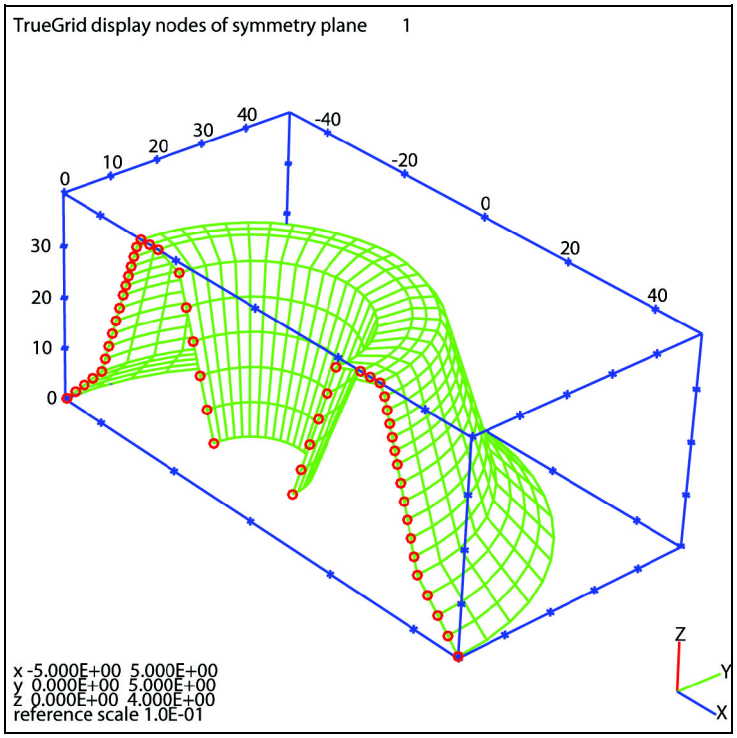

Figure 203 Boundary conditions from symmetry plane

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co ol **display outlets**

Remarks

The **co il** command works identically to this command.

Example

```
sd 1 cyli 2.5 0 2.5 0 1 0 1
block 1 5 11 17 23 27;
      1 5 9;
      1 5 11 17 23 27;
      1 2 2 3 3 4;
      1 3 5;
      1 2 2 3 3 4;
dei 1 2 0 5 6;2 3;;
dei ; 2 3;1 2 0 5 6;
dei 2 3 0 4 5;2 3;2 3 0 4 5;
dei 2 3 0 4 5;;1 3 0 4 6;
dei 1 2 0 5 6;;2 3 0 4 5;
mbi -1 0 -6;;-4 0 -5;z .25
mbi -1 0 -6;;-2 0 -3;z -.25
mbi -2 0 -3;;-1 0 -6;x -.25
mbi -4 0 -5;;-1 0 -6;x .25
sfi -2 -5;;-2 -5;sd 1
bb 2 1 4 3 3 4 1;
bb 3 1 4 3 3 5 1;
bb 4 1 4 4 3 5 2;
bb 4 1 4 5 3 4 2;
bb 4 1 3 5 3 3 3;
bb 4 1 2 4 3 3 3;
bb 3 1 2 3 3 3 4;
bb 2 1 3 3 3 3 4;
esm 3 3 2 4 3 5 & 2 3 3 3 3 4 & 4 3 3 5 3 4 100 .00001 1 .6 4
esm 2 1 3 5 1 4 & 3 1 4 4 1 5 & 3 1 2 4 1 3 100 .00001 1 .6 4
unifm 3 1 2 4 3 5 & 2 1 3 3 3 4 & 4 1 3 5 3 4 200 .00001 1
il 1 3 1 6 3 6
ol 1 1 1 6 1 6
endpart
merge
```

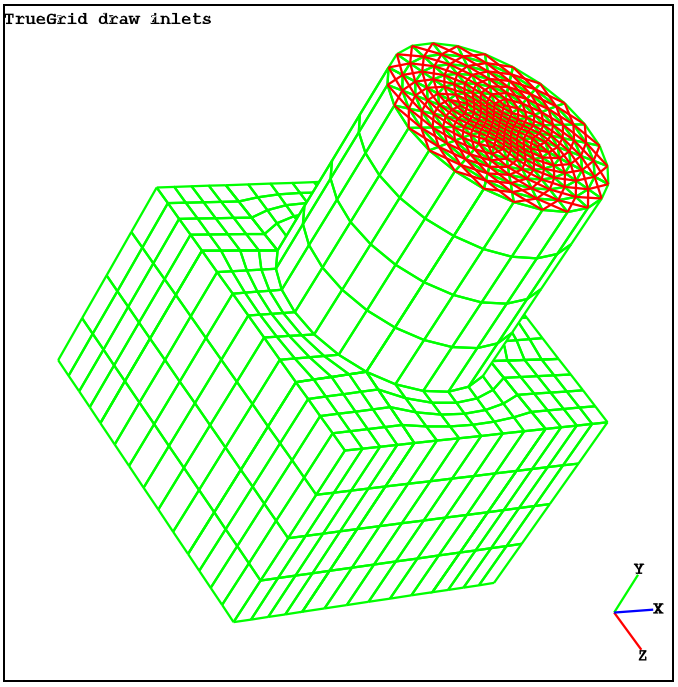

Figure 204 Inlet marked in hide mode

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co spw **display numbered spot welds**

Example

```
sd 1 cy 1.1 1.1 0 0 0 1 1
lmi 1
block -1 6 21;-1 6;1 21;
      .1 1.1 5 .1 1.1 0 5
sfi -1 2; -1 2; 1 2;sd 1
y=y+.1*sin((k-1)*36)
lct 1 ryz; lrep 0 1;
mate 1
endpart
block -1;1 11 21;1 21;
       0 -2.9 1.1 5.1 0 5
dom 1 2 1 1 2 2
y=y+.1*sin((k-1)*36)
mate 3
endpart
merge
spw 21 spwd 1 node 672 node 1323 node 126 ;;
```

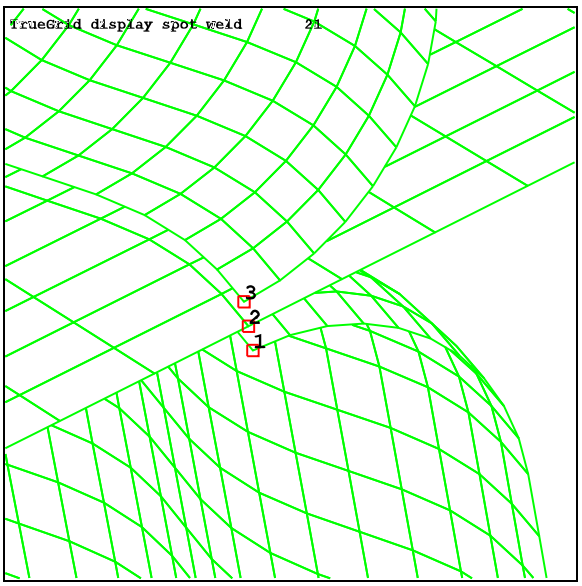

Figure 205 nodes of a numbered spotweld

co spwf **display facial spot welds**

Remarks

If the node defining the spot weld is not visible, a yellow box is draw.

Example

These commands are added to the previous example.

```
spwf rt -.1 1 4.875 3 mat 1 mat 2
spwf rt .1 .93 2.15 3 mat 1 mat 2
```

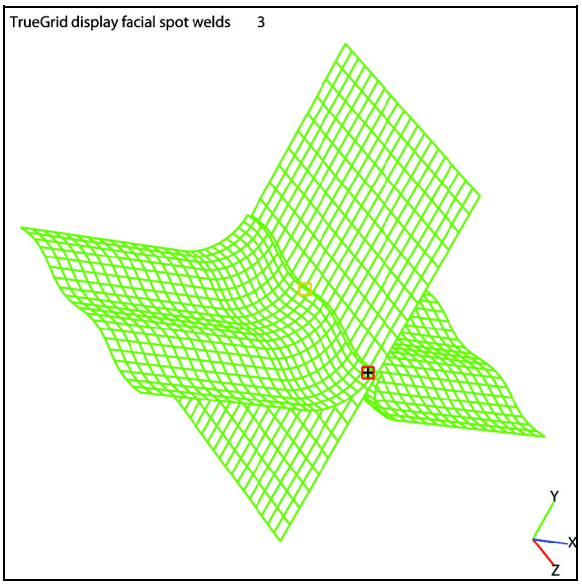

Figure 206 2 Facial spot welds

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co jt **display a numbered joint**

Example

```
cylinder 1 3;1 13;1 6;
         1 2 0 360 0 5
cylinder 1 3;1 13;1 6;
         1 2 0 360 5.5 10.5
merge
bm n1 42 n2 6 n3 180
   mate 2 nbms 2 ;
bm n1 289 n2 253 n3 235
   mate 2 nbms 2 ;
jd 1 uj pnlt .33 ;
jt 1 1 n 470 ;
jt 1 2 n 469 ;
jt 1 3 n 253 ;
jt 1 4 n 42 ;
```

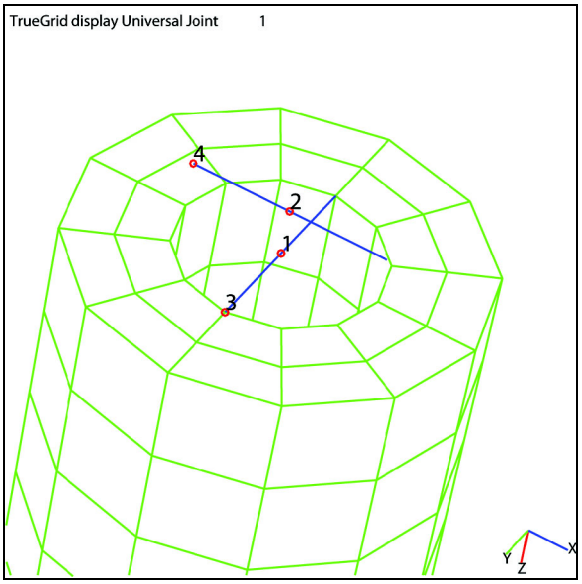

Figure 207 Nodes forming a joint

co dem **display discrete element spheres**

```
demicle
block 1 3;1 3;1 3;1 3;1 3;1 3;
merge
```

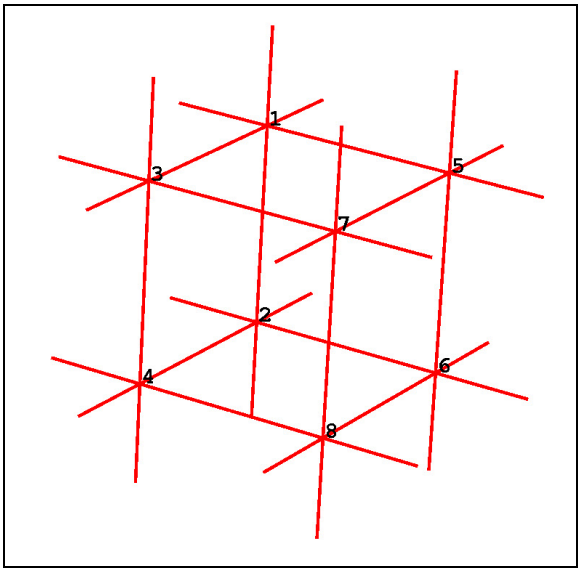

Figure 208 Discrete Element Spheres

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co mpce **multiple point constraint equations**

```
block 1 6 11 16;1 6 11 16;1 3;
      -3 -1 1 3 -3 -1 1 3 0 1
dei 2 3; 2 3;;
sd 1 cy 0 0 0 0 0 1 1
sfi -2 -3; -2 -3;;sd 1
merge
curd 1 arc3 whole rt 1 0 0
     rt 0 1 0 rt -1 0 0 ;
curd 2 arc3 whole rt 1 0 .5
     rt 0 1 .5 rt -1 0 .5 ;
curd 3 arc3 whole rt 1 0 1
     rt 0 1 1 rt -1 0 1 ;
nset one = crv 1 .01
nset two = crv 2 .01
nset three = crv 3 .01
mpce 1 rt 0 0 0 ; dx nsets one dx
; ;
mpce 2 rt 0 0 .5 ; dx nsets two dx
; ;
mpce 3 rt 0 0 1 ; dx nsets three
dx ; ;
co mpce
```

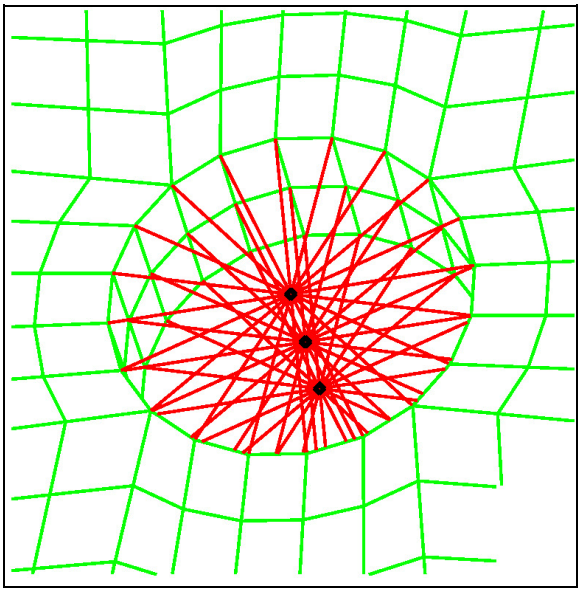

Figure 209 Multi Point Constraint Equations

labels specify type of label to be displayed

labels *option*
or
la *option*
where *option* can be any of:

off	to turn off labels display
size *scale*	to scale the size of the tokens displaying labels
angle θ	to change the cone angle for the tokens displaying labels
nodes	to label nodes by node number
1d	to label 1D beam elements by element number
2d	to label 2D shell elements by element number
2q	to label 2D quadratic shell elements by element number
3d	to label 3D brick element faces by element number

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3q	to label 3D quad'c brick element faces by element number
ijk1	for labeled and color coded reduced indices
ijk2	for labeled and color coded parts and reduced indices
ijk4	for color coded indices
locnd *node_number*	to locate a node by its number
loc1d *element_number*	to locate a 1D beam element by its number
loc2d *element_number*	to locate a 2D shell element by its number
loc3d *element_number*	to locate a 3D brick element by its number
loc2dq *element_number*	to locate a 2D quadratic shell element by its number
loc3dq *element_number*	to locate a 3D quadratic brick element by its number
sd	to display numbers of defined surfaces
crv	to display numbers of defined 3D curves
sdedge	to display surface edge identification numbers
sdpt	to display labels of points on defined surfaces
crvpt	to display labels of points on defined 3D curves
bb	to display the master block boundary interfaces
nodeset *set_name*	to display a node set
onset *set_name first last*	to display an ordered node set
faceset *set_name*	to display a face set
facesel *set_name*	to display and label a face set
elemset *set_name*	to display an element set
parts	to display parts
tol $part_1$ $part_2$	to display nodes merged between parts
where	
$part_1$	part sequence number (0 means all)
$part_2$	part sequence number (0 means all)
fraces	free faces (only actively displayed parts and materials)
fredges	free edges (only actively displayed parts and materials)
cracks θ	cracks below a specified subtended angle

Remarks

Labels do not overlap. If an object is not labeled, zoom in on it so there is more room for the label. Also try rotating the picture. The tokens used for a feature in the model may be different in the wire, hide, and fill graphics mode. These differences can used to your advantage. For example. when nodes are labeled in the wire mode, then nodes in the back of the mesh may also be labeled. Or, if the mesh is drawn in hide, only visible nodes are labeled. In another example, when using the fraces (free faces) option in wire mode, all free faces will be drawn in red and will stand out even in a large mesh. In hide mode, the hidden edges of free faces will appear as dotted lines. In fill graphics, they appear as red faces.

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labels nodes **label nodes by node number**

Example

Part 1 is composed of linear shells and Part 2 is composed of quadratic shells. The nodes are labeled in **Figure 210**. There are coincident nodes between parts, which are not merged.

```
linear
block 1 5;1 5;-1;1 5;1 5;0;
endpart
quadratic
block 1 5;1 5;-1;5 10;1 5;0;
merge
labels nodes
```

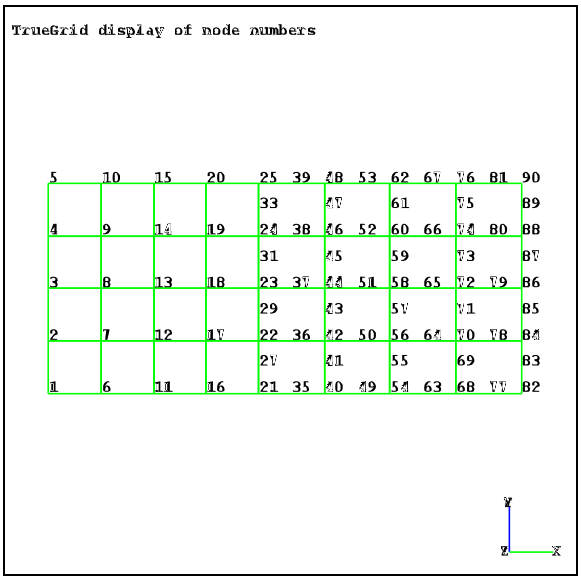

Figure 210 labeled nodes

labels 1d **label linear 1D beam elements by element number**

Example

20 beam elements are created along a curve. The elements are labeled. The simplified command file follows:

```
curd 1 arc3 seqnc rt -1 0 0 rt 0 1
0 rt 1 0 0 ;
curd 2 lp3 -1 -1 0 -1 0 0; ;
curd 3 lp3 1 -1 0 1 0 0; ;
curd 4 cpcds 2 1 3; ;
bm rt1 -1 -1 0 ; rt2 1 -1 0 ;
orient 0 0 0 mate 1 cs .1 nbms 20
cur 4 ;
labels 1d center
```

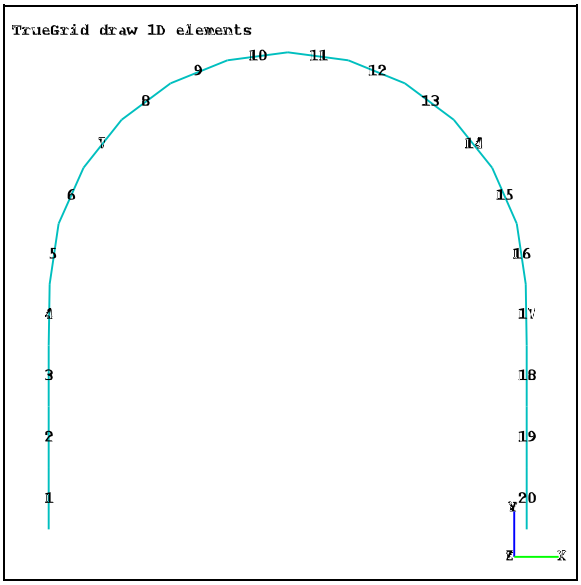

Figure 211 labeled beams

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labels 1q **label quadratic 1D beam elements by element number**

labels 2d **label 2D shell elements by element number**

Example

This example uses the mesh from the example for the **labels nodes** command. The numbering of elements always starts with 1 for each type of element.

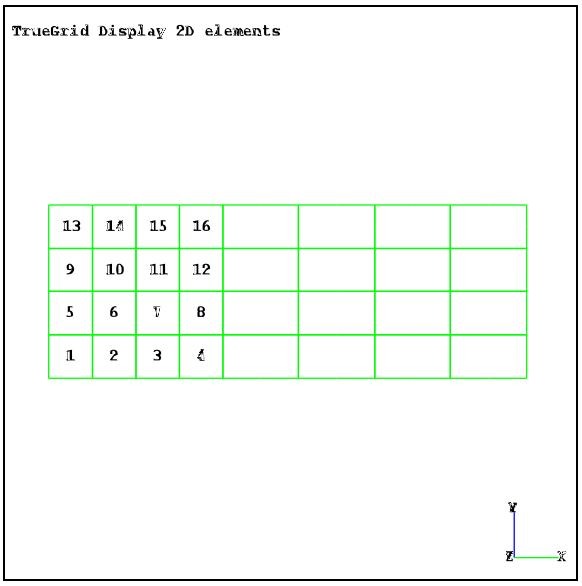

Figure 212 labeled linear shells

labels 2q **label 2D quadratic shell elements by element number**

Example

This example uses the mesh from the example for the **labels nodes** command. The numbering of elements always starts with 1 for each type of element.

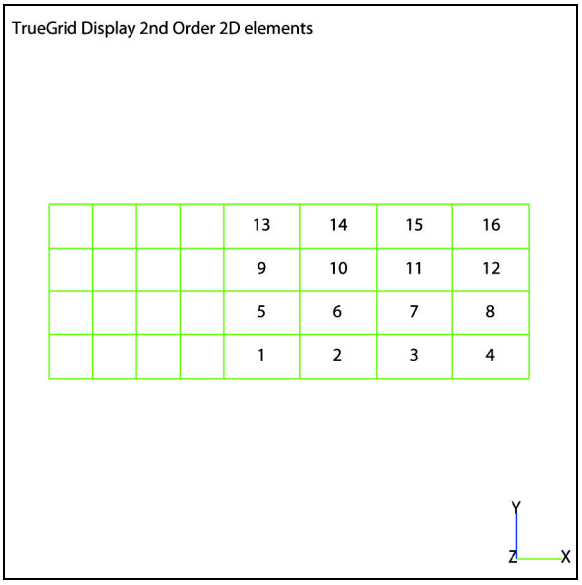

Figure 213 labeled quadratic shells

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labels 3d **label 3D brick faces by element number**

Example

Part 1 is composed of linear bricks and Part 2 is composed of quadratic bricks. The numbering of elements always starts with 1 for each type of element. The linear bricks are labeled by the element numbers. The simplified command file follows:

```
linear
block 1 5;1 5;1 5;1 5;1 5;1 5;
endpart
quadratic
block 1 5;1 5;1 5;5 12;1 5;1 5;
merge
labels 3d
```

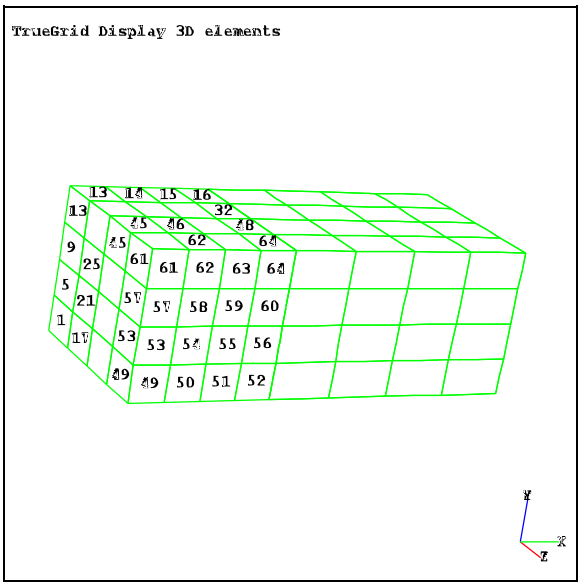

Figure 214 labeled bricks

labels 3q **label 3D quadratic brick faces by element number**

Example

This example uses the mesh from the example for the **labels 3d** command. The numbering of elements always starts with 1 for each type of element. The quadratic bricks are labeled by the element numbers.

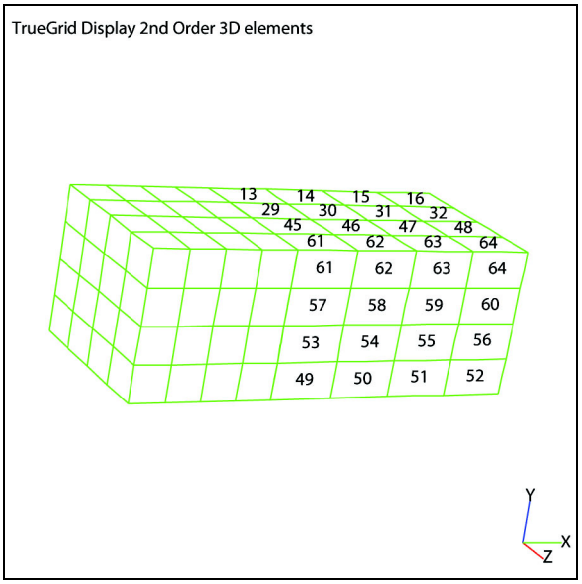

Figure 215 labeled quadratic bricks

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labels ijk1 **display color coded reduced indices**

Remarks

labels ijk1 can be used after issuing the Hide option in the Merge Phase. (I.e., this works when looking at a hidden line drawing.)

You can view the part information in the Merge Phase in a similar manner as in the Part Phase.

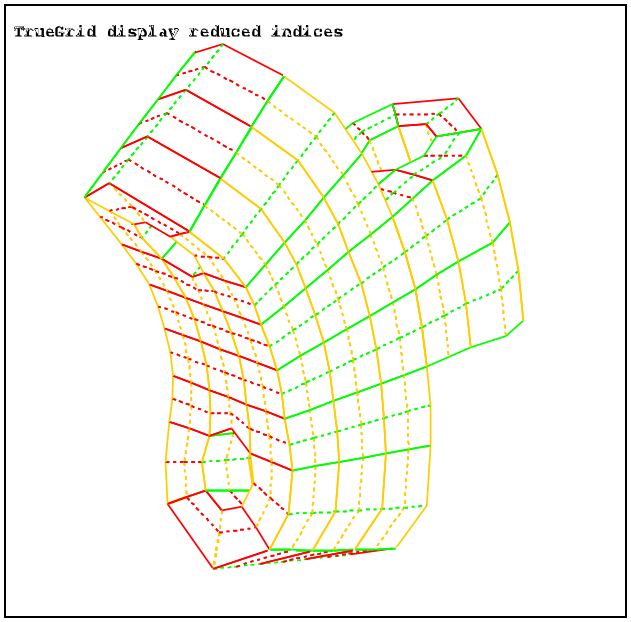

Figure 216 color coded reduced indices

labels ijk2 **display color coded and numbered reduced indices**

Remarks

Labels ijk2 can be used after issuing the Hide option in the Merge Phase.

You can view the part information in the Merge Phase in a similar manner as in the Part Phase.

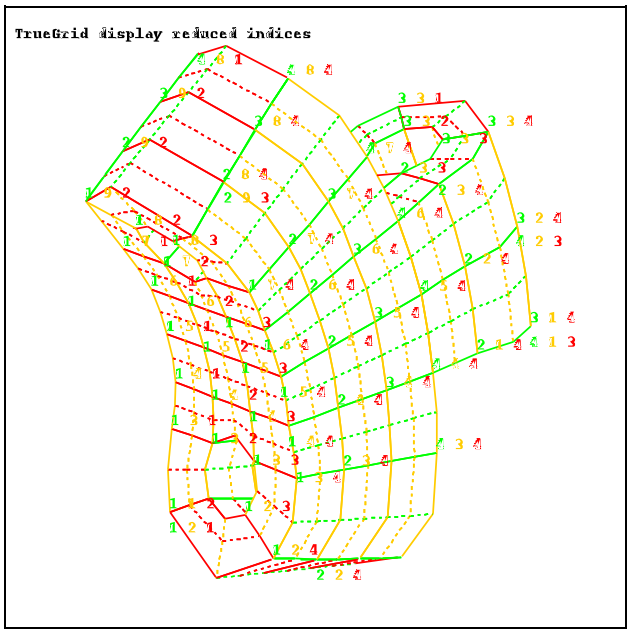

Figure 217 numbered reduced indices

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labels ijk4 **display color coded/numbered reduced indices/parts**

Remarks

labels ijk4 can be used after issuing the Hide option in the Merge Phase.

You can view the part information in the Merge Phase in a similar manner as in the Part Phase.

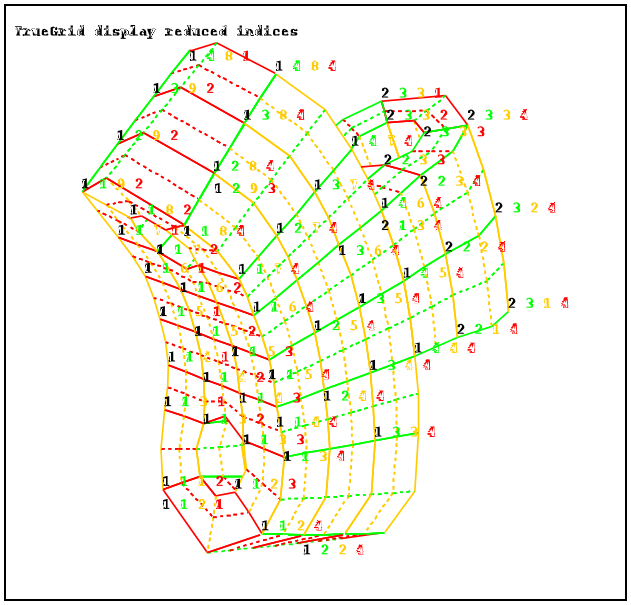

Figure 218 reduced indices and parts

labels locnd **locate a node by its number**

locnd *node_number*

Example

This example uses the mesh from the example for the **labels nodes** command. The node 20 is located. The located node is in the center of the picture by default.

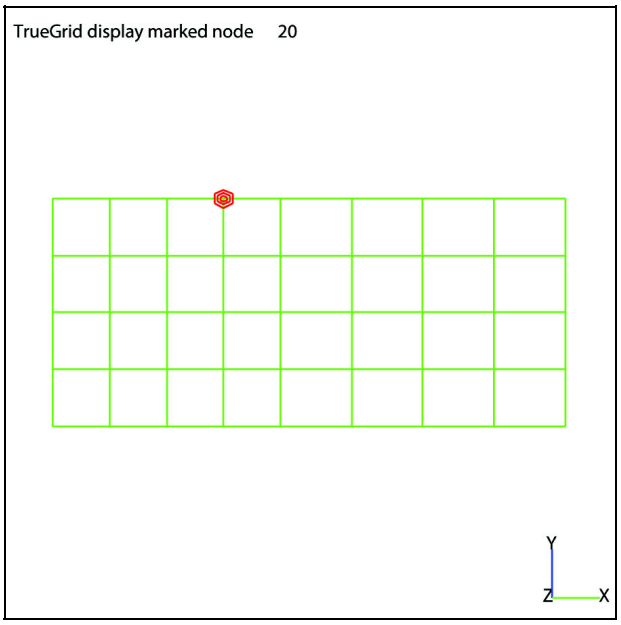

Figure 219 located node

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labels loc1d **locate a linear 1D beam element by its number**

loc1d *element_number*

Example

This example uses the mesh from the example for the **labels 1d** command. The beam 5 is located. The located beam is in the center of the picture by default.

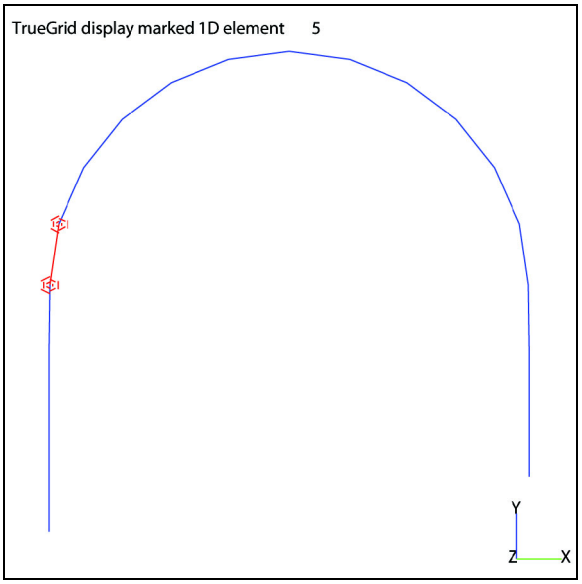

Figure 220 located beam

labels loc1dq **locate a quadratic 1D beam element by its number**

labels loc2d **locate a 2D shell element by its number**

loc2d *element_number*

Remarks

The diagonals of a shell are displayed to help check distorted elements.

Example

This example uses the mesh from the example for the **labels 2d** command. The shell number 10 is located. The located shell is in the center of the picture by default.

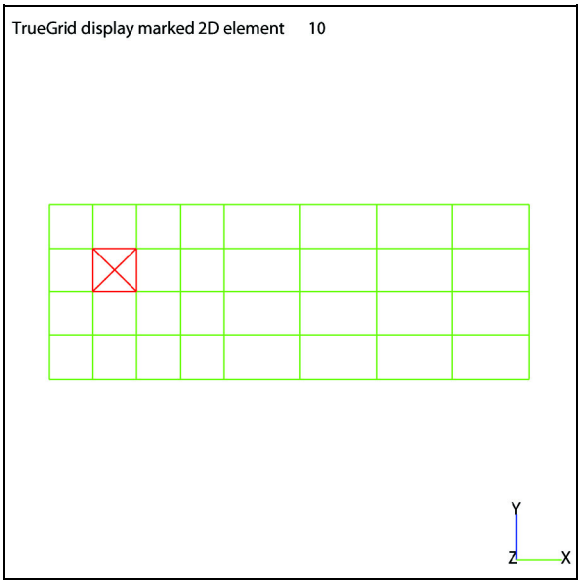

Figure 221 located linear shell

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labels loc2dq **locate a 2D quadratic shell element by its number**

loc2dq *element_number*

Remarks

The diagonals of a shell are displayed to help check distorted elements.

Example

This example uses the mesh from the example for the **labels 2d** command. The quadratic shell number 10 is located. The located shell is in the center of the picture by default.

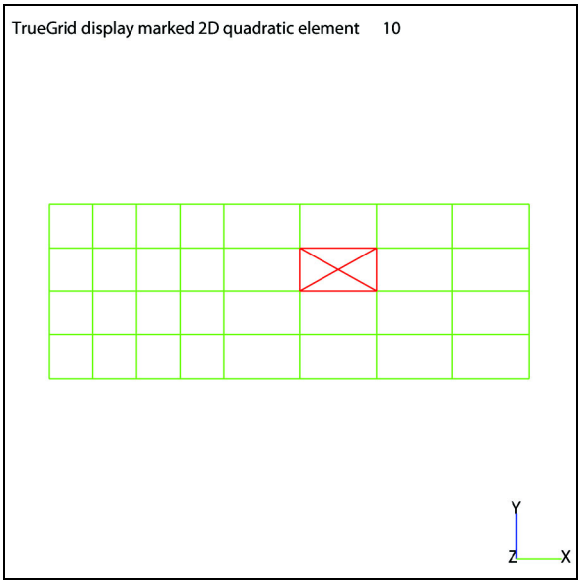

Figure 222 located quadratic shell

labels loc3d **locate a 3D brick element by its number**

loc3d *element_number*

Remarks

The diagonals of brick faces are displayed to help check distorted elements.

Example

This example uses the mesh from the example for the **labels 3d** command. The brick number 10 is located. The located brick is in the center of the picture by default.

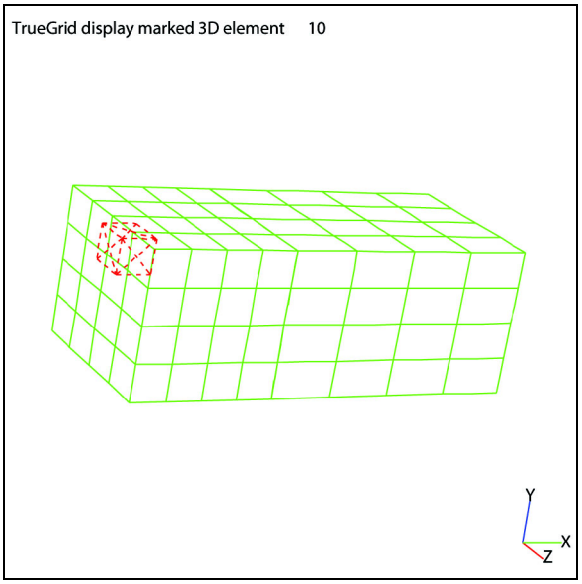

Figure 223 located brick

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labels loc3dq **locate a 3D quadratic brick element by its number**

loc3dq *element_number*

Remarks

The diagonals of brick faces are displayed to help check distorted elements.

Example

This example uses the mesh from the example for the **labels 3d** command. The quadratic brick number 10 is located. The located brick is in the center of the picture by default.

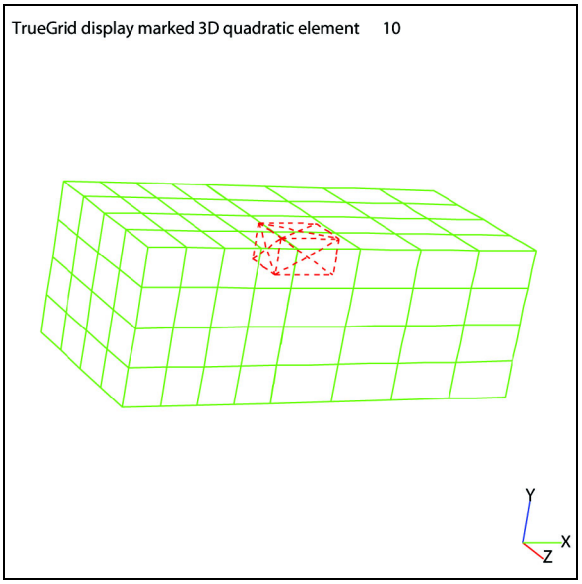

Figure 224 located quadratic brick

labels sd **label surfaces**

Example

```
curd 1 arc3 seqnc rt -1 0 0 rt 0 1 0
rt 1 0 0 mz -2;
curd 2 arc3 seqnc rt -1 0 0 rt 0 1 0
rt 1 0 0 mz 2;
curd 3 lp3 -1 0 -2 -1 0 2;;
curd 4 lp3 1 0 -2 1 0 2;;
curd 5 lp3 2 0 -2 2 0 2;;
curd 6 lp3 -3 0 -2 -3 0 2;;
curd 7 lp3 -3 -2 -2 -3 -2 2;;
curd 8 lp3 1 0 -2 2 0 -2;;
curd 9 lp3 1 0 2 2 0 2;;
curd 10 lp3 -1 0 2 -3 0 2;;
curd 11 lp3 -1 0 -2 -3 0 -2;;
curd 12 lp3 -3 0 -2 -3 -2 -2;;
curd 13 lp3 -3 0 2 -3 -2 2;;
sd 1 blend4 1 3 2 4;
sd 2 blend4 9 5 8 4;
sd 3 blend4 10 3 11 6; sd 4 blend4 13 7 12 6;dasd labels sd
```

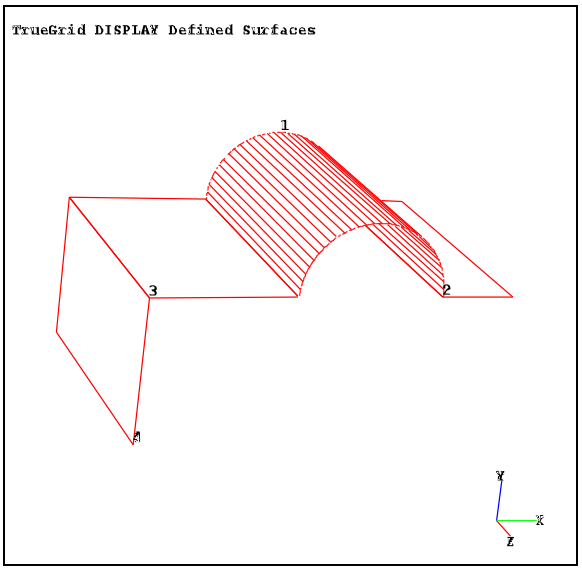

Figure 225 labeled surfaces

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labels crv **label curves**

Remarks

The **labels crv** command labels curves. It displays labels of curves in a non-overlapping way. To display more details use the **zf** or **zb** commands (zoom forward, zoom back).

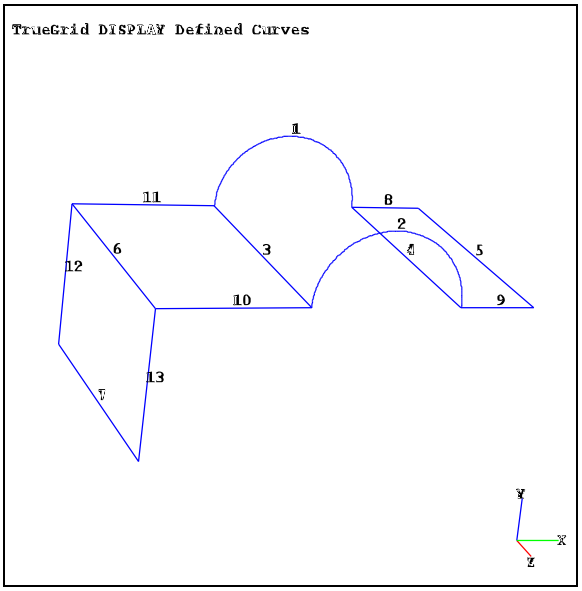

Figure 226 labeled curves in hide graphics

Example

Curves and surfaces are defined. The curves are labeled.

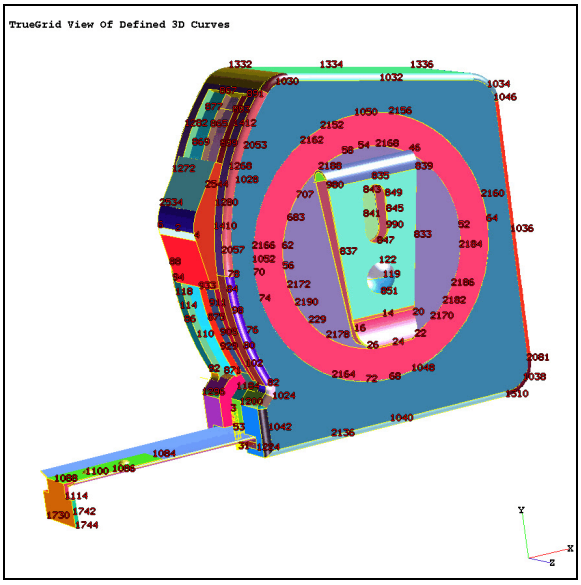

Figure 227 labeled curves in fill graphics

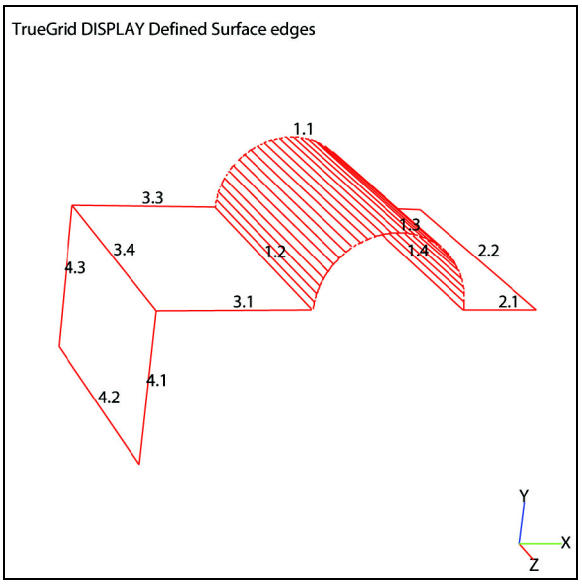

Figure 228 surface edges are labeled

labels sdedge **label surface edges**

Remarks

This command's functionality can be reproduced by using the ***Labels*** button in the Environment Window.

The **labels sdedge** command labels edges of surfaces. It displays labels of surface edges in a non-overlapping way. To display more details use **zf** or **zb** command (zoom).

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The label has the notation: *surface-#.edge-#*.

Example

Curves and surfaces are defined. The surface edges are labeled in 228.

labels sdpt **label surface points**

Remarks

The **labels sdpt** command labels surface definition points. It displays labels of surface definition points in a non-overlapping way. To display more details use **zf** or **zb** command (zoom).

The label has notation *surface_#.i_contour_#.j_contour_#* for most surfaces except for triangles from the IGES, VPOINT and STL files. For triangles the notation is *surface_#.triangle_#.point_#*, where *point_#* can be 1, 2 or 3.

See pg.? for a discussion on tessellation.

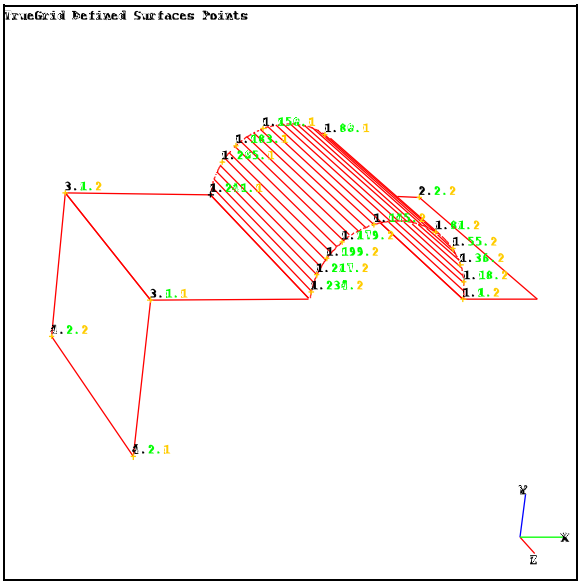

Figure 229 labeled surface definition points

labels crvpt **label curve points**

Remarks

The **labels crvpt** command labels curve definition points. It displays labels of curve definition points in a non-overlapping way. To display more details use the **zf** or **zb** commands (zoom).

The label has notation *curve.i.j* .

Example

Curves and surfaces are defined. The curve definition points are labeled in **Figure 230**.

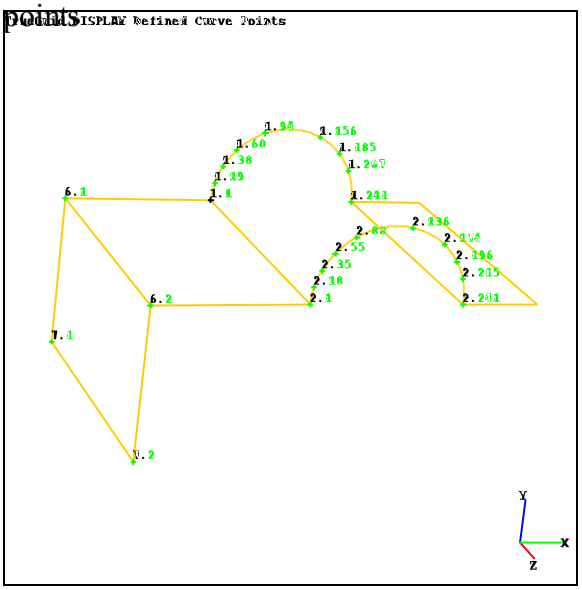

Figure 230 labeled curve definition points

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labels faceset **display face set**

Remarks

The orientation of a face in a face set can be confirmed using the "**la faceset**" command. Arrows are placed at the center of each face of the face set to indicate orientation. This is true only in the "hide" graphics mode. The face labels will only be visible if there is room for the label and if all corners of the face are visible. Orientation of a face set is controlled using the **orpt** command.

See the introduction to elements to see how the faces of brick elements are identified. **labels faceset** command works similarly as the **labels facesel** command, but without labeling.

labels facesel **display and label face set**

The faces in the face sets are labeled in a consistent fashion:

b-*number_of_element.number_of_face*	linear brick element
qb-*number_of_element.number_of_face*	quadratic brick element
s-*number_of_element.number_of_face*	linear shell element
qs-*number_of_element.number_of_face*	quadratic shell element
where	
number_of_element	is the number of the element
number_of_face	is the number of the face (1-6)

Remarks

Faces in a face set are oriented. If you use fill graphics to view the face set after issuing the **"la facesel"** command, the faces in the face set will be drawn as red. A subtle feature in the graphics will indicate if the face is oriented in towards the center of the parent element or if it is pointing outward. An inward face will be darkened as if the light sources in the picture are not reflecting off of the face. If it is pointing outward, the face will look much brighter. You need to rotate the mesh so that the faces are facing you to see this difference. Orientation of a face set is controlled using the **orpt** command.

See the introduction to elements to see how the faces of brick elements are identified. See also pg. ? for the face sets definition in the Part Phase.

Example

The face set **s1** is initialized by region, then it is enhanced 3 times by regions from various parts. The elements in the parts are linear bricks, quadratic bricks, linear shells and quadratic shells. The face

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set **s1** is displayed and labeled in **Figure 231**. The simplified command file follows:

```
linear          c set the type of element to linear

c mesh definition - linear bricks
block 1 3 5 7;1 3 5 7;1 3;1 2 3 4 1 2 3 4 -1 1

fset 1 1 2 4 4 2 = s1          c face set initialization
```

... mesh manipulations ...

```
c mesh definition - linear shells
block 1 3 5 7;1 3 5 7;-1;1 2 3 4 1 2 3 4 1


fset 1 1 1 4 4 1 u s1          c face set union
```

... mesh manipulations ...

```
endpart

quadratic       c set the type of element to quadratic

c mesh definition - quadratic bricks
block 1 3 5 7;1 3 5 7;1 3;1 2 3 4 1 2 3 4 -1 1

fset 1 1 2 4 4 2 u s1          c face set union
```

... mesh manipulations ...

```
c mesh definition for quadratic shells
block 1 3 5 7;1 3 5 7;-1;1 2 3 4 1 2 3 4 1

fset 1 1 1 4 4 1 u s1          c face set union
```

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... mesh manipulations ...

```
merge
labels facesel s1
```

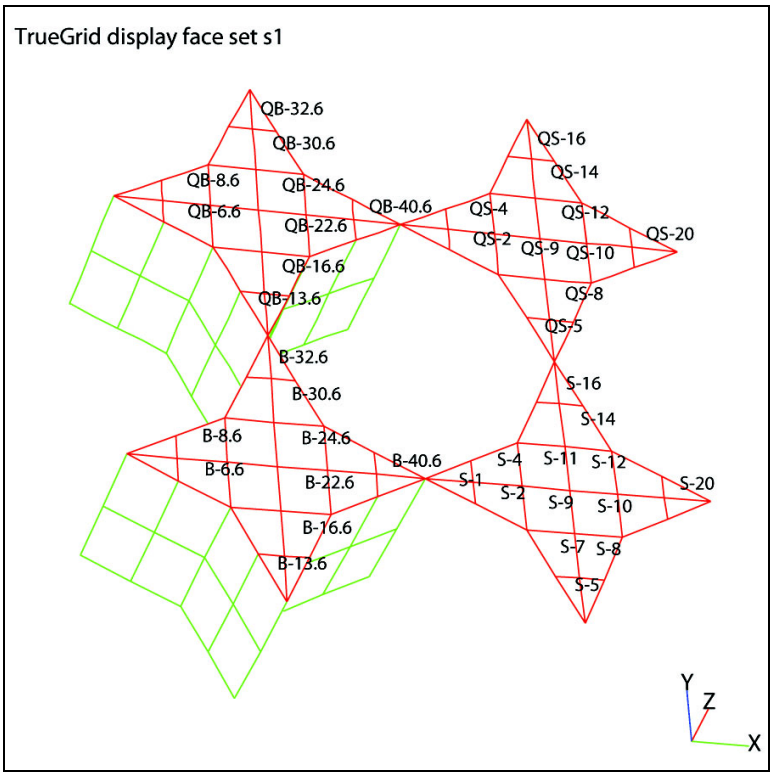

Figure 231 results of the **facesel** command

labels parts **label parts**

Remarks

Clicking the left mouse button on the part label displays a skeleton of the part.

Example

Two parts are created and labeled.

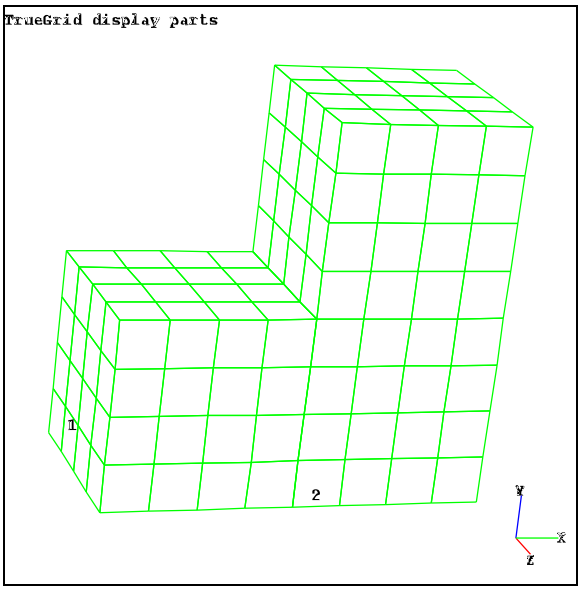

Figure 232 labeled parts

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labels tol **display nodes merged between parts**

tol $part_1$ $part_2$

where

$part_1$	part number (0 means all)
$part_2$	part number (0 means all)

See pg. 215, 470 for the rules which govern the merging of the nodes.

Example

Two parts are created. The common nodes are merged and displayed. The simplified command file follows:

```
stp .001 c merge nodes
labels tol 1 2
```

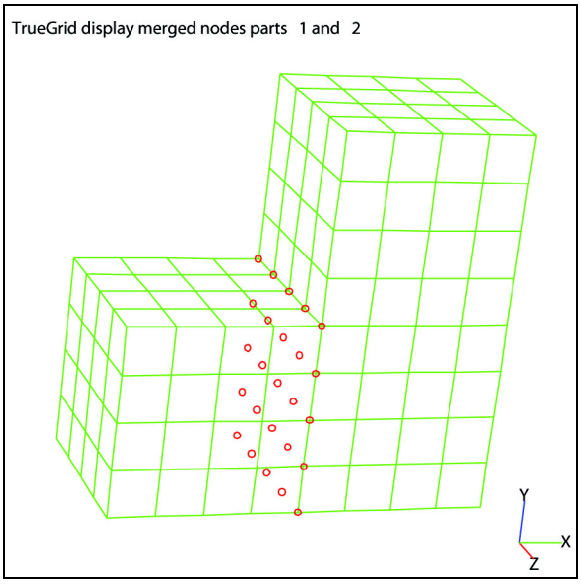

Figure 233 merged nodes

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labels fraces **display free faces**

By definition, at least one node is not merged on the free face. Free faces are constructed from the faces of the brick elements. Free faces are displayed only for active parts and materials in the picture. It is advantageous to label free faces only for a small number of parts. See pg.215, 470 for merging of the nodes. After invoking this option, you can remove or add parts and materials and the red faces will remain in the picture.

Example

The mesh from the **labels parts** example is used. The common nodes are merged. Free faces on the Part 2 are displayed.

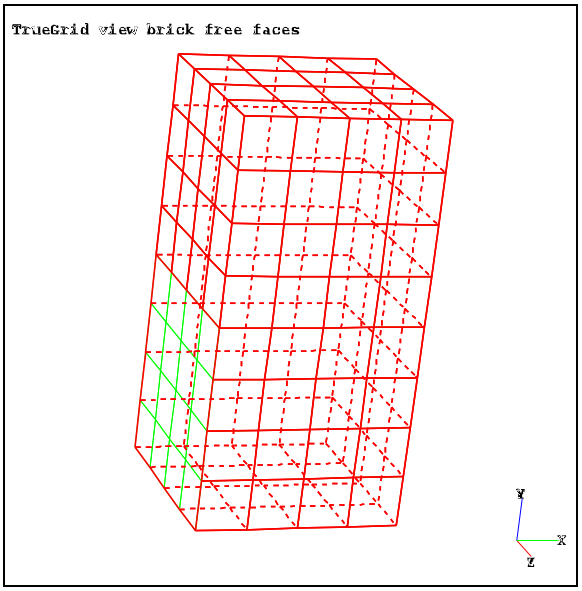

Figure 234 displayed free faces on Part 2

labels fredges **display free edges**

Remarks

By definition, at least one node is not merged on a free edge. Free edges are constructed from the edges of the shell elements. Free edges are displayed only for active parts in the picture. It is advantageous to label free edges only for a small number of parts. See pg. 215, 470 for the rules which govern the merging of the nodes. Only parts in the picture are considered.

Example

The mesh is created. The common nodes are merged. Parts 7 8 9 10 11 26 are displayed. Free edges are displayed.

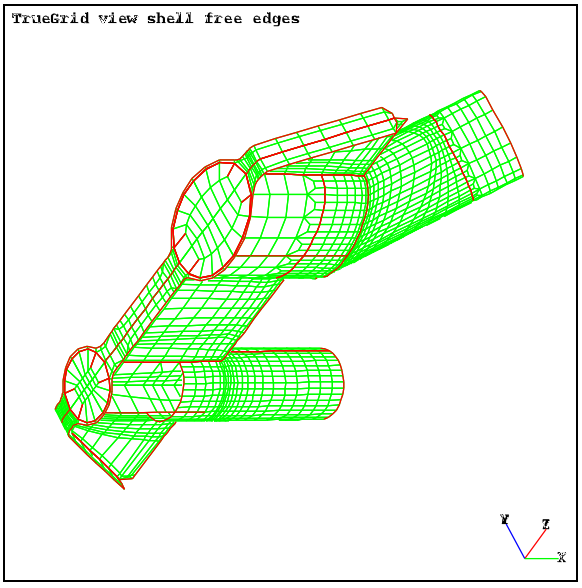

Figure 235 displayed free edges on shells

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labels cracks θ **display cracks in the mesh**

Remarks

The angle is used to measure between the exterior polygons. If the angle is below the specified angle, it is labeled as a crack. This is many times the minimum data needed to detect a problem in the node merging.

If a shell element is sandwiched between two bricks and the bricks do not form a crack, then no crack will be found, event if some of the nodes in the shell have not merged with the bricks. Use the "labels fraces" when trying to detect unmerged nodes between embedded shells and bricks.

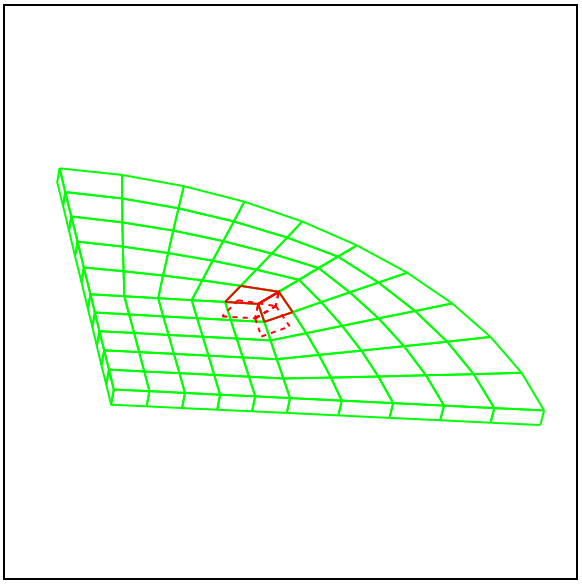

Figure 236 Crack in the mesh

Example

The butterfly mesh has not been merged, so this algorithm finds the critical spot where the merging first failed.

5. Animation Commands

You define a set of images and then generate additional images between these keyframe images in order to provide a smooth animation. This can give the appearance of flying around and even through a model.

To create an animation, you first establish up a few keyframe pictures. You to establish up to 30 keyframes. There is no limit to the number of interpolated images that may be generated.

The **sv** command will save a keyframe or view. You can go back to look at a single picture with **shv**. These two commands alone are useful for demonstrations and debugging. But with **avc** or **av** you can get an animated view of your model. The interpolated pictures allow you to produce a smooth animation while having to save a minimal number of images by hand.

With the **postscript** command you can save all these pictures as PostScript files for more processing.

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av animate views - linear interpolation

av *graphics #views* $view_1$ $\#frames_{12}$ $view_2$ $\#frames_{23}$ ...

where

graphics can be

poor	for a wireframe image,
disp	for a hidden line image, or
tvv	for a fill mode image

view1, *view2*,...	are the view numbers of views saved with **sv**, and
$\#frames_{ij}$	is the number of frames to interpolate between views numbered $view_i$ and $view_j$.

Remarks

Linear interpolation is used to form the specified number of intermediate pictures. This means that all of the parameters controlling the orientation of the picture are interpolated equally to produce a constant change in the orientation from one frame to another.

You can sequence up to 30 saved views. There is no limit on the number of frames interpolated between them.

avc animate views - cosine interpolation

avc *graphics #views* $view_1$ *#frames* $view_2$ *#frames* ...

where

graphics can be

poor	for a wireframe image,
disp	for a hidden line image, or
tvv	for a fill mode image

view1, *view2*,...	are the view numbers of views saved with **sv**, and
$\#frames_{ij}$	is the number of frames to interpolate between views numbered $view_i$ and $view_j$.

Remarks

Cosine interpolation is used to form the specified number of intermediate pictures. This means that all of the parameters controlling the orientation of the picture are interpolated in a manner that gives more frames near the saved views than midway between saved views.

You can sequence up to 30 saved keyframess. There is no limit on the number of frames interpolated between them.

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shv show a saved view

shv *view_number*
where *view_number* is the number of a view saved with the **sv** command.
Remarks

This feature restores the picture orientation to a previously saved orientation. This can be useful for demonstrations, debugging, or designing an animation.

sv save view

sv *view_number*
where *view_number* is a unique number to identify the view

Remarks

You can save up to 30 keyframes, or views. You can then use these views in the other commands in this section.

6. Exploded Views

These commands give you exploded views of your model. They affect only the graphics, not the actual model.

To make an exploded view of your model, you issue the **pexp** command to move some or all of the parts, with options of moving some parts farther than others, or in different directions. You can also issue the **mexp** command to move some or all of the materials, with options of moving some materials farther than others, or in different directions. The affect of these two commands issued multiple times is cumulative.

The following two explosion commands made the exploded view of the spaceship in the Introduction.

```
pexp 0 0 2   3 ;;
pexp 0 2 0   2 6; 1 5 ;;
```

The first command moved part 3 (front of secondary hull) to the right. The second command moved parts 2 and 6 up two units. Part 2 is the truss connected to the primary hull, and part 6 is the power unit supports - there are two supports modeled with one part. The second command also moved parts 1 and 5 up four units. Part 1 is the primary hull and part 5 is the power units - there are two power units modeled with one part. Part 4, the secondary hull, was not moved.

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exp reactivate exploded views

Display an exploded view that you had previously turned off with an **expoff** command.

exp (no arguments)

Remarks

The exploded view will be based on the translations you specified in the **pexp** and **mexp** commands.

expoff turn off exploded views

expoff (no arguments)

iniexp initialize explode offset to zero

Restore every part and material to an offset of zero.

iniexp (no argument)

mexp offset each subset of materials past the previous subset

mexp *x y z list_of_material_# ; list_of_material_# ; ... ;*
where (*x*,*y*,*z*) is the first vector offset.

Remarks

Elements with material numbers are in the N-th list will be offset in the picture by N times the specified vector.

There may be up to 50 different lists of material numbers.

If you issue both a **mexp** and a **pexp** command, the offsets you specify will be added to each other.

For a 3-D exploded view, use more than one **mexp** or **pexp** command to translate different sets of parts in different directions.

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pexp offset each subset of parts past the previous subset

pexp *x y z list_of_part_# ; list_of_part_# ; ... ;*

Remarks

Parts in the N-th list will be offset in the picture by N times the specified vector.

There may be up to 50 different lists of part numbers.

If you issue both a **mexp** and a **pexp** command, the offsets you specify will be simply added to each other.

For a 3-D exploded view, use more than one **mexp** or **pexp** command to translate different sets of parts in different directions.

sclexp explode scale factor

Multiply all the offsets for exploded views by a single scale factor.

sclexp *scale_factor*

7. Material Commands

Every material is identified with a number. Use the automatic parameter %nextmat for the next available material number. Material properties are defined by various commands associated with an output option. Refer to the various output manuals for further information.

tmm specify the total mass of a material

tmm *material_# total_mas*s

Remarks

When the Merge Phase begins or when this command is issued in the Merge Phase, then the volumes of the specified materials are calculated. The density of the material is determined in order to satisfy the total mass requirement. This feature is only available for the DYNA3D, NIKE3D, NASTRAN, TOPAZ3D, ES3D, ALE3D, ABAQUS, ANSYS, and NEUTRAL output options. One of these options must be selected before density can be determined. This feature applies to linear shells, bricks, and beams. The shell thickness and beam thickness must be set for shells and beams respectively.

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buoy specify a buoyancy condition for a list of materials

buoy *material_list ; options ;*

where

options can be

bform *type*

where *type* can be

1	displaced volume method
2	distributed pressure (head) method (brick, shell only)

styp *type parameters*

where *type* and *parameters* can be

1 *xpt ypt zpt xvec yvec zvec*	point on surface, downward vector
2 *node1 node2*	node1 on surface, node2 directly beneath

lcid *load_id*

where

load_id = 0	specifies a stationary surface
load_id > 0	specifies the load curve for surface motion vs. time

rhofg *density*

where

density > 0	means a constant density
density < 0	is the negative of a load curve for density vs. time

mgth *thickness*

rhomg *density*

where

density > 0	means a constant density
density < 0	is the negative of a load curve for density vs. time

snid *type parameters*

where *type* and *parameters* can be

1	RH rule applied to local node numbering
2 *xpt ypt zpt*	toward a point P
3 *xyt ypt zpt*	away from a point P
4 *node*	toward node N
5 *node*	away from node N

broach *flag*

where

flag can be

1	neglect beam thickness as beam broaches surface
2	include beam thickness as beam broaches surface

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am add material to the picture

am *material_number*

Remarks

The specified material is added to the list of materials to be displayed.

ams add a list of materials to the picture

ams *material_list ;*

Remarks
The specified list of materials are added to the list of materials to be displayed.

dam display all materials in the picture

dam *(no arguments)*

Remarks

All materials are added to the materials display list.

dms display a set of materials in the picture

dms *material_list* ;

Remarks

The specified list of materials becomes the list of materials to be displayed.

dm display one material in the picture

dm *material_number*

Remarks

The material display list is set to the specified material.

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ram remove all materials from the picture

ram

Remarks

No materials displayed.

rm remove one material from the picture

rm *material_number*

Remarks

The specified material is removed from the list of materials to be displayed.

rms remove a list of materials from the picture

rms *material_list ;*

Remarks

The specified materials are removed from the list of materials to be displayed.

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8. Interface Commands

chz cohesive shell elements

chz *face_set mat*

where

face_set	name of a face set
mat	material number for the cohesive

Remarks

All shells found in the region are included in the cohesion calculation. Every edge shared by two shells will be separated by a zero thick cohesion element and assigned material number *mat*. This applies to transitional regions and to edges with more than two shells.

This has not yet been generalized to triangular shells.

If an edge is selected, it will cause the material for the cohesive element to be changed along that edge. If the cohesive elements in a region or an edge have the material specified more than once, the last material number assigned will be used.

iss save interface segments

iss fset *face_set*

Interface element commands (sliding lines and surfaces, etc.) that are available during the Merge Phase.

si select nodes or faces for a sliding interface

si *type interface_# boundary options ;*

where

type can be one of:

n *node_numbe*r	to select a single node
rt *x y z*	to select a node close to a Cartesian point
cy *rho theta z*	to select a node close to a cylindrical point
sp *rho theta phi*	to select a node close to a spherical point
nset *name_of_set*	to select an entire node set

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fset *face_set* to select a face set

boundary can be one of

m master side of the interface

s slave side of the interface

options can be

fail *norm_failure_force shear_failure_force*

exp *norm_failure_exp shear_failure_exp*

fsf *coulomb_friction_scale viscous_friction_scale*

Remarks

The global properties of a sliding interface are defined using the **sid** command. Some interface types allow for nodes on the slave side. Most require face sets for both the master and slave sides. The dummy sliding interface type, which is used to control the merging without the side effect of causing a sliding interface definition in the output, allows for nodes on both the master and slave side.

9. Springs, Dampers, and Point Masses

npm creates a new node and assigns a point mass to it

npm *mp_node_# x y z mass options ;*

where

an *option* can be:

dx	no nodal displacement in the x-direction
dy	no nodal displacement in the y-direction
dz	no nodal displacement in the z-direction
rx	no nodal rotations about the x-axis
ry	no nodal rotations about the y-axis
rz	no nodal rotations about the z-axis
mdx	no mass displacement in the x-direction
mdy	no mass displacement in the y-direction
mdz	no mass displacement in the z-direction
mrx	no mass rotations about the x-axis
mry	no mass rotations about the y-axis
mrz	no mass rotations about the z-axis
ixx *mom*	specify the moment of inertia about the x-axis
iyy *mom*	specify the moment of inertia about the y-axis
izz *mom*	specify the moment of inertia about the z-axis
pdamp *alpha*	proportional damping factor (ABAQUS)
cdamp *fraction*	fraction of critical damping (ABAQUS)

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Remarks

This newly created node is separate from the existing mesh and can be attached by generating a beam or spring using this new node (see **bm** or **spring**). It can also be attached to the rest of the mesh by merging it to a neighboring node (see **t**, **tp**, **stp**, **bptol**, and **ptol)**. This is distinguished from assigning a mass to an existing node of the mesh. The latter can be done using the **pm** command. To create a new node and assign it a point mass such that it is replicated or transformed along with the part, then use the **npm** command in the Part Phase (see **lrep**, **grep**, and **pslv**). To assign a point mass to a vertex of a part such that it is replicated or transformed along with the part, use the **pm** command in the Part Phase. All of the options are not needed by all output options.

pm assigns a point mass to a node of the mesh

pm *node_# mass otions ;*

where

an *option* can be:

mdx	no mass displacement in the x-direction
mdy	no mass displacement in the y-direction
mdz	no mass displacement in the z-direction
mrx	no mass rotations about the x-axis
mry	no mass rotations about the y-axis
mrz	no mass rotations about the z-axis
ixx *mom*	specify the moment of inertia about the x-axis
iyy *mom*	specify the moment of inertia about the y-axis
izz *mom*	specify the moment of inertia about the z-axis
pdamp *alpha*	proportional damping factor (ABAQUS)
cdamp *fraction*	fraction of critical damping (ABAQUS)

Remarks

This is distinguished from creating a new node separate from the mesh and assigning a mass to it. The latter can be done using the **npm** command. To assign a point mass to a vertex within a part such that it is replicated or transformed along with the part, use the **pm** command in the Part Phase (see **lrep**, **grep**, and **pslv**). In order to create a new node and assign it a point mass such that it is replicated or transformed along with a part, then use the **npm** command in the Part Phase. All of the options are not needed by all output options.

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pminfo table of point masses information

pminfo (no arguments)

Example

The following example of the **pminfo** command was preceded by two point mass commands, the first (**npm** command) found in a part which was duplicated once, and the second (**pm** command) found in the Merge Phase.

```
npm 2 0 0 0 1.2 inc 2 dx rx mdx mrx iyy .1 izz .2 pdamp 2.3
    cdamp 1.2;
pm 377 2.1 mdx ixx .2 iyy .4 mrz pdamp .3 cdamp 2.3 ;

   STRUCTURED POINT MASSES TABLE
    node=     377, mass= 2.100E+00
     xxi= 2.000E-01, yyi= 4.000E-01, zzi= 0.000E+00
     proportional damping= 3.000E-01, critical damping= 2.300E+00
     dx=1, dy=0, dz=0, rx=0, ry=0, rz=1
   UNSTRUCTURED POINT MASSES TABLE
    number=   2 node=     126, mass= 1.200E+00
     xxi= 0.000E+00, yyi= 1.000E-01, zzi= 2.000E-01
     proportional damping= 2.300E+00, critical damping= 1.200E+00
     dx=1, dy=0, dz=0, rx=1, ry=0, rz=0
    number=   4 node=     252, mass= 1.200E+00
     xxi= 0.000E+00, yyi= 1.000E-01, zzi= 2.000E-01
     proportional damping= 2.300E+00, critical damping= 1.200E+00
     dx=1, dy=0, dz=0, rx=1, ry=0, rz=0
```

spring create/modify a numbered spring

spring *spring_# options ;*

where

an *option* can be

n1 *node_#*	assign a structure node as the first node
pm1 *pointmass_#*	assign a point mass as the first node
dx1	constrain spring in the x-direction at the first nod
dy1	constrain spring in the y-direction at the first node
dz1	constrain spring in the z-direction at the first node
rx1	constrain spring about the x-axis at the first node
ry1	constrain spring about the y-axis at the first node

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rz1 constrain spring about the z-axis at the first node
n2 *node_#* assign a structure node as the second node
pm2 assign a point mass as the second node
dx2 constrain spring in the x-direction at the second node
dy2 constrain spring in the y-direction at the second node
dz2 constrain spring in the z-direction at the second node
rx2 constrain spring about the x-axis at the second node
ry2 constrain spring about the y-axis at the second node
rz2 constrain spring about the z-axis at the second node
sddn *spd_#* specify the material properties
orop *flag* assign an orientation option

where

flag can be:

0 spring/damper acts along the axis
1 deflection/rotations are measured and force/moments applied along the following vector
2 deflection/rotations are measured and force/moments applied along the projection of the spring/damper onto the plane with the following normal
3 deflection/rotations are measured and force/moments applied along the vector between the following two nodes

prflg *flag* assign a print flag

where

flag can be:

0 forces are printed in DEFORC file
1 forces are not printed in DEFORC file

ofsi *offset* assign an initial offset
xco *x-component* assign a x-component of the orientation vector
yco *y-component* assign a y-component of the orientation vector
zco *z-component* assign a z-component of the orientation vector
n3 *node_#* assign a third node for orientation
n4 *node_#* assign a fourth node for orientation

Remarks

This command specifies the direction of the spring and the material. The nodes defining the ends of the spring can be from a node in the mesh or a point mass (see **npm** and **pm**). The **spring** command is usually invoked two times to generate a single spring, once for each node of the spring. This can be done across several parts or in the Merge Phase. All of the options are not needed by all output options.

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delspd Delete a numbered spring/damper

delspd *spring/damper-#*

Remarks

The spring/damper may have been defined using the **spring** command or it may have been read into the **True*Grid*®** internal data base using the **readmesh** command.

delspds Delete a list of numbered springs/dampers

delspds *spring/damper_list ;*

10. Element Commands

The recommended method of beam element generation in the merge phase uses the **bm** command. This method of beam generation has the advantage that one can connect any existing noe with any other existing node using a curve or a line. This method has the disadvantage that if a part is changed in a subsequent running of the session file, any reference in the **bm** command to an existing node may be incorrect and will have to be changed. This is the usual, non-parametric, nature of the merge phase. This command is fully interactive. Beams are strung along a 3D curve, interpolated along a line to connect existing nodes.

The second method of beam element generation extracts the needed nodes from an existing shell or brick part. This is only available within the Part Phase and has the usual advantage of the part phase of being parametric. If any parts are changed and the session file rerun, these beam commands will automatically adjust to the changes. The **ibm** and **ibmi** commands create beams along i-lines of the mesh, the **jbm** and **jbmi** along j-lines, and the **kbm** and **kbmi** along k-lines. This is a way to embed beam elements within a shell of a brick structure. Alternatively, the material of the parent shell of brick part can be set to 0 so that the part can be generated as usual, but so that the shell or brick elements will not be saved. The nodes that are used in any of these beam commands will be saved along with the beam elements.

The denigrated **beam** and **cbeam** commands create beams by connecting a pair of nodes with a string of beam elements. Intermediate nodes along the string are linearly interpolated. This is the INGRID method. It is not interactive. It is easy to change an INGRID beam part into this part.

Beam properties are defined using **bsd** and **bind**.

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Cross sectional properties, and in particular thicknesses, are not scaled by the **xsca**, **ysca**, **zsca**, and **csca** commands.

bm create a string of beam elements

bm *options ;*

where

option can be:

(Selection of the first node)

n1 *node_#*	to make an existing node the first node of the beams.
pm1 *point_mass_#*	to make a point mass node the first node of the beams.
rt1 *x y z const ;*	to create the first node of the beams in Cartesian coordinates.
cy1 *ρ θ z const ;*	to create the first node of the beams in cylindrical coordinates.
sp1 *ρ θ φ const ;*	to create the first node of the beams in spherical coordinates.

(Selection of the second node)

n2 *node_#*	to make an existing node the last node of the beams.
pm2 *point_mass_#*	to make a point mass the last node of the beams.
rt2 *x y z const ;*	to create the last node of the beams in Cartesian coordinates.
cy2 *ρ θ z const ;*	to create the last node of the beams in cylindrical coordinates.
sp2 *ρ θ φ const ;*	to create the last node of the beams in spherical coordinates.

(Selection of the orientation)

n3 *node_#*	to make an existing node the last node of the beams.
pm3 *point_mass_#*	to make a point mass the last node of the beams.
rt3 *x y z const ;*	to create the last node of the beams in Cartesian coordinates.
cy3 *ρ θ z const ;*	to create the last node of the beams in cylindrical coordinates.
sp3 *ρ θ φ const ;*	to create the last node of the beams in spherical coordinates.
n4 *node_#*	to make an existing node the intermediate quadratic node of a single beam.
pm4 *point_mass_#*	to make a point mass the intermediate quadratic node of a single beam.

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rt4 *x y z const ;* to create the intermediate quadratic node of a single beam in Cartesian coordinates.

cy4 *ρ θ z const ;* to create the intermediate quadratic node of a single beam in cylindrical coordinates.

sp4 *ρ θ φ const ;* to create the intermediate quadratic node of a single beam in spherical coordinates.

orient *x y z* to specify a coordinate triple to orient the beams.

sd *surface_#* to orient beam axes in the orientation of the normal of the surface

v *x y z* to orient beam axes in the direction of the vector

(Misc. options)

mate *material_#* to specify the material number.

cs *cross_section_#* to specify the cross section number (see **bsd**).

nbms *number_of_beams* to specify the number of beams in the string (default is 1).

indc *const ;* to specify the constraints on the intermediate nodes.

bb *n* to specify a 1D block boundary as the path

nd1c *const ;* to specify the constraints on the first end node along a block boundary.

nd2c *const ;* to specify the constraints on the last end node along a block boundary.

cur *3d_curve_#* to interpolate the string of beams along a 3D curve.

(Selection of the nodal spacing)

res *geometricratio* for relative spacing of nodes (default is equal spacing).

drs *first_geometricratio second_geometricratio* for double relative spacing of nodes.

nds *nodal_distribution_function_#* for nodal distribution by a function.

as 0 *first_thickness* first element length

as 1 *last_thickness* last element length

das *first_element_thickness last_element_thickness* first and last element length

sthi *sthi* for thickness in the y-direction.

sthi1 *sthi1* for thickness in the y-direction at the first end point.

sthi2 *sthi2* for thickness in the y-direction at the last end point.

tthi *tthi* for thickness in the z-direction.

tthi1 *tthi1* for thickness in the z-direction at the first end point.

tthi2 *tthi2* for thickness in the z-direction at the last end point.

csarea *csarea* for the cross section area

sharea *sharea* shear area

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inertia *Iss Itt Irr*	inertia moments
vold *volume*	volume of Discrete Beam
lump *inertia*	lumped inertia
cablcid *system_#*	local coordinate system id number defined by the **lsys**
cabarea *area*	cable area
caboff *offset*	cable offset

(Selection of the nodal offsets)

noint	for no interior node offset interpolation
roff1 *roff1*	for x-component of offset vector for first node
soff1 *soff1*	for y-component of offset vector for first node
toff1 *toff1*	for z-component of offset vector for first node
roff2 *roff2*	for x-component of offset vector for last node
soff2 *soff2*	for y-component of offset vector for last node
toff2 *toff2*	for z-component of offset vector for last node

(Selection of the pin flags)

ldr1	release the x-translation constraint at first node
lds1	release the y-translation constraint at first node
ldt1	release the z-translation constraint at first node
lrr1	release the x-rotation constraint at first node
lrs1	release the y-rotation constraint at first node
lrt1	release the z-rotation constraint at first node
ldr2	release the x-translation constraint at last node
lds2	release the y-translation constraint at last node
ldt2	release the z-translation constraint at last node
lrr2	release the x-rotation constraint at last node
lrs2	release the y-rotation constraint at last node
lrt2	release the z-rotation constraint at last node
ldr3	release the x-translation constraint at intermediate nodes
lds3	release the y-translation constraint at intermediate nodes
ldt3	release the z-translation constraint at intermediate nodes
lrr3	release the x-rotation constraint at intermediate nodes
lrs3	release the y-rotation constraint at intermediate nodes
lrt3	release the z-rotation constraint at intermediate nodes
ldp *displacement*	for the initial longitudinal displacement.
theta *angle*	for the orientation angle for the cross section.

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warpage *first_warpage_node second_warpage_node*
for two nodes used to determine warpage in the beam.

geom *option* for the method of determining curvature for the NASTRAN CBEND element.

where *option* can be

1 Center of curvature
2 Tangent of Centroid Arc
3 Bend Radius
4 Arc Angle

where *const* can be any of

dx	to constrain the x-displacement
dy	to constrain the y-displacement
dz	to constrain the z-displacement
rx	to constrain the x-axis rotation
ry	to constrain the y-axis rotation
rz	to constrain the z-axis rotation

Remarks

Use the **linear** (default) or **quadratic** command prior to this command to determine the geometric type of beam element to be generated.

There are many options to this command. However, many of the options are specific to a single simulation code. There is some overlap, but there is little consistency among the simulation codes on beam element properties. Care must be taken in selecting the options by knowing the options needed for the target simulation code. The dialogue box makes these selections easier.

This command is functional in the Merge Phase, and it is designed to create a general collection of beams or a single beam. We recommend that you use the dialogue box for **bm**.

You can use an existing node of the mesh for a beam, specify coordinates to create a new node for a beam, or you can use a point mass as a node for a beam. Coordinates can be specified in Cartesian, cylindrical, or spherical coordinates.

The cur option should not be used with **bb**, **nd1c**, **nd2c**, **n4**, **pm4**, **rt4**, **cy4**, and **sp4**.

Beam orientation can be defined using a third node, using a point mass, or by creating another node in Cartesian, cylindrical, or spherical coordinates. Use the output-code specific options in the MATERIAL Menu of the Control Phase to define materials for the beams.

Use the **bsd** to define a beam cross-section type.

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Nodes are automatically created if the number of beams specified is greater than 1.

You can define beam elements that follow a 3D curve, and specify the number of such elements, along with a spacing rule for the intermediate nodes.

Optional thickness parameters may be specified for the first and last beams when creating multiple beams. Intermediate beams will have thicknesses that are interpolated from the end beams. You may specify offsets for the first and last nodes, and optionally interpolate these offsets to intermediate nodes.

Constraints which couple the beams to the existing mesh can be eliminated. This may be done separately for the first, last, and intermediate nodes. An initial longitudinal displacement can be specified. An optional orientation angle can be specified. Warpage nodes can be defined for codes which support such options. **Bend** geometry options can be specified for codes which support such options.

Example

A simple frame structure composed of MARC beam elements is created. At first, the basic framework of beam elements is defined in the Part Phase with use of the **ibmi**, **jbmi** and **kbmi** commands. Then all 4 linear bricks are deleted by the **delem** command (or alternatively you can set the material of bricks to 0). The beams `1 7 8 2 27 26 25` are also deleted to get the desired model. Four new beams `27 28 29 30` are created by the **bm** command using existing nodes. Nodes are labeled in **Figure 237**. Beams are labeled in **Figure 238**. The simplified command file follows:

```
block 1 2 3;1 2 3;1 2;0 5 10;0 5 10;0 5;
     c strucured block mesh definition

ibmi 1 3; ; ;3 2 1 j 1 ;
 c creation of beams in the i-direction 3 columns in the
 c j-direction and 2 columns in the k-direction

jbmi 1 3; ; ;3 2 1 i 1 ;
 c creation of beams in the j-direction 3 columns in the
 c i-direction and 2 columns in the k-direction

kbmi 1 3; ; ;3 3 1 i 1 ;
 c creation of beams in the k-direction 3 columns in the
 c i-direction and 3 columns in the j-direction

merge
```

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```
delem lb 1 2 3 4;   c linear bricks 1 2 3 4 are deleted

delem lbm 1 7 8 2 27 26 25 ;
     c linear beams 1 7 8 2 27 26 25 are deleted

bsd 1 marc52 area .1 iyy .053 izz .039 ixx .0444
   etsay 100 etsaz 100  ; ;
 c cross section 1 definition
 c for marc52 beam - elastic beam
 c ... cross sectional properties

bm n1 4 n2 12 mate 1 cs 1 ;
 c beam definition by first node (n1) 4 and second node (n2) 12
 c with material 1 and cross section 1

bm n1 12 n2 16 mate 1 cs 1 ;
 c beam definition by first node (n1) 12 and second node (n2) 16
 c with material 1 and cross section 1

bm n1 3 n2 11 mate 1 cs 1 ;
 c beam definition by first node (n1) 3 and second node (n2) 11
 c with material 1 and cross section 1
```

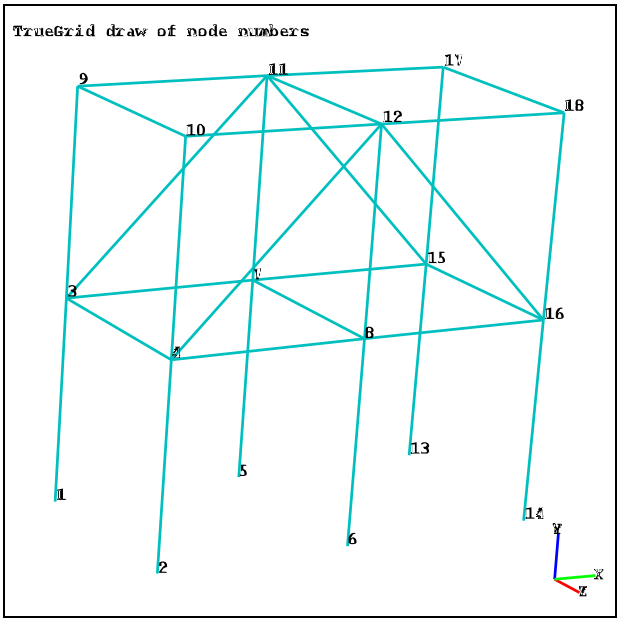

Figure 237 Beam Structure

```
bm n1 11 n2 15 mate 1 cs 1 ;
 c beam definition by first node
(n1) 11 and second node (n2) 15
  c with material 1 and cross
section 1

labels nodes    c   nodes are
                    labeled

labels 1d       c   beams are
                    labeled
```

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Example

The circular arc 3d curve is created by the **curd** command. Ten beam elements are generated along the curve by the **bm** command (**Figure 239**). The first and last node coordinates do not have to be coincident with the first and last point of the 3d curve.

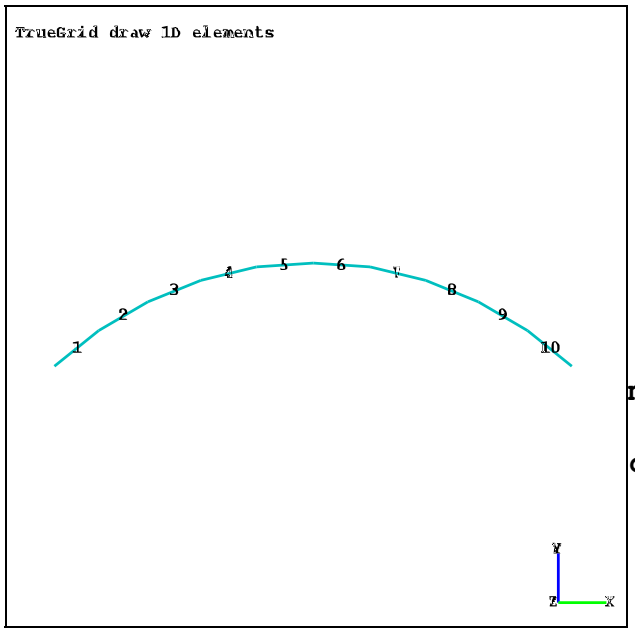

Figure 239 beams by bm

```
bm rt1 -10 0 0 ; rt2 10 0 0 ;
  mate 1 cs 1 nbms 10 cur 1 ;
 c beam definition
 c -10 0 0 and 10 0 0
 c are coordinates of the
 c first and last point
 c on arc
 c material 1 and cross
 c section 1 are assigned
 c to all 10 beams
 c 3d curve  number 1 is used

labels 1d      beams are labeled
```

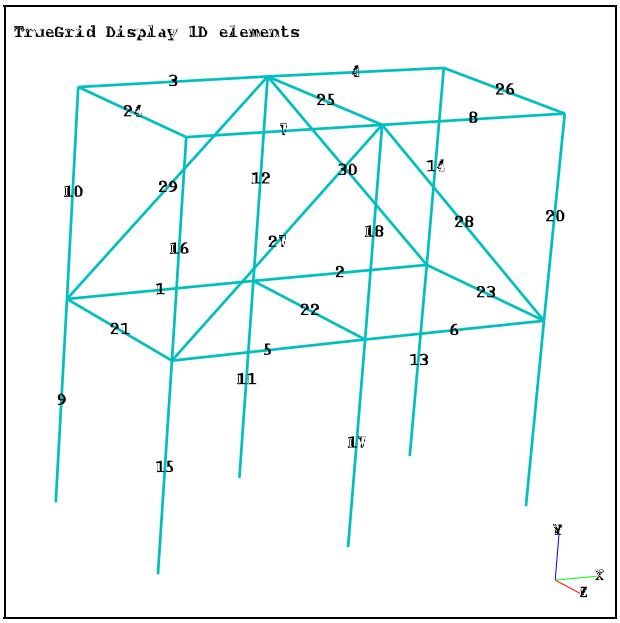

Figure 238 Beam Structure

```
merge

curd 1 arc3
 seqnc rt -10 0 0
       rt 0 4 0
       rt 10 0 0 ;
 c 3d curve definition
 c circular arc by sequence
 c of 3 points
```

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Example

The circular arc 3d curve is created by the **curd** command. Ten beam elements are generated along the curve by the **bm** command (**Figure 240**) by geometrical spacing with a factor of 0.9

```
merge
curd 1 arc3
 seqnc rt -10 0 0
       rt 0 4 0
       rt 10 0 0 ;
 c 3d curve definition
 c circular arc by sequence
 c of 3 points
bm rt1 -10 0 0 ; rt2 10 0 0 ;
  mate 1 cs 1 nbms 10 cur 1
  res .9;
 c Beam definition
 c -10 0 0 and 10 0 0
 c are coordinates of the
 c first and last node.
 c Material 1 and cross
 c section 1 are assigned
 c to all 10 beams
 c 3d curve  number 1 uses
 c the spacing factor .9

labels 1d      beams are labeled
```

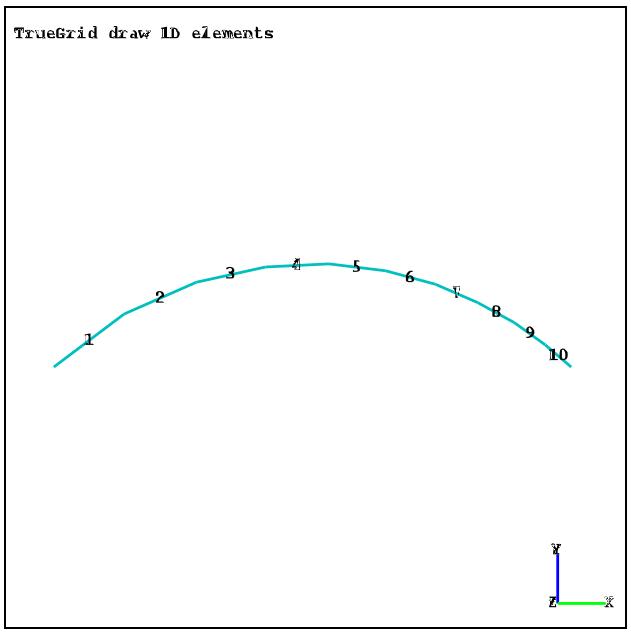

Figure 240 beams by bm

Example

Twenty beams are created by 2 **bm** commands from coordinates of the start and end nodes. The start and end nodes have assigned boundary conditions in x, y and z- directions. The local axes of all beams are pointing towards the same orientation node with coordinates `(.5,.5,0.)`. The intermediate nodes (marked by red circles in **Figure 241**) have constrained displacements in the z-direction. The simplified command file follows:

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```
merge
bm rt1 0 0 0 dx dy dz ;
   rt2 1 0 0  ; rt3 .5 .5 0. ;
   mate 1 cs 2 nbms 10
   indc dz ;;
 c beam definition
 c 0 0 0 and 1 0 0
 c are coordinates of the
 c first and last node
 c .5 .5 0 are coordinates
 c of the orientation node
 c material 1 and cross
 c section 2 are assigned
 c to all 10 beams
 c the intermediate nodes have
 c constrained displacements
 c dz
bm rt1 1 0 0  ;
   rt2 1 1 0 dx dy dz ;
   rt3 .5 .5 0. ;
  mate 1 cs 2 nbms 10
  indc dz ; ;
 c beam definition
 c 1 0 0 and 1 1 0
 c are coordinates of the
 c first and last node
 c .5 .5 0 are coordinates
 c of the orientation node
 c material 1 and cross
 c section 2 are assigned
 c to all 10 beams
 c the intermediate nodes have
 c constrained displacements
 c dz
```

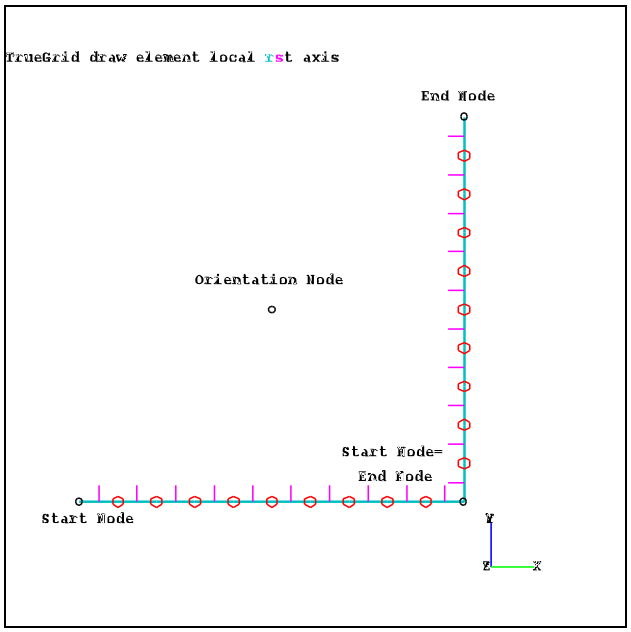

Figure 241 orientation of beams

Example

A beam structure is created by the **bm** command. A parametric 3d curve representing a parabola is created by the **curd** command. Twenty beams are generated along the 3d curve. The vertical beams (numbers `21` through `53`) are generated from the start nodes numbered `4, 6, 8, 10, 12, 14, 16, 18,` and `20` respectively. The coordinates of the last node in each vertical string of beams are formed from the coordinates of the first node, respectively, by zeroing the y-coordinate. It is useful to use ***Pick*** by ***Label*** in the Environment Window and F7 to fill the coordinates of a node

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into a dialogue box. Beams of the base (numbers 54 through 73) are generated from the pairs of existing nodes consecutively (**Figure 242** and **Figure 243**). The command file follows:

```
merge

curd 1 3dfunc -10 10 u ; -u*u/25+4 ;0; ;
 c 2D curve parametric definition - parabola

bm rt1 -10 0 0 ; rt2 10 0 0 ; mate 3 cs 2 nbms 20 cur 1 ;
 c 20 beams are generated on the curve 1 - parabola

bm n1 4  rt2 -8.2257433e+00 0.0 0.0 ; mate 2 cs 1 nbms 2 ;
bm n1 6  rt2 -6.3288207e+00 0.0 0.0 ; mate 2 cs 1 nbms 3 ;
bm n1 8  rt2 -4.3093190e+00 0.0 0.0 ; mate 2 cs 1 nbms 4 ;
bm n1 10 rt2 -2.1856163e+00 0.0 0.0 ; mate 2 cs 1 nbms 5 ;
bm n1 12 rt2 -2.3855813e-07 0.0 0.0 ; mate 2 cs 1 nbms 5 ;
bm n1 14 rt2  2.1856163e+00 0.0 0.0 ; mate 2 cs 1 nbms 5 ;
bm n1 16 rt2  4.3093190e+00 0.0 0.0 ; mate 2 cs 1 nbms 4 ;
bm n1 18 rt2  6.3288207e+00 0.0 0.0 ; mate 2 cs 1 nbms 3 ;
bm n1 20 rt2  8.2257433e+00 0.0 0.0 ; mate 2 cs 1 nbms 2 ;
 c The vertical beams (numbers 21 through 53) are generated
 c from the start node numbers
 c 4, 6, 8, 10, 12, 14, 16, 18, 20 and
 c last node coordinates,
 c which are derived from the start ones by zeroing
 c the y-coordinate. Number of beams is from 2 to 5

bm n1  1 n2 22 mate 2 cs 3 nbms 2 ;
bm n1 22 n2 24 mate 2 cs 3 nbms 2 ;
bm n1 24 n2 27 mate 2 cs 3 nbms 2 ;
bm n1 27 n2 31 mate 2 cs 3 nbms 2 ;
bm n1 31 n2 36 mate 2 cs 3 nbms 2 ;
bm n1 36 n2 41 mate 2 cs 3 nbms 2 ;
bm n1 41 n2 46 mate 2 cs 3 nbms 2 ;
bm n1 46 n2 50 mate 2 cs 3 nbms 2 ;
bm n1 50 n2 53 mate 2 cs 3 nbms 2 ;
bm n1 53 n2  2 mate 2 cs 3 nbms 2 ;
 c Beams of the base (numbers 54 through 73) are generated
 c from the pairs of existing nodes, consecutively

labels 1d      c beams are labeled
```

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```
labels nodes          c nodes are labeled
```

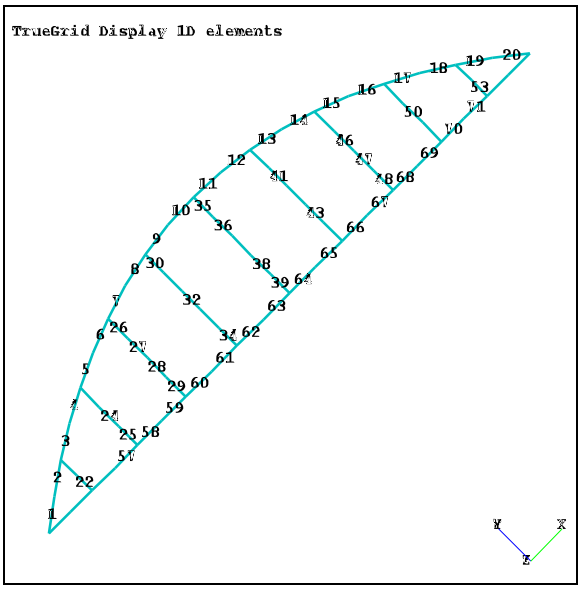

Figure 242 beams by bm

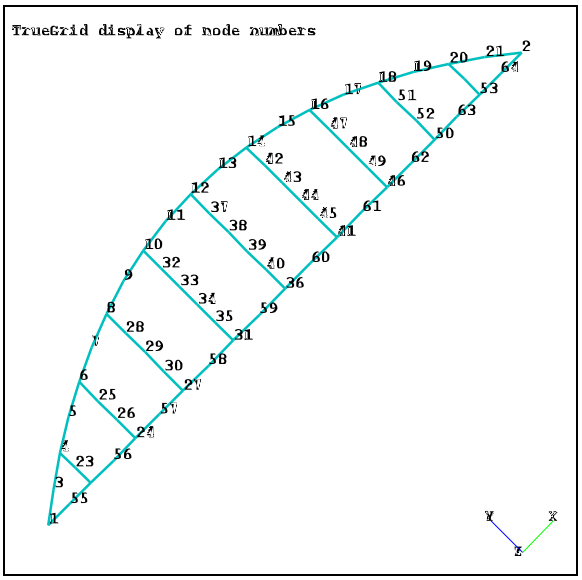

Figure 243 beams by bm

bms change the section properties of a set of beams

bms *selection options ;*

where

selection can be

set element_set	select an element set containing beam elements
lbm element_list ;	select a list of beam elements by number

option can be:

orient *x y z*	specify a coordinate triple to orient the beams.
v *x y z*	orient beam axes in the direction of the vector
cs *cross_section_#*	specify the cross section number (see **bsd**).
sthi *sthi*	thickness in the y-direction.
sthi1 *sthi1*	thickness in the y-direction at the first node
sthi2 *sthi2*	thickness in the y-direction at the last node
tthi *tthi*	thickness in the z-direction.
tthi1 *tthi1*	thickness in the z-direction at the first node
tthi2 *tthi2*	thickness in the z-direction at the last node
csarea *csarea*	the cross section area
sharea *sharea*	shear area

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inertia *Iss Itt Irr*	inertia moments
vold *volume*	volume of Discrete Beam
lump *inertia*	lumped inertia
cablcid *system_#*	local coordinate system id number defined by **lsys**
cabarea *area*	cable area
caboff *offset*	cable offset
noint	for no interior node offset interpolation
roff1 *roff1*	for x-component of offset vector for first node
soff1 *soff1*	for y-component of offset vector for first node
toff1 *toff1*	for z-component of offset vector for first node
roff2 *roff2*	for x-component of offset vector for last node
soff2 *soff2*	for y-component of offset vector for last node
toff2 *toff2*	for z-component of offset vector for last node
ldr1	release the x-translation constraint at 1st node
lds1	release the y-translation constraint at 1st node
ldt1	release the z-translation constraint at 1st node
lrr1	release the x-rotation constraint at 1st node
lrs1	release the y-rotation constraint at 1st node
lrt1	release the z-rotation constraint at 1st node
ldr2	release the x-translation constraint at 2nd node
lds2	release the y-translation constraint at 2nd node
ldt2	release the z-translation constraint at 2nd node
lrr2	release the x-rotation constraint at 2nd node
lrs2	release the y-rotation constraint at 2nd node
lrt2	release the z-rotation constraint at 2nd node
ldr1o	no release of the x-translation constraint at 1st node
lds1o	no release of the y-translation constraint at 1st node
ldt1o	no release of the z-translation constraint at 1st node
lrr1o	no release of the x-rotation constraint at 1st node
lrs1o	no release of the y-rotation constraint at 1st node
lrt1o	no release of the z-rotation constraint at 1st node
ldr2o	no release of the x-translation constraint at 2nd node
lds2o	no release of the y-translation constraint at 2nd node
ldt2o	no release of the z-translation constraint at 2nd node
lrr2o	no release of the x-rotation constraint at 2nd node
lrs2o	no release of the y-rotation constraint at 2nd node
lrt2o	no release of the z-rotation constraint at 2ns node
ldp *displacement*	for the initial longitudinal displacement.
theta *angle*	for the orientation angle for the cross section.

warpage *first_warpage_node second_warpage_node*

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two nodes used to determine warpage in the beam.

geom *option* for the method of determining curvature for the NASTRAN CBEND element.

where *option* can be

1 Center of curvature
2 Tangent of Centroid Arc
3 Bend Radius
4 Arc Angle

where *const* can be any of

dx	to constrain the x-displacement
dy	to constrain the y-displacement
dz	to constrain the z-displacement
rx	to constrain the x-axis rotation
ry	to constrain the y-axis rotation
rz	to constrain the z-axis rotation

Remarks

Refer to the bm command for more details on the meaning of these cross sectional properties.

delem delete a set of elements

delem *option list_of_elements ;*

where an

option must be one of the following:

lb	for 3D linear 8-node brick elements
ls	for 2D linear 4-node shell elements
lbm	for 1D linear 2-node beam elements
qb	for 3D quadratic 8-node brick elements
qs	for 2D quadratic 4-node shell elements
qbm	for 1D quadratic 3-node beam elements
eset *set_name*	to delete all elements in an element set

Remarks

All boundary constraints and conditions associated with these deleted elements are removed from the **True*Grid*®** internal database. Any nodes used to define these elements, which are not used in the definition of other elements, are removed. The sorting and data rearrangement due to element deletion can be extensive and time consuming. For this reason, it is recommended that a group of elements be deleted instead of one at a time. This command cannot be undone. When elements are deleted, they cannot be retrieved.

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etd specify the element types to be displayed in the graphics

etd options ;

where

an *option* can be any of the following:

1dl *switch*	for linear 1D beam elements
2dl *switch*	for linear 2D shell elements
3dl *switch*	for linear 3D brick elements
1dq switch	for quadratic 1D beam elements
2dq *switch*	for quadratic 2D shell elements
3dq *switch*	for quadratic 3D brick elements

where

switch can be

on

off

rbe rigid body elements

rbe *element_# type parameters*

where the *parameters* vary depending on the *type* from the following

rrod *node const ; node const ;*

where only one of the following constraints can be selected for *const*

mdx	dependent translation in x
mdy	dependent translation in y
mdz	dependent translation in z

where a *node* can be selected by it's number or by coordinates

node node_#	any exist node
rt *x y z*	new node in Cartesian coordinates
cy *rho theta z*	new node in cylindrical coordinates
sp *rho theta phi*	new node in spherical coordinates

rbar *node const ; node const ;*

where a constraint *const* can be formed using

ndx	independent translation in x
ndy	independent translation in y
ndz	independent translation in z
nrx	independent rotation in x
nry	independent rotation in y
nrz	independent rotation in z
mdx	dependent translation in x

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mdy dependent translation in y
mdz dependent translation in z
mrx dependent rotation in x
mry dependent rotation in y
mrz dependent rotation in z

where a *node* can be selected by it's number or by coordinates

node node_# any exist node
rt *x y z* new node in Cartesian coordinates
cy *rho theta z* new node in cylindrical coordinates
sp *rho theta phi* new node in spherical coordinates

rtrplt *node const ; node const ; node3 const ;*

where a constraint *const* can be formed using

ndx independent translation in x
ndy independent translation in y
ndz independent translation in z
nrx independent rotation in x
nry independent rotation in y
nrz independent rotation in z
mdx dependent translation in x
mdy dependent translation in y
mdz dependent translation in z
mrx dependent rotation in x
mry dependent rotation in y
mrz dependent rotation in z

where a *node* can be selected by it's number or by coordinates

node node_# any exist node
rt *x y z* new node in Cartesian coordinates
cy *rho theta z* new node in cylindrical coordinates
sp *rho theta phi* new node in spherical coordinates

rbe2 *node const ; list_nodes ;*

where a constraint *const* can be formed using

mdx dependent translation in x
mdy dependent translation in y
mdz dependent translation in z
mrx dependent rotation in x
mry dependent rotation in y
mrz dependent rotation in z

where a *node* can be selected by it's number or by coordinates

node node_# any exist node
rt *x y z* new node in Cartesian coordinates

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cy *rho theta z* new node in cylindrical coordinates
sp *rho theta phi* new node in spherical coordinates

rbe3 *node const ; [node const ; weight list_nodes ;];*

where a constraint *const* can be formed using

mdx dependent translation in x
mdy dependent translation in y
mdz dependent translation in z
mrx dependent rotation in x
mry dependent rotation in y
mrz dependent rotation in z

where a *node* can be selected by it's number or by coordinates

node node_# any exist node
rt *x y z* new node in Cartesian coordinates
cy *rho theta z* new node in cylindrical coordinates
sp *rho theta phi* new node in spherical coordinates

Remarks

This command generates rigid body elements for **NASTRAN** and **NE/NASTRAN**. The nodes used to form the rigid body are identified by node number or by coordinates in the neighborhood. This command supports 5 types of rigid bodies. They are:

1. **rrod** is for a rob rigid element formed with two nodes with one end pinned.
2. **rbar** is for a bar rigid element formed with two nodes and 6 DOF.
3. **rtrplt** is for a plate rigid element formed with three nodes.
4. **rbe2** is for a rigid body.
5. **rbe3** is for an interpolation constraint element.

The dependencies are best described in the **NASTRAN** or **NE/NASTRAN** manuals.

11. Commands Related to Parts

The are a number of element types that can be generated using the block or cylinder command. In addition, elements can be imported using the **readmesh** command. They are: linear 8 node hexa solid, linear 6 node penta (wedge) solid, linear 5 node pyramid solid, linear 4 node tetra solid, quadratic 20 node hexa solid, quadratic 15 node penta solid, quadratic 13 node pyramid solid, quadratic 10 node tetra solid, triquadratic 27 node hexa solid, linear 4 node quad shell, linear 3 node triangle shell, quadratic 8 node quad shell, quadratic 6 node triangle shell, triquadratic 9 node quad shell, linear 2 node beam, and quadratic 3 node beam. Both solids and shells can be generated using the **block** or **cylinder** command. Beams can also be generated using the **block** or **cylinder** command. They can also be generated using the **bm** command in the merge phase. In all cases, choose the order of element

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(linear, quadratic, or triquadratic) before creating parts. Once chosen, it cannot be changed. The default is linear.

Hexa elements are the easiest element types to generate. To generate a penta (wedge), pyramid, or tetra element, extra steps are needed to collapse some edges and/or faces of a hexa element. Alternatively, one can make a linear or triquadratic all hexa mesh. Then issue the **alltets** command in the merge phase. The hexa elements will be converted to tetra elements with 5 tretra elements derived from each hexa element. The **alltets** command can only be issued once at the end of generating parts.

linear specify elements to be linear

linear (no arguments)

Remarks

All brick elements with have 8 nodes, one at every corner of the brick. Shell elements will have 4 nodes, one at every corner. Beam elements will have 2 nodes, one at each end. This is the default.

After entering the merge phase and merging nodes, if an edge or face becomes degenerate, forming a different element type, such as a wedge or tetrahedron, the number of nodes are changed for that element.

This command has no affect on the models imported using the **readmesh** command.

quadratic specify elements to be quadratic

quadratic (no arguments)

Remarks

All brick elements with have 20 nodes, one at every corner of the brick and an additional mid-edge node for each of the 12 edges of the element. Shell elements will have 4 nodes, one at every corner and an additional mid-edge node for each of the 4 edges of the shell element. Beam elements will have 2 nodes, one at each end.

After entering the merge phase and merging nodes, if an edge or face becomes degenerate, forming a different element type, such as a wedge or tetrahedron, the number of nodes are changed for that element.

This command has no affect on the models imported using the **readmesh** command.

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triquadratic specify elements to be triquadratic

triquadratic (no arguments)

Remarks

All brick elements with have 27 nodes, one at every corner of the brick, an additional mid-edge node for each of the 12 edges of the element, an additional mid-face node for all 6 faces, and 1 mid-element node. Shell elements will have 4 nodes, one at every corner, an additional mid-edge node for each of the 4 edges of the shell element, and 1 midelement node. Beam elements will have 2 nodes, one at each end.

After entering the merge phase and merging nodes, if an edge or face becomes degenerate, forming a different element type, such as a wedge or tetrahedron, the number of nodes are changed for that element.

This command has no affect on the models imported using the **readmesh** command.

sparticle smooth particle hydrodynamic elements (SPH) for LSDYNA

sparticle (no arguments)

Remarks

For subsequent parts, all brick elements will be converted to smooth particle hydrodynamic elements. This is for LSDYNA output only. When building a part, the usual hex mesh graphics will be shown. Only after ending the part and entering the merge phase, will the representation of the SPH elements be shown. To continue using linear or quadratic bricks, issue the **linear**, **quadratic,** or **triquadratic** command before initiating a new part.

Use the **lsdymats** commands to assign properties for SPH elements. Be sure to specify the material density when defining this material. It is needed to complete the SPH elements.

Use the **plane** command to define the SPH symmetry plane.

Use the **vd** command to define a SPH box.

SPH control parameters can be set using the **lsdyopts** command.

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Example

```
sd 1 sp 0 0 0 3
sparticle
block 1 4 7 10;1 4 7 10;1 4 7 10;
         -1 -1 1 1 -1 -1 1 1 -1 -1 1 1
dei 1 2 0 3 4; 1 2 0 3 4;;
dei 1 2 0 3 4;; 1 2 0 3 4;
dei ; 1 2 0 3 4; 1 2 0 3 4;
sfi -1 -4; -1 -4; -1 -4;sd 1
```

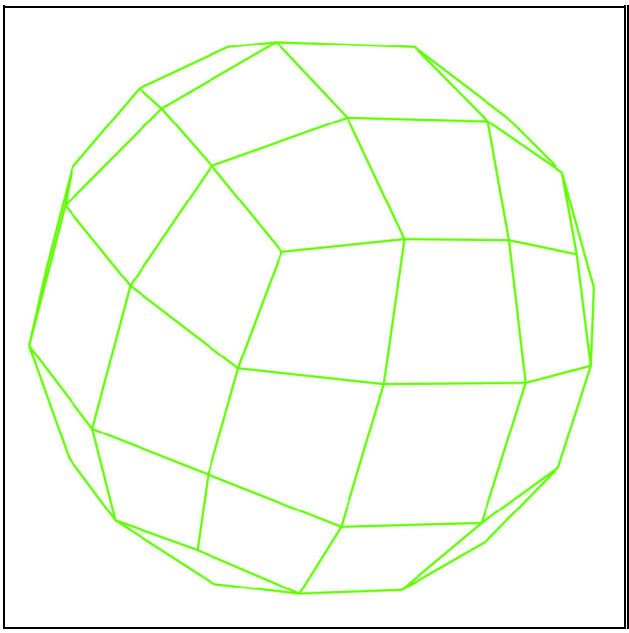

Figure 244 hex mesh in the part phase

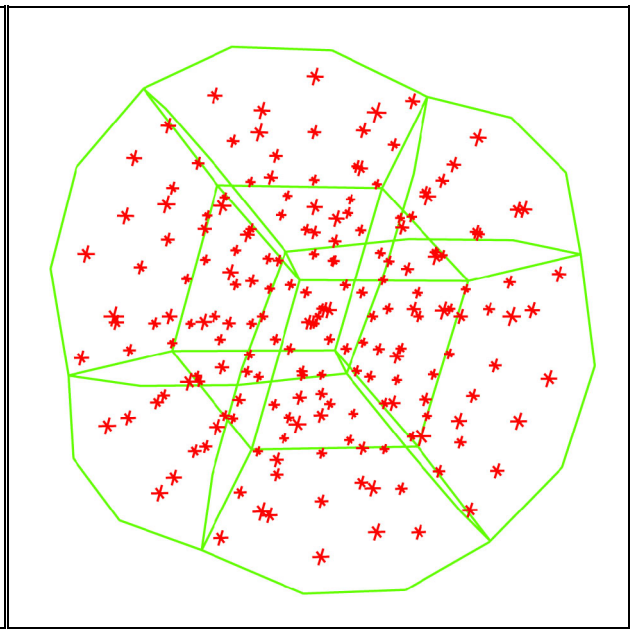

Figure 245 SPH centers in merge phase

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partmode part command indices format

partmode *mode*

where

mode can be

s	for standard (default)
i	for interval

Remarks

The standard method is described in the **block**, **cylinder**, and **blude** command remarks. The interval mode uses 3 lists of numbers, one for each index direction. The difference is that they are the number of elements for each region. This implies that there will be one less number in each list if the interval mode is specified. An analogy is a fence. The standard method indicates the position of each fence post from the start of the fence while the interval mode specifies gaps between the fence posts.

This command only applies to the **block**, **cylinder**, and **blude** commands.

The default is s.

The standard mode is more complex but has the advantage of creating both solids and shells. It also has the same format as the index progressions in the part phase commands. The interval mode is preferred because if the number of elements in one region is to be modified, only one number in the part command needs to be modified.

Examples

In these two examples, the result is the same.

```
parameters
    i1 1
    i2 [%i1+2]
    i3 [%i2+4]
    j1 1
    j2 [%j1+4]
    k1 1
    k2 [%k1+1]
    k3 [%k2+2] ;
partmode s c default
block %i1 %i2 %i3;%j1 %j2;%k1 %k2 %k3;0 .1 .2;1 1.2;2 2.3 2.5;
```

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partmode i
block 2 4;4;1 2;0 .1 .2;1 1.2;2 2.3 2.5;

trprt transform a part

trprt *part_# trans ;*

where *trans* is a product from left to right of the following operators

- **mx** *x_offset* to translate in the x direction
- **my** *y_offset* to translate in the y direction
- **mz** *z_offset* to translate in the z direction
- **v** *x_offset y_offset z_offset* to translate by a vector
- **scv** *xn yn zn offset* to translate a given distance along a vector
- **dv** $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$ to translate by a vector given by two points
- **rx** *theta* to rotate about the x axis
- **ry** *theta* to rotate about the y axis
- **rz** *theta* to rotate about the z axis
- **raxis** *angle x0 y0 z0 xn yn zn* axis of rotation
- **rxy** to reflect about the x-y plane
- **ryz** to reflect about the y-z plane
- **rzx** to reflect about the z-x plane
- **tf** *origin x-axis y-axis*

  where each argument consists of a coordinate type followed by coordinates:
  - **rt** *x y z* Cartesian coordinates
  - **cy** *rho theta z* cylindrical coordinates
  - **sp** *rho theta phi* spherical coordinates
  - **pt** *c.i* label of a labeled point from a 3D curve
  - **pt** *s.i.j* label of a labeled point from a surface
- **ftf** *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

  where each argument consists of a coordinate type followed by coordinates:
  - **rt** *x y z* Cartesian coordinates
  - **cy** *rho theta z* cylindrical coordinates
  - **sp** *rho theta phi* spherical coordinates
  - **pt** *c.i* label of a labeled point from a 3D curve
  - **pt** *s.i.j* label of a labeled point from a surface
- **inv** invert the present transformation
- **csca** *scale_factor* to scale all coordinates
- **xsca** *scale_factor* to scale the x coordinate
- **ysca** *scale_factor* to scale the y coordinate
- **zsca** *scale_factor* to scale the z coordinate

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Remarks

This command is a way to transform parts in the merge phase instead of transforming in the part phase.

```
sd 1 cy 0 0 0 0 1 0 1
block 1 11;1 11;1 11;-1 1 -1 1 -1 1
sfi -1 -2;; -1 -2;sd 1
block 1 11;1 11;1 11;-1 1 1 3 -1 1
sfi -1 -2;; -1 -2;sd 1
merge
stp .001
trprt 2 mx 1 my -1 rz 15 mx -1 my 1;
stp .001
```

In this example, 2 simple cylinder parts are generated, one on top of the other. In the merge phase, 121 nodes get merged. Then the **trprt** command is applied. The nodes are no longer merged. A new merge command is issued to merge nodes and only one node gets merged.

Notice that the rotation of the second part uses the point at (-1,1,0) to apply a rotation of 15 degrees about the z-axis. This is done by first translating the part so that the point that was at (-1,1,0) is translated to (0,0,0). Then the rotation is applied. As a last step, the part is translated so that the point that is now at (0,0,0) will be translated back to (-1,1,0).

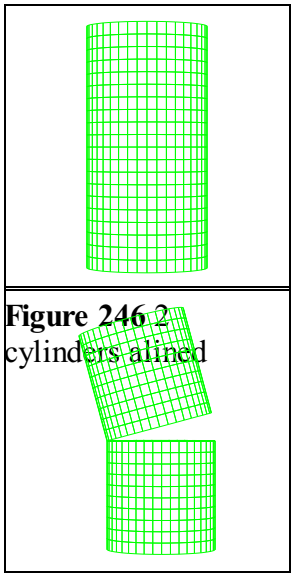

Figure 246 2 cylinders alined

Figure 247 2 cylinder not alined

alltets convert all hex elements to tet elements

alltets

Remarks

The **alltets** command can be used in the merge phase, just before writing the output file with the write command. This command converts hexahedron elements to tetrahedron elements. Care is needed in using this command. It can only be used once when generating the mesh and only after all elements have been generated. It only works on hex elements. No wedge, tet, or shell elements are allowed. There is one more issue. If you build a mesh which is equivalent to a torus (ring) and it has a twist of +/- 90 degrees or if there is an odd number of elements around the ring, then there be a disconnect someplace. This is because the hex element is being divided into 5 tet elements. There are other ways that a hex element can be divided into tets, but the quality of the mapping from hex to 5 tets was considered the best in quality. The purpose, primarily, is to allow for a fair comparison of the hex mesh to an equivalent tet mesh. However, by making this feature available to users, we expect that

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some will use it to get a high quality tet mesh to run a simulation. So user beware, this feature may not be useful as a general tool for building tet meshes.

Be sure not to use quadratic hexahedron elements. If you use linear hexahedron elements (default), you will get linear tetrahedron elements (4 nodes). If you use triquadratic hexahedron elements, you will get quadratic tetrahedron elements (10 nodes).

12. Displacements, Velocities, and Accelerations

These commands specify displacements, velocities, and accelerations, usually for initial or boundary conditions. In most of them, the arguments take the form

load_curve amplitude x y z

The displacement, velocity, etc. is applied in the direction given by the vector (*x*, *y*, *z*), which might be in Cartesian, cylindrical, or spherical coordinates. For some simulation codes like DYNA3D, the magnitude of the condition is the *product* of the *amplitude* and the *current value of the load curve*. In this case, the load curve is a time-dependent function given by the load curve number, *load_curve*. In some other simulation codes, such as ABAQUS, the load curve number is associated with a step, and in other simulation codes like NASTRAN, the load curve number is associated with a load case.

Also see the Displacement, Velocities, and Acceleration section in the Part Phase (Generation chapter).

acc Cartesian prescribed nodal boundary acceleration

acc *nodes load_curve ampl options x y z*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

load_curve	for a load curve or case number or zero
ampl	for the amplitude

where *options* can be any of

exclude	exclude normal directions (for Lsdyna)
birth *time*	specify starting time (for Lsdyna)
death *time*	specify ending time (for Lsdyna)

where

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x y z	for the x, y, and z components of the acceleration

deform assigns deformations to beam elements

deform *type object sid deformation*

where *type* and *object* are related

beam *beam_#*	beam number (rods or bars)
eset *element_set*	element set name, assumed to contain rods or bars
where *sid*	set identification number
where *deformation*	magnitude of deformation

Remarks

This command defines rod and bar element deformations for NASTRAN and NE/NASTRAN. It does not check to see if the beam element cross section type (**bsd**) is a rod or bar.

dis initial nodal displacement

dis *nodes option x y z*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

option can be

sid *id*	for a set identification number
r	to flag this command for rotational conditions
x y z	for the x, y, and z-components of the displacement

Remarks

When a node is assigned several displacements, the last one specified is the one that is used. The **r** flag converts this command to mean nodal rotational velocities. The **sid** is used by several output options to indicate a load case or step.

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fd displacement boundary condition

fd *nodes load_curve ampl options x y z*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

load_curve	for a load curve or case number or zero
ampl	for the amplitude

where *options* can be any of

exclude	exclude normal directions (for Lsdyna)
birth *time*	specify starting time (for Lsdyna)
death *time*	specify ending time (for Lsdyna)

where

x y z	are the x, y, and z-components of the displacement.

fv prescribed velocities

fv *nodes load_curve ampl options x y z*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

load_curve	for a load curve or case number or zero
ampl	for the amplitude

where *options* can be any of

exclude	exclude normal directions (for Lsdyna)
birth *time*	specify starting time (for Lsdyna)
death *time*	specify ending time (for Lsdyna)

where

x y z	for the x, y, and z-components of the velocity

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frb prescribed rotational boundary

frb *nodes load_# amplitude options condition direction*

where *nodes* can be

n *node_number*
rt *x y z*
cy ρ θ *z*
sp ρ θ φ
nset *set_name*

where *options* can be any of the following

birth *time*
death *time*
o**ffset** $offset_1$ $offset_2$

where *condition* must be one of the following

v	velocities
a	accelerations
d	displacements
dofv	nodal dof velocities
dofa	nodal dof accelerations
dofd	nodal dof displacements

where *direction* must be one of the following

x	about the x-axis
y	about the y-axis
z	about the z-axis
v x_0 y_0 z_0	about an arbitrary axis
ex	not about the x-axis
ey	not about the y-axis
ez	not about the z-axis
ev x_0 y_0 z_0	not about an arbitrary axis

Remarks

A condition can be a velocity, acceleration, displacement, or a nodal rotation. This is suited for Dyna3D (velocities, accelerations, and nodal rotations) and Lsdyna. In these codes, the selected nodes are prescribed this condition relative to an axis of rotation.

Use the **frb** option in the **co** merge phase command in the merge phase to display these conditions.

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fvv prescribed variable velocities

fvv *nodes load_curve amp_expr ; x_expr ; y_expr ; z_expr ;*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

load_curve	for a load curve or case number or zero
amp_expr	for the FORTRAN amplitude
x_expr	for the FORTRAN expression for the velocity x-component
y_expr	for the FORTRAN expression for the velocity y-component
z_expr	for the FORTRAN expression for the velocity z-component

Remarks

This command assigns velocities, allowing for the amplitude factor and the Cartesian vector components to be calculated using a FORTRAN expression. Each expression can reference the nodal coordinates x, y, and z.

vacc Cartesian prescribed variable nodal boundary acceleration

vacc *nodes load_curve amp_expr ; x_expr ; y_expr ; z_expr ;*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

load_curve	for a load curve or case number or zero
amp_expr	for the FORTRAN amplitude
x_expr	for the FORTRAN expression for the x-component of the acceleration
y_expr	for the FORTRAN expression for the y-component of the acceleration
z_expr	for the FORTRAN expression for the z-component of the acceleration

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Remarks

This command assigns accelerations, allowing for the amplitude factor and the Cartesian vector components to be calculated using a FORTRAN expression. Each expression can reference the nodal coordinates X, Y, and Z.

ve initial nodal velocity

ve *nodes option x y z*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

option can be

sid *id*	for a set identification number
r	to flag this command for rotational conditions
x y z	for the x, y, and z-components of the velocity

Remarks

This overrides the velocities specified by the **velocity** and **rotation** commands. When a node is assigned several velocities, the last one specified is the one that is used. The **r** flag converts this command to mean nodal rotational velocities. The **sid** is used by several output options to indicate a load case or step.

13. Force, Pressure, and Loads

These commands specify loads, displacements, or similar conditions. In most of them, the arguments include

load_curve amplitude x y z

The load, displacement, etc. is applied in the direction given by the vector (*x*, *y*, *z*), which might be in Cartesian, cylindrical, or spherical coordinates. For some simulation codes like DYNA3D, the magnitude of the load, displacement, etc. is the *product* of the *amplitude* and the *current value of the load curve*. In this case, the load curve is a time-dependent function given by the load curve number, *load_curve*. In some other simulation codes, such as ABAQUS, the load curve number is associated with a step, and in other simulation codes like NASTRAN, the load curve number is associated with a load case.

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dsload distributed surface loads for ABAQUS

dsload *face_set load_case [options] type parameters*

where *options* can be any of the following:

cr	to choose constant resultant
nf	to choose no follower forces
lcase *load_case_#*	for load case
or *local_sys_#*	for orientation
rn *node_#*	for reference node

where *type* and associated *parameters* must be one of

p *ampl*	for uniform pressure
pnu *ampl*	for nonuniform pressure
sp *ampl*	for stagnation pressure
vp *ampl*	for viscous pressure
hp *ampl z0 zp*	for hydrostatic pressure
trvec *ampl x y z*	for uniform general traction
trshr *ampl x y z*	for uniform shear traction
edld *ampl x y z*	for uniform general edge traction
edmom *ampl*	for unform edge moment
ednor *ampl*	for uniform normal edge traction
edshr *ampl*	for uniform shear edge traction
edtra *ampl*	for uniform transverse edge traction

Remarks

For the meaning of these options, types, and parameters, see the ABAQUS manual under the *DSLOAD command.

An ABAQUS *SURFACE card is formed in the output file for ABAQUS for every face set built and the name of the face set can be used as an argument in this command. If a load curve is formed using the **lcd** or **flcd** command, the load curve will be referenced in the ABAQUS *AMPLITUDE card and can be referenced in this **dsload** command using the **lcd** option. If the load curve is not defined, then no ABAQUS *AMPLITUDE card will be formed and the AMPLITUDE option will not show up in the ABAQUS *DSLOAD card. But it is still necessary to specify a load curve number in the **dsload** command so that an ABAQSTEP command can activate the **dsload**.

Use the **lsys** command to create a local coordinate system and associate this system with the **dsload** using the **or** option. The **abaqstep** command, used to associate conditions with different analysis steps, has an option to associate the loads defined in the **dsload** command with an analysis step.

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fa fixed nodal rotations

fa *nodes angle_x angle_y angle_z*

where

nodes can be

n *node_number*

rt *x y z*

cy *rho theta z*

sp *rho theta phi*

nset *name_of_set*

angle_x	is the rotation about the x axis,
angle_y	is the rotation about the y axis,
angle_z	is the rotation about the z axis

fc concentrated nodal loads

fc *nodes load_# amplitude fx fy fz*

where

nodes can be

n *node_number*

rt *x y z*

cy *rho theta z*

sp *rho theta phi*

nset *name_of_set*

load_#	is a load curve or case number or zero
amplitude	is a load curve multiplier,
fx	is the initial force in the x-direction
fy	is the initial force in the y-direction
fz	is the initial force in the z-direction

ffc concentrated nodal load with a follower force

ffc *nodes load_# amplitude* $node_1$ $node_2$ $node_3$

where *nodes* can be

n *node_number*

rt *x y z*

cy $\rho\ \theta\ z$

sp $\rho\ \theta\ \varphi$

nset *set_name*

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Remarks

A follower plane is defined using the three nodes. The normal of the plane defines the direction of the force. This is used for Dyna3D and Lsdyna.

Use the **co** merge phase command with the **fcc** option to view this condition.

fmom follower nodal moment

fmom *nodes load_# amplitude* $node_1$ $node_2$ $node_3$

where *nodes* can be

n *node_number*
rt *x y z*
cy ρ θ *z*
sp ρ θ φ
nset *set_name*

Remarks

A follower plane is defined using three nodes. The moment is about the normal of the follower plane. This is used for Lsdyna.

Use the **co** merge phase command with the **fmom** option to view this condition.

mom nodal moment about one global coordinate axis

mom *nodes load_curve_# moment direction*

where

load_curve_# load curve or case number or zero

nodes can be

n node_number
rt *x y z*
cy *rho theta z*
sp *rho theta phi*
nset *name_of_set*

direction can be

x,
y, or
z axis

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pr pressure load

pr fset *face_set load_curve amplitude*

where

face_set	face set name
load_curve_#	load curve or case number or zero
load_curve	is a load curve number or zero
amplitude	is an amplitude factor

Remarks

Pressure is a scalar quantity applied to a face of an element. A positive pressure acts on a face in the direction opposite the positive normal of the face.

All faces within the set are assigned the same pressure condition. When a load curve accompanies the condition, the pressure becomes time dependent. If the load curve number is zero, no load curve is specified and the pressure load is considered a constant.

The positive normal direction can be specified using the **orpt** command.

tract traction

tract fset *face_set load_curve amplitude amplitude* x_n y_n z_n

where

face_set	face set name
load_curve	is a load curve or case number or zero
amplitude	is an amplitude factor
x_n	x-component of displacement
y_n	y-component of displacement
z_n	z-component of displacement

Remarks

This command assigns traction to face sets.

14. Boundary and Constraint conditions

b nodal displacement and rotation constraints

b *nodes option_list ;*

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where
nodes can be

n *node_number*	for a node
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a pont in cylindrical coordinates
sp *rho theta phi*	for a point in spherical coordinates
nset *name_of_set*	for a node set

an *option* is

sid *n*	set identification number
dx	for x-displacement
dy	for y-displacement
dz	for z-displacement
rx	for rotation about the x-axis
ry	for rotation about the y-axis
rz	for rotation about the z-axis

followed by a value of

0	in order to initialize to no constraint
1	to constrain

Remarks

This command adds nodal constraints in the global coordinate system. Initially there are no constraints. Each **b** command modifies the constraints for the nodes of the region. Thus several commands may set different constraints for the same node. This has a cumulative effect. For example, you can remove degrees of freedom in the **x**-direction for nodes of an edge of the mesh. Then the constraint lists of all nodes along this edge are modified to reflect this constraint. Then you could place a displacement constraint in the **y**-direction on an adjoining edge of the mesh. The corner node where these two edges meet would then be supported in both the **x** and **y**-directions.

Several other commands can affect the constraints. For example, the **plan** command with the **symm** option for symmetry can add constraints if the symmetry plane is parallel to one of the coordinate planes.

Different nodes may be merged into one. The merged node inherits ALL of the constraints of the nodes which were merged into it. To view the different constraints in the model, use the **condition** command with the **dx**, **dy**, **dz**, **rx**, **ry**, or **rz** options.

Sid specifies a set identification number so that the nodal constraints are written to the NASTRAN and NE/NASTRAN output using the SPC1 and SPCADD keywords. For ABAQUS output, the set identification number becomes the load set number used in abcload option of the **abaqstep** to associate the boundary condition with a step in the analysis.

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cfc boundary conditions for the CF3D output option

cfc *id type parameters faces*

where

the *type* and *parameters* can be:

fv *vx vy vz*
ft *temperature*
fsp *species_# amplitude*
ol *pressure*
il *vx vy vz*
wall
ufl *amplitude*
vfl *amplitude*
wfl *amplitude*
tfl *temperature*
spf *species_# amplitude*
cb

faces can be a sequence of the following terminated with a ";"

lb1 *list* ; for linear brick face 1
lb2 *list* ; for linear brick face 2
lb3 *list* ; for linear brick face 3
lb4 *list ;* for linear brick face 4
lb5 *list ;* for linear brick face 5
lb6 *list ;* for linear brick face 6

list is a sequence of element numbers

or

faces can be one of the following

surface *surface_# tolerance #_nodes*

where

tolerance is the accepted distance from the surface
#_nodes is the required number of corner nodes within tolerance

set *face_set_name*

Remarks

A set of faces can be generated using options in the Part Phase with the **fset** command. Face sets can be modified or generated in the Merge Phase with either the **fset** command or the set window in the pick panel of the environment window. Only the linear brick element faces will be used by this command. It is usually better to define face sets in the Part Phase so that it is parametric. However,

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the **cfc** command is also available in the Part Phase so if you are going to define the set in the Part Phase to be used for the **cfc** command, it is simpler to use the **cfc** command in the Part Phase. In those cases when the regions in the part do not create the desired effect, use the set feature in the Merge Phase.

The surface option applies only to the nodes in the picture. This is all of the nodes in the active list of parts and materials.

Examples

```
cfc first fv 1.2 2.3 3.4 surface 33 .001 4
```

This example creates a fixed velocity boundary condition called first. All faces of the mesh with all 4 nodes within .001 distance of surface 33 will be assigned this condition.

```
cfc house ob set hs1
```

This example assigns the obstruction boundary condition called house to all of the faces in the hs1 face set.

```
cfc 3142 ol -2.34e+08 lb1 4 219 410:435; lb3 3:8 29:34;;
```

This command assigns the outlet pressure condition numbered 3142 to face number 1 of the linear brick elements 4, 219, and 410 thru 435. It assigns the same condition to face number 3 of the linear brick elements 3 thru 8 and 29 thru 34.

fbc FLUENT boundary conditions

fbc fset *face_set type zone*

where *type* can be

interior	
wall	
pr_inlet	pressure-inlet
inlet_ve	inlet-vent
intake_f	intake-fan
pr_outle	pressure-outlet
exhaust_	exhaust-fan
outlet_v	outlet-vent
symmetry	
per_shad	periodic-shadow

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pr_far_f	pressure-far-field
velocity	velocity-inlet
periodic	
fan	
porous_j	porous-jump
radiator	
mass_flo	mass-flow-inlet
interfac	interface
outflow	
axis	

il identifies an inlet for fluid flow.

il fset *face_set*

infol print nodal information with a specific load/condition

infol *nodes type*

where

nodes can be

n *node_number*	a node number
rt *x y z*	a Cartesian coordinate
cy *rho theta z*	a point in cylindrical coordinates
sp *rho theta phi*	a point in spherical coordinates
nset *name_of_set*	a node set

type can be

all *load_curve_#*	for all nodal conditions
fc *load_#*	for force
fd *load_curve_#*	for fixed displacement
fv *load_curve_#*	for fixed velocity
ft *load_curve_#*	for forced temperature
tm	for initial temperature
sw *stone_wall_#*	for stone wall
v	for electrostatic potential
acc *load_curve_#*	for acceleration
fa	for fixed angular displacement
mp	for magnetic potential
mom *load_#*	for moment

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lb *local_system_#*	for constraint in local coordinate system
ve	for velocity
te	for temperature
si *interface_#* ***S***	for nodal sliding interface
tepro *load_curve_# scale base_temp*	for Temperature

where load_curve_#, stone_wall_#, interface_#, or local_system_# can be 0 for all cases.

Remarks

This command will print a table of nodes with specific loads and conditions to the Text Window. The set of nodes can be selected by node number, location, or by containment in a set. The **b** command is not included since the **conditions** command can graphically display the nodes of a specific constraint.

Some of the types of loads include a load case or load curve number. If 0 is used, then all cases will be used in the command.

jt assign nodes to a joint

jt *joint_# local_node_# options* ;

where *local_node* depends on the type of joint (see **jd**)

where an *option* can be

n node_number	use an existing node
p *x y z rigid_body_material_#*	new node in Cartesian coordinates
c *rho theta z rigid_body_material_#*	new node in cylindrical coordinates
s *rho theta phi rigid_body_material_#*	new node in spherical coordinates
v *x_offset y_offset z_offset*	offset the new node
dx	x-displacement constraint on new node
dy	y-displacement constraint on new node
dz	z-displacement constraint on new node
rx	x-rotation constraint on new node
ry	y-rotation constraint on new node
rz	z-rotation constraint on new node

Remarks

This command assigns a node to a numbered joint defined by **jd**. Each node in a joint is assigned a node sequence number which is referred to as its local node number. Each type of joint requires a different number of nodes and the role a node plays in the joint depends on the joint type and the node local node number. The simplest example of a joint is when a set of nodes share constraints. In this

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case the ordering of the nodes are not important. Up to 16 nodes can be included in a shared constraint joint. The nodal constraints available in this command apply to new nodes defined using the p, c, or s options.

kivabc boundary conditions for kiva4 output format

kivabc *face_set type*

where *type* can be

moving	for moving piston
movingb1	for moving bottom face of 1st valve
movingt1	for moving top face of 1st valve
movingb2	for moving bottom face of 2nd valve
movingt2	for moving top face of 2nd valve
movingb3	for moving bottom face of 3rd valve
movingt3	for moving top face of 3rd valve
movingb4	for moving bottom face of 4th valve
movingt4	for moving top face of 4th valve
solid	for a face of a solid
solidh	for a solid face of a cylinder head
axis	for a face on the axis
fluid	for a fluid face (default)
inflow	for an inlet
outflow	for an outlet
presin	for pressure inflow
presout	for pressure outflow
userdef *value*	for user defined condition

lb local nodal displacement and rotation constraints

lb *nodes system_# option_list ;*

where a *system* is defined using the command **lsys**

where an *option* is

dx *flag*	for x-displacement
dy *flag*	for y-displacement
dz *flag*	for z-displacement
rx *flag*	for rotation about the x-axis
ry *flag*	for rotation about the y-axis
rz *flag*	for rotation about the z-axis

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where the *flag* can be

0	for initialize to no constraint
1	for constrain

where *nodes* must be one of

n *node*	for an existing node
rt *x y z*	for a new node in Cartesian coordinates
cy *rho theta z*	for a new node in cylindrical coordinates
sp *rho theta phi*	for a new node in spherical coordinates
nset *name_of_set*	for a node set

Remarks

Use this command to set constraints on nodes that cannot be set using the **b** or **bi** commands because they are restricted to the global coordinate system. Care is needed not to over specify the constraints on a node. No warnings are given if a node is over constrained. Use the **lsys** command to define the local coordinate system.

mpce multiple point constraint equation

mpce *id depend_node dof option independ_nodes* ;

where *id* is a positive integer uniquely identifying this equation

where *depend_node* can be

n *node_#*	existing node
pm *point_mass_#*	exiting point mas
rt *x y z constraints* ;	create a new node in Cartesian coordinates
cy *ρ θ z constraints* ;	create a new node in cylindrical coordinates
sp *ρ θ φ constraints* ;	create a new node in spherical coordinates

where *constraints* can be any of

dx	for x-displacement constrained
dy	for y-displacement constrained
dz	for z-displacement constrained
rx	for x-rotation constrained
ry	for y-rotation constrained
rz	for z-rotation constrained

where *option* can be

c0 *constant*	constant term in equation (zero default)

where *dof* can be any one of

dx	for x-displacement of dependent node

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dy	for y-displacement of dependent node
dz	for z-displacement of dependent node
rx	for x-rotation of dependent node
ry	for y-rotation of dependent node
rz	for z-rotation of dependent node

where *independ_nodes* must be one of

nsets [*node_set dof*]	for independent nodes in sets
nlists [*node_list ; dof wieght*]	for independent nodes in lists

Remarks

There are two ways to specify the independent nodes. When **nsets** is used, multiple node sets can be used. There is only one degree of freedom (*dof*) for each node set specified. Weights are all assumed to be 1.

Alternatively, if **nlists** is used, then lists of node numbers are required. There is only one degree of freedom (*dof*) and *weight* for each node list.

Suppose there are n independent nodes from either **nsets** or **nlists**. Also, let ND_i be the nodal component of node i indicated by the corresponding DOF with C_i is the corresponding weight. Then the multiple point constraint equation has the following explicit form:

$$ND_0 = C_0 + \sum_{1}^{n} ND_i \times C_i$$

To check the MPCEs, use the **co** command with the **mpce** option to display the relation between the dependent node and the independent nodes. They are shown by a line connecting the dependent node and each of the independent nodes. The dependent node has a special mark. This works for wire, hide, fill, and Fast Graphics.

In all three commands (**mpce**, **mpc**, and **joint**) that generate multiple point constraints, if any two nodes are in a MPC, they cannot be merged. All merge commands (**t**, **tp**, **st**, and **stp**) will avoid merging such pairs.

Example

```
block 1 6 11 16;1 6 11 16;1 3; -3 -1 1 3 -3 -1 1 3 0 1
dei 2 3; 2 3;;
sd 1 cy 0 0 0 0 0 1 1
```

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```
sfi -2 -3; -2 -3;;sd 1
merge
curd 1 arc3 whole rt 1 0 0 rt 0 1 0 rt -1 0 0 ;
curd 2 arc3 whole rt 1 0 .5 rt 0 1 .5 rt -1 0 .5 ;
curd 3 arc3 whole rt 1 0 1 rt 0 1 1 rt -1 0 1 ;
nset one = crv 1 .01
nset two = crv 2 .01
nset three = crv 3 .01
mpce 1 rt 0 0 0 ; dx nsets one dx ;;
mpce 2 rt 0 0 .5 ; dx nsets two dx ;;
mpce 3 rt 0 0 1 ; dx nsets three dx ;;
co mpce
```

In this example, a hole is formed in a brick mesh. Then three circles corresponding to the edges exposed by the hole are formed. Node sets are formed from these circles. Finally, MPCEs are formed from these three nodes sets and displayed.

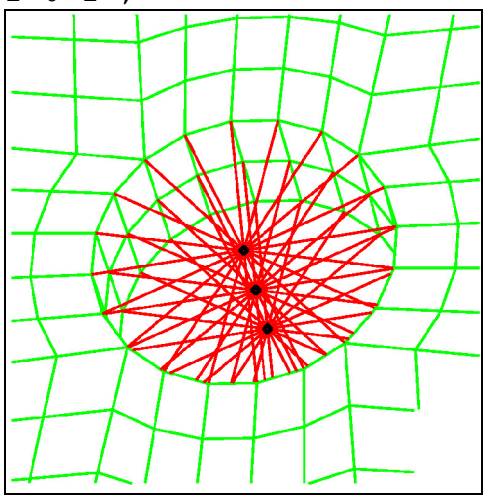

Figure 248 3 Spider Web MPCEs

nr non-reflecting boundary

nr fset *face_set*

ol identifies a face of the mesh as an outlet for fluid flow

ol fset *face_set*

rigid create a rigid body from a nodal set

rigid *node_set_name options ;*

where

options can be

cid *local_sys_id*

rgm *rigid_body_material_number* for NIKE3D

com *x_center y_center z_center*

trm *translational_mass*

lcid *local_coordinate_system_for_inertia* for inertia tensor

ixx *xx-component_of_inertia_tensor*

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ixy *xy-component_of_inertia_tensor*
ixz *xz-component_of_inertia_tensor*
iyy *yy-component_of_inertia_tensor*
iyz *yz-component_of_inertia_tensor*
izz *zz-component_of_inertia_tensor*
trv *x_velocity y_velocity z_velocity*
rtv *x_rot_velocity y_rot_velocity z_rot_velocity*

Remarks

This command allows you to append nodes to an existing rigid body for NIKE3D, or to create a nodal rigid body for LS-DYNA. The nodes appended to a NIKE3D rigid body inherit constraints, and so no constraints can be specified; the material number of the existing rigid body is required. For LS-DYNA, the nodal rigid body is created and all properties of the rigid body can be specified.

rml remove specific loads or conditions on a set of nodes

rml *nodes type*
where
nodes can be

n *node_number*	a node number
rt *x y z*	a Cartesian coordinate
cy *rho theta z*	a cylindrical coordinate
sp *rho theta phi*	a spherical coordinate
nset *name_of_set*	a node set

type can be

all *load_curve_#*	for all conditions for a load case (curve)
fc *load_#*	for Nodal force by load curve
fd *load_curve_#*	for Nodal displacement
fv *load_curve_#*	for Nodal velocity
ft *load_curve_#*	for Nodal velocity
tm	for Initial temperature
sw *stone_wall_#*	for Nodes impacting a stone wall
v	for Electrostatic temperature
acc *load_curve_#*	for Nodal acceleration
fa	for Fixed nodal angular acceleration
mp	for Magnetic potential
mom *load_#*	for Nodal moment about an axis
lb *local_system_#*	for Local coordinate nodal constraints

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ve for Velocity
te for Temperature
si *interface_#* **S** for Nodal sliding interface
tepro *load_curve_# scale base_temp* for Temperature
where load_curve_#, stone_wall_#, interface_#, or local_system_#
can be 0 for all cases

Remarks

There is no undo command for the Merge Phase. However, you can remove a condition placed on a node set using the **rml** command. The **rml** command is more general than an **undo** because it can be used to remove conditions from any node set, whether or not the condition was applied directly to the node set in the first place.

rsl restore specific loads or conditions on a set of nodes

rsl *nodes type*
where
nodes can be
n *node_number* a node number
rt *x y z* a Cartesian coordinate
cy *rho theta z* a cylindrical coordinate
sp *rho theta phi* a spherical coordinate
nset *name_of_set* a node set
type can be
all *load_curve_#* for all conditions for a load case (curve)
fc *load_#* for Nodal force by load
fd *load_curve_#* for Nodal displacement
fv *load_curve_#* for Nodal velocity
ft *load_curve_#* for Nodal velocity
tm for Initial temperature
sw *stone_wall_#* for Nodes impacting a stone wall
v for Electrostatic temperature
acc *load_curve_#* for Nodal acceleration
fa for Fixed nodal angular acceleration
mp for Magnetic potential
mom *load_#* for Nodal moment about an axis
lb *local_system_#* for Local coordinate nodal constraints
ve for Velocity
te for Temperature

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si *interface_#* **S** for Nodal sliding interface
tepro *load_curve_# scale base_temp* for Temperature
where load_curve_#, stone_wall_#, interface_#, or local_system_# can be 0

Remarks

This command is the inverse of **rml**.

spotweld interactive selection of spot welds

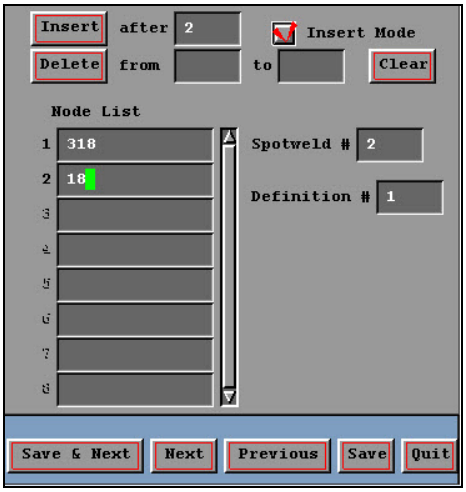

Figure 249 Spotweld dialogue box

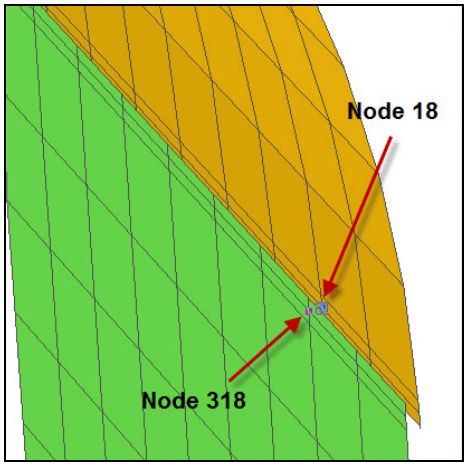

Figure 250 Selecting nodes

spotweld *(no arguments)*

Remarks

This feature is an easy and interactive way to create multiple spot welds using the mouse to click on the nodes. It is only available in the merge phase. It creates **spw** commands and writes them to the tsave (session) file.

The ***Spotweld #*** window should be set to the first spot weld number to be generated (see the **spw** command). Also select the appropriate spot weld property definition number (see the **spwd** command). Then click on nodes to be tied as a spotweld. Typically only two are needed, but NASTRAN allows for many nodes to be tied as a rigid body, so you can click on more than 2 nodes for NASTRAN. Then click on the ***Save & Next*** button. Repeat this for every spot weld.

An existing spot weld can be changed. Select the spot weld number by typing it into the ***Spotweld #***

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window and entry. The node numbers can be changed. Deselect the ***Insert Mode*** check mark. Click on the node number in the Node List that you wish to change and then select a new node in the picture of the mesh in the physical window. Nodes can be deleted (use the ***Delete*** button) or a node can be inserted (use the ***Insert*** button) after a node already in the Node List. Be sure to ***Save*** this new spot weld.

You can scroll through the existing spot welds by clicking on the ***Next*** or ***Previous*** button. The nodes in the physical window will be marked.

Use the **co spw** feature to view each spot weld.

spw create a single spot weld

spw *id option nodes* ;

where

id	positive integer identifying the spot weld

where an *option* can be

spwd *spwd_id*	see **spwd** for the property number

where a *node* can be

ann *node_#*	which can be any literal node number
node *node_#*	which must be an existing node number
loc *x0 y0 z0 material_#*	to locate closest node
locc *x0 y0 z0 material_#*	to locate closest node and change its location

Remarks

A literal node number will be output exactly as it is specified in this command. It does not have to refer to an existing node. This option makes it possible to spot weld between two separately generated meshes. Typically one or both meshes have node offsets (see **offset**).

This will generate the

*CONSTRAINED_SPOTWELD

and

*CONSTRAINED_SPOTWELD_FILTERED_FORCE

cards in LSDYNA. Use the **spwd** command to define the properties of the spot welds. When doing so, be sure to specify both the number of force vectors and the time window to activate the filtered force

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card.

For NASTRAN, the RBE2 card is generated.

Use the **co spw** feature to view each spot weld.

spwd spot weld property definitions

spwd *id options* ;

where

id	positive integer identifying the spot weld properties definition

where an *option* can be

dx	displacement in the x-direction for all nodes
dy	displacement in the y-direction for all nodes
dz	displacement in the z-direction for all nodes
rx	rotation about the local x-axis for all nodes
ry	rotation about the local y-axis for all nodes
rz	rotation about the local z-axis for all nodes
sn	nornal_force
ss	shear_force
n	normal_force_exponent
m	shear_force_exponent
tf	failure_time
ep	effective_plastic_strain
nf	#_force_vectors
tw	time_window

Remarks

This command defines the properties of a spotweld. These definitions are needed by the **spw** command and the associated interactive command **spotweld**. The shared degrees of freedom flags are used for NASTRAN. The other options are for LSDYNA.

spwf spot welds for LSDYNA material 100

spwf *points sw_material 1st_contact 2nd_contact*

where *points* can be

eqsp *3D_curve_# flag #_spotwelds*

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where *flag* can be

0	to include the last point
1	to not include the last point

pnts *3D_curve_# flag*

where *flag* can be

0	to include the last point
1	to not include the last point

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates

where *1st_contact* and *2nd_contact* can be

mat *material_number*

Remarks

Use the **co spwf** feature to view each spot weld.

sw select nodes that may impact a stone wall

sw *nodes stone_wall_#*

where

nodes can be:

n *node_number*	node number
rt *x y z*	a Cartesian coordinate
cy *rho theta z*	a cylindrical coordinate
sp *rho theta phi*	a spherical coordinate
nset *name_of_set*	a node set

Remarks

First use the **plane** command to define the stone wall.

syf assign faces to a numbered symmetry plane with failure

syf fset *face_set symmetry_plane_# failure*

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trp create tracer particles

trp *tracking list_options ;*

where

tracking can be **fixed** or **free**, and

an *option* is followed by some parameters

time *start_time*

point *x0 y0 z0*

lnpy *x1 y1 z1 x2 y2 z2 #_tracers*

Remarks

This defines the location and properties of Tracer Particles in Ls-dyna.

15. Radiation and Temperature Commands

These commands let you set various boundary conditions related to radiation and temperature. See also the radiation and temperature commands in the Part Phase.

bf bulk fluid

bf fset *face_set id_# load amplitude a b*

where

face_set	name of the face set
id_#	**bfd** bulk fluid identification number
load	load curve number
amplitude	multiplier of the load curve
a	exponent a
b	exponent b

Remarks

Use this command to identify those faces (surfaces) which are to be part of the bulk fluid (node) calculation.

Use the **orpt** command to orient the faces as desired.

Use the **co** command with the **bf** option to display the bulk fluid faces.

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cv boundary convection

cv fset *face_set load_curve$_1$ amplitude$_1$ load_curve$_2$ amplitude$_2$ exponent*

where

face_set	name of the face set
load_curve$_1$	first load curve number or zero,
amplitude$_1$	amplitude factor for the first load curve,
load_curve$_2$	second load curve number or zero,
amplitude$_2$	amplitude factor for the second load curve, and
exponent	exponent.

Remarks

First use the **orpt** command to specify the surface orientation; that is, how to orient the normal vector.

A zero load curve number means that the condition is constant in time. If a curve is specified which has not been defined, a warning message will be issued.

cvt convection thermal loads

cvt fset *face_set coefficient temperature*

where

face_set	name of the face set
coefficient	film coefficient
temperature	temperature near convection

Remarks

First use the **orpt** command to specify the surface orientation; that is, how to orient the outward normal vector.

This command is used to create convection thermal loads for ANSYS. The first parameter is the film *coefficient*. This is followed by the *temperature* near convection. This command will create the EP cards for the ANSYS input file.

fl prescribed boundary flux

fl fset *face_set load_curve_# amplitude*

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where

face_set	name of the face set
load_curve_#	a load curve number
amplitude	amplitude constant

Remarks

First use the **orpt** command to specify the surface orientation; that is, how to orient the outward normal vector.

ft prescribed temperature

ft *nodes load_# temperature*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point spherical coordinates
nset *name_of_set*	for a node set

Remarks

This specifies a time-dependent temperature boundary condition on the nodes. It is assumed that the temperature used by the appropriate simulation code will be the product of the temperature and the amplitude of the load curve at the appropriate time in the simulation.

rb prescribed radiation boundary condition

rb fset *face_set load_curve$_1$ amplitude$_1$ load_curve$_2$ amplitude$_2$*

where

face_set	name of the face set
load_curve$_1$	a load curve number
amplitude$_1$	amplitude constant
load_curve$_2$	a load curve number
amplitude$_2$	amplitude constant

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Remarks

First use the **orpt** command to specify the surface orientation; that is, how to orient the outward normal vector.

If a load curve number is specified as zero, then the condition is constant in time.

re radiation enclosure

re fset *face_set 0 temperature obstruction_flag*

where

face_set	name of the face set
temperature	constant
obstruction flag is:	
yes	to include surface obstruction calculations
no	to not include surface obstruction calculations

Remarks

First use the **orpt** command to specify the surface orientation; that is, how to orient the outward normal vector.

This command generates enclosure radiation data for TOPAZ3D. Use the **emissivity** and **rband** commands to specify the emissivity and wavelength breakpoint tables associated with the enclosure radiation data. For more details, see the TOPAZ3D manual.

te constant temperature for all nodes

te *nodes load_# temperature*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point spherical coordinates
nset *name_of_set*	for a node set

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Remarks

These temperatures will override the temperatures specified by **temp**. When this command is invoked, all nodal temperatures are assumed to be specified. See the **temp** commands.

tepro nodal temperature profile

tepro *nodes load_# amp_expr ; temp_expr ;*

where

nodes can be

n *node_number*	for a node number or zero
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point spherical coordinates
nset *name_of_set*	for a node set

amp_expr for a FORTRAN expression forming the scale factor

temp_expr for a FORTRAN expression forming the base temperature

Remarks

Both the base temperature and load curve scaling factor can be functions of the x, y, and z-coordinates of the node. This command is used for input to DYNA3D, LS-DYNA, and NIKE3D.

tm initial temperature

tm *nodes load_# temperature*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point spherical coordinates
nset *name_of_set*	for a node set

Remarks

This command does not invoke temperatures for all nodes like the **temp** commands and the **te** and **tei** commands.

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vvhg volumetric heat generation w/ functional amplitude

vvhg *set_name load_# expression ;*

where

set_name	element set name
load_#	load curve number
expression	amplitude expression

Remarks

The expression can be any valid FORTRAN like expression with x, y, and z. The coordinates used in this expression will be the centroid of the element.

16. Electric Conditions

efl electric flux boundary condition

efl fset *face_set value_of_flux*

Remarks

This command produces four-node polygons with an assigned constant flux, one polygon for each face within the region.

mp constant magnetic potential

mp *nodes potential*

where

nodes can be

n *node_number*	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
nset *name_of_set*	for a point spherical coordinates
sp *rho theta phi*	for a node set

v constant nodal electrostatic potential boundary condition

v *nodes potential*

where

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nodes can be

n *node_numbe*r	for a node number
rt *x y z*	for a point in Cartesian coordinates
cy *rho theta z*	for a point in cylindrical coordinates
sp *rho theta phi*	for a point spherical coordinates
nset *name_of_set*	

17. Sets

These commands are unique to the Merge phase. Named sets are a useful tool in the definition of boundary conditions and properties in the mesh and form an alternative to selecting regions in the Part phase. An arbitrary selection can be made and this is the advantage to using set functions in the Merge phase. The disadvantage is that the selection may no longer be parametric. If you go back and make a change to the mesh, you may have to redefine the set since the element and node numbers have changed.

The name of the set can be up to 8 alphanumeric characters long. Each name of the set must be unique. Since LSDYNA numbers its sets, one can use an integer as the name of the set.

In some of the set commands, the logical or Boolean set operators AND and OR are used to create new sets from existing sets. The AND operator between two sets means to take their intersection. This should not be confused with the common usage of and which might be interpreted to mean the addition of two sets. The OR operator does this function.

The **True*Grid*®** internal face IDs of elements depends on the element geometry after nodes have been properly merged so that **True*Grid*®** can determine the element geometry (i.e. hexa, penta /wedge, or tetra). Pyramid elements are treated as degenerate hexa. These face IDs can be seen when in the merge phase by using the **la facesel** command. The face IDs differ among the different simulation codes. You need to interpret what you see in **True*Grid*®** graphics to correspond with the simulation code. This interpretation of the face ID numbers is only needed when using the **la facesel** command in the merge phase. When **True*Grid*®** writes an output file for a particular simulation code, it does the translation of the face ID numbers appropriate for that simulation code.

adnset add nodes to an ordered node set

adnset *set_name insertion_node before/after list_of_nodes*

where

set_name	is the name of the node set
insertion_node	is the sequence number of the node for the insertion point
before/after	is either

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0	for the insertion before the insertion node, or
1	for the insertion after the insertion node
list_of_nodes	is the list of nodes, which will be inserted to the ordered node set.

Remarks

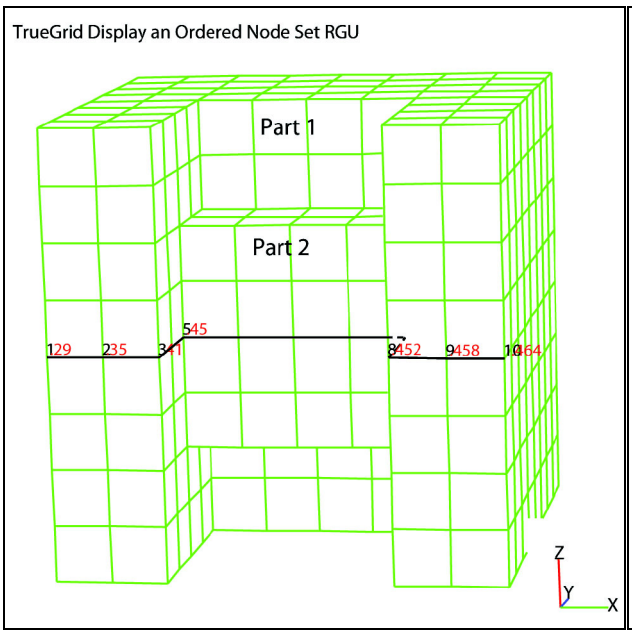

Figure 251 before **adnset**

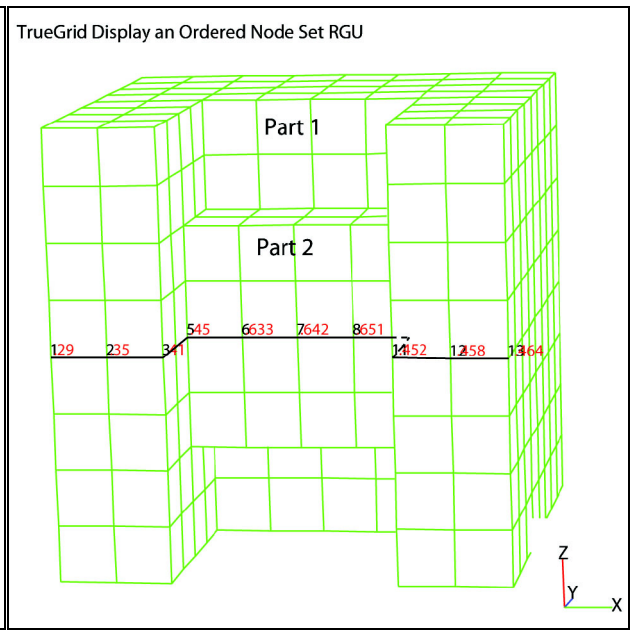

Figure 252 after **adnset**

This command adds or inserts a list of nodes into an ordered node set.

Example

Part 1 is created and a region is deleted. The node set named RGU is created using the index progression `1 2;-1;-3;`. The index progression `-2;1 2 ;-3;` is used to enhance the RGU node set by union. Part 2 is defined. Ordered node set RGU is labeled (**Figure 251**).

Then the nodes `633 642 651` are added to the ordered node set RGU, after sequence node number `5`. The result is displayed on **Figure 252**. The command file follows:

```
block 1 3 7 9; 1 3 5 7 9;1 3 5 7 9;1 3 7 9; 1 3 5 7 9;1 3 5 7 9;
dei 2 3;1 3;1 5;
nseti  1 2;-1;-3;= RGU
```

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```
nseti -2;1 2 ;-3;OR RGU
block 1 5;1 3;1 3 5;3 7;3 5;3 5 7;
merge
dap
nset RGU OR L 456 454 452 458 464 ;
labels onset RGU 1 0
adnset RGU 5 1 633 642 651 ;
labels onset RGU 1 0
```

crvnset order a segment of an ordered node set (3D curve)

crvnset *set_name first_node last_node 3D_curve-#*

where

set_name	is the name of the node set
first_node	is the sequence number of the node
last_node	is the sequence number of the node
3D_curve-#	is the number of the 3D curve used for reordering of the nodes in the set.

Remarks

The nodal sequence number, referred to above, can be from 0 to the total number of nodes in the set. The sequence number **1** is always the first node in the set and sequence number **0** is the last node number in the set. Each node is projected to a 3D curve and it is assigned the arc length of the projected point on the 3D curve. Then the nodes are ordered according to the arc length. The effect is that the ordered set of nodes will approximate the shape and order of the 3D curve. This can be useful if the mesh is to be parameterized. Direction of the 3D curve plays an important role. If it is reversed, the ordering of the nodes in the set will be reversed. Use a negative curve number to accomplish this.

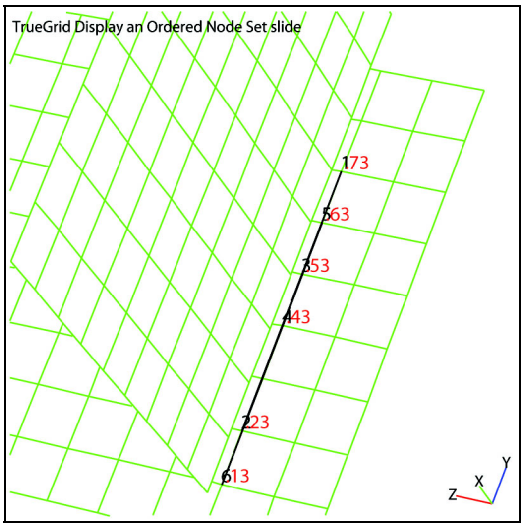

Figure 253 before **crvnset**

Example

The node set **slide** is at first defined by node numbers without any order (**Figure 253**). Then it is ordered according to curve 1 (**Figure 254**).

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```
block -1;1 10;1 10;1;1 10;1 10;
block 1 10;1 10;-1;1 10;1.8 11.1;3.2;
merge
nset slide = L 73 23 53 43 63 13;
labels onset slide 1 0
curd 1 lp3 1.3 2 2 1.5 10 2;;
crvnset slide 1 0 1
```

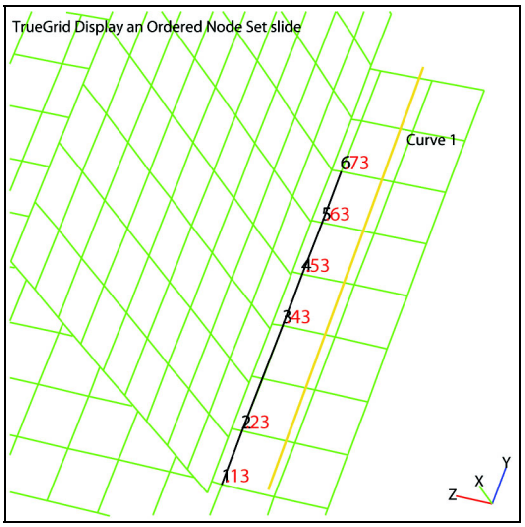

Figure 254 after **crvnset**

eset add/remove elements to/from a set of elements

eset *set_name operator option*;

where

set_name is the name of the element set

operator can be

=	for initial assignment,
AND	for intersection with element set,
OR	for union with the element set,
-	for removal from the element set
RPL	to replace one element type (only with **LS**, **LB**, **LBM**, **QS**, and **QB**)

option can be

LBM *element_list* ;	linear beam element list
QBM *element_list* ;	quadratic beam element list
LS *element_list* ;	linear shell element list
QS *element_list* ;	quadratic shell element list
LB *element_list* ;	linear brick element list
QB *element_list* ;	quadratic brick element list
S set_name	a set of elements
MT *material*	elements of a material
SRF *surface tolerance #_nodes*	elements within a tolerance of a surface
CRV *curve tolerance #_nodes*	elements within a tolerance of a curve

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Remarks

The initial assignment creates an element set. If the element set with the same name already existed, then it is deleted and recreated. The intersect operator redefines an element set to be only those elements which are found to be both in the original set and among the selected elements. Selected elements can be added by using the union operator. This causes any selected elements to be included in a set, if it is not already in that set. The minus operator removes all elements in a set which are among the selected elements.

The **SRF**, **CRV**, and **MT** options only consider those elements that are in the display set. If a part or material is removed from the picture, those elements will not be searched. **SRF** and **CRV** options have a number of nodes as a parameter which varies depending on the type of element. An element will be selected if the specified number of corner nodes of the element are within tolerance. 0 nodes means to use the element centroid.

esetc assign a comment to a element set

esetc *set_name comment*

Remarks

The esetc command can be used to set the title for an element set for LSDYNA.

fset add/remove faces to/from a set of faces

fset *set_name operator faces*

where *set_name* is the name of the face set

where *operator* can be

=	for initial assignment
AND	for intersection with face set
OR	for union with the face set
-	for removal from the face set

where *faces* can be

srf *surface_id tolerance #_nodes*	specified number of nodes fall within the specified distance from the surface
set *set_name*	set operation with this previously defined face set
enumerations ;	each type of face must be in a separate list

where an *enumeration* can be

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ls *list_elements* ;	list linear shell elements
lb1 *list_elements* ;	list linear bricks
lb2 *list_elements* ;	list linear bricks
lb3 *list_elements* ;	list linear bricks
lb4 *list_elements* ;	list linear bricks
lb5 *list_elements* ;	list linear bricks
lb6 *list_elements* ;	list linear bricks
qs *list_elements* ;	quadratic shell elements
qb1 *list_elements* ;	list quadratic bricks
qb2 *list_elements* ;	list quadratic bricks
qb3 *list_elements* ;	list quadratic bricks
qb4 *list_elements* ;	list quadratic bricks
qb5 *list_elements* ;	list quadratic bricks
qb6 *list_elements* ;	list quadratic bricks

Remarks

The initial assignment creates a face set. If the face set with the same name already existed, then it is deleted and recreated. The intersect operator redefines a face set to be only those faces which are found to be both in the original set and among the selected faces. Selected faces can be added by using the union operator. This causes any selected faces to be included in a set, if it is not already in that set. The minus operator removes all faces in a set which are among the selected faces. See also pg.277 for labeling of the face sets.

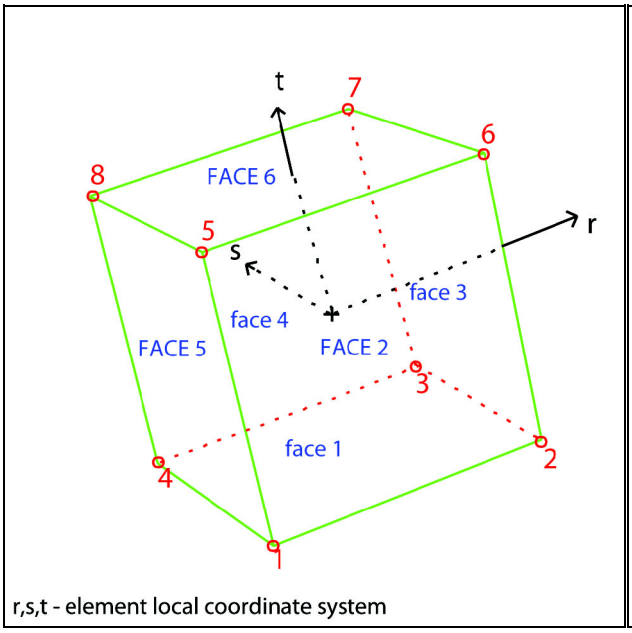

Figure 255 face order in brick element

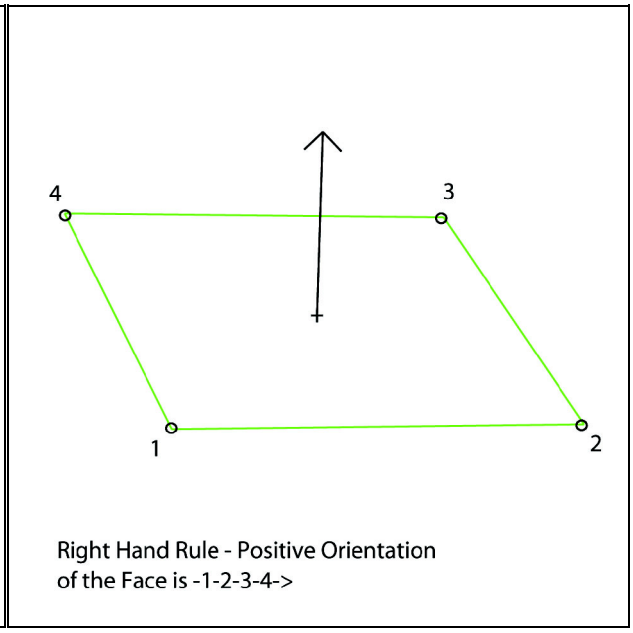

Figure 256 orientation of nodes in face

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Faces in face sets are identified by an element number and an order number of a face in the element. Face order in the linear hexa element is shown in 255. Nodes in the face are ordered by the right hand rule The vector in 256 is always oriented outward from the element. A similar method is used to identify the faces for every geometric element type. They are not enumerated in this manual because it is complicated details that are not necessary when using sets or applying loads. There are numerous geometric ways of selecting a face set which relieve the user from the tedium of face identification. If you need to know how a face is identified, make a face set and, using the **la** command in the merge face, label the face set.

The **SRF** option only considers those elements that are in the display set. If a part or material is removed from the picture, those elements will not be searched. The **SRF** option has a number of nodes as a parameter which varies depending on the type of element. An element will be selected if the specified number of corner nodes of the element are within tolerance. 0 nodes means to use the element centroid.

fsetc assign a comment to a face set

fsetc *set_name comment*

Remarks

The fsetc command can be used to set the title for a face set for LSDYNA.

infol information on nodal loads

infol *nodes type*

where

nodes can be

N *node_number*	node number
RT *x y z*	node closest to Cartesian point
CY *rho theta z*	node closest to cylindrical point
SP *rho theta phi*	node closest to spherical point
NSET *name_of_set*	node set by name

type can be

ALL *load_curve_#*	any type
FC *load_curve_#*	forces
FD *load_curve_#*	fixed displacements

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FV *load_curve_#*	fixed velocity
FT *load_curve_#*	forced temperature
TM	initial temperature
SW *stone_wall_#*	stone wall
V	electrostatic potential
ACC *load_curve_#*	acceleration
FA	fixed angular displacement
MP	magnetic potential
MOM *load_curve_#*	moment
LB *local_coordinate_system_#*	constraint in local coordinate system
VE	velocity
TE	temperature
SI *sliding_interface_#*	nodal sliding interface
NPB	nodal print blocks
TEPRO *load_curve_#*	temperature profile

where

load_curve_#, *stone_wall_#*, *local_coordinate_system_#* can be **0** for all cases.

Remarks

Command **infol** prints information about nodes with a specific load or condition. The results on the **infol** are printed to the dialog window and to the session file. See the **rml** command for example of use.

mvnset move a subset of nodes in an ordered node set

mvnset *set_name first_node last_node insertion_node before/after*

where

set_name	is the name of the node set
first_node	is the first node sequence number of the group of nodes to be moved
last_node	is the last node sequence number of the group of nodes to be moved
insertion_node	is the sequence number of the node for the insertion point
before/after	is either
0	for before the insertion node
1	for after the insertion node

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Remarks

Command **mvnset** moves a subset of nodes in an ordered nodal set. This can be used with **onset** to get the proper ordering of nodes. If the first sequence number is greater than the second sequence number, then the interval of nodes will have their order reversed before they are inserted back into the ordered set.

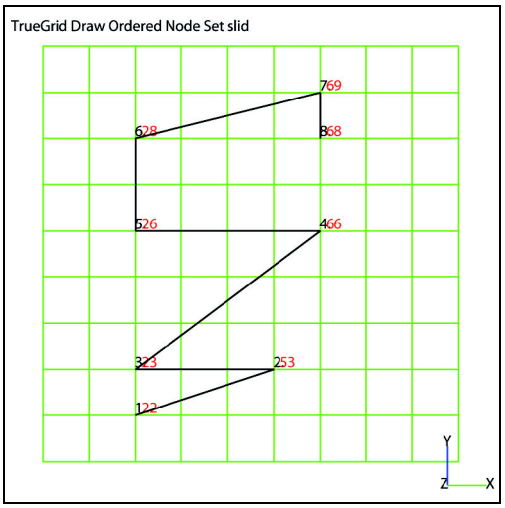

Figure 257 before **mvnset**

Example

The node set SLID is initialized by the list of nodes (**Figure 257**). The node set is reordered by **mvnset** command (**Figure 258**).

```
block 1 10;1 10;-1;1 10;1 10;0;
merge
nset SLID = L 22 53 23 66 26 28 69 68
;
labels onset SLID 1 0
mvnset SLID 2 2 3 1
mvnset SLID 7 7 8 1
labels onset SLID 1 0
```

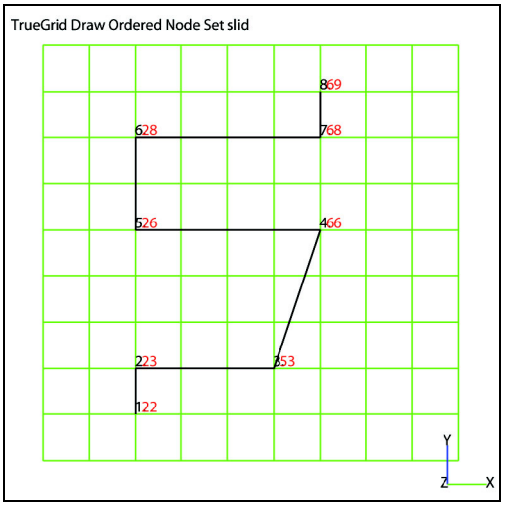

Figure 258 after **mvnset**

nset add/remove nodes to/from a set of nodes

nset *set_name operator node_selection*

where *set_name* is the name of the node set

where *operator* can be

=	for initial assignment
AND	for intersection with node set
OR	for union with the node set
+	to append the selected nodes to the node set

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- for removal from the node set

where *node_selection* can be

L *list;* list of node numbers

S *set_name* named set of nodes

SRF *surface_number tolerance* nodes near a surface

CRV *curve_number tolerance* nodes near a curve

MT *material_number*

LOAD *load* nodes with a specific load

where *load* can be

ALL *load_curve_#* any type

FC *load_curve_#* forces

FD *load_curve_#* fixed displacements

FV *load_curve_#* fixed velocity

FT *load_curve_#* forced temperature

TM initial temperature

SW *stone_wall_#* stone wall

V electrostatic potential

ACC *load_curve_#* acceleration

FA fixed angular displacement

MP magnetic potential

MOM *load_curve_#* moment

LB *local_coordinate_system_#* constraint in local coordinate system

VE velocity

TE temperature

SI *sliding_interface_#* nodal sliding Interface

NPB nodal Print blocks

TEPRO *load_curve_#* temperature Profile

B *constraint flag*

where *constraint* can be

dx constrain the x translation

dy constrain the y translation

dz constrain the z translation

rx constrain the rotation about the x axis

ry constrain the rotation about the y axis

rz constrain the rotation about the z axis

where *flag* can be

1 constrained

0 unconstrained

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Remarks

This command is probably the easiest and most versatile method for constructing a node set. The initial assignment creates a node set. If the node set with the same name already existed, then it is deleted and recreated. The intersect operator redefines a node set to be only those nodes which are found to be both in the original set and among the selected nodes. Selected nodes can be added by using the union operator. This causes any selected nodes to be included in a set, if it is not already in that set. The add operator will always append selected nodes to a set. This is used to create ordered node sets where duplicate nodes are allowed. The minus operator removes all nodes in a set which are among the selected nodes.

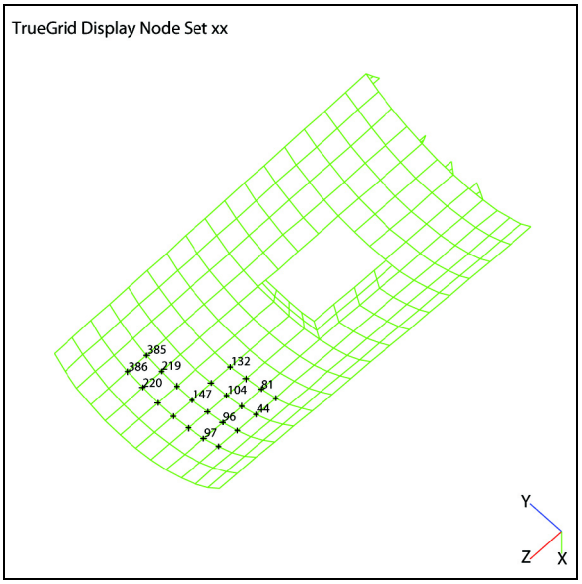

Figure 259 Node Set by **nset**

Example

The node set is defined by the list of node numbers in an arbitrary order (**Figure 259**).

```
cylinder 1 -3;1 -5 -7 -9 -13;1 -5 -9 -13 17;
          2.5 3;-45 -15 0 15 45;1 3 5 7 9;
dei -2; 2 4; 2 3;
dei 1 2; -3; 2 3;
merge
nset xx = L
 44 45 81 96 97 104 105 132 147
148 219 220 316 329
 330 339:341 351 352 385 386;
labels nodeset xx
```

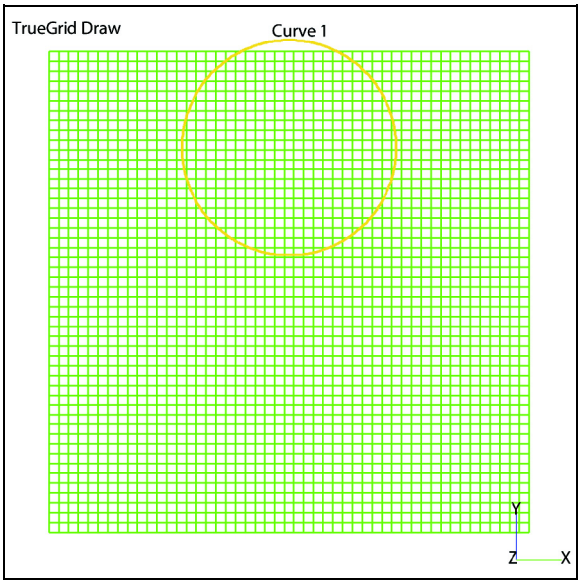

Figure 260 Mesh and Curve

Example

This example demonstrates the parametric capabilities of node sets. The mesh and the curve (circle) are shown (**Figure 260**). Node set Press_1 is defined by a distance .3 from curve number 1 (**Figure 261**). Node set Press_2 is defined by a distance of 1 from curve number 1 (**Figure 262**).

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```
block 1 50;1 50;-1;0 10;0 10; 0;
curd 1 arc3 whole rt 7 7 0
 rt 6 6 0 rt 4 6 0; merge
nset Press_1 = crv 1 .3
labels nodeset  Press_1
nset Press_2 = crv 1 1


labels nodeset  Press_2
```

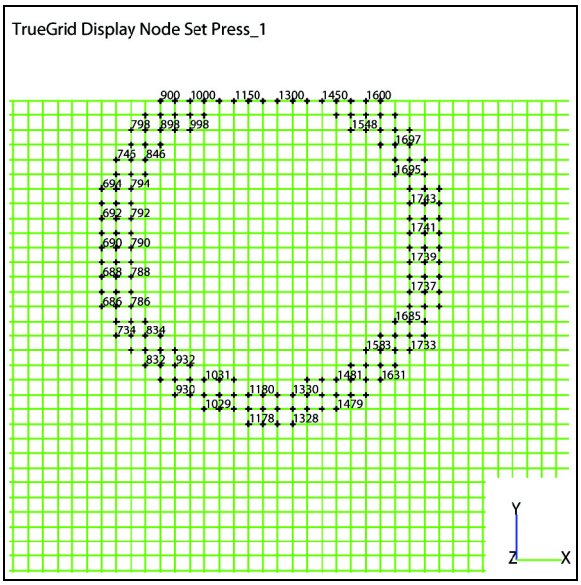

Figure 261 Node Set Press_1

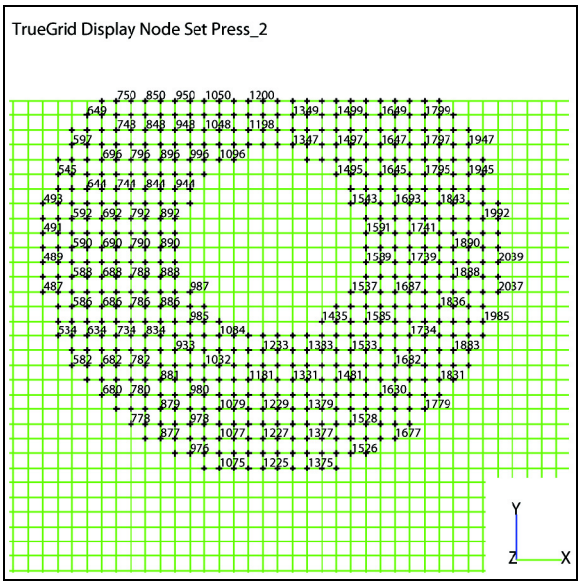

Figure 262 Node Set Press_2

Example

Another useful application of the **nset** command is the definition of a sliding interface (**Figure 263**). The node set RAD1 is composed of all nodes within a distance of 2 units from the cylinder (**Figure 263**). The result of the **nset** command is on the **Figure 264**.

```
nset RAD1 = srf 6 2
```

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labels nodeset RAD1

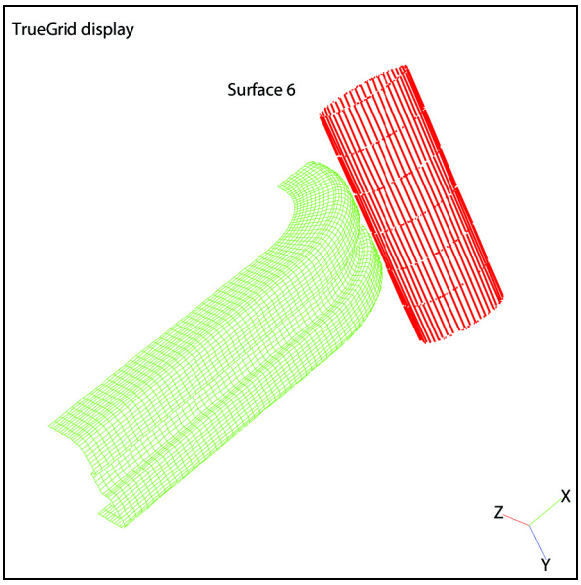

Figure 263 Cylinder and mesh

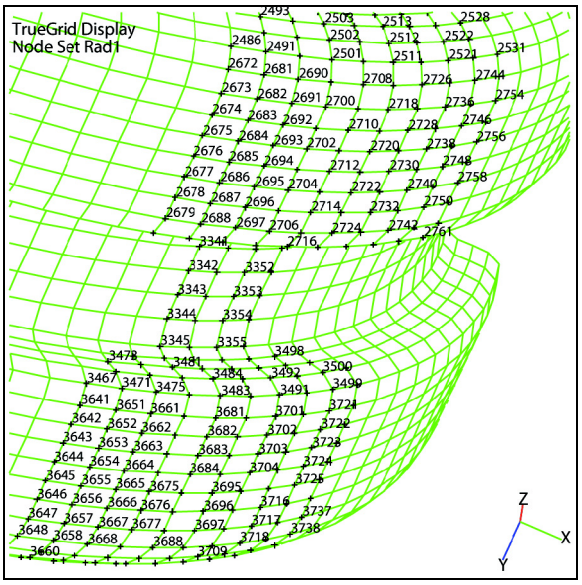

Figure 264 Node Set Rad1

nsetc assign a comment to a node set

nsetc *set_name comment*

Remarks

The nsetc command can be used to set the title for a node set for LSDYNA.

onset order a segment of a nodal set

onset *set_name first_node last_node*

where

set_name	name of the node set
first_node	first node sequence number of the group of nodes to be moved
last_node	last node sequence number of the group of nodes to be moved.

Remarks

This command is useful when building a 1D slide line, for example, where the nodes must be in a

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certain order. The nodes can be included in the node set as each part in generated using the **nset** command. Additional nodes can be included in a set using the **nset** command in the Merge Phase. Use the **labels** command under graphics to view the order of the nodes in the set. Use this command to reorder a segment of the nodes. Select the beginning and ending nodes of the sequence to be reordered. Use the sequence numbers which are in white. See also **mvnset** command.

pset create or modify a polygon set

pset *set_name operator selection*

where *set_name* is the name of the polygon set

where *operator* can be

=	for initial assignment
AND	for intersection with node set
OR	for union with the node set
+	to append the selected nodes to the node set
-	for removal from the node set

where the *selection* can be

l s_1 p_1 s_2 p_2 ... ;	for a list of polygons
where	
s_i	surface number
p_i	polygon number of that surface
s *polygon_set_name*	for another polygon set

Remarks

Associated with this command is an interactive feature to select or modify the polygons in a set with the ***Sets*** window from the Environment Window.

Polygon sets can then be turned into surfaces using (see the **sd** command):

sd *surface#* **poly** *polygon_set trans* ;

These features along with the **wrsd** can be used to sort out complex polygon surfaces and split them into multiple surfaces or remove features. This can also be used to create an normal offset by using the normal option as one of the transformation primitives.

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rml remove specific loads or conditions on a set of nodes

rml *nodes type*

where

nodes can be

N *node_number*	node number
RT *x y z*	node closest to Cartesian point
CY *rho theta z*	node closest to cylindrical point
SP *rho theta phi*	node closest to spherical point
NSET *name_of_set*	node set by name

type can be

ALL *load_curve_#*	any type
FC *load_curve_#*	forces
FD *load_curve_#*	fixed displacements
FV *load_curve_#*	fixed velocity
FT *load_curve_#*	forced temperature
TM	initial temperature
SW *stone_wall_#*	stone wall
V	electrostatic potential
ACC *load_curve_#*	acceleration
FA	fixed angular displacement
MP	magnetic potential
MOM *load_curve_#*	moment
LB *local_coordinate_system_#*	constraint in local coordinate system
VE	velocity
TE	temperature
SI *sliding_interface_#*	nodal sliding interface
NPB	nodal print blocks
TEPRO *load_curve_#*	temperature profile

and where

load_curve_#, *stone_wall_#*, *local_coordinate_system_#* can be **0** for all cases

Remarks

The **rml** command is very useful in conjunction with the **readmesh** command to alter boundary conditions. The **rml** can be used also to edit the forces. To change of boundary condition, first remove it and them issue a new boundary condition.

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Example

```
block 1 20;1 20; -1;0 20;0 20;0;        c mesh definition
merge
nset force = l 155 154 153 173 174 175 ;
                                  c node set force initialization
lcd 1 0 1 1 1;                    c load curve definition
fc nset force 1 1 1 0 0           c concentrated force definition
                                  c for the node set force
infol nset force all 1
     c information print for the node set force
     c for all conditions (stat is either on or off):
                        LOAD TABLE
NODE STAT  LOAD   CASE     X          Y          Z       AMPLITUDE
155  on    fc       1     1.0        0.0        0.0        1.0
154  on    fc       1     1.0        0.0        0.0        1.0
153  on    fc       1     1.0        0.0        0.0        1.0
173  on    fc       1     1.0        0.0        0.0        1.0
174  on    fc       1     1.0        0.0        0.0        1.0
175  on    fc       1     1.0        0.0        0.0        1.0
mom nset force 1 2 x     c moment around x definition
                                  c for the node set force
infol nset force all 1
     c information print for the node set force
     c for all conditions  (stat is either on or off):
                       LOAD TABLE
NODE STAT  LOAD   CASE    X          Y          Z       AMPLITUDE
 155  on    fc       1    1.0        0.0        0.0        1.0
 154  on    fc       1    1.0        0.0        0.0        1.0
 153  on    fc       1    1.0        0.0        0.0        1.0
 173  on    fc       1    1.0        0.0        0.0        1.0
 174  on    fc       1    1.0        0.0        0.0        1.0
 175  on    fc       1    1.0        0.0        0.0        1.0
 155  on    mom      1     1                               2.0
 154  on    mom      1     1                               2.0
 153  on    mom      1     1                               2.0
 173  on    mom      1     1                               2.0
 174  on    mom      1     1                               2.0
 175  on    mom      1     1                               2.0
rml nset force fc 1               c removal of the concentrated
               c forces of the node set force from the model
infol nset force all 1
     c information print for the node set force
```

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```
     c for all conditions  (stat is either on or off):
                          LOAD TABLE
NODE STAT  LOAD   CASE     X          Y          Z       AMPLITUDE
155  off   fc       1     1.0        0.0        0.0        1.0
154  off   fc       1     1.0        0.0        0.0        1.0
153  off   fc       1     1.0        0.0        0.0        1.0
173  off   fc       1     1.0        0.0        0.0        1.0
174  off   fc       1     1.0        0.0        0.0        1.0
175  off   fc       1     1.0        0.0        0.0        1.0
155  on    mom      1     1                                2.0
154  on    mom      1     1                                2.0
153  on    mom      1     1                                2.0
173  on    mom      1     1                                2.0
174  on    mom      1     1                                2.0
175  on    mom      1     1                                2.0
```

rsl restore specific loads or conditions on a set of nodes

rsl *nodes type*

where

nodes can be

N *node_number*	node number
RT *x y z*	node closest to Cartesian point
CY *rho theta z*	node closest to cylindrical point
SP *rho theta phi*	node closest to spherical point
NSET *name_of_set*	node set by name

type can be

ALL *load_curve_#*	any type
FC *load_curve_#*	forces
FD *load_curve_#*	fixed displacements
FV *load_curve_#*	fixed velocity
FT *load_curve_#*	forced temperature
TM	Initial temperature
SW *stone_wall_#*	stone wall
V	electrostatic potential
ACC *load_curve_#*	acceleration
FA	fixed angular displacement
MP	magnetic potential
MOM *load_curve_#*	moment
LB *local_coordinate_system_#*	constrain in local coordinate system

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VE	velocity
TE	temperature
SI *sliding_interface_#*	nodal sliding interface
NPB	nodal print blocks
TEPRO *load_curve_#*	temperature profile

and where

load_curve_#, *stone_wall_#*, *local_coordinate_system_#* can be **0** for all cases

Remarks

The **rsl** command is the complement to the **rml** command. It can be used only after **rml** is issued. You can use the **infol** command to make sure which loads are turned off.

Example

```
infol nset force all 1
     c information print for the node set force
     c for all conditions  (stat is either on or off):
                         LOAD TABLE
NODE STAT  LOAD   CASE     X          Y          Z       AMPLITUDE
155  off   fc       1     1.0        0.0        0.0        1.0
154  off   fc       1     1.0        0.0        0.0        1.0
153  off   fc       1     1.0        0.0        0.0        1.0
173  off   fc       1     1.0        0.0        0.0        1.0
174  off   fc       1     1.0        0.0        0.0        1.0
175  off   fc       1     1.0        0.0        0.0        1.0
155  on    mom      1     1                                2.0
154  on    mom      1     1                                2.0
153  on    mom      1     1                                2.0
173  on    mom      1     1                                2.0
174  on    mom      1     1                                2.0
175  on    mom      1     1                                2.0
rsl nset force fc 1  c restore load case 1for the node set force
infol nset force all 1
     c information print for the node set force
     c for all conditions  (stat is either on or off):
                         LOAD TABLE
NODE STAT  LOAD   CASE     X          Y          Z       AMPLITUDE
155  on    fc       1     1.0        0.0        0.0        1.0
154  on    fc       1     1.0        0.0        0.0        1.0
153  on    fc       1     1.0        0.0        0.0        1.0
173  on    fc       1     1.0        0.0        0.0        1.0
174  on    fc       1     1.0        0.0        0.0        1.0
```

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```
175   on     fc        1       1.0          0.0          0.0          1.0
155   on     mom       1       1                                      2.0
154   on     mom       1       1                                      2.0
153   on     mom       1       1                                      2.0
173   on     mom       1       1                                      2.0
174   on     mom       1       1                                      2.0
175   on     mom       1       1                                      2.0
```

rvnset remove a node subset

rvnset *set_name first_node last_node*

where

set_name	name of the node set
first_node	first node sequence number of the group of nodes to be moved
last_node	last node sequence number of the group of nodes to be moved.

Example

The node set SLID is defined, reordered and displayed (**Figure 265**). The nodes with sequence numbers `2,3,4,5` are removed from the node set SLID. The node set SLID is displayed in **Figure 266**.

```
labels onset SLID 1 0
c display labeled ordered set SLID
rvnset SLID 2 5
c remove node sequence position
c from 2 to 5 from the set SLID
labels onset SLID 1 0
c display labeled ordered set SLID
```

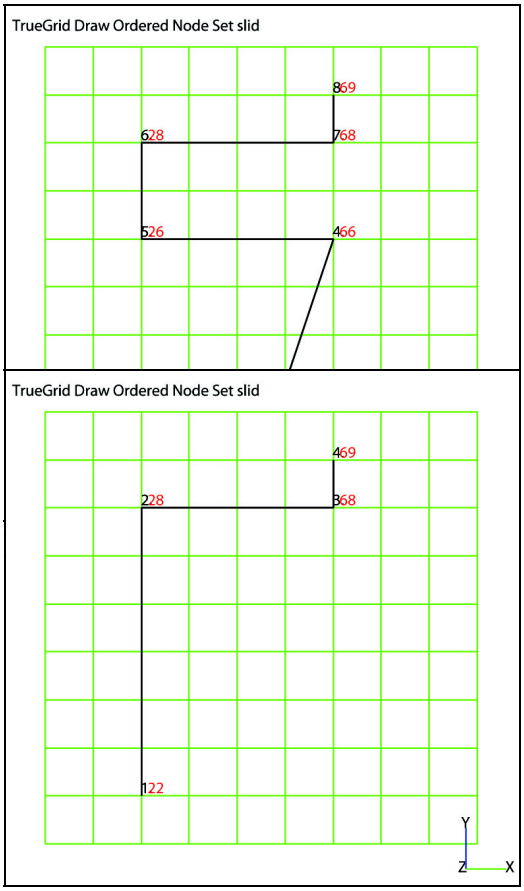

Figure 266 after **rvnset**

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III. Global Commands

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1. Materials

Every material is identified with a number. Use the automatic parameter %nextmat for the next available material number. Material properties are defined by various commands associated with an output option. Refer to the various output manuals for further information.

delmats delete a material definition

delmats *list* ;
where
list a list of material numbers of material models to be removed from the data base

Remarks

This command is useful when reading a model using the **readmesh** command. There may be material models for one FEM code which will not be used when the model is translated to a new format. Deleting the material models is the safest way to avoid complications.

2. Parts

There are several ways to create a part. The **block** command is recommended for most problems. Only rarely is a **cylinder** part justify because of its unusual behavior near the local z-axis. The **readmesh** creates several parts by reading a mesh file with a specific format. Each material within the mesh file becomes a part. The **blude** command is even more rarely used and was designed to build a fluid mesh around an existing shell mesh. The **beam** and **cbeam** commands are included to stay compatible with old versions of Ingrid, **True*Grid*®**'s predecessor. The **bm** command in the merge phase is the preferred method for generating a string of beams.

If the **partmode** command precedes the **block**, **cylinder**, or **blude** commands, it affects the meaning of those commands. If the **linear** or **quadratic** command precedes the **block**, **cylinder**, **blude**, or **bm** commands, it affects the meaning of those commands. These three commands cannot be issued within the scope of a part command. For that reason, they are no discussed in this section. They can be found under assembly commands.

There are commands to display selected parts of your model. When you enter the merge phase, all parts for graphics are automatically activated. When you enter into the part phase, only that part will be active in the graphics. When a sequence of parts is needed, you only need to give the first and last numbers separated by a colon.

block create a brick-shaped part

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block *i_indices ; j_indices ; k_indices ; x_coordinates y_coordinates z_coordinates*

where

i_indices, *j_indices*, and *k_indices* are lists of indices (not reduced indices). Each list of indices takes the form i1 i2 ... in, where i1=1 and |i1| < |i2| < ... < |in|. Normally each of these indices is a positive integer. But a negative integer can be used to create a shell surface, and a zero is used to separate regions of the mesh.

x_coordinates, *y_coordinates*, and *z_coordinates* are lists of values of physical coordinates, one for each index value: an x coordinate for each i index, a y coordinate for each j index, and a z coordinate for each k index. These define the physical locations of every region interface.

Example

The block part is defined by a list of i-indices and x-coordinates, a list of j-indices and y-coordinates, and a list of k-indices and z-coordinates. An example command file follows:

```
c Block part definition.
c  Indices in the x-direction (i-dir.) are:1 5 7 8 11.
c    x-coordinates are:-1.3 5.2 8 8.7 11.3 units
c  Indices in the y-direction (j-dir.) are: 1 3 5.
c    y-coordinates are:1.2 3.4 5.2 units.
c  Indices in the z direction (k-dir.) are:1 10.
c    Z-coordinates are:-5 5 units.

block 1 5 7 8 11;   1 3 5;   1 10;
      -1.3  5.2  8  8.7  11.3;
        1.2  3.4  5.2;
        -5  5;
```

This block is illustrated in the following 6 pictures. The faces that can be selected using the index bars correspond to the indices in the index lists. These faces partition the part into blocks and are sometimes referred to as i-, j-, and k-partitions for descriptive purposes.

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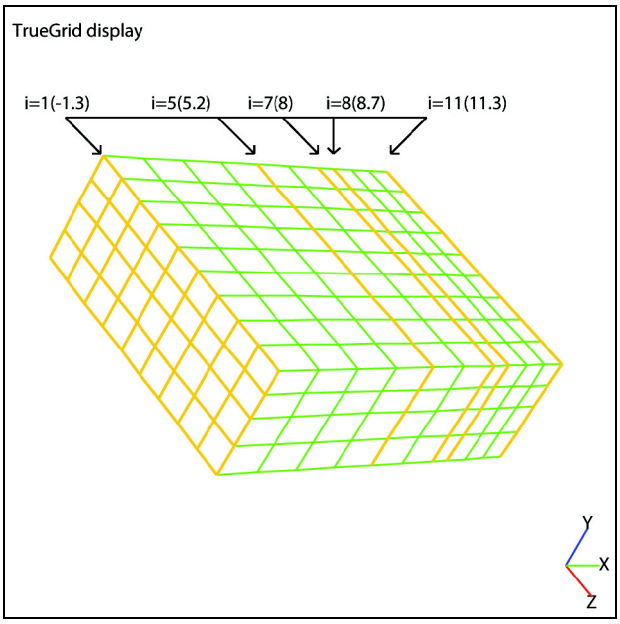

Figure 267 i-partitions and x-coordinates

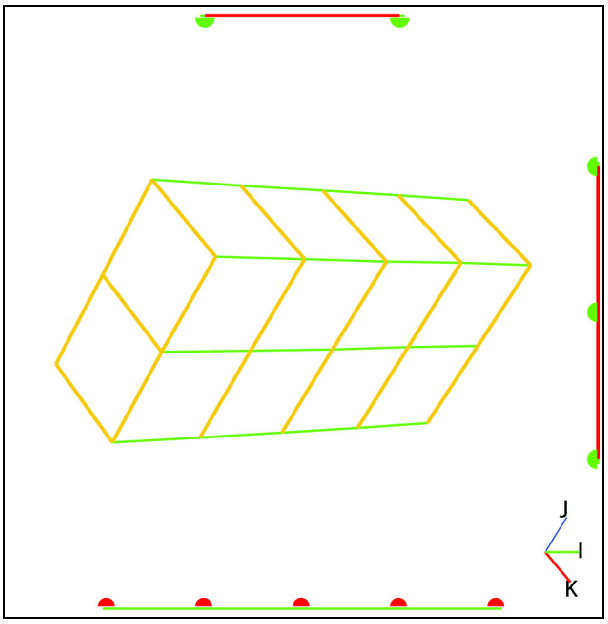

Figure 268 i-partitions

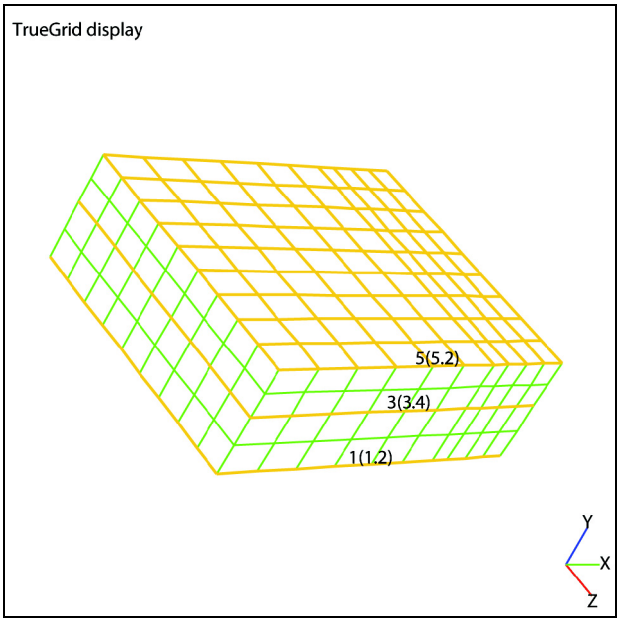

Figure 269 j-partitions and y-coordinates

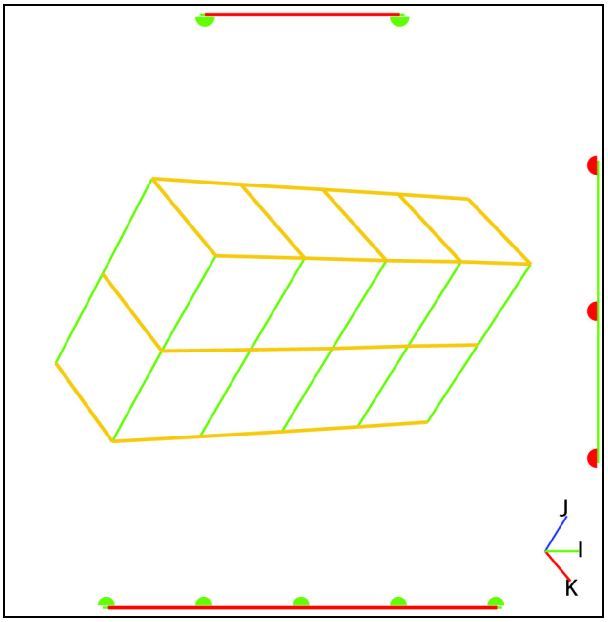

Figure 270 j-partitions

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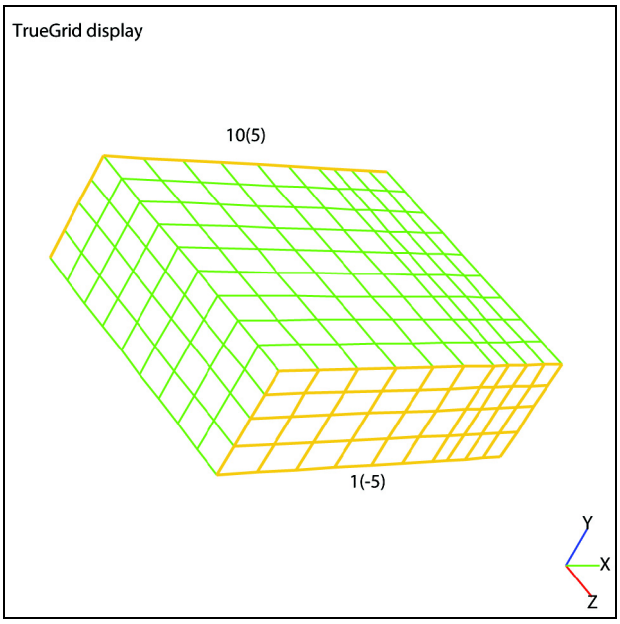

Figure 271 k-partitions and z-coordinates

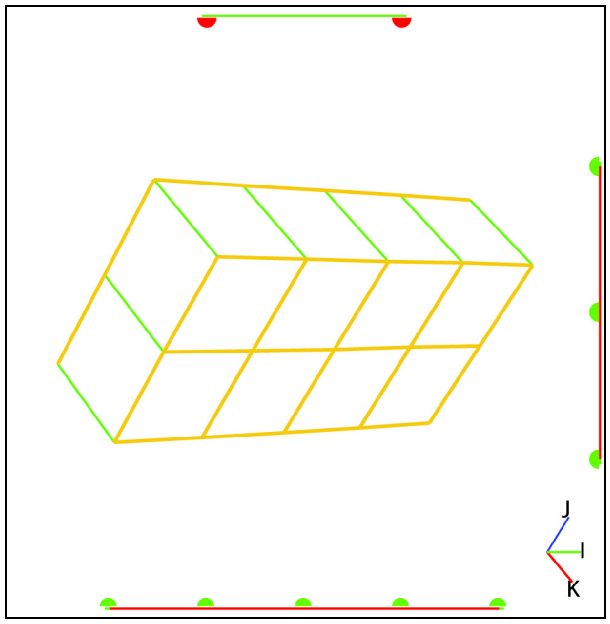

Figure 272 k-partitions

Remarks

This command is also discussed in the Introduction.

This command transitions the code to the Part Phase. This is the standard way to generate parts. In order to complete the Part Phase and return to the Control Phase, use the **endpart** or **control** command. In order to complete the Part Phase and return to the Merge Phase, use the **merge** command. In order to abort the part and return to the Control Phase, use the **abort** command.

When this command is issued, the previous part (if any) is ended as if the **endpart** command had been issued.

Six lists of numbers follow the **block** command. The first three lists consist of integers, each list terminated with a semi-colon. The second three lists consist of real numbers, each list optionally terminated by a semi-colon. The first list of integers must start with a 1 or -1. The integers that follow must be zero or have absolute value greater than the absolute values of the integers that preceded it in that list. These are the number of nodes to create in the first dimension of the computational mesh. A positive integer indicates that there will be a partition at that nodal index in the first dimension of the computational mesh, which can be referenced by most commands in the Part Phase. These partitions are used to break the part into multiple structured blocks. Since most part commands reference only the partition sequence numbers and not the nodal indices, it is a trivial matter to change the density of the mesh by only changing this list of integers. Moreover, one may use

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parameter values here in order to be able to later alter the mesh density. The partition sequence numbers corresponding to the integers in this list are symbolic references to the structure of the mesh. When positive integers are used, solid elements are created. If you are generating solid elements exclusively, you may prefer the alternative method of specifying the number of elements in each list by first issuing the **partmode** command. A negative integer in the list also produces a partition in the mesh with a nodal index corresponding to the absolute value of the integer, with shell elements created along that partition in the computational mesh. A zero in the list means there is a gap in the mesh, i.e. the part is not connected. The second and third lists have the same meaning applied to the second and third dimensions in the computational mesh, respectively. The fourth list is a list of x-coordinates, one coordinate for each integer in the first list. This initializes each partition in the first dimension of the computational mesh. The fifth list consists of y-coordinates corresponding in like manner to the second list of partitions in the second dimension of the computational mesh. The sixth list consists of z-coordinates, corresponding in like manner to the third list of partitions in the third dimension of the computational mesh.

cylinder create a cylindrical part

cylinder *i_indices ; j_indices ; k_indices ; r_coordinates θ_coordinates z_coordinates*

where

i_indices, *j_indices*, and *k_indices*

are lists of indices (not reduced indices). Each list of indices takes the form i1 i2 ... in, where i1=1 and |i1| < |i2| < ... < |in|. Normally each of these indices is a positive integer. But a negative integer can be used to create a shell surface, and a zero can be used to separate parts of the mesh.

r_coordinates, *θ_coordinates*, and *z_coordinates*

are lists of values of physical coordinates, one for each index value: an r coordinate for each i index, a θ coordinate for each j index, and a z coordinate for each k index.
These define the physical locations of every region interface.

Example

The cylinder part is defined by a list of i-indices and radii, a list of j-indices and angles, and a list of k-indices and z-coordinates. A sample command file follows:

```
c Cylinder part definition.
c  Indices in the radial direction (i-dir.) are:1 3 5 7 11.
c    Radii are:6 9 12 15 18 units
c  Indices in the angular direction (j-dir.) are:
c                                1 5 9 13 17 21 25 29 33 37.
c    Angles are:0 50 80 120 160 200 240 280 310 360 degrees.
```

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```
c  Indices in the z direction (k-dir.) are:1 3 5 7 9.
c    Z-coordinates are:-10 0 10 20 30 units.

cylinder 1 3 5 7 11; 1 5 9 13 17 21 25 29 33 37; 1 3 5 7 9;
          6 9 12 15 18;
           0 50 80 120 160 200 240 280 310 360;
            -10 0 10 20 30;
```

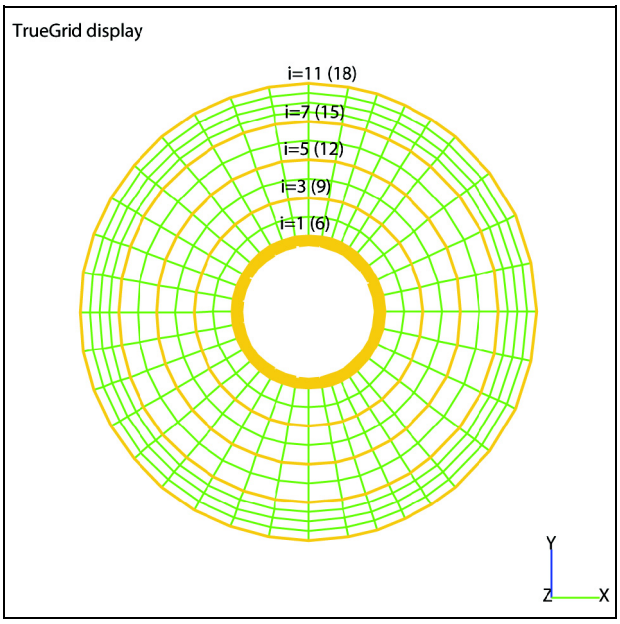

Figure 273 i-partitions and radiuses

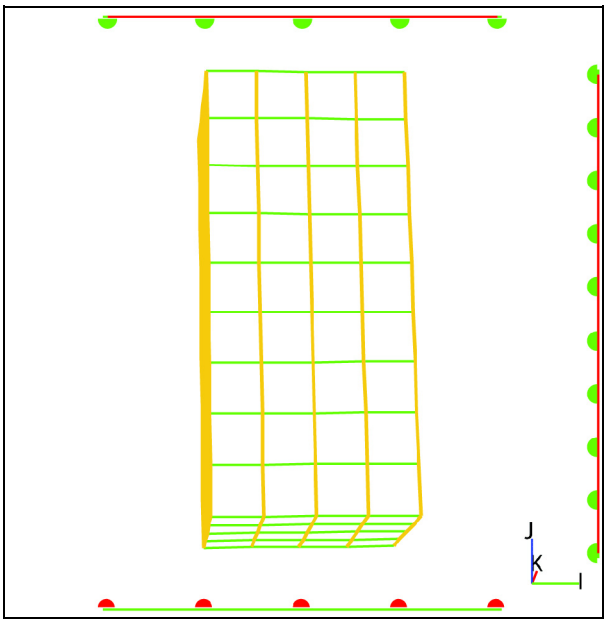

Figure 274 i-partitions

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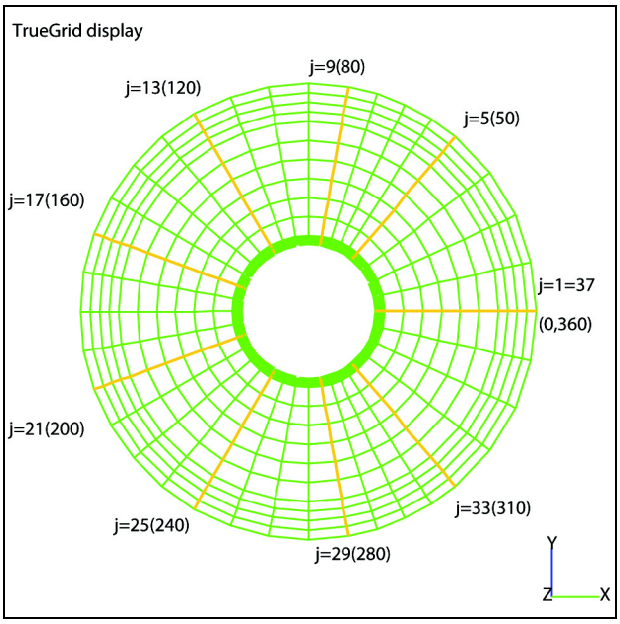

Figure 275 j-partitions and angles

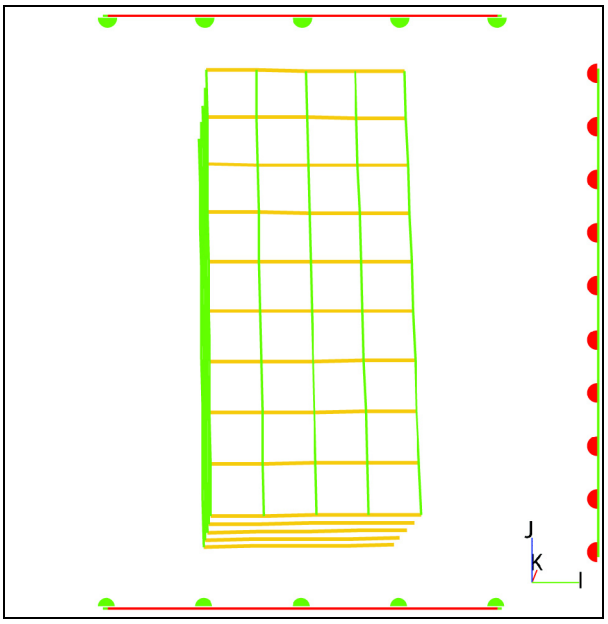

Figure 276 j-partitions

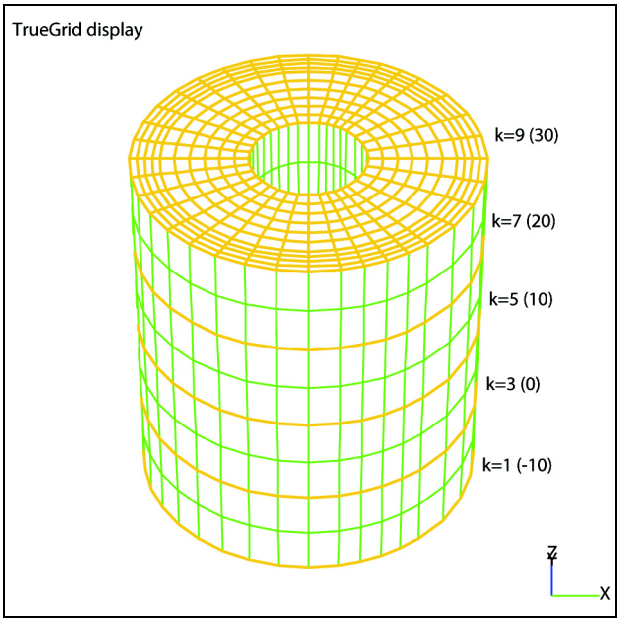

Figure 277 k-partitions and z-coordinates

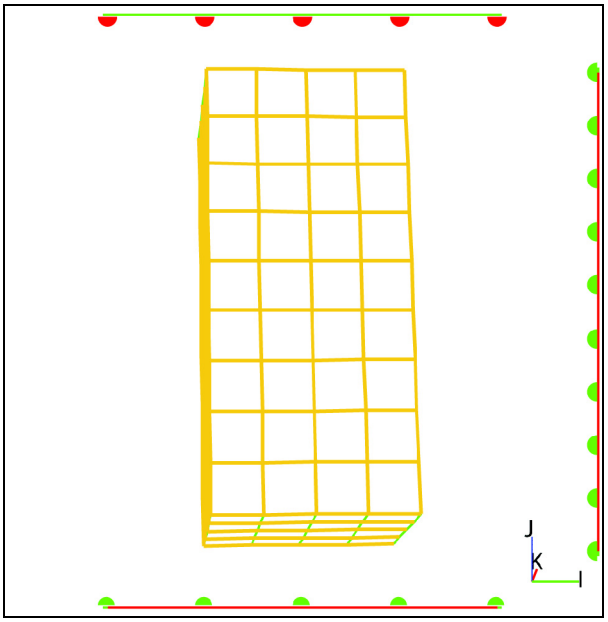

Figure 278 k-partitions

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Remarks

When this command is issued, the present part is ended as if the **endpart** command was issued. Then this new part is immediately initialized.

Like **block**, this command transitions the code to the Part Phase. This is the standard way to generate a part. In order to complete the Part Phase and return to the Control Phase, use the **endpart** or **control** command. In order to complete the Part Phase and return to the Merge Phase, use the **merge** command. In order to abort the part and return to the Control Phase, use the **abort** command.

Six lists of numbers follow the **cylinder** command. The first three lists consist of integers, each list terminated with a semi-colon. The second three lists consist of real numbers, each list optionally terminated by a semi-colon. The first list of integers must start with a `1` or `-1`. The integers that follow must be zero or have absolute value greater than the absolute values of the integers that preceded it in that list. These are the number of nodes to create in the first dimension of the computational mesh. A positive integer indicates that there will be a partition at that nodal index in the first dimension of the computational mesh, which can be referenced by most commands in the Part Phase. These partitions are used to break the part into multiple structured blocks. Since most part commands reference only the reduced indices and not the nodal indices, it is a trivial matter to change the density of the mesh by only changing lists of nodal indices in the part definition. When positive nodal indices are used, solid elements are created. If you are generating solid elements exclusively, you may prefer the alternative method of specifying the number of elements in each list by first issuing the **partmode** command. A negative nodal index in the list also produces a partition in the mesh with a nodal index corresponding to the absolute value of the integer, but with shell elements created along that partition in the computational mesh. A zero in the list means there is a gap in the mesh, i.e. the part is not connected. The second and third lists have the same meaning applied to the second and third dimensions in the computational mesh, respectively.

The fourth list is a list of radial coordinates, one coordinate for each integer in the first list. This initializes each partition in the first dimension of the computation mesh. The fifth list consists of angular coordinates, corresponding in like manner to the second list of partitions in the second dimension of the computational mesh. The sixth list consists of z-coordinates, corresponding in like manner to the third list of partitions in the third dimension of the computational mesh.

The cylindrical coordinates relate to Cartesian coordinates by

$$x = r\cos(\theta) \qquad y = r\sin(\theta) \qquad z = z\,.$$

When the part is completed, **True*Grid*®** will apply this transformation to put the part in the global Cartesian coordinate system. By default, this means that the axis of the part's cylindrical coordinate system will be the same as the z-axis of the global coordinate system. You can select a local frame

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of reference by choosing the z-axis of this coordinate system to be any vector in the global coordinate system by issuing the **cycorsy** command. Of course, after you finish making the part you can replicate or transform it into another part with whatever axis you like.

But for some uses of the **sf** and **sfi** commands, it is more convenient to define your part at its final location in the global coordinate system. If you define a part with projections and then transform it to its final location, the projection surfaces will have to be transformed from the final location to the place where the part is defined. If the **cylinder** command does not let you define your part at a convenient location, then you should use the **block** command. You can project a **block** part to make the same shape that a **cylinder** part would make.

Many computations that apply to the part will use cylindrical coordinates. An example of this is interpolations: an interpolation that will put points along a line for a **block** part will put points along an arc for a **cylinder** part if the angular coordinates at the end nodes are different.

Some part commands, such as **fd**, use physical coordinates in their arguments. For simplicity this manual names such command arguments in terms of Cartesian coordinates. For example, *x y z* would be arguments defining the physical coordinates of a point. However, the coordinate system of such arguments is always the same as the coordinate system of the relevant part. Thus if the part is set up by a **cylinder** command, you would give the physical coordinates of a point as its r-, θ-, and z-coordinates.

readmesh read a file containing a mesh

readmesh *format filename cmds* **endpart**

where

format can be:

nastran	is partially supported at this time
neutral	is partially supported at this time
dyna3d	is nearly fully supported
lsdyna	reads only nodes and elements
dynain	this is a file written by LS-DYNA
iges	reads a FEM model from an IGES file

filename is the path and file name of the formatted data

cmds can be

cvtab	to write the conversion tables from NASTRAN to **True*Grid*®**
exclude	exclude the NASTRAN model when writing the LS-DYNA
mate *mat*	used only for IGES

where *mat* is the material number for all elements in the file.

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mt *type mat* used only for IGES. *Type mat* can be repeated any number of times.
where *type* can be any or all of

beams	
lshells	for linear shells
qshells	for quadratic shells
lbricks	for linear bricks
qbricks	for quadratic bricks
springs	
dampers	

where *mat* is the material number for all elements of type in the file. Use 0 to ignore that type of element from the IGES file.

ptmass *mass* used only for IGES
where *mass* is the mass at all point masses. Use 0 to ignore the point masses in the IGES file.

rigid *switch* used only for IGES
where *switch* can be
on to include rigid beams (default)
off to ignore rigid beams

cond *switch* used only for IGES
where *switch* can be
on to include nodal conditions (default)
off to ignore nodal conditions

ndcons *load_case_list* ; used only for IGES
where *load_case_list* is a list of the load cases to be combined

maplabel *list_maps* ; used only for IGES to ANSYS translations
where a single map is formed by a label, file name, and material Id

Remarks

This is a part, and it must be ended with the **endpart** command. This command will read in a formatted NASTRAN, PATRAN, DYNA3D, LS-DYNA, and IGES file. The data in any of these files will be added to the **True*Grid*®** internal data base. It can be written out in another format. Only some of the data is understood at this time. The list of features for each format is found below.

Material models, cross sectional properties for beams and shells, spring properties, and other specialized elements do not convert directly from one format to another. For this reason, you should redefine these features within **True*Grid*®**.

The loads, boundary conditions, etc. can be viewed using the **condition** command in the graphics menu while in the Merge Phase. They can be modified as well - see the command in the Merge Phase for that specific condition. Use the node, face, and element set features to select objects from these parts

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and assign boundary conditions and properties while in the merge phase.

When a NASTRAN file is imported, **True*Grid*®** will report the number of features it translates into the **True*Grid*®** data base. That report may look like:

```
 NASTRAN objects included in the TRUEGRID data base:
      4153 node points
        32 beam elements
       333 linear shell elements
        21 quadratic shell elements
        11 linear brick elements - including tets and wedges
       357 quadratic brick elements - including tets and wedges
         2 beam element cross section definitions
         6 material models
         2 local coordinate systems defined
         4 load sets
                  45     324       3     655
         spring material properties
        22 springs
 There were       12 warning conditions from the NASTRAN deck
```

Load sets are similar to load curves in many of the dynamics codes and it will be necessary to define the appropriate load curves when porting to these codes. In the above example, if the output format is DYNA3D, then the load curves 45, 324, 3, and 655 should be defined.

The NASTRAN features which are supported in the READMESH command are:

CBAR	saved as a 2D element - cross sectional properties can be changed
CBEAM	saved as a 2D element - cross sectional properties can be changed
CBEND	saved as a 2D element - cross sectional properties can be changed
CELAS1	scalar and ground points are not supported, a PELAS must be referenced
CELAS2	scalar and ground points are not supported, this also produces a spring material definition
CHEXA	
CHEXA27	
CORD	
CPENTA	
CPTRAN	
CPYRAM	
CQUAD4	
CQUAD8	
CQUADR	

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CROD	saved as a 2D element - cross sectional properties can be changed
CSHEAR	
CTETRA	
CTRIA3	
CTRIA4	
CTRIA6	
CTRIA8	
CTRIAR	
CTUBE	saved as a 2D element - cross sectional properties can be changed
FORCE	
FORCE1	
FORCE2	
GRDSET	
GRID	
MAT1	must be redefined if using another output format
MAT2	must be redefined if using another output format
MAT4	must be redefined if using another output format
MAT5	must be redefined if using another output format
MAT8	must be redefined if using another output format
MAT9	must be redefined if using another output format
MAT10	must be redefined if using another output format
MATT9	must be redefined if using another output format
MOMENT	
MOMENT1	
MOMENT2	
PBAR	must be redefined if using another output format
PBCOMP	must be redefined if using another output format
PBEAM	must be redefined if using another output format
PBEND	must be redefined if using another output format
PCOMP	must be redefined if using another output format
PELAS	must be redefined if using another output format
PLOAD	
PROD	must be redefined if using another output format
PSHEAR	must be redefined if using another output format
PSHELL	must be redefined if using another output format
PSOLID	must be redefined if using another output format
PTUBE	must be redefined if using another output format
RBE2	

When a LSDYNA file is imported, **True*Grid*®** will report the number of features it translates into the **True*Grid*®** data base. That report may look like:

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```
Nodes from         1 to      8131
Materials from      1 to      2
Bricks and thick shells from         1 to      1000
Shells from         1 to      6720
```

The LSDYNA features which are supported in the READMESH command are:
nodes, elements, materials, and sets.

The node set numbers are offset depending on what numbered node sets already existing when the **readmesh** command is issued. The segment sets numbers are offset depending on what numbered segment sets already exist when the **readmesh** command is issued. The element sets are also offset and with a prefix to the number: BM for beam, SH for shell, BR for brick, and TS for thick shell.

An experimental feature with the **readmesh** command for NASTRAN files can be used in a limited way to import an existing NASTRAN model and add parts which can be merged to the nodes of the NASTRAN model (preserving the original node numbers) such that when the model is written out to an LS-DYNA keyword format, only the new nodes and elements are written out. This is the **exclude** option when reading a NASTRAN file. The **exclude** option means exclude objects from the **readmesh** when writing the output file. The is useful when building an add on to a model that already exists as a NASTRAN model that has been converted to LS-DYNA. If the **readmesh** command were able to read an LS-DYNA input and write it out to produce the exact model that was raed in, then this exclude feature would not be useful.

No new nodes or elements will be given numbers that are already in use in the original NASTRAN model. This feature is to meet a specific need in a timely fashion. However, this feature may have a more general use if it is extended to all keyword input and output. This would make it possible to modify a model incrementally by creating add-on files which can be concatenated. It is not clear that such an improvement will be made. Comments on this feature will be appreciated.

There are two limitations to keep in mind. The **readmesh** command should be issued before any parts with nodes that should be merged to nodes from the NASTRAN input file. This will probably remain a limitation or feature. Secondly, only the new nodes and elements are protected from colliding with the objects in the NASTRAN input file. Materials, springs and dampers, sets, etc. must still be managed by the user.

The IGES option reads in FEM elements, properties, and loads from an IGES file. Currently, entity 134 (NODE), most element types for entity 136 (element type 1, 2, 3, 5, 6, 12, 13, 14, 15, 17, 18, 27, 29, 31, 32, 33, and 35), 406 (tabular data, form 11, property type 12 nodal loads/constraint, and 418 (nodal load/constraint) are supported. The material number may be set for all elements in the IGES file or individually for beams, linear shells, quadratic shells, linear bricks, quadratic bricks, springs and dampers. (Setting the material for an element type to zero causes TrueGrid to skip all elements

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of that type in the file.) Also, the mass for the point mass (entity 136 element type 31) may be set (a point mass of zero means skip). Rigid beams (entity 136 element type 32) and nodal loads/constraints (entity 418) may be skipped using the **rigid off** and **cond off** options, respectively. Nodal load cases may be selected using the **ndcons** option followed by the list of nodal load cases to be combined. Note that **cond off** will cause all nodal conditions to be ignored even if **ndcons** is set.

The **maplabel** option is used in conjunction with an ANSYS output to reference beam cross sections data from a file that the ANSYS code references. If the entity 406, form 11, ptype 5008 is found in the IGES file, then the referenced beam elements will be given a beam type of 188 and the appropriate file for the cross section data will be referenced. The **maplabel** option can be repeated any number of times.

Related to the **dynain** option in the **readmesh** command is the **dynain** output format. One will usually read a **dynain** file using the **readmesh** command, modify the data using the **smags** and **delem** commands, and write the results back to a file using the **dynain** and **write** commands. Node and element numbers are preserved.

blude extrude a set of polygons

blude *direction face_set_name*
i_indices ; j_indices ; k_indices ; x_coordinates y_coordinates z_coordinates
where

direction is the face of the block, where the extrusion begins and can be one of:

- **1** for the minimum i-face
- **2** for the maximum i-face
- **3** for the minimum j-face
- **4** for the maximum j-face
- **5** for the minimum k-face
- **6** for the maximum k-face

face_set_name is the name of the face set to be extruded

i_indices, *j_indices*, and *k_indices* are lists of indices (not reduced indices). Each list of indices takes the form i1 i2 ... in, where i1=1 and |i1|< |i2| < ... < |in|. Normally each of these indices is a positive integer. But a negative integer can be used to create a shell surface, and a zero to separate parts of the mesh.

x,y,z_coordinates are lists of values of physical coordinates, one for each index value: an x coordinate for each i index, a y coordinate for each j index, and a z coordinate for each k index. These define the physical locations of every

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region interface.

Remarks

This command extrudes a set of polygons. It is based upon the **block** command and an understanding of that command is essential to understanding this one. This extrusion is done by the construction of a block part with no deleted regions. The polygons in the face set are extruded, one polygon at a time, by following the mesh lines formed by the block part. See also the **block** command. The set of polygons is defined using a face set. See also the **fset** command. The block part is discarded. Its only purpose is to define the flow lines for the extrusion of the polygons. Be sure, when you are building the block part, that one of its 6 outer faces completely covers the face set of polygons.

This command transitions the code to the Part Phase. In order to complete the Part Phase and return to the Control Phase, use the **endpart** or **control** command. In order to complete the Part Phase and return to the Merge Phase, use the **merge** command. In order to abort the part and return to the Control Phase, use the **abort** command.

This part is difficult to use. Its application is to a small class of problems and many conditions must be near perfect for it to work properly. There is no guarantee that it will work for a problem. This part is not recommended for the casual user.

Example

In this example, a simple NASTRAN ship model is read using the **readmesh** command. This mesh is unstructured because there is a small region of the hull with triangles. A face set is constructed using the mouse to select the faces on the starboard side. This face set is then used to construct a

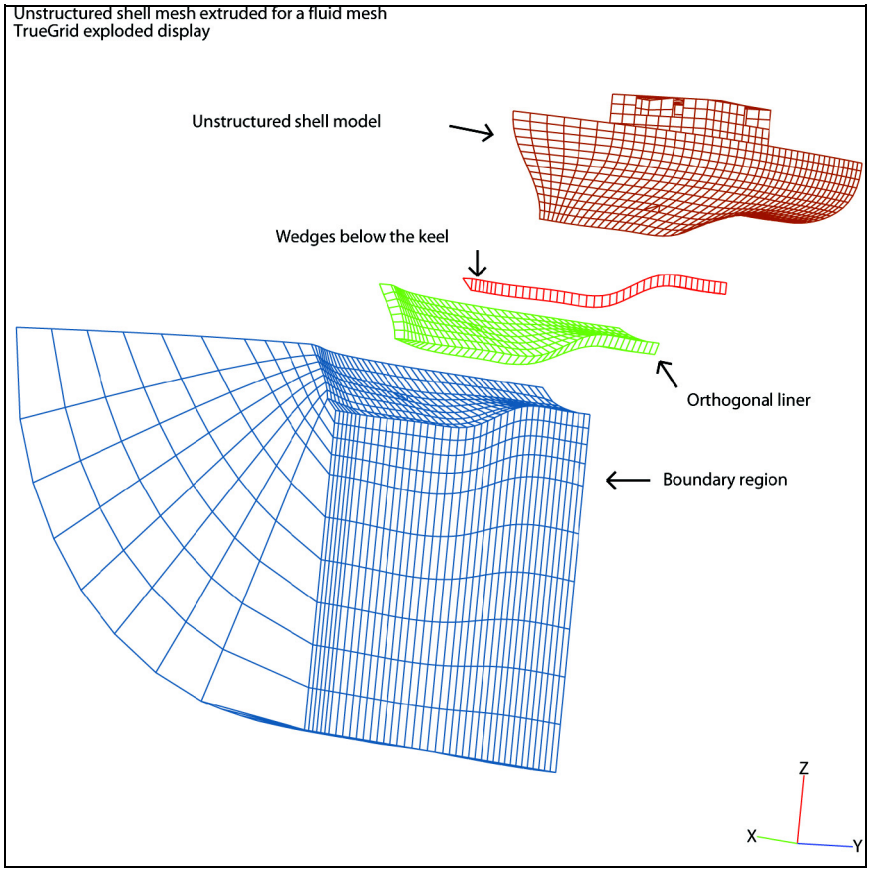

Figure 279 Shell structure extruded forming a solid

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face surface for projection and for the **blude** command. An intra-part **bb** command with a normal offset is used in the construction of the liner to form an orthogonal part. This liner is needed because a wedge part just below the keel is needed to round off the keel. Only then can a good fluid mesh be extruded to the outer cylinder wall of the region of interest. It would be possible to use the intra-part **bb** command with an offset to form the boundary layer, however in this example, for graphical purposes, the boundary region was made too large for the normal offset feature in the **bb** command to work effectively. The normal offset feature is only good for short distances relative to the changes in curvature on the master surface. This entire model is too complicated to have the input files included in this text. The input files are available.

meshscal scale up the mesh density for all parts

meshscal *scale*

where *scale* is a positive integer

Remarks

This will increase the mesh density uniformly in all 3 directions by an integer scale factor. Typically a factor of 2 (8 times as many bricks, 4 times as many shells, 2 times as many embedded beams are the result) or 3 is sufficient. This should come first before any parts are made. This cannot be used if the **update** command is used. The default is 1.

beam initialize a beam part in Cartesian coordinates

beam *options* **endpart**

where

options are any of the following (this may be continued across many lines)

abort

bm $node_1$ $node_2$ *number_of_elements material cross-section* $node_3$ *;*

to create a sequence of beam elements, where:

$node_1$ and $node_2$	form the end nodes of the sequence of beams,
number_of_elements	is the number of beam elements,
material	is the material number,
cross-section	is a cross-section number, and
$node_3$	is the local coordinate system orientation node.

cy *radius angle z-coordinate constraints ;* for cylindrical coordinates

where:

dx	for x-displacement,

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dy for y-displacement,
dz for z-displacement,
rx for rotation about the x-axis,
ry for rotation about the y-axis,
rz for rotation about the z-axis,
followed by a value of
0 for initialization to no constraint
1 for constraint

jt *joint_# joint_node_# beam_node_# joint_increment ;* for joint nodes
joint_# for the **True*Grid*®** joint definition number,
joint_node_# each joint type consists of a certain number of nodes, which are numbered in a specified way for **True*Grid*®** (consult the user's manual). You must specify which node of the joint this beam node must become,
beam_node_# this is the **True*Grid*®** beam node number for the current beam part,
joint_increment this is used when there are replications and indicates how the joint_no should be adjusted for replications

lct *#_transforms ; first_transforms ; ... ; last_transforms ;* transformations

lrep *list_local_transformation_# ;* for replication of the beams

rt *x-coordinate y-coordinate z-coordinate constraints ;* for Cartesian coord's
where the constraint list is composed of
dx for x-displacement
dy for y-displacement
dz for z-displacement
rx for rotation about the x-axis
ry for rotation about the y-axis
rz for rotation about the z-axis
followed by a value of
0 for initialize to no constraint
1 for constrain

si *interface_# option_list ;* for sliding interface nodes
where an option can be
ffn normal_force
ffs shear_force
enf e_normal_force
esf e_shear_force

sp *radius angle_1 angle_2 constraints ;* for spherical coordinates
where the constraint list is composed of:
dx for x-displacement

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dy for y-displacement
dz for z-displacement
rx for rotation about the x-axis
ry for rotation about the y-axis
rz for rotation about the z-axis
followed by a value of
0 for initialize to no constraint
1 for constrain

Remarks

This command may be issued in any phase. If it is issued in the block or cylinder Part Phase, this command will end that part and create this new part. The implementation of the beam part is not interactive, although it is fully functional. It is provided to be nearly compatible with the INGRID beam part. Please note the differences in the two parts.

The intent here is to first create a set of nodes, one node at a time using either **rt**, **cy**, or **sp**. Each node is assigned a set of nodal constraints as part of the definition of the node. No constraints are needed. Usually beam nodes are merged to other nodes in the model, so that they will then inherit the constraints of the other nodes that they are merged to. These nodes are then tied together to form beam elements using the **bm** command. The nodes are referenced by their sequence number within the beam command. Some simulation codes, such as DYNA3D, require a third node in a beam element definition to define the orientation of the beam. Use the **bsd** command to define the cross sectional properties of the beam. Use the appropriate material definition command to assign properties to the numbered material referenced in the **bm** command.

Coordinates are referenced in their respective systems but all interpolation is done in Cartesian coordinates.

See the **jd** and **jt** command within the block part to understand the beam **jt** command. Joints require at least two nodes. The **jd** command is used to define joint properties. The nodes of a joint are numbered in a set way. You can flag nodes created in a beam part for use in a joint using the **jt** command

cbeam initialize a beam part in cylindrical coordinates

cbeam *options* **endpart**
where
options are the same as in **beam** except that the interpolation of node points in the **bm**

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option is done in a cylindrical coordinate system.

Remarks

See the **beam** command.

demicle generate discrete element spheres

demicle (*no arguments*)

Remarks

Discrete element spheres can be generated using the **demicle** command prior to creating a part. This will cause all brick elements of the next part to be converted to discrete element spheres. This is for LSDYNA output.

Each sphere will have a center coincident with the solid element. The radius will be the largest radius of a sphere that fits within the solid element.

The **demp**/**dempi** commands in the part phase will assign properties to the spheres

Use the **condition** (**co**) **dem** command to view the discrete element numbers. The default length of the lines that cross to represent a discrete Element will be the radius of that element. The size can be scaled using the **co size** option

ap add a part to the picture

ap *part_number*

Remarks

The display list determines what objects are in the picture. A part in the display list is referred to as being active. Part number 0 refers to the present part. If the selection is the present part, then it can be viewed in full. If it is any other part, it can only be viewed as a wire frame.

aps add a list of parts to the picture

aps *part_list ;*

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Remarks

The display list determines what objects are in the picture. A part in the display list is referred to as being active. Part number 0 refers to the present part. If one of the active parts is the present part, then it can be viewed in full. All other parts are only viewed as a wire frame. A list of part numbers can include a sequence of part numbers with only the first and last numbers separated by a colon (e.g. aps 2:5;).

dap display all parts

dap *(no arguments)*

Remarks

The display list determines what objects are in the picture. A part in the display list is referred to as being active. The present part will entirely be shown. The previously generated parts will be wire frame only.

dp display one part in the picture

dp *part_number*

This command will display one, and only one, part.

dps display a set of parts in the picture

dps *part_list* ;
where
part_list is a blank-delimited list of part numbers

Remarks

The display list determines what objects are in the picture. Part number 0 refers to the present part. A part in the display list is referred to as being active. If a selection is the present part, then it can be viewed in full. If it is any other part, it can only be viewed as a wire frame. A list of part numbers can include a sequence of part numbers with only the first and last numbers separated by a colon.

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pinfo part information

pinfo *(no arguments)*

rap remove all parts from the picture

rap *(no arguments)*

Remarks

The display list determines what objects are in the picture. A part in the display list is referred to as being active.

rp remove one part from the picture

rp *part_number*
where
part_number is a part number.

Remarks

The display list determines what objects are in the picture. A part in the display list is referred to as being active. Part number 0 means the present part.

rps remove a set of parts from the picture

rps *part_list ;*

Remarks

The display list determines what objects are in the picture. A part in the display list is referred to as being active. Part number 0 refers to the present part. A list of part numbers can include a sequence of part numbers with only the first and last numbers separated by a colon.

3. Motion

These commands apply initial velocities to all parts. The **rotation** or **velocity** commands in the Part

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Phase can override these commands for just one part. The **ve** or **vei** commands in the Part Phase can override these commands for selected regions of the part.

rotation global initial velocities as a rigid body rotation

rotation x_0 y_0 z_0 *x_rotation y_rotation z_rotation*

where

(x_0,y_0,z_0)	is any point on the axis of rotation, and
(x_rotation,y_rotation,z_rotation)	is the rotation vector in radians per unit time.

Remarks

By default, an initial body rotation is assigned to all parts defined after this command. Use the units expected by the simulation code, as **True*Grid*®** does no conversions. (**True*Grid*®** is dimension less.)

Examples

The following example assigns initial velocities which rotate the shell cylinder mesh about its axis of symmetry. The velocity magnitude is 0.2 because the radius of the cylinder is 2.

```
sd 1 cy 0 2 2 1 0 0 2
rotation 0 2 2 .1 0 0
block 1 11;-1 -11;-1 -11;-1 1 1 3 1 3
sfi ; -1 -2; -1 -2;sd 1
endpart
```

velocity global initial velocities as a rigid body translation

velocity *x_velocity y_velocity z_velocity*

where

(x_velocity,y_velocity,z_velocity)	is a velocity vector.

Remarks

An initial rigid body velocity is assigned to all parts defined after this command. Use the units expected by the target simulation code as **True*Grid*®** does no conversions.

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4. Boundary Conditions and Constraints

detp detonation points or lines

detp *material options ;*

where

material is the material number and

options is a list of any of the following:

time *detonation_time*

point *x y z*

lnpt x_1 y_1 z_1 x_2 y_2 z_2 *#_detonators*

Remarks

This creates detonation points and lighting times for high explosives. Use the **point** option if there is only one detonation point. Otherwise, use the **lnpt** option for a line of detonation points equally spaced.

jd joint definition

jd *joint_# type options ;*

where *joint_#* must be between 1 and 300

where *type* can be (default is shared constraints)

sj	for spherical joint
rj	for revolute joint
cj	for cylindrical joint
pj	for planar joint
uj	for universal joint
tj	for translational joint
sw	to constrain nodes as a spotweld

where *option* can be

pnlt *joint_penalty*	for joint penalty
repe *#_joint_replications*	for joint replications
dx	to specify shared x-displacement between the nodes
dy	to specify shared y-displacement between the nodes
dz	to specify shared z-displacement between the nodes
rx	to specify shared x-axis rotation between the nodes
ry	to specify shared y-axis rotation between the nodes
rz	to specify shared z-axis rotation between the nodes

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Remarks

Use this command first to define the properties of a numbered joint. Then use the **jt** command to identify the nodes used in the joint.

There are two types of joints. **Sj** , **rj**, **cj**, **pj**, **uj**, and **tj** form the first type and "nodes constrained together" form the second type. See the diagrams below for the relationship between the numbered nodes of the first type.

The **pnlt** option is used for the **DYNA3D**, **LSDYNA** and **NIKE3D** output only in the first type of joint definition. The **repe** option makes it possible to create a sequence of duplicate joint definitions starting with the one specified by the **jd** command.

Up to 16 nodes may be assigned to a joint. If the joint is of a type that requires fewer nodes, **True*Grid*®** will ignore any superfluous nodes assigned to that joint. If there are more than 16 nodes to share constraints, use the **mpc** command.

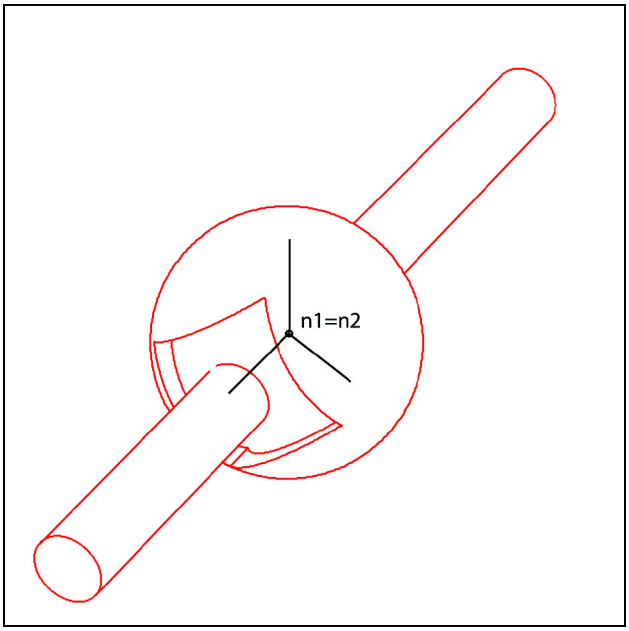

Figure 280 Spherical Joint

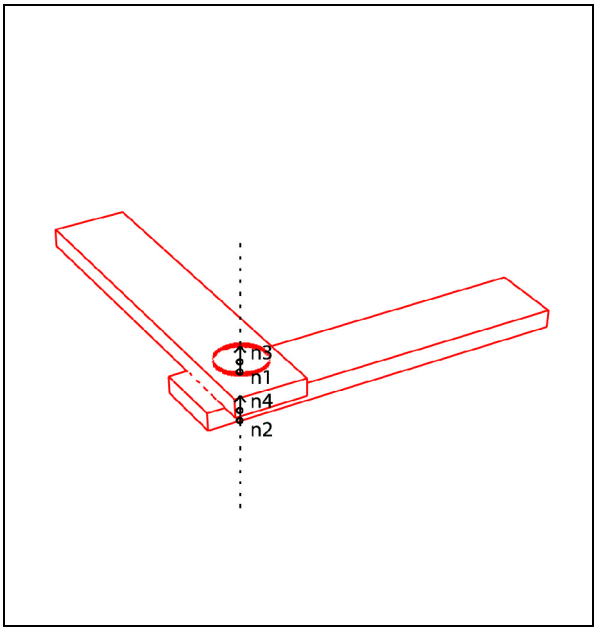

Figure 281 Revolute Joint

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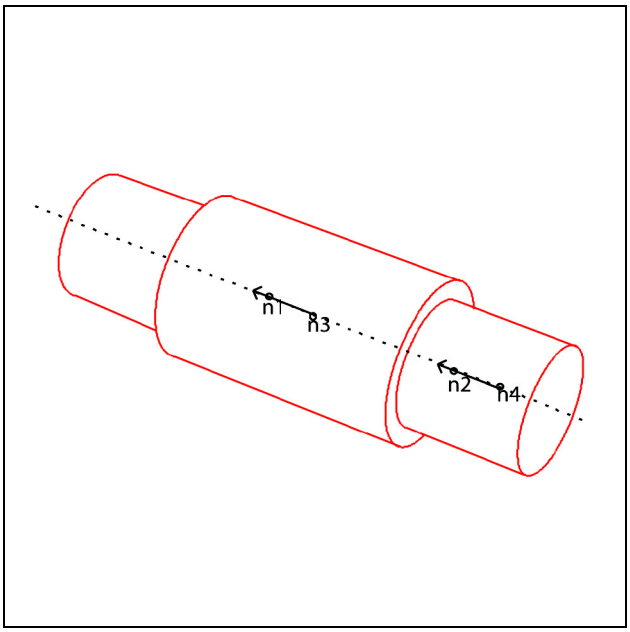

Figure 282 Cylindrical Joint

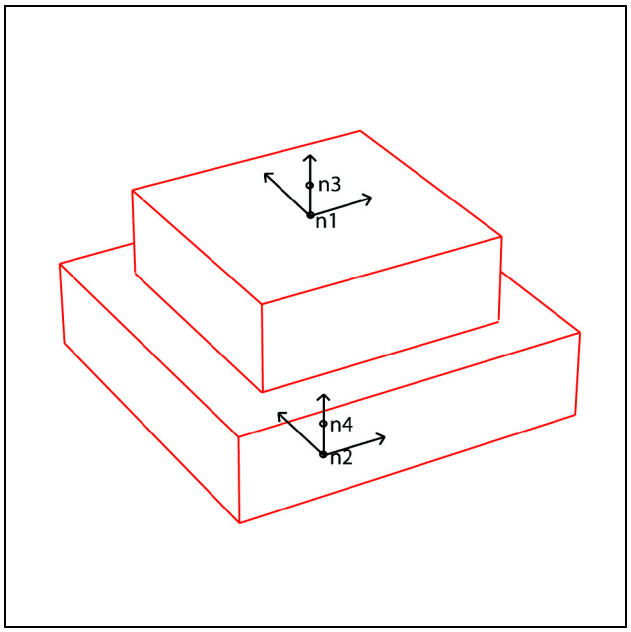

Figure 283 Planar Joint

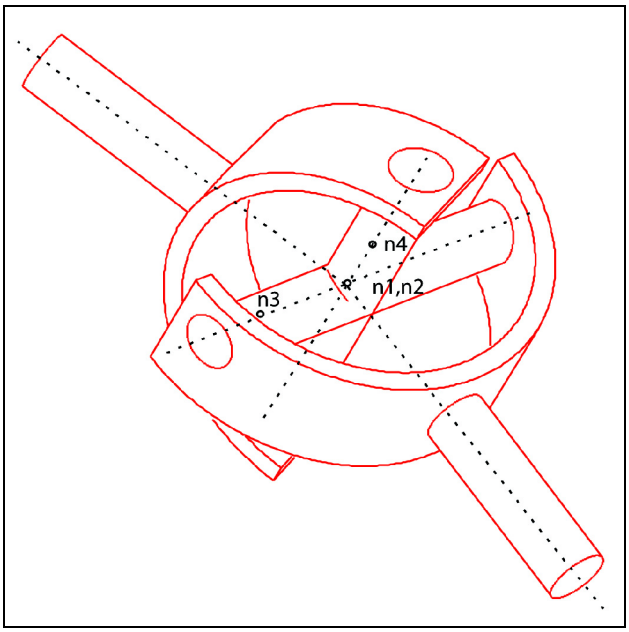

Figure 284 Universal Joint

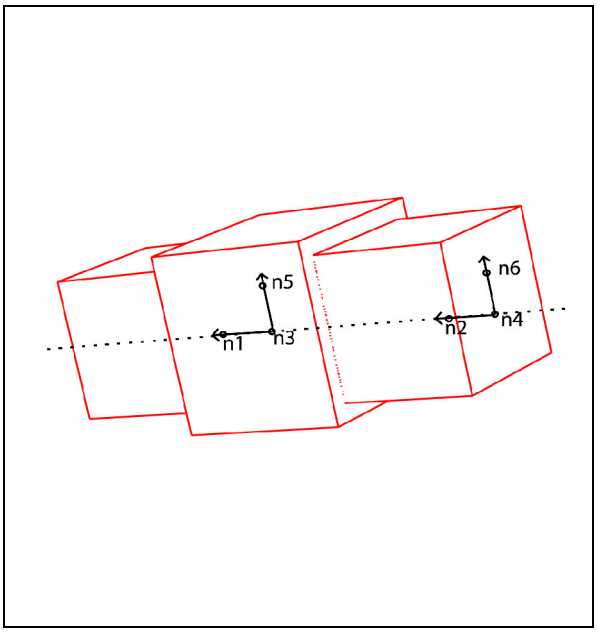

Figure 285 Translational Joint

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jtinfo write information about joints

jtinfo *<no arguments>*

Remarks

This command lists all of the joint definitions in the model.

lsys define a local coordinate system for the lb command

lsys *system transformations ;*

where

system is the "system number" which **lb** will use to identify this local coordinate system, and

transformations is a list of transformations that define the local coordinate system, chosen from the following:

mx *x_offset* for translation in x

my *y_offset* for translation in y

mz *z_offset* for translation in z

v *x_offset y_offset z_offset* for translation in an arbitrary direction

rx θ for rotation about the x axis

ry θ for rotation about the y axis

rz θ for rotation about the z axis

tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* for Cartesian coordinates

cy *rho theta z* for cylindrical coordinates

sp *rho theta phi* for spherical coordinates

pt *c.i* for a label of a labeled point from a 3D curve

pt *s.i.j* for a label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* for Cartesian coordinates

cy *rho theta z* for cylindrical coordinates

sp *rho theta phi* for spherical coordinates

pt *c.*i for a label of a labeled point from a 3D curve

pt *s.i.j* for a label of a labeled point from a surface

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inv to invert the present transformation

Remarks

This command is used before using the **lb** command. This command defines the local coordinate system to constrain a set of nodes in a coordinate system different from the global coordinate system.

lsysinfo list all of the local coordinate systems

lsysinfo <*no arguments*>

Remarks

This command lists all of the local coordinate systems defined using the **lsys** command.

mpc shared nodal (multiple point) constraints for a nodal set

mpc *node_set_name constraints ;*
where
constraints can be any combination of:
dx for constrained displacement in the x-direction,
dy for constrained displacement in the y-direction,
dz for constrained displacement in the z-direction,
rx for constrained rotations about the x-axis,
ry for constrained rotations about the y-axis, and
rz for constrained rotations about the z-axis.

Remarks

The **mpc** command assigns constraints to be shared by a set of nodes. This set of nodes is defined by the **nset** command. The nodes in the set share a specified degree of freedom. The first node is the master node for those codes requiring a master node.

Example

```
block 1 3 5 7;1 3 5 7;1 3 5 7;1 3 5 7;1 3 5 7;1 3 5 7;
 c definition of the mesh
nset 2 1 1 3 1 4 = narxs
 c definition of node set narxs
```

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```
 c from nodes of region
 c 2 1 1 3 1 4
mpc narxs dy ; ;
 c multiple point constraint
 c in dy direction
 c is assigned to nodes
 c of the node set narxs
merge
```

mpcinfo multiple point constraint sets information

mpcinfo *<no arguments>*

Remarks

Mpcinfo lists all of the MPC (multiple point constraints). These conditions are assigned to the nodes of node sets using the **mpc** command. There are several ways to view the nodes in a node set in the merge phase. The easiest way is to click on the Sets button in the Pick Panel of the environment window. Alternatively, use the **Labels Nodeset** command found in the Graphics menu.

plane define a boundary plane

plane *plane_#* x_0 y_0 z_0 x_n y_n z_n *tolerance type*

where *type* can be

- **symm** for symmetry plane
- **syf** for symmetry plane with failure
- **ston** *features* for stonewall with optional features
  - where a *feature* can be
    - **stick** *stick_condition*
    - **limit** x_c y_c z_c x_1 y_1
    - **move** *mass initial_velocity*
    - **penalty** *penalty_stiffness*
    - **dlcv** *velocity_ld_curve_#*
- **sph** for SPH element symmetry plane

Remarks

The point (x_0, y_0, z_0) is on the plane with a normal vector (x_n, y_n, z_n).

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This command is used to define nodal constraints for nodes on a symmetry plane. Some simulation codes also have a symmetry with failure and a stone wall feature which is supported by this command.

Nodes are automatically selected for the symmetry plane constraint if they are within the specified tolerance of the plane. The tolerance is not used for the symmetry plane with failure and the stone wall. Use the **syf** and **syfi** commands to select nodes on the symmetry plane with failure. Use the **sw** and **swi** commands to select nodes on the stone wall. These planes are numbered so that there can be multiple stone wall and symmetry plane with failure conditions.

The symmetry feature is complicated, depending on the type of plane and the simulation code. If the symmetry plane is parallel to one of the planes where x=0, y=0, or z=0, then the nodes on the symmetry plane are assigned constraints in the global coordinate system. These types of symmetry planes are referred to as canonical symmetry planes and are equivalent to the following constraints:

plane parallel to x=0: x-displacement, y-rotation, z-rotation
plane parallel to y=0: y-displacement, x-rotation, z-rotation
plane parallel to z=0: z-displacement, x-rotation, y-rotation

Some simulation codes support symmetry planes other than these canonical forms, such as a symmetry plane where x=y (45 degrees). In this case, a different option is used when the model is output. A complication occurs when a set on nodes fall on more than one of these types of symmetry planes.

For DYNA3D, when a node is found on 2 non-canonical distinct symmetry planes, then a sliding boundary plane of vector type is automatically created for this node. If a node is found on 3 or more distinct symmetry planes, it has all of its DOFs constrained. If a node was assigned nodal boundary constraints orthogonal to the sliding boundary plane to which it is assigned, those nodal boundary constraints will be ignored and a warning message will be written.

For LS-DYNA, the canonical (x,y or z) planes are handled as global constrains in the *NODES section. The non-canonical symmetry planes are handled with the *BOUNDARY_SPC_SETS and *BOUNDARY_SPC_NODE keywords along with the *DEFINE_COORDINATE_SYSTEM.

For ABAQUS, ANSYS, NASTRAN and NE/NASTRAN, nodes on non-canonical symmetry planes are constrainted in local coordinate systems.

plinfo write information about defined boundary planes

plinfo *<no arguments>*

Remarks

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This command lists all of the planes defined in the **plane** command.

5. Radiation and Temperatures

temp global default constant temperature

temp *temperature*

Remarks

A temperature is assigned to all parts defined after this command. Use the units expected by the target simulation code as **True*Grid*®** does no conversion.

For the NASTRAN and NE/NASTRAN output, the last value is used for the TEMPD command.

For DYNA3D output, this produces the temperatures for the temperature option 2.

For LS-DYNA output, this produces the *LOAD_THERMAL_CONSTANT_NODE.

Use the **te** or **tei** command to change the default constant temperature for a specific region of the mesh.

bfd bulk fluid definition

bfd *id_# material_# volume*

where

id_#	identification number of this definition to be used by the **bf** command
material_#	material number associated with the bulk fluid
volume	total volume of the material

Remarks

Use **bfd** to define the material number and total volume of a bulk node.

6. Springs, Dampers, and Point Masses

There are two ways to generate springs or dampers. The **spdp** command is used to create an array of springs or dampers between two disconnected faces of the mesh, one spring between each pair of corresponding nodes. The **spring** command is used to create a single spring between two nodes. The **spdp** command is only found in the part phase. The **spring** command is found in both the part and

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merge phase, but the syntax differs. Both of these commands refer to a spring definition number. The **spd** command is used to define the properties of a spring and assign a definition number to these properties.

There are two ways to generate a point mass. The **npm** command creates a new node with a point mass. Since this node is not connected to the rest of the model, you must take additional steps to make sure it is connected through merging or building a beam or spring using this node. The **pm** command can also be used to assign a mass to an existing node. Although the **npm** command may appear to be a global command, it causes point mass replications when found in the part phase when the part is replicated, distinguishing it from the **npm** command found in the merge phase. The **pm** command is found in both the part and merge phase, but their syntax differs.

Springs, dampers, and point masses are replicated when they are specified within a part and the part is replicated. When the spring definition number is incremented in the spring command, make sure that you use the **spd** command to define each spring property.

spd define the properties of a set of springs or dampers

spd *spring/damper_# option type parameters*
where
option can be
dro *flag*
where flag can be:
0 for a linear spring/damper
1 for a torsional spring/damper
dmf *dynamic_factor*
tv *test_velocity*
cl *clearance*
fd *failure_deflection*
dlc *limit_compression*
dlt *limit_tension*
where
type is the spring or damper's material model and
parameters is a list of corresponding parameters, as in the following:
le *stiffness* for linear elastic
lv *damping* for linear viscous
iep *elastic tangent yield* for isotropic elastic
ne *ld_curve_#* for nonlinear elastic
nv *ld_curve_#* for nonlinear viscous
gn *loading_# unloading_# hardening tension compression*

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for general nonlinear

dhpt for a dashpot

mv for a three parameter maxwell viscoelastic

itc for a inelastic tension or compression only

se *elastic_value damping stress* for scalar elastic

mus *l0 vmax sv a fmax tl tv fpe lmax ksh* for muscle

where

l0	for initial muscle length
vmax	for maximum CE shortening velocity
sv	for scale factor for Vmax vs. activs state
a	for activation level vs. time function
fmax	for peak isometric force
tl	for active tension vs. length function
tv	for active tension vs. velocity function
fpe	for force vs. length function
lmax	for relative length
ksh	for exponential rise constant

Remarks

A spring or damper is defined using either the **spdp** command forming a set of springs/dampers between two surfaces, or using the **spring** command to create a single spring at a time. In each case, the definition of a spring includes a reference to a material definition **spd** number.

ANSYS, use linear elastic, linear viscous (damper), isotropic elastoplastic, or the dashpot.

DYNA3D, use linear elastic, linear viscous (damper), isotropic elastoplastic, nonlinear elastic, nonlinear viscous, general tabulated nonlinear, and dashpot.

LS-DYNA, use linear elastic, linear viscous (damper), isotropic elastoplastic, nonlinear elastic, nonlinear viscous, general tabulated nonlinear, dashpot, three parameter Maxwell viscoelastic, inelastic tension or compression only, and muscle. LS-DYNA3D options also include **dro**, **dmf**, **tv**, **cl**, **fd**, **dlc**, and **dlt**.

LS-NIKE3D, use linear elastic, linear viscous (damper), isotropic elastoplastic, nonlinear elastic, and nonlinear viscous.

MARC, use linear elastic, linear viscous (damper), isotropic elastoplastic, or the dashpot.

NASTRAN, use scalar elastic.

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NE/NASTRAN, use scalar elastic.

NIKE3D, use linear elastic, linear viscous (damper), isotropic elastoplastic, nonlinear elastic, nonlinear viscous, general tabulated nonlinear, and dashpot.

If the output option has been selected prior to using the dialogue box to make a selection, only the options available to that output option will be displayed in the dialogue box.

Example

```
nastran
spd 1 se 1 .1 2.1
block 1 6;1 6;1 6;-1 1 -1 1 -1 1
spdp 1 1 2 2 2 2 1 m 1 ;
block 1 6; 1 6; 1 6;-1 1 -1 1 1.1 3.1
spdp 1 1 1 2 2 1 1 s ; ;
merge
write
```

spinfo write information about springs and dampers

spinfo (no arguments)

7. Interfaces

bbinfo block boundary interface information

bbinfo

Remarks

Only the master side of the interface appears in this table. Additional information is printed when first entering the merge phase and performing nodal merging. Master interfaces defined within a part are not included in the table while in that part generation phase. Only after that part is completed will its master block boundary interfaces appear in the table.

Example

The following commands form 7 parts. The first part is a simple cylinder formed with a single block. Each of the faces of this part forms master side of another block boundary. Each of the subsequent parts are glued to this center part using the block boundary interface command.

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```
sd 1 sp 0 0 0 1                          c sphere surface
block 1 11; 1 11; 1 11;-1 1 -1 1 -1 1 c part 1
sfi -1 -2; -1 -2; -1 -2;sd 1             c project to sphere
bb 1 1 1 1 2 2 1;                        c master BB 1
bb 2 1 1 2 2 2 2;                        c master BB 2
bb 1 1 1 2 1 2 3;                        c master BB 3
bb 1 2 1 2 2 2 4;                        c master BB 4
bb 1 1 1 2 2 1 5;                        c master BB 5
bb 1 1 2 2 2 2 6;                        c master BB 6
block 1 11;1 11;1 11;-3 -1 -3 3 -3 3  c part 2
bb 2 1 1 2 2 2 1;                        c slave BB 1
block 1 11;1 11;1 11;1 3 -3 3 -3 3    c part 3
bb 1 1 1 1 2 2 2;                        c slave BB 2
block 1 11;1 11;1 11;-3 3 -3 -1 -3 3  c part 4
bb 1 2 1 2 2 2 3;                        c slave BB 3
block 1 11;1 11;1 11;-3 3 1 3 -3 3    c part 5
bb 1 1 1 2 1 2 4;                        c slave BB 4
block 1 11;1 11;1 11;-3 3 -3 3 -3 -1  c part 6
bb 1 1 2 2 2 2 5;                        c slave BB 5
block 1 11;1 11;1 11;-3 3 -3 3 1 3    c part 7
bb 1 1 1 2 2 1 6;                        c slave BB 6
merge
```

The **bbinfo** command then produced this table of master block boundaries.

```
 Master Block Boundary Interface Table
 BB number      1 is     11 by     11 nodes
 BB number      2 is     11 by     11 nodes
 BB number      3 is     11 by     11 nodes
 BB number      4 is     11 by     11 nodes
 BB number      5 is     11 by     11 nodes
 BB number      6 is     11 by     11 nodes
```

getbb retrieve a block boundary from a part file

getbb *filename source_BB_# TG_alias_BB_#*

where

filename	name of the part file
source_BB_#	block boundary identification number in the part file
TG_alias_BB_#	new identification number for this block boundary

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Remarks

Use the **savepart** command to save the part data including the master block boundary information to a file. The **getbb** command makes it possible to retrieve the block boundary information from this file without having to re-generate the entire part used to create the master side of the block boundary interface.

This feature can be used to define the connections between portions of a model so that a large model can be built by more than one individual. A shell part can be used to form such an interface by forming a cross section of the model. This splits the model. Then each individual must match this cross section as they build their portion of the mesh. Care must be taken to assure that the interface can be matched by both sides of the model.

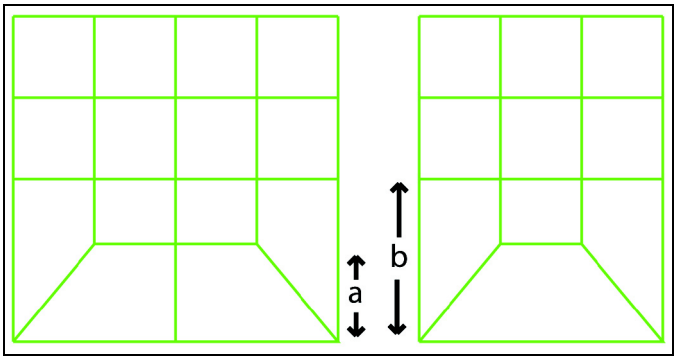
inttr block boundary transition element interpolation factor

inttr α

Remarks

The interior elements of a transition block are interpolated using a parameter between 0 and 1. The default for this parameter is 0.5. Referring to the diagram,

$$\alpha = a / b$$



Transition parameter

Example

```
inttr .6
block 1 5 0 6 9;1 3;-1;1 5 0 6 9 1 3 0 c part 1
bb 1 1 1 2 1 1 1;bb 4 1 1 5 1 1 2; c master BBs
block 1 3 0 4 5;1 2;-1;1 5 0 6 9 -1 1 0 c part 2
trbb 1 2 1 2 2 1 1;trbb 4 2 1 5 2 1 2; c slave transitions
merge
```

mbb master block boundary from point data

mbb *n #_rows #_columns x1 y1 z1 ... xn yn zn trans ;*
where

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n	is the block boundary interface number
#_rows	is the number of rows in the table of coordinates
#_columns	is the number of columns in the table of coordinates
xi yi zi	are the coordinate triples for each node point
trans	is a coordinate transformation

Remarks

This creates a master block boundary from a table of coordinates. This feature is useful when a mesh is imported with the **readmesh** command. A master block boundary can be extracted from this part by picking nodes and using the ***F7*** function key to print out the coordinates. Care must be taken to select a block region and to pick the nodes in the correct order.

Example

```
mbb 10 3 3
 1 3  0  2  3.1 .25 3   3 0
.9 2 .25 2  2  -.25 3.1 2 .25
 1 1  0  2 .9   .25 3   1 0 ;
```

prtrnd project intermediate transition nodes

prtrnd *switch*

where *switch* can be

on

off

Remarks

A transition node is an intermediate node in a transition block. The transition block is the region were the transition from coarse to fine or fine to coarse mesh is located. This is always on the slave side of a transition interface. By default, these intermediate nodes are not projected. They are interpolated from the position of the surrounding nodes..

Care must be taken when using this feature when there are several parts or regions on the slave side of the transition (such as bricks and embedded shells), so that all slave parts or regions involved in the transition have the appropriate faces projected. If only some of the faces are projected to surfaces, then the nodes may not merge properly.

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sid sliding interface definition

sid *slide_# option_list* ;

where *option_list* consists of

- **tied** for tied sliding surface (DYNA3D,NIKE3D)
- **sl** for sliding only (DYNA3D,NIKE3D)
- **sv** for sliding with voids (DYNA3D,NIKE3D)
- **single** for single sided slide surface (DYNA3D,NIKE3D)
- **dni** for discrete nodes impacting surface (DYNA3D)
- **dnt** for discrete nodes tied to surface (DYNA3D)
- **sets** for shell element edge tied to shell element surface (DYNA3D)
- **nsw** for nodes spot welded (DYNA3D)
- **break** for tie-break interface (DYNA3D)
- **owsv** for one way sliding with voids (DYNA3D)
- **dummy** is only used to insure that nodes in this interface will not be merged
- **sand** *options* for the Slide Surface with Adaptive New Definitions (DYNA3D)
  where *options* can be
  - **sms** *slave_material_list* ;
  - **mms** *master_material_list* ;
- **auto** for automatic contact (DYNA3D)
- **inter** for Interface elements (ABAQUS)
- **tcrs** *thermal_contact_resistance* for thermal sliding with thermal contact resistance (TOPAZ3D)
- **rebar** *options* to define properties of REBAR 1D sliding interface (DYNA3D)
  where *options* can be any of the following:
  - **rbrad** radius
  - **rbstr** strength
  - **rbshr** modulus
  - **rbumax** displacement
  - **rbexp** exponent
  - **rbibond** non-negative_number
- **pnlt** *penalty_factor* for sliding penalty (NIKE3D)
- **fric** *friction_factor* for static coefficient of friction (DYNA3D,NIKE3D,ABAQUS)
- **kfric** *kinetic_coefficient_of_friction* for kinetic coefficient of friction (DYNA3D,NIKE3D)
- **decay** *exponential_decay_coefficient* for exponential decay coefficient (DYNA3D,NIKE3D)
- **pen** for small penetration flag (DYNA3D)
- **sfif** for slave to be printed in force file (DYNA3D)

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mfif for master to be printed in force file (DYNA3D)
pnlts *slave_penalty_factor* for slave penalty factor (DYNA3D)
pnltm *master_penalty_factor* for master penalty factor (DYNA3D)
bwmrad *#_facets* to set the bandwidth minimization radius (NIKE3D)
fric2 *friction_factor* for the anisotropic friction coefficient (ABAQUS)
stif *stiffness* for stiffness in stick (ABAQUS)
essl *stress* for the equivalent shear stress limit (ABAQUS)
penmax *distance* to set the small penetration search distance
iaug *flag* to set the augmentation flag
where the *flag* can be
1 to augment until convergence tolerance is satisfied
0 for no augmentations (penalty method)
-n for the number of augmentations per step
altoln *tolerance* to set the normal direction convergence tolerance
altolt *tolerance* to set the tangential direction convergence tolerance
tkmult *multiplier* to set the tangent stiffness multiplier
dtime *time* to set the interface death time
bury *time* to set the interface burial time
concon *conductance* contact conductance
radcon *conductance* radiation conductance
lsdsi *type options* ; for LS-DYNA sliding interfaces (LS-DYNA)
where *type* can be:
1 for Sliding without penalties
p1 for Symmetric sliding with penalties
2 for Tied
3 for Sliding, impact, friction
a3 for Sliding, impact, friction, no segmentation orientation
4 for Single surface contact
5 for Discrete nodes impacting surface
a5 for Discrete nodes impacting surface, no segmentation orientation
6 for Discrete nodes tied to surface
7 for Shell edge tied to shell surface
8 for Nodes spot welded to surface
9 for Tiebreak interface
10 for One way treatment of sliding, impact, friction
a10 for One way treatment, no segmentation orientation
13 for Automatic single surface with beams and arbitrary orientations
a13 for Automatic single surface with beams and arbitrary

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orientations with extra search for airbag contact

14 for Surface to surface eroding contact

15 for Single surface eroding contact

16 for Node to surface eroding contact

17 for Surface to surface symmetric/asymmetric constraint method

18 for Node to surface constraint method (Taylor and Flanagan 1989)

19 for Rigid body to rigid body contact with arbitrary force/deflection curve

20 for Rigid nodes to rigid body contact with arbitrary force/deflection curve

21 for Rigid body to rigid body contact with arbitrary force/deflection curve (one way treatment)

22 for Single edge treatment for shell surface edge to edge treatment

23 for Simulated draw bead

24 for Automatic surface to surface tiebreak

25 for Automatic one way surface to surface tiebreak

34 for Automatic general

35 for Automatic general interior

36 for Force transducer constraint

37 for Force transducer penalty

38 for Forming node to surface

39 for Forming one way surface to surface

40 for Forming surface to surface

41 for Discrete nodes impacting surface w/ interference

42 for One way treatment of sliding, impact, friction w/ interference

43 for Spotweld

44 for Spotweld with torsion

45 for Sliding, impact, friction w/ interference

46 for Tiebreak nodes only

47 for Tied with failure

rebar for rebar in concrete 1D sliding

and an *option* can be:

lcrsgo *load_curve_#* for optional load curve defining the resisting stress vs. gap opening

isrch *flag* small penetration in contact search

where *flag* can be

0 for check is off

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1 for check is on
2 for check is on, shortest diagonal used

visdam *percent* for viscous damping coefficient in percent of critical

kpf *flag* kinematic partition factor for constraint

where

flag can be

0 for fully automatic treatment
1 for one way treatment with slave nodes constrained to master surface
-1 for one way treatment with master nodes constrained to slave surface

lcair *load_curve_#* load curve defining airbag thickness

penmax *penetration* maximum penetration

thkopt f*lag* thickness option

where

flag can be

0 for default from the control cards
1 for thickness is not considered
2 for thickness is considered but rigid bodies are excluded
3 for thickness is considered including rigid bodies
4 for thickness effects are not included

lcfpb *load_curve_#* force vs. penetration behavior load curve

fcm *flag* force calculation method

where

flag can be

1 for total normal force on surface vs. max. penetration of any node
2 for normal force on each node vs. penetration of node through the surface
3 for normal pressure vs. penetration of node into surface
4 for total normal force vs. max. soft penetration

unstf unloading stiffness

lcbcrf *load_curve_#* load curve giving the bending component of the retaining force

lcnflen *load_curve_#* load curve giving the normal force per unit draw bead length as a function of displacement

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dbd *depth* draw bead depth
scllc *factor* scale factor for load curve
nitdb *#_iterations* number of integration points laong the draw bead
slvmat *material_list* ; automatic slave segment materials
sypl slave, do not include faces with normal boundary constraints
serin slave erosion/interior node option
sadjmat slave storage is allocated so that eroding contact can occur
scoufsf *factor* oulomb friction scale factor
svfsf *factor* viscous friction scale factor
snffs *stress_or_force* slave normal stress at failure
ssffs *stress_or_force* slave shear stress at failure
senf *exponent* slave exponent for normal force
sesf *exponent* slave exponent for shear force
mstmat *material_list* ; automatic master segment materials
mypl master, do not include faces with normal boundary constraints
merin master erosion/interior node option
madjmat master storage is allocated so that eroding contact can occur
mcoufsf *factor*
mvfsf *factor*
mnffs *stress_or_force* master normal stress at failure
msffs *stress_or_force* master shear stress at failure
menf *exponent* master exponent for normal force
mesf *exponent* master exponent for shear force
scoef *coefficient* static coefficient of friction
dcoef *coefficient* dynamic coefficient of friction
decay *coefficient* dynamic decay coefficient
incslv include slave side in printed and binary force interface file
incmst include master side in printed and binary force interface file
sfsps *factor* scale factor on default slave penalty stiffness
sfmps *factor* scale factor on default master penalty stiffness
vfcoef *coefficient* coefficient for viscous friction
thss *thickness* optional thickness for slave surface
thms *thickness* optional thickness for master surface

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sthss *factor* scale factor for slave surface thickness
sthms *factor* scale factor for master surface thickness
btime *time* birth time
dtime *time* death time
softc *flag* soft constraint option
where
flag can be:
0 for penalty formulation
1 for soft constraint formulation
ssoftc *factor* scale factor for constraint forces of soft constraint option
maxpcss *coordinate* maximum parametric coordinate in segment search
srchdp *depth* search depth in automatic contact
ncybs *cycles* number of cycles between bucket sorts
ncyup *cycles* number of cycles between contact force updates for penalty formulations
diskoc disable logic in thickness offset contact to avoid shooting nodes
concon *conductance* contact conductance
radcon *conductance* radiation conductance
gapcs *size* gap critical size
ctofst
atbo *args*
where *args* can be
1 for Slave nodes in contact and which come into contact will permanently stick. Tangential motion is inhibited.
2 normal_stress shear_stress
for Tiebreak is active for nodes which are initially in contact. Until failure, tangential motion is inhibited.
3 normal_stress shear_stress
for Same as 1st option but with failure after sticking
4 for Tiebreak is active for nodes which are initially in contact but tangential motion with frictional sliding is permitted.
5 plastic_stress load_curve
for Tiebreak is active for nodes which are initially in contact. Damage is defined by a

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load curve.

6 distance
for Tiebreak is active for brick and thick shell nodes which are initially in contact. damage is a linear function between points.

lcid1 *load_curve*	load curve dynamic interface stiffness
lcid2 *load_curve*	load curve transient interface stiffness
isym *option_#*	symmetric plane option
i2d3d *option_#*	segment searching option
sldthk *thickness*	solid element thickness
sldstf *thickness*	solid element stiffness
igap *option_#*	flag implicit convergence behavior
ignore *option_#*	ignore initial penetration in automatic interfaces
edge *distance*	edge to edge penetration check
rbrad *radius*	
rbstr *strength*	
rbshr *modulus*	
rbumax *strain*	
rbexp *exponent*	

Remarks

Sliding interfaces or contact surfaces are constructed in 3 steps. These steps can be done in any order.

1. define the properties
2. select the slave side
3. select the master side, if applicable

The **sid** command is used to define the properties. The **si** and **sii** commands are used in the part phase or the merge phase to select the nodes or faces that form the master and slave sides of the interface.

The options for LS-DYNA are large and unique so they have been singled out and fall under the **lsdsi** option of the **sid** command. Some of these sliding interfaces require nodes on the slave side while others require only a set of faces on the slave side. This definition is required so that the proper data can be written to the output file.

With some simulation code formats, one can construct a node set or a face set. This will be written to the output file as a set. Then it is a simple matter to add the keyword command to the output file using a text editor to transform that set into a contact surface or sliding interface. This approach has the problem that nodes may be merged across the two sides because they are not defined as sliding

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interfaces.

When nodes are merged, nodes across a sliding interface will not be merged. When a merge command is first issued in the merge phase, a table is written listing the number of nodes and faces associated with each sliding interface.

The **dummy** type interface is actually used to avoid merging of nodes. A sliding interface of this type is not written to the output file.

The nodes and faces of a sliding interface or contact surface can be viewed in the merge phase using the **si** option of the **co** command.

If the output option has been selected prior to using the dialogue box to make a selection, only the options available to that output option will be displayed in the dialogue box.

Examples

```
sid 1 tied pnlt 10;
sid 10 rebar rbrad .01;;
sid 13 lsdsi 9 lcrsgo 1 visdam .34 isrch 1 ; ; ;
sid 54 dummy;
```

siinfo list sliding interface definitions

siinfo <*no arguements*>

8. Elements

The default face IDs of brick elements depends on the element geometry after nodes have been properly merged so that **True*Grid*®** can determine the element geometry (i.e. hex, wedge, or tet). The face IDs can be seen when in the merge phase by using the "**la facesel**" command. The face IDs differ among the different simulation codes. When you use the graphics to help you inspect and verify your model, you may need to convert the face IDs you see in **True*Grid*®** graphics to the face IDs defined for your simulation code. This interpretation of the face ID numbers is only needed when using the "**la facesel**" command in the merge phase. When **True*Grid*®** writes an output file for a particular simulation code, it does the translation of the face ID numbers appropriate for that simulation code.

Almost all simulation codes that use face IDs agree on how to identify faces of a hex element. But there is significant variation among the different simulation codes on how to identify faces of a wedge or tet element. The diagrams below show how **True*Grid*®** identifies faces for all three element

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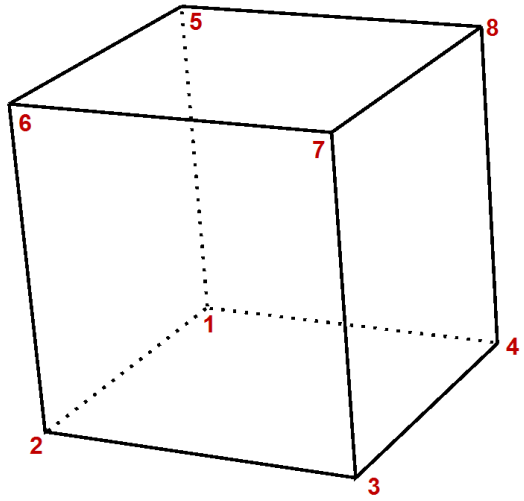
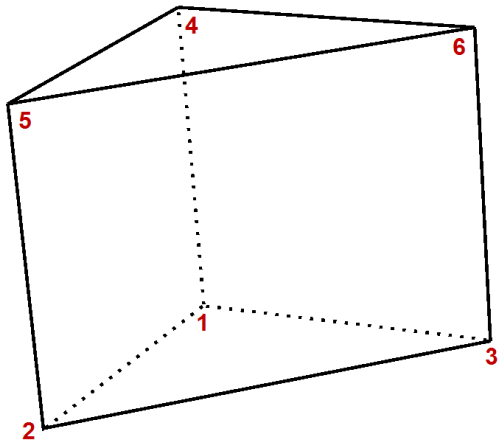
geometries. They are all based on the order of nodes that define the element. The sequence of nodes are identified on the left diagrams. The corresponding faces are identified in the table to the right of the diagram. These rules apply for **linear**, **quadratic**, and **triquadratic** elements.

This is a change from the past versions of **True*Grid***®. In the past, when a wedge or tet element was formed from a degenerate hex, the face ID numbers from the "**la facesel**" command showed the face ID numbers that were inherited from the hex element that formed the wedge or tet.

NOTE: Dotted line means it is hidden behind faces of the element.



Face ID	Node 1	Node 2	Node 3	Node 4
1	1	2	3	4
2	1	5	6	2
3	2	6	7	3
4	4	3	7	8
5	1	4	8	5
6	5	8	7	6



Face ID	Node 1	Node 2	Node 3	Node 4
1	1	2	3	-
2	4	5	6	-
3	1	4	5	2
4	1	3	6	4

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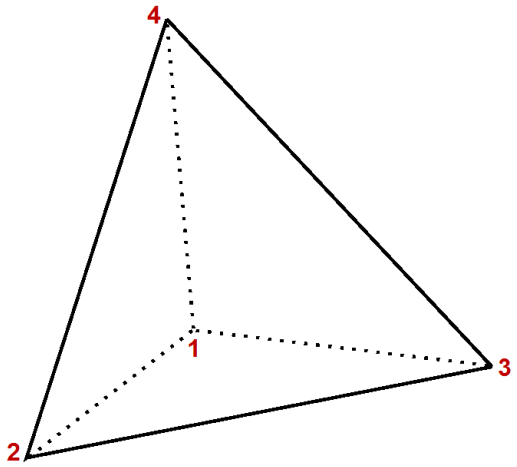
5	2	5	6	3



Face ID	Node 1	Node 2	Node 3
1	1	2	3
2	1	4	2
3	2	4	3
4	3	4	1

Beam elements can be created in three ways. The primary method of beam element generation extracts the needed nodes from an existing shell of brick part. This is only available within the block or cylinder part phase. The **ibm** and **ibmi** commands create beams along i-lines of the mesh, the **jbm** and **jbmi** along j-lines, and the **kbm** and **kbmi** along k-lines. This is a way to embed beam elements within a shell of brick structure. Alternatively, the material of the parent shell of brick part can be set to 0 so that the part can be generated as usual, but so that the shell or brick elements will not be saved when the part generation is ended. The nodes that are used in any of these beam commands will be saved along with the beam elements.

The second method of beam element generation uses the **bm** command. This is fully interactive. Beams are strung along a 3D curve.

The **beam** and **cbeam** parts create beams by connecting a pair of nodes with a line of beam elements. This is an old method included to be backwards compatible, but it is not interactive.

Beam cross section properties are defined using **bsd** and **bind**. Shell cross section properties are defined with the **sind** command.

Elements and nodes can have their numbers offset using the **offset** command. This makes it easier to glue several models together by concatenating two files.

Cross sectional properties, and in particular thicknesses, are not scaled by the **xsca**, **ysca**, **zsca**, and

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csca commands.

bsd global beam cross section definition

bsd *cross_section_# option_list ;*

where

option_list consists of some of the following:

sthi *thickness*	for both thicknesses in s-direction
sthi1 *thickness*	for first thickness in s-direction
sthi2 *thickness*	for second thickness in s-direction
tthi *thickness*	for both thicknesses in t-direction
tthi1 *thickness*	for first thickness in t-direction
tthi2 thickness	for second thickness in t-direction
ldp *distance*	for NEUTRAL file beams

ABAQUS beams

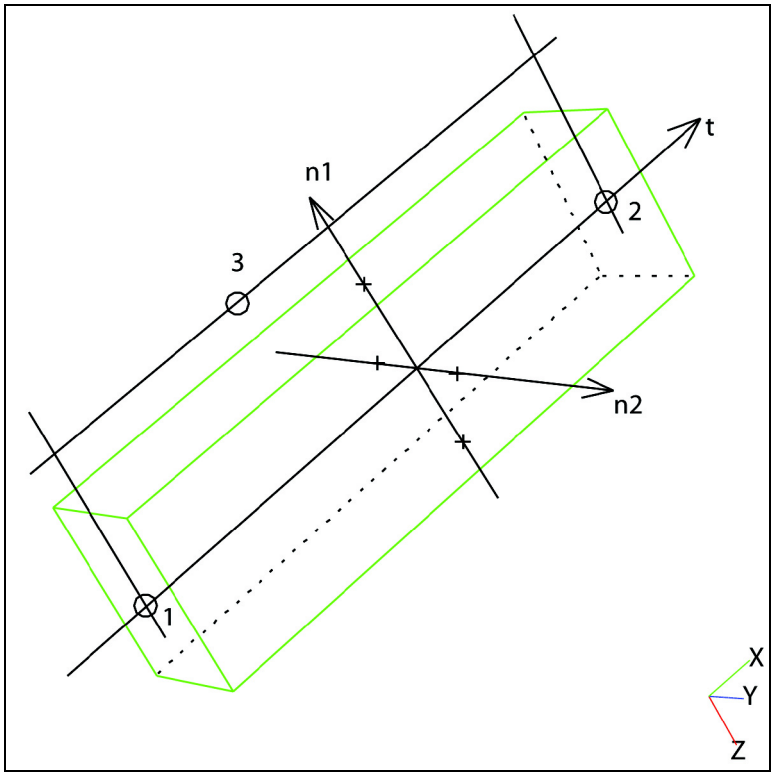

Figure 290 Beam Local Coordinate System for ABAQUS

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cstype *type t_options* ;

where

type and *t_options* can be:

7 *t_options ;* for PIPE (ABAQUS)

where

t_options can be

ABCS1 *radius*
ABCS2 *thickness*
NABIP1 *#_integrations*
TRSS *stiffness*
ABTEMP *Temperature*

8 *t_options ;* for BOX (ABAQUS)

where

t_options can be

ABCS1 *width*
ABCS2 *height*
ABCS3 *thickness*
ABCS4 *thickness*
ABCS5 *thickness*
ABCS6 *thickness*
NABIP1 *#_integrations*
NABIP2 *#_integrations*
TRSS *stiffness*
ABTEMP *Temperature*

9 *t_options ;* for CIRCLE (ABAQUS)

where

t_options can be

ABCS1 *radius*
NABIP1 *#_integrations*
NABIP2 *#_integrations*
TRSS *stiffness*
ABTEMP *temperature*

10 *t_options ;* for I-BEAM (ABAQUS)

where

t_options can be

ABCS1 *depth*
ABCS2 *height*
ABCS3 *width*
ABCS4 *width*
ABCS5 *thickness*

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ABCS6 *thickness*
ABCS7 *thickness*
NABIP1 *#_integrations*
NABIP3 *#_integrations*
TRSS *stiffness*
ABTEMP *temperature*

11 *t_options ;* for RECTANGLE (ABAQUS)

where

t_options can be

ABCS1 *width*
ABCS2 *height*
NABIP1 *#_integrations*
NABIP2 *#_integrations*
TRSS *stiffness*
ABTEMP *temperature*

12 *t_options ;* for HEXAGON (ABAQUS)

where

t_options can be

ABCS1 *thickness*
ABCS2 *thickness*
NABIP1 *#_integrations*
NABIP2 *#_integrations*
TRSS *stiffness*
ABTEMP *temperature*

13 *t_options ;* for ELBOW (ABAQUS)

where

t_options can be

ABCS1 *radius*
ABCS2 *thickness*
ABCS3 *radius*
NABIP1 *#_integrations*
NABIP2 *#_integrations*
NABIP3 *#_integrations*
TRSS *stiffness*
ABTEMP *temperature*

14 *t_options ;* for TRAPEZOID (ABAQUS)

where

t_options can be

ABCS1 *width*

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ABCS2 *height*
ABCS3 *width*
ABCS4 *depth*
NABIP1 *#_integrations*
NABIP2 #_integrations
RSS *stiffness*
ABTEMP *temperature*

15 *t_options ;* for I-SECTION (ABAQUS)

where

t_options can be

ABCS1 *width*
ABCS2 *height*
ABCS3 *thickness*
ABCS4 *thickness*
NABIP1 *#_integrations*
NABIP2 *#_integrations*
TRSS *stiffness*
ABTEMP *temperature*

16 *t_options ;* for ARBITRARY (ABAQUS)

where

t_options can be

CSCRV *y1 z1 ... yn zn* ;
CSSTH thick1 ... *thickn* ;
TRSS *stiffness*
ABTEMP *temperature*

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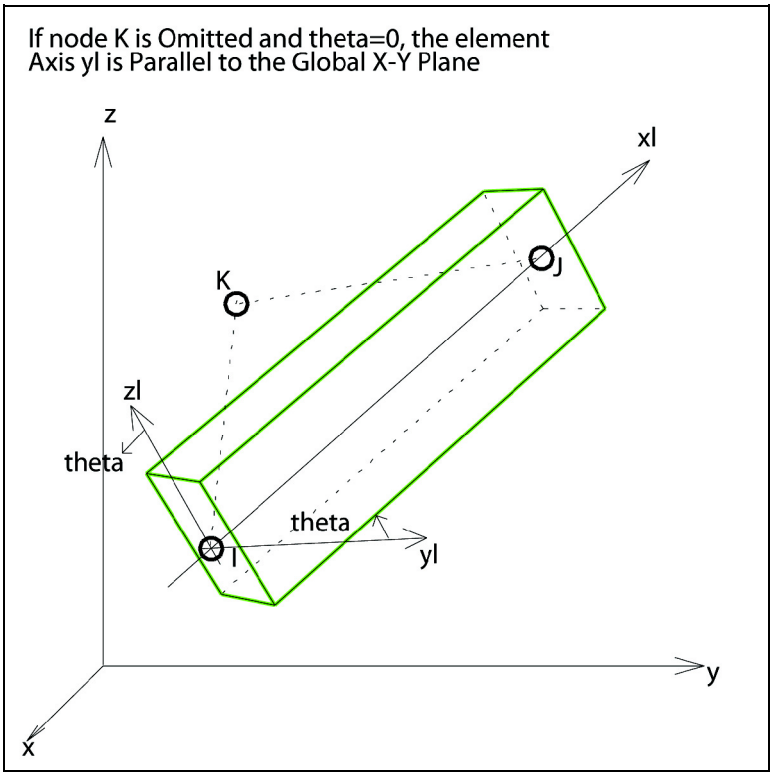

Figure 291 Beam Local Coordinate System for ANSYS

ban4 *t_options* ; for ELASTIC BEAMS (ANSYS)

where

t_options can be

AREA *area*
IXX moment
IYY *moment*
IZZ *moment*
HEIGHT *height*
WIDTH *width*
THETA *angle*
INSTR *strain*
SHEARY *constant*
SHEARZ *constant*

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ban8 *t_options* ; for SPARS (ANSYS)
where
t_options can be
AREA *area*
INSTR *strain*
ban10 *t_options* ; for TENSION/COMPRESSION SPARS (ANSYS)
where
t_options can be
AREA *area*
INSTR *strain*

ban24 *t_options* ; for THIN WALLED PLASTIC BEAMS (ANSYS)
where
t_options can be
CSCRV *y1 z1 ... yn zn* ;
CSSTH *thick1 ... thickn* ;
RXOFF1 *x-offset*
RXOFF2 *x-offset*
SHEARY *constant*
SHEARZ *constant*
ban33 *t_options* ; for THERMAL BARS (ANSYS)
where
t_options can be
AREA *area*
ban44 *t_options* ; for TAPPERED UNSYMMETRICAL BEAMS (ANSYS)
where
t_options can be
AREA *area*
AREA1 *area*
AREA2 *area*
IXX *moment*
IXX1 *moment*
IXX2 *moment*
IYY *moment*
IYY1 *moment*
IYY2 *moment*
IZZ *moment*
IZZ1 *moment*
IZZ2 *moment*
YB *width*

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YB1 *width*
YB2 *width*
YA *width*
YA1 *width*
YA2 *width*
ZB *height*
ZB1 *height*
ZB2 *height*
ZA *height*
ZA1 *height*
ZA2 *height*
XOFF1 *x-component*
YOFF1 *y-component*
ZOFF1 *z-component*
XOFF2 *x-component*
YOFF2 *y-component*
ZOFF2 *z-component*
SHEARY *constant*
SHEARZ *constant*

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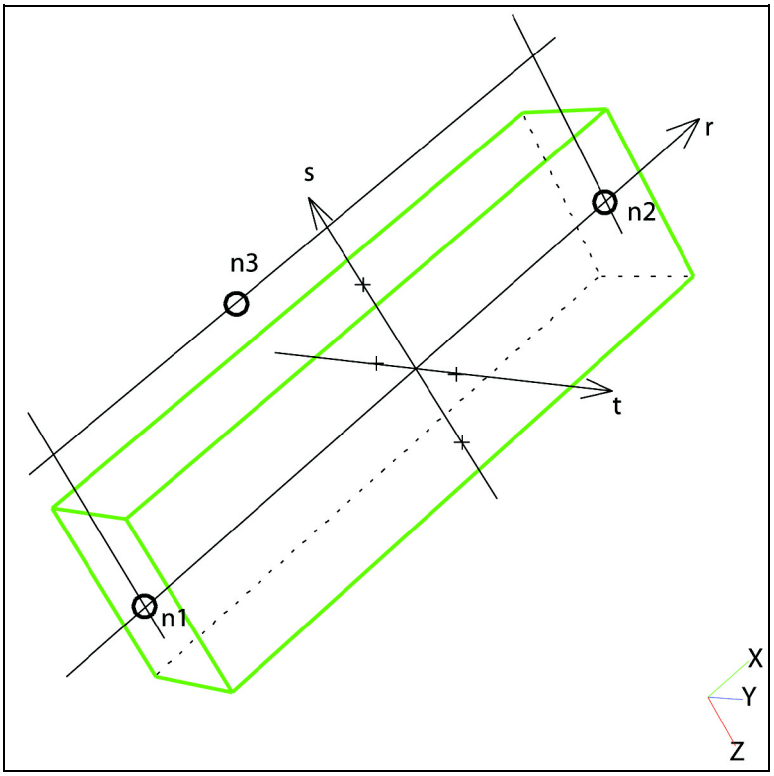

Figure 292 Beam Local Coordinate System for DYNA3D

Parameters

where

parameters can be

(Hughes-Liu beam, constant thickness)

STHI *thickness* (s-thickness at both ends)

TTHI *thickness* (t-thickness at both ends)

or

(Hughes-Liu beam, variable thickness)

STHI1 *thickness* (s-thickness at beginning)

STHI2 *thickness* (s-thickness at ending)

TTHI1 *thickness* (t-thickness at beginning)

TTHI2 *thickness* (t-thickness at ending)

or

(Belytschko-Schwer beam)

carea *area* cross section area

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iss *iss* area moment of inertia about s-axis
itt *itt* area moment of inertia about t-axis
irr *irr* area moment of inertia about r-axis
sarea *area* shear area of cross section
or
(Truss)
carea *area* cross section area

LS-DYNA beams

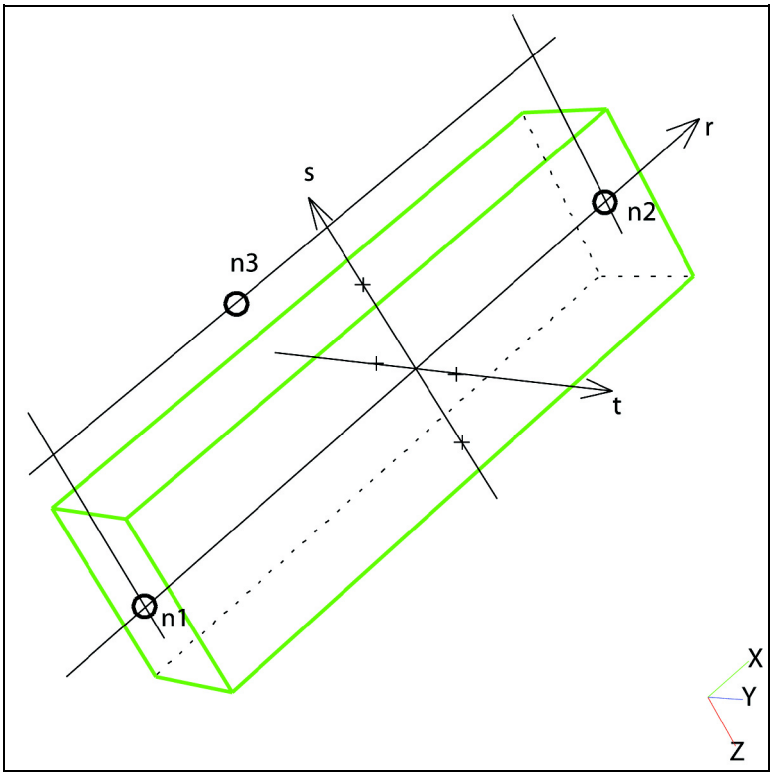

Figure 293 Beam Local Coordinate System for LS-DYNA

Parameters ; for LS-DYNA beams
(Dimensions of Standard Cross Sections are in Global Coordinates)
where
parameters can be either:

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(Hughes-Liu Standard Sections)
lsd 1 *flange_width flange_thickness depth web_thickness* (W-section)
lsd 2 *flange_width flange_thickness depth web_thickness* (C-section)
lsd 3 *flange_width flange_thickness depth web_thickness* (angle)
lsd 4 *flange_width flange_thickness depth web_thickness* (T-section)
lsd 5 *flange_width flange_thickness depth web_thickness* (rectangular)
lsd 6 *flange_width flange_thickness depth web_thickness* (Z-section)
lsd 7 *flange_width depth web_thickness* (trapezoidal)
SREF *location*
where
location can be:
1 meaning the side where s is 1
0 meaning centered
-1 meaning the side where s is -1
TREF *location*
where
location can be:
1 meaning the side where t is 1
0 meaning centered
-1 meaning the side where t is -1
or
(Hughes-Liu Constant Thickness or Diameter)
STHI *thickness* (s-thickness or outer diameter at both ends)
TTHI *thickness* (t-thickness or inner diameter at both ends)
SREF *location*
where
location can be:
1 meaning the side where s is 1
0 meaning centered
-1 meaning the side where s is -1
TREF *location*
where
location can be:
1 meaning the side where t is 1
0 meaning centered
-1 meaning the side where t is -1
or
(Hughes-Liu Variable Thicknesses or Diameters)
STHI1 *thickness* (s-thickness or outer diameter at beginning)
STHI2 *thickness* (s-thickness or outer diameter at ending)

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TTHI1 *thickness* (t-thickness or inner diameter at beginning)
TTHI2 *thickness* (t-thickness or inner diameter at ending)
SREF *location*
where
location can be:
1 meaning the side where s is 1
0 meaning centered
-1 meaning the side where s is -1
TREF *location*
where
location can be:
1 meaning the side where t is 1
0 meaning centered
-1 meaning the side where t is -1
or
(Belytschko-Schwer beam)
carea *area* cross section area
iss *iss* area moment of inertia about s-axis
itt *itt* area moment of inertia about t-axis
irr *irr* area moment of inertia about r-axis
sarea *area* shear area of cross section
or
(Truss)
carea *area* cross section area
or
(Belytschko-Schwer Full Integration Beam Standart Sections)
lsd 1 *flange_width flange_thickness depth web_thickness* (W-section)
lsd 2 *flange_width flange_thickness depth web_thickness* (C-section)
lsd 3 *flange_width flange_thickness depth web_thickness* (angle)
lsd 4 *flange_width flange_thickness depth web_thickness* (T-section)
lsd 5 *flange_width flange_thickness depth web_thickness* (rectangular)
lsd 6 *flange_width flange_thickness depth web_thickness* (Z-section)
lsd 7 *flange_width depth web_thickness* (trapezoidal)
or
(Belytschko-Schwer Full Integration Beam Constant Thickness or Diameters)
STHI *thickness* (s-thickness or outer diameter at both ends)
TTHI *thickness* (t-thickness or inner diameter at both ends)
or
(Belytschko-Schwer Full Integration Beam Variable Thicknesses or

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Diameters)
STHI1 *thickness* (s-thickness or outer diameter at beginning)
STHI2 *thickness* (s-thickness or outer diameter at ending)
TTHI1 *thickness* (t-thickness or inner diameter at beginning)
TTHI2 *thickness* (t-thickness or inner diameter at ending)

or
(Belytschko-Schwer Tubular Beam Constant Diameter)
STHI *outer_diameter* (outer diameter at both ends)
TTH *inner_diameter* (inner diameter at both ends)
or
(Belytschko-Schwer Tubular Beam Variable Diameter)
STHI1 *first_outer_diameter* (outer diameter at beginning)
STHI2 *last_outer_diameter* (outer diameter at ending)
TTHI1 *first_inner_diameter* (inner diameter at beginning)
TTHI2 *last_inner_diameter* (innter diameter at ending)
or
(Discrete 3D Beam)
vold *volume*
cabarea *cable_area*
lump *inertia* (lumped geometric inertia)
cablcid *local_coordinate_system_#* (defined by the **lsys** command)
caboff *cable_offset*
or
(Spot Weld Beam Standart Sections)
lsd 1 *flange_width flange_thickness depth web_thickness* (W-section)
lsd 2 *flange_width flange_thickness depth web_thickness* (C-section)
lsd 3 *flange_width flange_thickness depth web_thickness* (angle)
lsd 4 *flange_width flange_thickness depth web_thickness* (T-section)
lsd 5 *flange_width flange_thickness depth web_thickness* (rectangular)
lsd 6 *flange_width flange_thickness depth web_thickness* (Z-section)
lsd 7 *flange_width depth web_thickness* (trapezoidal)
or
(Spot Weld Beam Constant Thickness)
STHI *thickness* (s-thickness at both ends)
TTHI *thickness* (t-thickness at both ends)
or
(Spot Weld Beam Variable Thicknesses)
STHI1 *thickness* (s-thickness at beginning)
STHI2 *thickness* (s-thickness at ending)

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TTHI1 *thickness* (t-thickness at beginning)
TTHI2 *thickness* (t-thickness at ending)

MARC beams

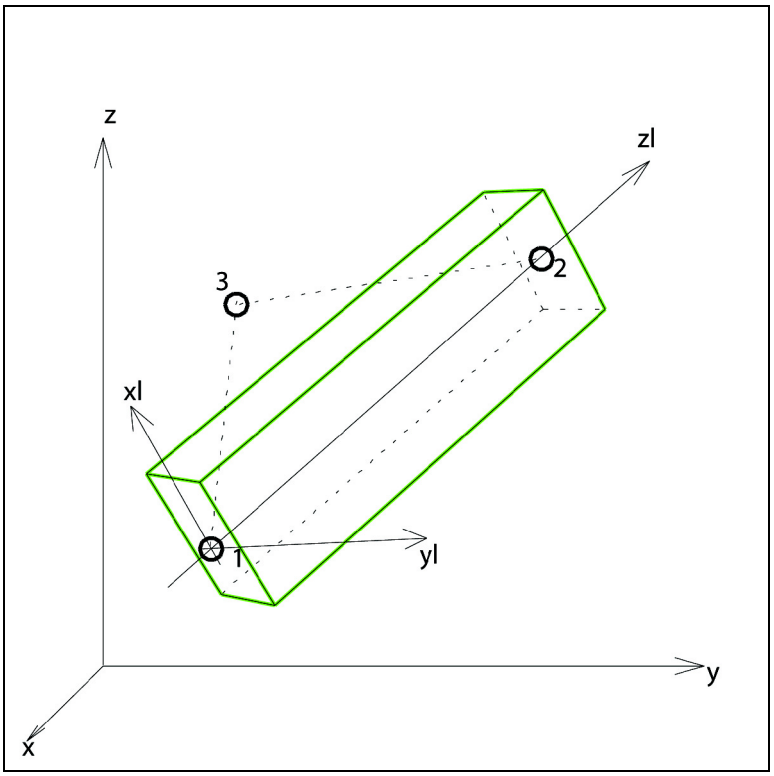

Figure 294 Beam Local Coordinate System for MARC

marc13 *t_options* ; for OPEN SECTION THIN WALLED BEAMS (MARC)
where
t_options can be:
CSCRV *y(1) z(1) ... y(n+1) z(n+1) ;*
CSSTH *thick(1) ... thickn* ;
VCSSTH *first_thick(1) last_thick(1)... first_thick(n) last_thick(n)* ;
CSSLEN *length(1) ... length(n)* ;
CSSSLP *first_dy(1) first_dz(1) last_dy(1) last_dz(1) ... first_dy(n) first_dz(n) last_dy(n) last_dz(n) ;*
CSDIV *#_divisions(1) ... #_divisions(n)* ;

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marc14 *t_options* ; for THIN WALLED BEAMS W/O WARPING (MARC)

where

t_options can be:

CSCRV *y(1) z(1) ... y(n+1) z(n+1) ;*

CSSTH *thick(1) ... thickn ;*

VCSSTH *first_thick(1) last_thick(1) ...first_thick(n) last_thick(n) ;*

CSSLEN *length(1) ... length(n)* ;

CSSSLP *first_dy(1) first_dz(1) last_dy(1) last_dz(1) ... first_dy(n) first_dz(n) last_dy(n) last_dz(n)* ;

CSDIV *#_divisions(1) ... #_divisions(n)* ;

THICK *thickness*

RADIUS *radius*

marc25 *t_options* ; for THIN WALLED BEAMS (MARC)

where

t_options can be:

CSCRV *y(1) z(1) ... y(n+1) z(n+1) ;*

CSSTH *thick(1) ... thickn* ;

VCSSTH *first_thick(1) last_thick(1) ...first_thick(n) last_thick(n)* ;

CSSLEN *length(1) ... length(n)* ;

CSSSLP *first_dy(1) first_dz(1) last_dy(1) last_dz(1) ... first_dy(n) first_dz(n) last_dy(n) last_dz(n)* ;

CSDIV *#_divisions(1) ... #_divisions(n)* ;

THICK *thickness*

RADIUS *radius*

marc31 *t_options* ; for ELASTIC CURVED PIPE (MARC)

where

t_options can be:

AREA *area*

IYY *moment*

IZZ *moment*

IXX *moment*

ETSAY *area*

ETSAZ *area*

THICK *thickness*

RADIUS *radius*

BEND *radius*

marc52 *t_options* ; for ELASTIC BEAMS (MARC)

where

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t_options can be
AREA *area*
IYY *moment*
IZZ *moment*
IXX *moment*
ETSAY *area*
ETSAZ *area*

marc76 *t_options* ; for THIN WALLED BEAMS W/O WARPING (MARC)
where
t_options can be:
CSCRV *y(1)* z(1) ... *y(n+1) z(n+1)* ;
CSSTH *thick(1) ... thickn* ;
VCSSTH *first_thick(1) last_thick(1)... first_thick(n) last_thick(n)* ;
CSSLEN *length(1) ... length(n)* ;
CSSSLP *first_dy(1) first_dz(1) last_dy(1) last_dz(1) ... first_dy(n) first_dz(n) last_dy(n) last_dz(n)* ;
CSDIV *#_divisions(1) ... #_divisions(n)* ;
THICK *thickness*
RADIUS *radius*

marc79 *t_options* ; for THIN WALLED BEAMS WITH WARPING (MARC)
where
t_options can be:
CSCRV *y(1) z(1) ... y(n+1) z(n+1)* ;
CSSTH *thick(1) ... thickn* ;
VCSSTH *first_thick(1) last_thick(1)... first_thick(n) last_thick(n)* ;
CSSLEN *length(1) ... length(n)* ;
CSSSLP *first_dy(1) first_dz(1) last_dy(1) last_dz(1) ... first_dy(n) first_dz(n) last_dy(n) last_dz(n)* ;
CSDIV *#_divisions(1) ... #_divisions(n)* ;

marc98 *t_options* ; for ELASTIC BEAMS WITH TRANSVERSE SHEAR (MARC)
where
t_options can be:
AREA *area*
IYY *moment*
IZZ *moment*
IXX *moment*
ETSAY *area*
ETSAZ *area*

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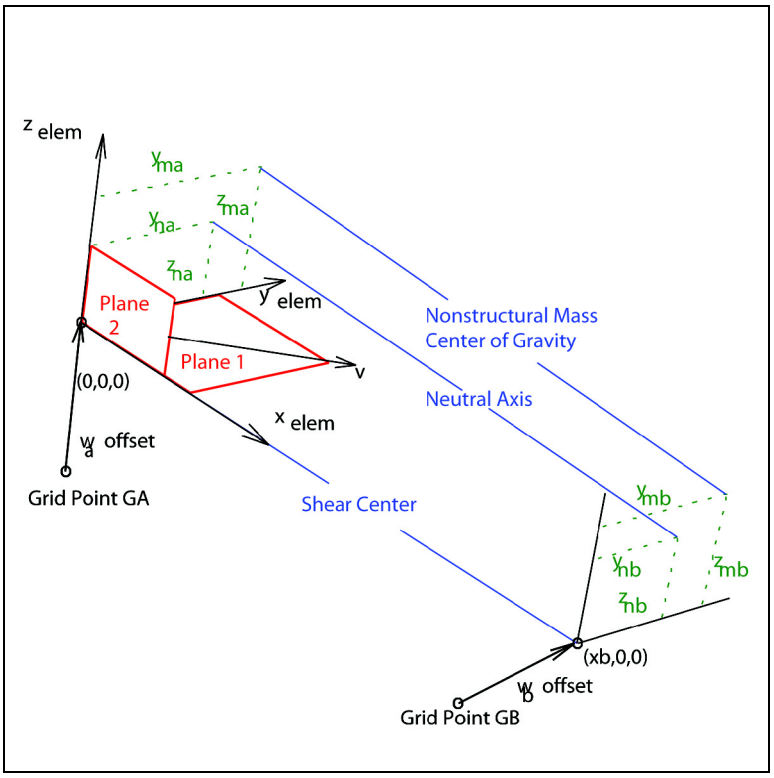

Figure 295 Beam Local Coordinate System for NASTRAN Beam

bna1 *t_options ;*

where

t_options can be:

SHSTF *y-shear* z-shear

SHRLF *y-shear* z-shear

NSMMI *moment*

NSMMI1 *moment*

NSMMI2 *moment*

WARP *coefficient*

WARP1 *coefficient*

WARP2 *coefficient*

CENGRAV *y1 z1 y2 z2*

NEUAXIS *y1 z1 y2 z2*

SPCSD *position c_options* ;

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where
c_options can be
AREA *area*
IYY *moment*
IZZ *moment*
IYZ *moment*
IXX *moment*
MASS *mass*
SDR1 *y1 z1*
SDR2 *y2 z2*
SDR3 *y3 z3*
SDR4 *y4 z4*
DX1
DZ1
RX1
RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

bna2 *t_options* ; for OFFSET RODS (NASTRAN)
where
t_options can be:
AREA *area*
IYY *moment*
IZZ *moment*
IYZ *moment*
IXX *moment*
MASS *mass*
CENGRAV y *z*
NEUAXIS y *z*
CSTYPE *type c_options* ;
where
type can be
1 for the default elliptic
2 for symmetry about y and z
3 for symmetry about y

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4 for symmetry about z
5 for symmetry about y=z=0
6 for arbitrary

c_options can be
CSCRV *y1 z1 ... yn zn* ;
CSSAR *area1 ... arean* ;
CSSMAT *material1 ... materialn* ;

DX1
DY1
DZ1
RX1
RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

bna3 *t_options* ; for CURVED BEAMS (NASTRAN)
where
t_options can be:
AREA *area*
IYY *moment*
IZZ *moment*
IXX *moment*
RB *radius*
THETAB *angle*
SHSTF *y-shear z-shear*
SDR1 *y1 z1*
SDR2 *y2 z2*
SDR3 *y3 z3*
SDR4 *y4 z4*
RC *offset*
ZC *offset*
DELTAN *offset*
DX1
DY1
DZ1
RX1

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RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

bna4 *t_options* ; for ELBOW & CURVED PIPE (NASTRAN)

where

t_options can be:

FSI *intensification*
MCSR *radius*
WTH *thickness*
IP *pressure*
RB *radius*
THETAB *angle*
MASS *mass*
RC *offset*
ZC *offset*
DX1
DY1
DZ1
RX1
RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

bna5 *t_options* ; for SIMPLE BEAM (BAR) (NASTRAN)

where

t_options can be:

AREA *area*
IYY *moment*
IZZ *moment*
IYZ *moment*
IXX *moment*

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MASS *mass*
SHSTF *y-shear* z-shear
SDR1 *y1* z1
SDR2 *y2* z2
SDR3 *y3* z3
SDR4 *y4* z4
DX1
DY1
DZ1
RX1
RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

bna6 *t_options* ; for ROD (NASTRAN)

where

t_options can be:

AREA *area*
IXX *moment*
TSC *coefficient*
MASS *mass*
DX1
DY1
DZ1
RX1
RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

bna7 *t_options* ; for TUBE (NASTRAN)

where

t_options can be:

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OD *diameter*
WTH *thickness*
MASS *mass*
OD2 *diameter*
DX1
DY1
DZ1
RX1
RY1
RZ1
DX2
DY2
DZ2
RX2
RY2
RZ2

NE/NASTRAN Beams (in addition to all of the beam cross sections for NASTRAN)

bna8 *t_options* ; for CABLE (NE/NASTRAN)
where
t_options can be:
islack *slack*
itens *tension*
area *area*
iyz *moment*
ats *stress*

Remarks

Choose a positive integer for the cross section number. This number is used as reference in the **ibm**, **ibmi**, **jbm**, **jbmi**, **kbm**, **kbmi**, and bm commands. It is also used in reference in the definition of materials **lsdymats** and **dyanmats**.

For detailed information on definition of the cross-section see the manual of the specific simulation code.

If the output option has been selected prior to using the dialogue box to make a selection, only the options available to that output option will be displayed in the dialogue box.

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bsinfo write information about defined beam cross sections

bsinfo *bsd_#*

where

bsd_# is the one assigned to the definition in the **bsd** command.

bind Hughes-Liu beam user-defined integration points

bind *rule_#* s_1 t_1 w_1 s_2 t_2 w_2 s_3 t_3 w_3 ... ;

where

rule_# can be any positive integer used to refer to this rule followed by a list of local coordinates of integration points s_i and t_i and corresponding weights w_i .

Remarks

This command is used to define the integration points for the Hughes-Liu beam in **DYNA3D** and **LS-DYNA**. The coordinates s_i and t_i are parametric coordinates of integration points from interval <-1,1> (Denoted by crosses). The weights w_i are determined from the term:

$$w_i = A_i / A$$

where A_i is the area corresponding to the *i-th* integration node. A is the total area of the cross section determined by:

$$A = \sum A_i.$$

The t_t and s_t dimensions are used for scaling from parametric to real coordinates. The t_t and s_t dimensions are specified using the **bsd**, **bm**, **ibm**, **ibmi**, **jbm**, **jbmi**, **kbm**, **kbmi**, **dynamats**, or **lsdymats** commands.

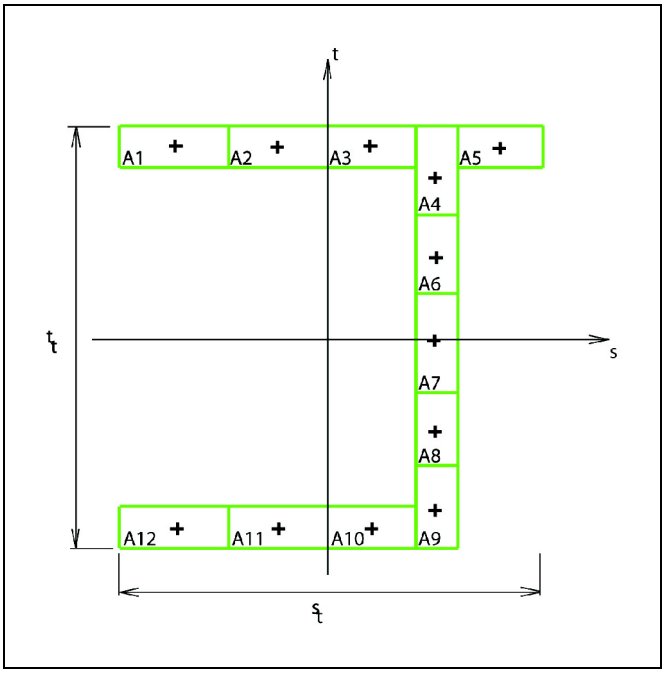

Figure 296 Cross Section with Integration Points

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lsbsd list the defined beam cross sections.

lsbsd

Remarks

Beam cross section properties are listed by number and type. Use **bsinfo** to get full information about a specific beam cross section definition.

offset add offset to numbered entities in the output

offset [*type offset*] ;

where *type* can be

nodes	node numbers
bricks	brick elements (or all elements)
shells	shell elements (if numbered independently)
beams	beam elements (if numbered independently)
tshells	thick shells
nsets	node sets if they are automatically numbered (not named)
fsets	face sets if they are automatically numbered (not named)
esets	element sets if automatically numbered (not named)
parts	parts
lsys	local coordinate systems
vecids	vector ID

Remarks

Only keyword outputs can use this number offset feature. For example, **Ls-dyna**, **Abaqus**, **Ansys**, **Marc**, **Nastran**, **NE/Nastran**, and **Neutral**.

Abaqus uses **nodes**, **bricks**, **nsetoff**, **esetoff**, **partoff,** and **nsetoff** affects the automatically numbered node sets as a result of the **fc**, **fd**, **fv**, **ft**, **acc**, and **mom** nodal boundary conditions. **Esetoff** affects the automatically numbered element sets as a result of the **pr** condition. **Partoff** affects the automatic numbering of element sets based on the part number.

Ansys uses the **bricks** and **nodes** offsets.

Lsdyna uses all of the offsets.

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Marc uses the **bricks** and **nodes** offsets.

Nastran and **NE/Nastran** uses the **bricks**, **nodes**, **nsetoff**, and **esetoff** offsets.

Neutral uses the **bricks** and **nodes** offsets.

If the output option has been selected prior to using the dialogue box to make a selection, only the options available to that output option will be displayed in the dialogue box.

sind shell user-defined integration rules

sind *rule* s_1 t_1 w_1 s_2 t_2 w_2 ... s_n t_n w_n ;

where

rule	is the integration rule number, 0 or 1. If 1, then all other arguments are ignored.
t_i w_i m_i	are local coordinate, integration weight, and material number

9. Sets

These commands are found in all phases. Named sets are a useful tool in the definition of boundary conditions and properties in the mesh and form an alternative to selecting regions in the Part phase. An arbitrary selection can be made and this is the advantage to using set functions in the Merge phase. The disadvantage is that the selection may no longer be parametric. If you go back and make a change to the mesh, you may have to redefine the set since the element and node numbers have changed.

The name of the set can be up to 8 alphanumeric characters long. Each name of the set must be unique.

In some of the set commands, the logical or Boolean set operators AND and OR are used to create new sets from existing sets. The AND operator between two sets means to take their intersection. This should not be confused with the common usage of and which might be interpreted to mean the addition of two sets. The OR operator does this function.

delset delete a set

delset *type set_name*

where *type* is the type of set which can be one of the following

node	node set
face	face set
element	element set

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where *set_name* is the name of the set.

Remarks

If a node set was constructed but is no longer needed, then it is best to delete it with this command. This can be important if an output file is going to be written which automatically writes all sets. When deleted, the set will not be written to the output file. In addition, it will not be using memory. This deletion has no affect on the nodes and elements of the mesh. All that is deleted is the list of nodes, faces, or elements that formed the set.

esetc element set comment

esetc *set_name text*

where

set_name	is the name of the element set
text	a text comment

Remarks

It is necessary to specify a comment using **esetc** for each node set to be written to ALE3D. This command can also be used to specify a title to an element set for LSDYNA.

esetinfo report the element set names and number of elements

esetinfo (no arguments)

Remarks

Command **esetinfo** reports the element set names and number of elements.

fsetc face set comment

fsetc *set_name text*

where

set_name	is the name of the face set
text	a text comment

Remarks

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It is necessary to specify a comment using **fsetc** for each node set to be written to ALE3D. This command can also be used to specify a title for a face set for LSDYNA.

fsetinfo report the face set names and number of faces

fsetinfo (no arguments)

Remarks

Command **fsetinfo** reports the face set names and number of faces.

nsetc attach a comment to a node set

nsetc *set_name text*

where

set_name	is the name of the node set
text	a text comment

Remarks

It is necessary to specify a comment using **nsetc** for each node to be written to ALE3D. This command can also be used to specify a title for a node set.

nsetinfo report the node set names and number of nodes

nsetinfo (no arguments)

Remarks

Command **nsetinfo** reports the node set names and number of nodes.

10. Coordinate Transformations and Part Replication

Coordinate transformations are used to translate, scale, reflect, and rotate an object from a local coordinate system to the global coordinate system. A local coordinate system is a frame of reference in which to generate a part. A local coordinate system is almost always a matter of convenience. The global coordinate system refers to the frame of reference used to create the complete model. In many

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cases, a part is generated in the global coordinate system, so there is no need to transform it. When you take advantage of symmetry by generating a section of the model, you will need some of the commands in this section to replicate the part.

The **lct** (local coordinate system) command should not be confused with the notion of a local coordinate system and the **gct** (global coordinate transformation) command should not be confused with the notion of a global coordinate system. This unfortunate naming cannot be repaired because such a change would cause many old input files to fail to generate the expected models.

In some cases, a component is duplicated many times. Each duplicate component must be transformed or placed into its proper location within the global coordinate system. The part duplication commands depend on the coordinate system transformations. Coordinate transformations can be defined in the **lct**, **gct**, and **lev** commands. Parts can be replicated with the **lrep**, **grep**, and **pslv** commands which refer back to the transformations defined in the **lct**, **gct**, and **lev** commands, respectively. The **lrep** and **grep** commands are discussed in the part phase section.

Properties are replicated along with the parts, such as material numbers and sliding interfaces (contact surface). Some of these properties are numbered and the number can be increased for each replication. These are replication increment commands found in this section.

In all cases, a coordinate transformation in **True*Grid*®** is composed of a sequence of basic operations. Each basic operation is given by a keyword possibly followed by some parameters. Each basic operation is performed in order from left to right. This ordering of the basic operations is sometimes referred to as a product or composition of basic operations. The composition of the basic operations is referred to as a single coordinate transformation.

It can be difficult to think of a complex transformation in three dimensions. You can simplify this by thinking of the object already in the global coordinate system. Then build the transformation, one operation at a time until you have moved, rotated, reflected, and scaled it to the proper position, orientation, and size.

Some coordinate transformations are applied to all parts, such as **csca**, **xcsa**, **ycsa**, **zcsa**, **xoff**, **yoff**, **zoff**, **exch**, and **gexch**. These transformations are referred to as final transformations because they are the last transformations that are applied to the parts. These commands are usually issued at the beginning of a command file to change dimensions.

There are four types of transformations that can be applied to a part. They are: local (**lct**), global (**gct**), levels (**lev**), and final transformations and they are applied in the order they are listed. Since the **lev** and the associated **pslv**/**pplv** commands allow for many nested levels of transformations, the ordering is based on the ordering of the nested levels. The inner nested level is performed first, with the outer level performed last. This ordering is like nested loops in a programming language.

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Element cross sectional properties, and in particular thicknesses, are not scaled by these transformations.

lct define local coordinate transformations

lct *n* $trans_1$ *; ... ;* $trans_n$ *;*

where $trans_i$ is a left-to-right product of the following basic operations:

mx *x_offset* to translate in the x direction
my *y_offset* to translate in the y direction
mz *z_offset* to translate in the z direction
v *x_offset y_offset z_offset* to translate by a vector
scv *xn yn zn offset* to translate a given distance along a vector
dv $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$ to translate by a vector given by two points
rx *theta* to rotate about the x axis
ry *theta* to rotate about the y axis
rz *theta* to rotate about the z axis
raxis *angle x0 y0 z0 xn yn zn* axis of rotation
rxy to reflect about the x-y plane
ryz to reflect about the y-z plane
rzx to reflect about the z-x plane
tf *origin x-axis y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface
ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*
where each of the arguments consist of a coordinate type followed by coordinate information:
rt *x y z* Cartesian coordinates
cy *rho theta z* cylindrical coordinates
sp *rho theta phi* spherical coordinates
pt *c.i* label of a labeled point from a 3D curve
pt *s.i.j* label of a labeled point from a surface
inv invert the present transformation)
csca *scale_factor* to scale all coordinates
xsca *scale_factor* to scale the x coordinate

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ysca *scale_factor*	to scale the y coordinate
zsca *scale_factor*	to scale the z coordinate
repe *#_repetitions*	to repeat powers of the current transformation
save *transform_#*	to apply a previous transformation
last	to use the last transformation

Remarks

This command defines a sequence of *n* local coordinate transformations; once you issue an **lct** command the results of any previous **lct** commands are no longer available. Each transformation is a product of primitive operators. The product is formed from the left to the right. One copy of the part is created and transformed for each transformation defined in the **lct** command and referenced in the **lrep** command. There is a limit of 500 transformations in the **lct** table of transformations.

You may wish to build a part in a convenient local coordinate system, and then transform the part to the proper position in the global coordinate system. The **lct** command, in conjunction with the **lrep** command, is used to move, scale, reflect, and rotate the part to put it in its place in the model relative to the other parts.

Examples

lct 3 **rz** 45 ; **rz** 45 **mz** 10 ; **rz** 45 **mz** 10 **mx** 25 ;

This command defines 3 local coordinate transformations numbered from 1 to 3. Once you have issued this command, you can no longer reference any previously defined local coordinate transformations.

The first transformation is a rotation about the z-axis. The second transformation is defined in two operations. First the object is rotated about the z-axis 45 degrees. Then the object is moved 10 units in the z-direction. The third transformation is defined by three operations. The first and second operations are the same as the second transformation. In addition, the third operation then moves the object in the x-direction 25 units. This command is equivalent to

lct 3 **rz** 45 ; **last mz** 10 ; **last mx** 25 ;

which is also equivalent to

lct 3 ; **rz** 45 ; **save** 1 **mz** 10 ; **save** 2 **mx** 25 ;

gct define global coordinate transformations

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gct *n trans$_1$; ... ; trans$_n$;*

where *trans$_i$* is one or more of:

mx *x_offset* to translate in the x direction

my *y_offset* to translate in the y direction

mz *z_offset* to translate in the z direction

v *x_offset y_offset z_offset* to translate by a vector

scv *xn yn zn offset* to translate a given distance along a vector

dv $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$ to translate by a vector given by two points

rx *theta* to rotate about the x axis

ry *theta* to rotate about the y axis

rz *theta* to rotate about the z axis

raxis *angle x0 y0 z0 xn yn zn* axis of rotation

rxy to reflect about the x-y plane

ryz to reflect about the y-z plane

rzx to reflect about the z-x plane

tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* Cartesian coordinates

cy *rho theta z* cylindrical coordinates

sp *rho theta phi* spherical coordinates

pt *c.i* label of a labeled point from a 3D curve

pt *s.i.j* label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z* Cartesian coordinates

cy *rho theta z* cylindrical coordinates

sp *rho theta phi* spherical coordinates

pt *c.i* label of a labeled point from a 3D curve

pt *s.i.j* label of a labeled point from a surface

inv invert the present transformation

csca *scale_factor* to scale all coordinates

xsca *scale_factor* to scale the x coordinate

ysca *scale_factor* to scale the y coordinate

zsca *scale_factor* to scale the z coordinate

repe *#_repetitions* to repeat powers of the current transformation

save *transform_#* to apply a previous transformation

last to use the last transformation

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Remarks

This command defines the entire sequence of global coordinate transformations; once you issue a **gct** command the results of any previous **gct** commands are no longer available. Each transformation is a product of primitive operators. The product is formed from the left to the right.. One copy of the part is created and transformed for each transformation defined in the **gct** command and referenced in the **grep** command. There is a limit of 500 transformations in the **gct** table of transformations.

You may wish to build a part in a convenient local coordinate system, and then transform the part to the proper position in the global coordinate system. The **gct** command, in conjunction with the **grep** command, is used to move, scale, reflect, and rotate the part to put it in its place in the model relative to the other parts.

Examples

```
gct 7 rxy; ryz; rzx; ryz rzx; rxy ryz; rxy rzx; rxy ryz rzx;
```

This set of transformations can be used to reflect a part from the first octant into all of the other 7 octants. The next example shows these transformations in use.

```
gct 7 rxy;ryz;rzx;ryz rzx;
rxy ryz;rxy rzx;rxy ryz rzx;
block 1 2 3;1 2 3 4;1 2 3 4;
     .25   .975  1
     -.2  -.025 .025 .22
     -.2   -.025 .025 .2
dei 1 2;1 2 0 3 4;;
dei 1 2;;1 2 0 3 4;
dei 2 3;1 2 0 3 4;1 2 0 3 4;
mb 2 2 2 3 3 3 x .25
tr 1 1 1 3 4 4 rz 45 ry -45;
grep 0 1 2 3 4 5 6 7;
endpart
```

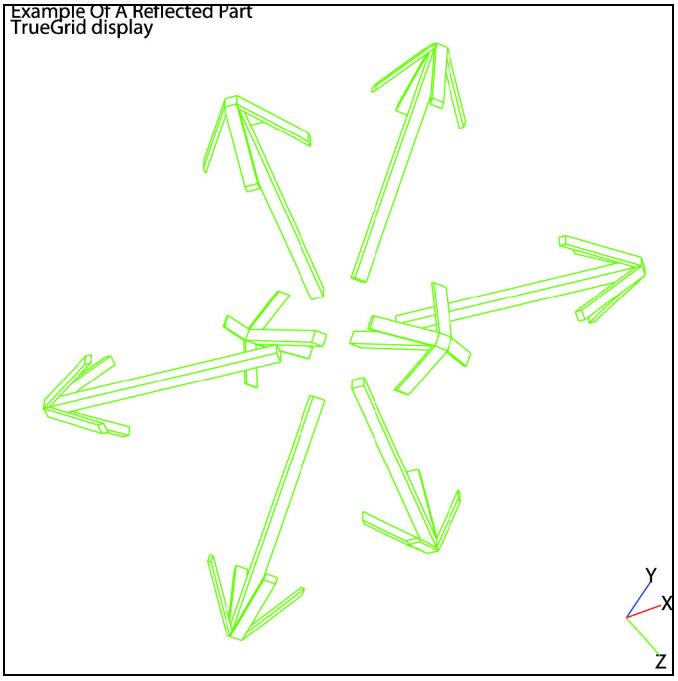

Figure 297 A Part Reflected Into 7 Octants

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lev define a set of transformations to replicate a set of parts

lev *level_# options ;*

level_# is a small positive number to uniquely identify the level

where an *option* can be any of:

grep *list_global_transform_# ;* a list of global coordinate transformations numbers (see the **gct** command)

add *level_#* include all transforms from another numbered **lev** command

prod *first_level_# second_level_#* to include a product of transforms of two numbered **lev** commands

levct *n* $trans_1$ *;* $trans_2$ *; ... ;* $trans_n$ *;*

where $trans_i$ is one or more of:

mx *x_offset*	to translate in the x direction
my *y_offset*	to translate in the y direction
mz *z_offset*	to translate in the z direction
v *x_offset y_offset z_offset*	to translate by a vector
scv *xn yn zn offset*	to translate a given distance along a vector
dv $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$	to translate by a vector given by two points
rx *theta*	to rotate about the x axis
ry *theta*	to rotate about the z axis
rz *theta*	to rotate about the z axis
raxis *angle x0 y0 z0 xn yn zn*	axis of rotation
rxy	to reflect about the x-y plane
ryz	to reflect about the y-z plane
rzx	to reflect about the z-x plane

tf *origin x-axis y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates
pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface

ftf *1st_origin 1st_x-axis 1st_y-axis 2nd_origin 2nd_x-axis 2nd_y-axis*

where each of the arguments consist of a coordinate type followed by coordinate information:

rt *x y z*	Cartesian coordinates
cy *rho theta z*	cylindrical coordinates
sp *rho theta phi*	spherical coordinates

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pt *c.i*	label of a labeled point from a 3D curve
pt *s.i.j*	label of a labeled point from a surface
inv	invert the present transformation
csca *scale_factor*	to scale all coordinates
xsca *scale_factor*	to scale the x coordinate
ysca *scale_factor*	to scale the y coordinate
zsca *scale_factor*	to scale the z coordinate
repe *#_repetitions*	to repeat powers of the current transformation
save *transform_#*	to apply a previous transformation
last	to use the last transformation

Remarks

This command defines a set of coordinate transformations associated with a level number. After defining a level, you can apply it to several parts at once, with the **pslv** and **pplv** commands. These two commands bracket the set of parts that the level is applied to. **True*Grid*®** will apply the level transformations to every part defined after a **pslv** and before the matching **pplv**. The way you first define and then apply a level is analogous to the way you first define a local or global coordinate transformation, **lct** or **gct**, and then apply them with **lrep** or **grep**.

You can define up to 20 levels of transformations. And each level can have an unlimited number of transformations. Once you define a level, you can reference it with **pslv** to replicate a set of parts. A level of transformations can be referenced by **pslv** commands many times. A level can be redefined.

The **levct** option defines a list of coordinate transformations in the same fashion as the **lct** and **gct** commands. That is, this option defines a transformation as a composition of a number of basic operations.

The **grep** option includes in the level a subset of previously defined global coordinate transformations. You identify them by their sequence numbers in the last **gct** command. A warning message is issued if a referenced global transformation was not previously defined by a **gct** command. Do not confuse this option with the part command **grep**. The **grep** command directly uses global transformations to replicate and transform a part. This **grep** option only uses global transformations to help define the coordinate transformations that constitute the level.

The **add** option will include in the present level definition all of the transformations associated with a previously defined level.

With the **prod** option, you can form the products of all transformations in two previously defined levels, and include all those products as transformations of your new level. This means that for every

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coordinate transformation in the first level and every coordinate transformation in the second level, **True*Grid*®** forms the product transformation and includes that in the new level's list of coordinate transformations. A product of two coordinate transformations is defined to be the result of first applying the coordinate transformation from the first level and then applying the coordinate transformation from the second level.

Examples

```
gct 3 mx 10 ; repe 3 ;
lev 1 levct 3 rx 30 ; rx 30 mz 10 ; rx 30 mz 10 my 10 ;;
lev 2 grep 0 2 3 ; ;
lev 3 add 2 levct 1 mx -10 ; grep 1 ; ;
lev 4 prod 1 3 ; ;
```

In the following example, a small part, in the shape of a sphere is created and then replicated using the **lev** command. The picture is below. The point to this example is that the second **lev** is nested in the first **lev** command using the **pslv** and **pplv** pair of commands, much like the way loop statements can be nested in a programming language.

```
title Example Of The LEV Command
gct 1 mx 4 my 4 rz 15;
lev 1 grep 0;
      levct 11 rz 30;repe 11;;
lev 2 grep 0 1;;
pslv 1
pslv 2
block 1 4;1 4;1 4;5 7 5 7 5 7
sfi -1 -2;-1 -2;-1 -2;sp 6 6 6 1
endpart
pplv
pplv
merge
```

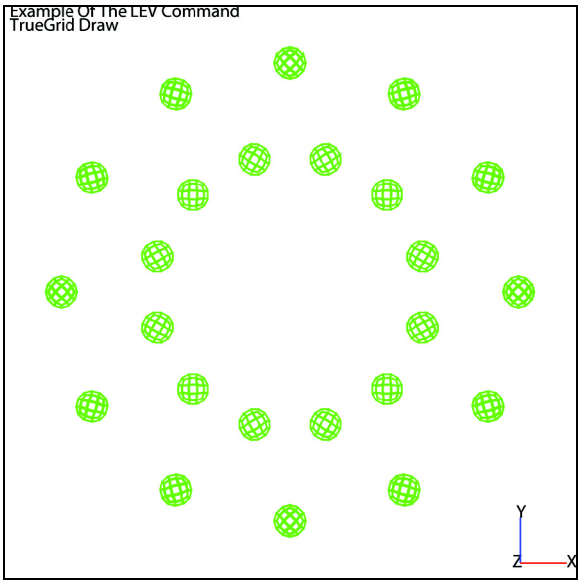

Figure 298 Nested LEV Replications

pslv begin replicating multiple parts

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pslv *level_#*

where *level_#* is the level definition number assigned in a **lev** command.

Remarks

This command is usually not issued interactively, although there is no compelling reason for this. It usually requires good planning, since it applies to multiple parts. Typically a set of parts are generated and the commands saved in a batch command file for regeneration. Then this command is inserted in the batch command file to produce the desired replications.

Before using the **pslv** and the related **pplv** commands, define a numbered set of transformations with the **lev** command.

The **pslv** and **pplv** commands will replicate a set of parts. Each use of the **pslv** command is paired with the use of a **pplv** command. The **pslv** marks the beginning and **pplv** marks the ending. All parts in between these two commands are replicated, one replication for each transformation found in the corresponding **lev** command. Neither the **pslv** nor the **pplv** command can be issued in the part phase.

For each **pslv** command there must follow somewhere a **pplv** command. The **pslv** and **pplv** pair of commands can be nested. When you nest one **pslv**/**pplv** pair within another, the effect is that of taking the product of the transformations in the two corresponding levels of transformations. A stacking technique is used to handle the products of replications. Each nesting of replications adds another level to the replication stack. The command **pslv** stands for "**pu**sh a **l**e**v**el onto the stack" and the command **pplv** stands for "**p**o**p** the top **l**e**v**el off of the stack".

If there are nested **pslv**/**pplv** commands, care is needed to determine the pairing of the **pslv** command with its corresponding **pplv** commands. This pairing is necessary to determine the scope of each **pslv** command (i.e. the parts that are replicated). The best way to determine which **pplv** is paired to which **pslv** command is to use a stacking technique. As you inspect the sequence of parts being generated, mentally place each **pslv** command encountered onto a stack. For each **pplv** that is encountered, pop the top **pslv** off of the stack which is then paired with the **pplv** command.

Pslv/**pplv**s pairs can be nested up to 20 deep. Local or global replication commands (**lrep** or **grep**) are also pushed onto and later popped off of the replication stack as a part is replicated. So when either of these commands are used, the total number of possible nested **pslv**/**pplv**s that can be used is reduced.

Examples

```
pslv 1
pslv 2
```

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```
c part number 1
   block ...
    ...
   endpart
  pplv   c   this pplv is paired with the pslv 2 command above
  pslv 3
c part number 2
   block ...
    ...
   endpart
  pplv   c   this pplv is paired with the pslv 3 command above
 pplv   c this pplv is paired with the pslv 1 command above
 pslv 4
c part number 3
  block ...
   ...
   grep 0 1 2 3 4 5 6;
   lrep 0 1 2 3 4 5;
  endpart
 pplv   c   this pplv is paired with the pslv 1 command above
```

In this example, the first part has two nested levels of part replications from levels numbered 1 and 2. The second part has two nested levels of part replications from levels numbered 1 and 3. The third part has three levels of part replications, one each from level number 4 and global and local coordinate transformations.

pplv end replicating multiple parts

pplv *(no arguments)*

Remarks

See the remarks on **pslv**. **Pplv** marks the end of the scope of part replications begun by a **pslv** command.

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csca scale all coordinates of all following parts

csca *scale*

Remarks

This command scales all parts that follow it.

xsca scale all x-coordinates of all following parts

xsca *scale*

Remarks

This command scales the x-coordinates of all parts that follow it.

ysca scale all y-coordinates of all following parts

ysca *scale*

Remarks

This command scales the y-coordinates of all parts that follow it.

zsca scale all z-coordinates of all following parts

zsca *scale*

Remarks

This command scales the z-coordinates of all parts that follow it.

xoff translate all x-coordinates of all following parts

xoff *offset*

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Remarks

This command translates the x-coordinates of all parts that follow it.

yoff translate all y-coordinates of all following parts

yoff *offset*

Remarks

This command translates the y-coordinates of all parts that follow it.

zoff translate all z-coordinates of all following parts

zoff *offset*

Remarks

This command translates the z-coordinates of all parts that follow it.

gexch permute the coordinates of all following parts

gexch *new_x_# new_y_#*
where

new_x_#	identifies the old coordinate which will become x: 1 for x, 2 for y, 3 for z
new_y_#	identifies the old coordinate which will become y: 1 for x, 2 for y, 3 for z

Remarks

See the remarks on **exch** below. **Gexch** is the same as **exch**. If both the **exch** and **gexch** are issued, **exch** is performed first.

exch permute the coordinates of all following parts

exch *new_x_# new_y_#*
where

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new_x_# identifies the old coordinate which will become x: 1 for x, 2 for y, 3 for z
new_y_# identifies the old coordinate which will become y: 1 for x, 2 for y, 3 for z

Remarks

This command permutes the x, y, and z coordinates. It applies to all parts defined after this command is issued.

A permutation of (1,2,3) is normally denoted by three numbers, e.g. (3,2,1). There are only two arguments to this command with the third number implied.

There is no way to undo this command except by issuing another **exch** command that applies the inverse permutation.

This permutation is applied before any permutations from **gexch**.

Examples

```
exch 2 1
```

For every part after this command is issued, the x and y coordinates will be interchanged.

```
exch 1 3
```

For every part after this command is issued, the y and z coordinates will be interchanged.

nerl rule used to number the nodes and brick elements

nerl *flag*

where the *flag* can be

0	for default numbering method
1	for ordered by increasing i-index last
-1	for ordered by decreasing i-index last
2	for ordered by increasing j-index last
-2	for ordered by decreasing j-index last
3	for ordered by increasing k-index last
-3	for ordered by decreasing k-index last

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Remarks

The default numbering is complicated to describe and is good for most problems. Nodes and brick elements can be numbered by specifying the last index so that the part can be split into layers. This way, the first sequence of nodes and elements will be found in the first layer, the second sequence of nodes and elements will be found in the second layer, and so on.

gmi material number increment for global replication

gmi *material_#_increment*

Remarks

This command sets the increment used to change the material numbers in a part that is being replicated using the **grep** command. For each replication of a part, the material numbers are increased from the previous replication. For example, if materials 1 and 6 are used in the creation of a part and if the global material increment is 3, then a replication of this part using **grep** would be assigned materials 4 and 9, respectively.

The default is 0.

Example

```
gmi 5 c set the material number increment for grep replications
lmi 1 c set the material number increment for lrep replications
block 1 2;1 2;1 2;1 2;1 2;1 2; c single element part
mate 1 c set the material number for the original part
lct 4 mx 1;repe 4; lrep 0:4;  c 5 local replications
gct 4 my 1;repe 4; grep 0:4;  c 5 global replications
endpart
```

In this example, 25 replications of a single element part are created with each element having a different material number. The material numbers are incremented according to the order of the transformations in the **lrep** and **grep** commands.

lmi material number increment for local replication

lmi *material_#_increment*
Remarks

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This command sets the increment used to change the material numbers in a part that is being replicated using the **lrep** command. For each replication of a part, the material numbers are increased from the previous replication. For example, if materials 1 and 6 are used in the creation of a part and if the local material increment is 1, then a replication of this part using **lrep** would be assigned materials 2 and 7, respectively.

The default is 0. See the example for the **gmi** command.

gsii sliding interface number increment for global replication

gsii *sliding_increment*

Remarks

This command sets the increment used to change the sliding interface numbers in a part that is being replicated using the **grep** command. For each replication of a part, the sliding interface numbers are increased from the previous replication. For example, if sliding interface numbers 1 and 6 are used in the creation of a part and if the global sliding interface increment is 3, then a replication of this part using **grep** would be assigned materials 4 and 9, respectively.

The default is 0.

Example

```
gsii 5
lsii 1
sid  1 sv; sid  2 sv; sid  3 sv; sid  4 sv; sid  5 sv;
sid  6 sv; sid  7 sv; sid  8 sv; sid  9 sv; sid 10 sv;
sid 11 sv; sid 12 sv; sid 13 sv; sid 14 sv; sid 15 sv;
sid 16 sv; sid 17 sv; sid 18 sv; sid 19 sv; sid 20 sv;
sid 21 sv; sid 22 sv; sid 23 sv; sid 24 sv; sid 25 sv;
block 1 2;1 2;1 2;1 2;1 2;1 2;
si 1 1 1 2 2 1 1 m
lct 4 mx 1;repe 4; lrep 0:4;
gct 4 my 1;repe 4; grep 0:4;
endpart
```

In this example, there are 25 replications. The first, with no transformation applied, has a sliding interface numbered 1. Every subsequent replication has a sliding interface that is one greater.

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lsii sliding interface number increment for local replication

lsii *slide_#_increment*

Remarks

This command sets the increment used to change the sliding interface numbers in a part that is being replicated using the **lrep** command. For each replication of a part, the sliding interface numbers are increased from the previous replication. For example, if sliding interface numbers 1 and 6 are used in the creation of a part and if the local sliding interface increment is 1, then a replication of this part using **lrep** would be assigned materials 2 and 7, respectively.

The default is 0. See the **gsii** command for an example.

11. Control Statements

These statements are much like the control statements in a programming language. They are useful when creating template command files to automate mesh generation in **True*Grid*®**. A set of parameters are assigned values prior to entering these commands. The control statements make it possible to execute a different set of commands in **True*Grid*®**, depending on the parameter values.

In order to remain backwards compatible, the **if/elseif/else/endif** group of commands are incompatible with the **when/elsewhen/else/endwhen**, **while/endwhile**, and the **for/endfor** commands. When using either of these command groups, use the **when/elsewhen/else/endwhen** command group when a conditional statement is required.

If you use a combination of the **while**, **for**, and **when** group of statements, their scopes must obey the embedding rule. The embedding rule is that the scope of one group of statements must be fully contained within the scope of the other group of statements, or the scopes of the two group of commands must be disjoint. The scope of the **if** group of statements must be disjoint from the scope of all of the **while**, **for**, and **when** group of statements.

For example, the following shows a **for**/**endfor** group of statements fully embedded within the **when**/**endwhen** group of statements:

```
when(%x.gt.1.0) then
  for idx 1 10 1
      .....
  endfor
endwhile
```

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The following is not allowed and has no meaning:

```
when(%x.gt.1.0) then
  for idx 1 10 1
    ...
endwhen
  endfor
```

break jump out of a for or while loop

break (no arguments)

Remarks

The **break** command is used to jump out of a **for** or **while** loop.

for iterate through a set of commands

for *p_name start end increment*

where

p_name	is the index variable name
start	is the starting index
end	is the ending index
increment	is the increment to be added to the index variable each iteration

Remarks

The **for** command is a loop similar in functionality to the DO statement in FORTRAN and the FOR statement in C. The **endfor** command flags the end of the **for** block. All commands between these two are repeated the designated number of times.

The **for** and **while** commands are similar in functionality and have the same combine limitations. They can be nested 20 deep. Every **for** and **endfor** statement should be on a separate line. Use the **break** command to jump out of a **for/endfor** loop. An **interrupt** inside a **for** loop will be ignored.

Example

```
curd 1 lp3
for idx 1 100 1
```

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```
   [sin(%idx*5)] [cos(%idx*6)] %idx
endfor
;;;
```

endfor terminate the for loop

endfor (no arguments)

Remarks

An **endfor** command is required to terminate a **for** loop. There must be one **endfor** statement **for** every **for** statement.

if begin an if... elseif ... else ... endif

if (*expression* **) then**

where *expression* can be relational or logical with the following features:

Integer, floating point, and exponential numbers and parameters can be used as operands and arguments to functions.

Unitary and binary arithmetic operators + and -.

Binary arithmetic operators *, /, **, and ^.

Binary relational operators:
.gt. and > for greater than,
.ge. and >= for greater than or equal,
.lt. and < for less than,
.le. and <= for less than or equal,
.eq. and == for equal,
.ne. and != for not equal.

Binary logical operators:
.and. and & and && for the logical *and* operator,
.or. for the logical *or* operator.
.eqv. for the logical *equivalent* operator,
.neqv. for the logical *not equivalent* (exclusive or) operator.

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Unitary logical operators .not. and ! .

Parenthesis to specify order of operations, where the default
order is given to exponentiation ** and ^ first,
then multiplication * and division / ,
addition + and subtraction - ,
relational operators, logical operators, and negation.

All calculations are done in floating point.
All illegal operations cause warnings.
All trigonometric operations are in degrees.
INT(x): truncates x to an integer.
NINT(x): rounds x to the nearest integer.
ABS(x): absolute value of x.
MOD(a,b): a modulo b.
SIGN(a,b): transfer the sign of b to a.
MAX(x1,x2,...,xn): maximum value of a list of numbers.
MIN(x1,x2,...,xn): minimum value of a list of numbers.
SQRT(x): square root of x where x most be positive.
EXP(x): exponential of x where x must not exceed 85.19.
LOG(x): natural logarithm where x must be positive.
LOG10(x): common logarithm base 10 where x must be positive.
SIN(x): trigonometric sine of x.
COS(x): trigonometric cosine of x.
TAN(x): trigonometric tangent of x where abs(x) can not be 90.
ASIN(x): trigonometric arcsine of x where abs(x) can not exceed 1.
ACOS(x): trigonometric arccosine of x where abs(x) can not exceed 1.
ATAN(x): trigonometric arctangent.
ATAN2(y,x): trigonometric arctangent with two arguments.
SINH(x): hyperbolic sine.
COSH(x): hyperbolic cosine.
RAND: uniform random number of unit length with default *mea*n 0 w/ forms
RAND,
RAND(seed),
RAND(seed, mean)
where *seed* is the random number seed, and
mean is the mean of the distribution.

NORM: normal random number w/ default standard deviation 1 & mean 0, w/ forms
NORM,
NORM(seed),
NORM(seed, mean),

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NORM(seed, mean, sig)
where *seed* is the random number seed,
mean is the mean of the distribution, and
sig is the standard deviation from the mean and must be positive.

Remarks

The **if** statement must be on a line by itself and it cannot be wrapped around to another line. This is like the block **if** statement in FORTRAN. Every **if** statement must be ended with an **endif**. If there is another **if** statement that follows before the **endif** statement, then the second **if** statement must be ended with an **endif** before the first one can be ended. This is called embedding of **if** statements. 19 embeddings of **if** statements are allowed. Optionally, multiple **elseif** statements can be used within the **if** statement. An **else** statement can optionally be used as the final option in the **if** statement. The **elseif** and **else** can also have up to 19 embedded **if** statements following it. The relational or logical expression found in the **if** statement is evaluated. This expression can have any of the features found in expressions. Relational and logical operators are also available. The order of precedence matches that of FORTRAN. If the expression is true, then all statements and commands found between the **if** and any associated **elseif**, **else**, or **endif** statement will be processed. This set of statements is referred to as the scope of the **if** statement. If the expression is false, these same statements and commands, those within the scope of the **if** statement, will be ignored. The **if** statement is a command and it cannot be used to select one set of arguments from another by embedding the **if** statement within another command.

When two numbers, which may be from the evaluations of expressions, are compared, possible roundoff errors are estimated and are used to determine equality.

This command is typically added to a session file and rerun. It is more difficult to plan a set of commands such that you can enter the **if-elseif-else-endif** statement interactively.

Example

```
if(%rebar.eq.1)then
 if(%k.ge.2*%k)then
  insprt 1 5 1 [%k]
  parameter k7 [%k1+1] k2 3;
  if(%sk.eq.1)then
   sd [%ns+1] plan %xx 0 [%zz+%k*%ze] [-sin(%p)] 0 [cos(%p)]
  elseif(%sk.le.0)then
   echo Error in parameter sk
   end
  else
```

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```
   sd [%ns+1] plan %xx 0 [%zz+%k*%ze] 0 1 0
  endif
  sfi ;;-2;sd [%ns+1]
 else
  parameter k 2;
 endif
endif
```

elseif add an option to an if statement

elseif (*expression* **) then**

where *expression* is defined in the **if** statement above.

Remarks

The **elseif** statement must be on a line by itself and it cannot be wrapped around to another line. This is like the block **elseif** statement in FORTRAN. Every **elseif** belongs to an **if** statement which must be ended with an **endif.** If there is another **if** statement that follows before the **endif** statement, then that **if** statement must be ended with an **endif** before the first **if**..**elseif** one can be ended. This is called embedding of **if** statements. **True*Grid*®** allows for 19 embeddings of **if** statements. Optionally, multiple **elseif** statements can be used within the **if** statement. An **else** statement can optionally be used as the final option in the **if** statement. The **elseif** and **else** can also have up to 19 embedded **if** statements following it. The relational or logical expression found in the **elseif** statement is evaluated. This expression can have any of the features found in expressions. Relational and logical operators are also available. The order of precedence matches that of FORTRAN. If the preceding associated **if** statement expression is false and all preceding associated **elseif** statement expressions are also false, then the expression in this **elseif** statement is evaluated. If it is true, then all statements and commands found between the **elseif** and any subsequent associated **elseif**, **else**, or **endif** statement will be processed. The statements and commands found between this **elseif** statement and the next associated **elseif**, **else**, or **endif** statement are referred to as the scope of the **elseif** statement. If the expression is false, these same statements and commands, the commands in the scope of this **elseif** statement, will be ignored.

When two numbers, which may be from the evaluations of expressions, are compared, possible roundoff errors are estimated and are used to determine equality.

This command is typically added to a session file and rerun. It is more difficult to plan a set of commands such that you can enter the **if-elseif-else-endif** statement interactively.

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else final option in an if or when statement

else

Remarks

The **else** statement must be on a line by itself. This is like the block **else** statement in FORTRAN. Every **else** belongs to an **if** or **when** statement which must be ended with an **endif** or **endwhen**, respectively. If there is another **if** or **when** statement that follows before the **endif** or **endwhen** statement, then that **if** or **when** statement must be ended with an **endif** or endwhen before the first **if** or **when** statement is ended. This is called embedding of **if** or **when** statements. **True*Grid*®** allows for 19 embeddings of **if** and **when** statements. An **else** statement can optionally be used as the final option in the **if** or **when** statement. The **else** can also have up to 19 embedded **if** or **when** statements following it.

This command is typically added to a session file and rerun. It is more difficult to plan a set of commands such that you can enter the **if-elseif-else-endif** or **when-elsewhen-else-endwhen** statement interactively.

endif end an if statement

endif

Remarks

The **endif** statement must be on a line by itself. This is like the block **if**..**elseif**..**else**..**endif** statement in FORTRAN. Every **endif** ends an **if** statement.

endwhile end a while statement

endwhile

Remarks

The **endwhile** statement must be on a line by itself. Every **endwhile** ends or closes a **while** statement.

when begin a when... elsewhen ... else ... endwhen

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when (*expression* **) then**

where *expression* can be relational or logical with the following features:

Integer, floating point, and exponential numbers and parameters can be used as operands and arguments to functions.

Unitary and binary arithmetic operators + and -.

Binary arithmetic operators *, /, **, and ^.

Binary relational operators:
.gt. and > for greater than,
.ge. and >= for greater than or equal,
.lt. and < for less than,
.le. and <= for less than or equal,
.eq. and == for equal,
.ne. and != for not equal.

Binary logical operators:
.and. and & and && for the logical *and* operator,
.or. for the logical *or* operator.
.eqv. for the logical *equivalent* operator,
.neqv. for the logical *not equivalent* (exclusive or) operator.

Unitary logical operators .not. and ! .

Parenthesis to specify order of operations, where the default
order is given to exponentiation ** and ^ first,
then multiplication * and division / ,
addition + and subtraction - ,
relational operators, logical operators, and negation.

All calculations are done in floating point.
All illegal operations cause warnings.
All trigonometric operations are in degrees.
INT(x): truncates x to an integer.
NINT(x): rounds x to the nearest integer.
ABS(x): absolute value of x.
MOD(a,b): a modulo b.
SIGN(a,b): transfer the sign of b to a.
MAX(x1,x2,...,xn): maximum value of a list of numbers.
MIN(x1,x2,...,xn): minimum value of a list of numbers.

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SQRT(x): square root of x where x most be positive.
EXP(x): exponential of x where x must not exceed 85.19.
LOG(x): natural logarithm where x must be positive.
LOG10(x): common logarithm base 10 where x must be positive.
SIN(x): trigonometric sine of x.
COS(x): trigonometric cosine of x.
TAN(x): trigonometric tangent of x where abs(x) can not be 90.
ASIN(x): trigonometric arcsine of x where abs(x) can not exceed 1.
ACOS(x): trigonometric arccosine of x where abs(x) can not exceed 1.
ATAN(x): trigonometric arctangent.
ATAN2(y,x): trigonometric arctangent with two arguments.
SINH(x): hyperbolic sine.
COSH(x): hyperbolic cosine.
RAND: uniform random number of unit length with default *mea*n 0 w/ forms
RAND,
RAND(seed),
RAND(seed, mean)
where *seed* is the random number seed, and
mean is the mean of the distribution.

NORM: normal random number w/ default standard deviation 1 & mean 0, w/ forms
NORM,
NORM(seed),
NORM(seed, mean),
NORM(seed, mean, sig)
where *seed* is the random number seed,
mean is the mean of the distribution, and
sig is the standard deviation from the mean and must be positive.

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Remarks

The **when** statement must be on a line by itself and it cannot be wrapped around to another line. This is like the block **if** statement in FORTRAN. Every **when** statement must be ended with an **endwhen**. If there is another **when** statement that follows before the **endwhen** statement, then the second **when** statement must be ended with an **endif** before the first one can be ended. This is called embedding of **when** statements. 19 embeddings of **when** statements are allowed. Optionally, multiple **elsewhen** statements can be used within the **when** statement. An **else** statement can optionally be used as the final option in the **when** statement. The **elsewhen** and **else** can also have up to 19 embedded **when** statements following it. The relational or logical expression found in the **when** statement is evaluated. This expression can have any of the features found in expressions. Relational and logical operators are also available. The order of precedence matches that of FORTRAN. If the expression is true, then all statements and commands found between the **when** and any associated **elsewhen**, **else**, or **endwhen** statement will be processed. This set of statements is referred to as the scope of the **when** statement. If the expression is false, these same statements and commands, those within the scope of the **when** statement, will be ignored. The **when** statement is a preprocessed command and it can be used to select one set of arguments from another by embedding the **when** statement within another command. For example:

```
block
  when(%x1.gt.%x2) then
    1 3;1 3;1 3;
  else
    1 2;1 2;1 2;
  endwhen
 1 2 1 2 1 2
```

This feature of the **when**, **while**, and **for** group of statements is the reason these commands are preprocessed and incompatible with the **if** group of statements.

When two numbers, which may be from the evaluations of expressions, are compared, possible roundoff errors are estimated and are used to determine equality.

This command is typically added to a session file and rerun. It is more difficult to plan a set of commands such that you can enter the **when-elsewhen-else-endwhen** statements interactively.

Example

```
when(%rebar.eq.1)then
 when(%k.ge.2*%k)then
  insprt 1 5 1 [%k]
```

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```
  parameter k7 [%k1+1] k2 3;
 when(%sk.eq.1)then
   sd [%ns+1] plan %xx 0 [%zz+%k*%ze] [-sin(%p)] 0 [cos(%p)]
  elsewhen(%sk.le.0)then
   echo Error in parameter sk
   end
  else
   sd [%ns+1] plan %xx 0 [%zz+%k*%ze] 0 1 0
  endwhen
  sfi ;;-2;sd [%ns+1]
 else
  parameter k 2;
 endwhen
endwhen
```

elsewhen add an option to a when statement

elsewhen (*expression* **) then**

where *expression* is defined in the **when** statement above.

Remarks

The **elsewhen** statement must be on a line by itself and it cannot be wrapped around to another line. This is like the block **elseif** statement in FORTRAN. Every **elsewhen** belongs to a **when** statement which must be ended with an **endwhen.** If there is another **when** statement that follows before the **endwhen** statement, then that **when** statement must be ended with an **endwhen** before the first **when**..**elsewhen** one can be ended. This is called embedding of **when** statements. **True*Grid*®** allows for 19 embeddings of **when** statements. Optionally, multiple **elsewhen** statements can be used within the **when** statement. An **else** statement can optionally be used as the final option in the **when** statement. The **elsewhen** and **else** can also have up to 19 embedded **when** statements following it. The relational or logical expression found in the **elsewhen** statement is evaluated. This expression can have any of the features found in expressions. Relational and logical operators are also available. The order of precedence matches that of FORTRAN. If the preceding associated **when** statement expression is false and all preceding associated **elsewhen** statement expressions are also false, then the expression in this **elsewhen** statement is evaluated. If it is true, then all statements and commands found between the **elsewhen** and any subsequent associated **elsewhen**, **else**, or **endwhen** statement will be processed. The statements and commands found between this **elsewhen** statement and the next associated **elsewhen**, **else**, or **endwhen** statement are referred to as the scope of the **elsewhen** statement. If the expression is false, these same statements and commands, the commands in the scope of this **elsewhen** statement, will be ignored.

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When two numbers, which may be from the evaluations of expressions, are compared, possible roundoff errors are estimated and are used to determine equality.

This command is typically added to a session file and rerun. It is more difficult to plan a set of commands such that you can enter the **when-elsewhen-else-endwhen** statement interactively.

endwhen end a when statement

endwhen (no arguments)

Remarks

The **endwhen** statement must be on a line by itself. This is like the block **if**..**elseif**..**else**..**endif** statement in FORTRAN. Every **endwhen** ends a **when** statement.

while begin a loop iteration

while (*expression* **)**

where *expression* can be relational or logical with the following features:

Integer, floating point, and exponential numbers and parameters can be used as operands and arguments to functions.

Unitary and binary arithmetic operators + and -.

Binary arithmetic operators *, /, **, and ^.

Binary relational operators:
.gt. and > for greater than,
.ge. and >= for greater than or equal,
.lt. and < for less than,
.le. and <= for less than or equal,
.eq. and == for equal,
.ne. and != for not equal.

Binary logical operators:
.and. and & and && for the logical *and* operator,
.or. for the logical *or* operator.
.eqv. for the logical *equivalent* operator,

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.neqv. for the logical *not equivalent* (exclusive or) operator.

Unitary logical operators .not. and ! .

Parenthesis to specify order of operations, where the default
order is given to exponentiation ** and ^ first,
then multiplication * and division / ,
addition + and subtraction - ,
relational operators, logical operators, and negation.

All calculations are done in floating point.
All illegal operations cause warnings.
All trigonometric operations are in degrees.
INT(x): truncates x to an integer.
NINT(x): rounds x to the nearest integer.
ABS(x): absolute value of x.
MOD(a,b): a modulo b.
SIGN(a,b): transfer the sign of b to a.
MAX(x1,x2,...,xn): maximum value of a list of numbers.
MIN(x1,x2,...,xn): minimum value of a list of numbers.
SQRT(x): square root of x where x most be positive.
EXP(x): exponential of x where x must not exceed 85.19.
LOG(x): natural logarithm where x must be positive.
LOG10(x): common logarithm base 10 where x must be positive.
SIN(x): trigonometric sine of x.
COS(x): trigonometric cosine of x.
TAN(x): trigonometric tangent of x where abs(x) can not be 90.
ASIN(x): trigonometric arcsine of x where abs(x) can not exceed 1.
ACOS(x): trigonometric arccosine of x where abs(x) can not exceed 1.
ATAN(x): trigonometric arctangent.
ATAN2(y,x): trigonometric arctangent with two arguments.
SINH(x): hyperbolic sine.
COSH(x): hyperbolic cosine.
RAND: uniform random number of unit length and default *mean* 0 /w forms
RAND,
RAND(seed),
RAND(seed,mean)
where *seed* is the random number seed, and
mean is the mean of the distribution.

NORM: normal random number w/ default standard deviation 1 & mean 0, w/ forms
NORM,

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NORM(seed),
NORM(seed,mean),
NORM(seed,mean,sig)
where *seed* is the random number seed,
mean is the mean of the distribution, and
sig is the standard deviation from the mean and must be positive.

Remarks

The **while** statement must be on a line by itself and it cannot be wrapped around to another line. Every **while** statement must be ended with a closing **endwhile** statement. The statements and commands that fall between the **while** and closing **endwhile** statements is referred to as the scope of the command. If there is another **while** statement that follows before the closing **endwhile** statement, then the second **while** statement must be ended with an **endwhile** before the first one can be ended. This is called embedding of **while** statements. 19 embeddings of **while** statements are allowed. The relational or logical expression found in the **while** statement is evaluated. This expression can have any of the features found in expressions. Relational and logical operators are also available. The order of precedence matches that of FORTRAN. If the expression is true, then all statements and commands found between the **while** and **endwhile** statement will be processed and, upon completion, this whole process is repeated until the expression is false. When the expression becomes false, these same statements and commands will be ignored and the next statement or command to be executed will be the one that follows the closing **endwhile** statement. Use the **break** feature to jump out of a while statement. The **while** statement is a command and it cannot be embedded within another command.

Use the **break** command to jump out of a **while/endwhile** loop. An **interrupt** inside a **while** loop will be ignored.

When two numbers, which may be from the evaluations of expressions, are compared, possible roundoff errors are estimated and are used to determine equality.

This command is typically added to a session file and rerun. It is more difficult to plan a set of commands such that you can enter the **while** statement interactively.

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Example

```
para n 0;
while(%n.lt.10)
 block 1 2;1 2;1 2;0 1 0 1 0 1
 tri ;;; v %n %n %n;
 endpart
 para n [%n+1];
endwhile
```

This example produces 10 unit brick elements along a diagonal.

12. Merging Nodes

In the Merge phase, nodes that are close to one another are merged into a single node. Tolerance commands allow you to define how close is close. All tolerances are in absolute distances. There are commands for specifying tolerances for the general merging of all nodes over all parts or just nodes on the exterior faces of the mesh. There are commands for specifying the tolerances for the special merging of nodes between parts or within a part. These special tolerances override the general ones. If no tolerance commands are specified, then no merging is done. However, the Merge phase must be entered in order to build the node map which is used to generate output for a simulation code.

Invocation of a tolerance command (**t**, **tp**, **st**, **stp**) within the Merge phase causes an immediate merging of nodes. These commands can also be invoked any time; when the Merge phase is entered, those tolerance commands are immediately executed or re-executed as the case may be. The merge process is always performed on the nodes in their original (prior to any merging) state. Merging is not cumulative. If you leave the Merge phase and reenter it, all merging is recalculated with what ever new parts that have been added. This lets you interactively experiment with merging and tolerances. Setting a tolerance to a negative value is an easy way to restore the nodes to their original states. Graphical displays of the mesh in the Merge phase always reflect the results of any merging.

Nodes are merged depending on the distance between them. If a node lies within a tolerance distance of more than one other node, then it is merged with the closest one. When merging several nodes into one node, the first-defined node survives. This can be overridden by the **bptol** command. Nodes within a joint, spring, or weld spot and across the two sides of a sliding surface are not merged. When the first merging of nodes occurs, a sliding interface table is calculated which is used in the merging process. This table is written to the screen and to the save file and is intended as diagnostics. The following is a sample of that table:

```
SLIDING INTERFACE SUMMARY
```

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Surf	S-node	S-lseg	S-qseg	M-node	M-lseg	M-qseg
1	105	84	0	468	418	0
2	232	0	52	468	418	0
3	221	0	0	390	304	0
4	221	0	0	390	304	0
5	158	0	0	120	88	0
6	158	30	30	120	88	0
7	204	102	0	204	102	0
8	232	0	52	90	52	0
9	101	18	18	161	132	0
10	101	18	18	161	132	0
11	548	120	120	3216	0	1056
12	133	84	0	161	132	0
13	133	84	0	161	132	0
14	308	240	0	3216	0	1056

This table is organized by the sliding interface number on the right. Columns 2, 3, and 4 are datum pertaining to the slave side on the interface; columns 5, 6, and 7 to the master side. Columns 2 and 5 (`S-node` and `M-node`) are node counts. Columns 3 and 6 (`S-lseg` and `M-lseg`) are linear face counts, and columns 4 and 7 (`S-qseg` and `M-qseg`) are quadratic face counts.

A table of merged nodes is always written after the **tp** or **stp** commands are executed.

```
                MERGED NODES SUMMARY
  12 nodes merged between parts        1 and        2
  16 nodes merged between parts        2 and        2
  12 nodes merged between parts        1 and        3
  16 nodes merged between parts        3 and        3
 216 nodes merged between parts        4 and        4
  30 nodes merged between parts        7 and        7
  88 nodes merged between parts        8 and        8
 390 nodes were deleted by tolerancing
```

Up through 4000 parts can be merged under general tolerancing (i.e. no use of the **ptol** or **bptol** commands). 300 parts can be merged under special tolerancing (**ptol** and **bptol**).

The following is a common error to avoid. Suppose you create three parts that meet as shown in the figures below. Then define a sliding interface between parts 1 and 2 and also between parts 1 and 3. No nodes will be merged between parts 1 and 2 and between parts 1 and 3. However, nodes can be merged between parts 2 and three. Sometimes you need to look closely in the graphics or carefully check the Merged Nodes Summary to detect this error. To fix this error, if indeed it is an error, use a dummy sliding interface between parts 2 and 3 to force no merging between those parts.

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Alternatively, use the **bptol** command with a negative number to avoid merging between those parts. You should also consider extending both interfaces 1 and 2 across to parts 3 and 2, respectively, because they may come in contact. This is an ambiguous situation since there are equally plausible situations where parts 2 and 3 should be merged together.

```
sid 1 sv;
sid 2 sv;
block 1 3;1 3;1 3;
     1 2 1 2 1 2
sii -2;;;1 s;
sii ;-2;;2 s;
block 1 3;1 3;1 3;
     2.1 3 1 2 1 2
sii -1;;;1 m;
block 1 3;1 3;1 3;
     1 2 2.1 3 1 2
sii ;-1;;2 m;
merge
stp .2
```

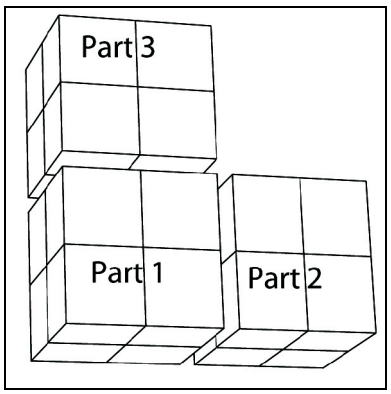

Figure 299 Before **stp**

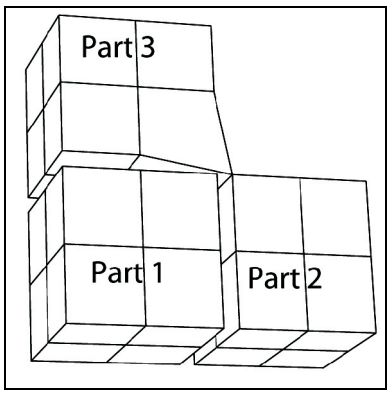

Figure 300 After **stp**

bnstol between node set tolerance

bnstol *node_set*$_1$ *node_set*$_2$ *tolerance*

where

node_set$_1$	first set of nodes to be compared
node_set$_2$	second set of nodes to be compared
tolerance	tolerance that must be met for merging between node

Remarks

A pair of nodes may be merged into one node based on this new command. If one node of the pair is in one of these two node sets and the other node is in the other node set, they will be merged into one node if the distance between them is less than the *tolerance*. This is checked for all possible pairs of nodes. This command works similarly to the **ptol** and **bptol** commands. These commands only set the tolerance to be used when a merge command, such as **stp** or **tp**, is issued.

It is possible with the **ptol**, **bptol**, and the **bnstol** commands for a pair of nodes to have several tolerances, making the merging process ambiguous. To resolve this issue, when the merging of a pair

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of nodes are ruled by the issuing of several of these commands, then the one that was issued last will be the command to determine if the pair of nodes are to be merged. The key to this is that the order that these commands are issued can be important.

For example, one may wish to merge nodes in a specific region and rule out merging in all other regions. This can be done by creating two node sets, each set containing the nodes on opposite faces to be merged and issuing the **bnstol** command. Then an **stp** command with a tolerance of -1 could be issued to rule out all merging except for the region of the two node sets.

This feature has the opposite effect of the sliding interface (**sid** command) of type dummy.

merge switch to the merge (assembly) phase

merge *(no arguments)*

Remarks

This command does not work in the merge phase since you will already be in the merge phase. If this command is issued while in the part phase, it will cause the part to end. This command is useful in the control phase to switch to the merge phase where there is a 3D graphics window.

mns merge node sets

mns *mns# node_set_1 node_set_2*

where

mns#	mns identification number
node_set_1	name of the first node set
node_set_2	name of the second node set

Remarks

Two node sets are specified and any pair of nodes, one in each set, are allowed to be merged if they are found on opposite sides of a sliding interface. A merge command in the Merge phase is still required for the nodes to be merged, and only those nodes within the specified distance between them will be merged. Up to 15 pairs of node sets are allowed.

This feature can be undone by using a ";" for one of the sets in order to prematurely end the command.

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Example

Two shell parts are generated with a small gap between them. The first part is assigned to be the master side of a sliding interface, with the faces oriented towards the second part. The second part plays the opposite role as the slave side, facing the first side. Two node sets are created, one from the first part along an edge, and the second node set along the corresponding edge of the second part. These nodes along the corresponding edges are allowed to merge.

```
sid 1 sv;
block 1 11;1 11;-1;-1 1 -1 1 0
orpt + 0 0 1
si 1 1 1 2 2 1 1 m;
nset 1 1 1 1 2 1 = ifc1
block 1 11;1 11;-1;-1 1 -1 1 .001
orpt + 0 0 -1
si 1 1 1 2 2 1 1 s;
nset 1 1 1 1 2 1 = ifc2
merge
mns 1 ifc1 ifc2
stp .0011
```

st set tolerance and merge surface nodes

st *tolerance*

Remarks

Only nodes that lie on the surfaces or exterior faces of parts are considered for merging. Exterior faces include faces of the mesh that are physically matching but are logically distinct faces of the mesh. No table of results is printed. If the sliding interface table has not yet been printed, it will be printed at this time.

stp set tolerance and merge surface nodes, with diagnostics

stp *tolerance*

Remarks

This command performs as **st** except that the results of the merge are reported. Exterior faces include

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faces of the mesh that are physically matching but are logically distinct faces of the mesh. A table of the merged nodes is printed to the screen and to the save file. If there are any sliding interfaces, then the first merge command will cause a sliding interface report to be written first.

Example

In the example above for the **mns** command, the **stp** command would cause the following tables to be written.

```
stp .002
                SLIDING INTERFACE SUMMARY
   Surf   S-node   S-lseg   S-qseg   M-node   M-lseg   M-qseg
      1      121      100        0      121      100        0
                  MERGED NODES SUMMARY
         11 nodes merged between parts        1 and        2
         11 nodes were deleted by tolerancing
```

t set tolerance and merge nodes

t *tolerance*

Remarks

All nodes of all parts are considered for merging. If the sliding interface table has not yet been printed, it will be printed at this time.

tp set tolerance and merge nodes, with diagnostics

tp *tolerance*

Remarks

This command performs as **t** except that the results of the merge are reported. See the **stp** command for an example. If the sliding interface table has not yet been printed, it will be printed at this time.

ztol minimum non-zero absolute coordinate

ztol *tolerance*

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Remarks

This command causes each coordinate of each node, whose absolute value is less than *tolerance*, to be set to zero prior to merging and prior to the generation of output. Each invocation of **ztol** causes the **original** coordinates to be compared against *tolerance*. This lets you interactively experiment with **ztol**. This command does not change the data base. It only affects the merging process and the output.

When a node is picked in the merge phase, the environment window displays the coordinates. **Ztol** affects these coordinates displayed in the pick panel.

bptol between parts tolerance specification

bptol *part$_1$ part$_2$ tolerance*

where

part$_1$ and *part$_2$*	are part numbers, and
tolerance	is the between part tolerance.

Remarks

The *tolerance* is used as the between-part tolerance for merging nodes between parts *part$_1$* and *part$_2$*. This command does not initiate the merge procedure. The *tolerance* overrides the default tolerance specified by the **t**, **tp**, **st**, and **stp** tolerance commands just between these two parts. When nodes are merged, the nodes belonging to part *part$_1$* survive. This is also a way to change the ordering of the merging procedure. Use a negative number to avoid merging between two parts. There is no limit to the number of **bptol** and **ptol** commands.

This command can be easily abused. Most meshes do not need this command to force the merging. If the nodes between parts do not merge with a reasonable tolerance using, for example, the **stp** command, then perhaps the interface between the two parts were not generated identically. This may warrant some investigation into the cause. Many of these problems can be avoided by using the **bb** command.

ptol part tolerance specification

ptol *part_number tolerance*

Remarks

The *tolerance* is used as the absolute tolerance for merging nodes within part *part_number*. This

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command does not initiate the merge procedure. The *tolerance* overrides the default tolerance specified by the **t**, **tp**, **st**, and **stp** tolerance commands just within this part. Use a negative number to avoid merging within the part. There is a maximum of 1000 **bptol** and **ptol** commands.
This command can be easily abused. Most meshes do not need this command to force the merging. If the nodes within a part do not merge with a reasonable tolerance using, for example, the **stp** command, then perhaps the interfaces within the part were not generated identically. This may warrant some investigation into the cause.

rigbm identify two rigid bodies to be merged

rigbm $material_1$ $material_2$
where
$material_1$ $material_2$ must be rigid materials in DYNA3D or LS-DYNA.

13. Global Graphics Commands

The following commands are available at any time.

ibzone control the computational window frame

ibzone *option*
where the *option* can be
on *size* turn on the frame and set the size of the index zone
off turn off the frame

Remarks

This command activates or deactivates the drawing of the boundary between part and index-bar on the computational window. The outer part of the window is the zone of influence of the index-bar. This is where the mouse can move to cause highlighting of the mesh in both the physical and computational window. This is also the area where a mouse click or a click-and-drag can make selections in the mesh. The inner box of the zone becomes the area where the computational window can be seen and where a click-and-drag can select a region directly from the picture, instead of from the index bar. One can also change the size of the zone. The default is **on 1.0**.

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noplot turn all graphics off

noplot *(no arguments)*

Remarks

This command is useful when you have a batch file with some graphics commands. When running this batch file, you may wish to suppress the graphics. It will speed the process. If you wish to avoid all graphics including the creation of the graphics windows, use the **nogui** option on the execute line. If you use the **nogui** option, the **noplot** command is not needed to suppress the graphics.

plot turn graphics back on

plot *(no arguments)*

Remarks

This command re-activates the graphics after issuing the **noplot** command. Graphics are active by default.

14. Miscellaneous Commands

apainfo list all automatic parameters

apainfo (no arguements)

Remarks

Apainfo lists all of the automatic parameters and their values. Below is an example. Note that the automatic parameters associated with the part phase (the first 6 below) will only appear in the part phase. The same information can be gotten, one automatic parameter at a time, using the dc command (desk calculator).

Examples

```
List of full i-indices of the part
     idxlist =    1    3    5    7
List of full j-indices of the part
     jdxlist =    1    3    5    7    9   11   13
List of full k-indices of the part
     kdxlist =    1    3    5    7    9   11   13   15   17
```

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```
                19   21   23   25   27   29   31   33   35
                37   39
Maximun reduced index in the i-direction
     maxrudi =       4
Maximun reduced index in the j-direction
     maxrudj =       7
Maximun reduced index in the k-direction
     maxrudk =      20
Next surface number
     nextsrf =       1
Next 3D curve number
     nextcrv =       1
Next load curve number
     nextlc  =       1
Next 2D curve number
     nextln  =       1
Next material number
     nextmat =       2
Next Block Boundary ID number from BB command
     nextbb  =       1
Next part number
     nextprt =       2
Node number result from ajnp/sajnp and PICK>NODE functions
     node    =       0
Element number result from ajel/sagel functions
     element =       0
The functions ipil, trapt, tricent,
bulc, circent, project, ptcor, pptcd, and nodcor/snodcor
record their results in the following three parameters:
X-component of projection
     xprj    = 0.000000E+00
Y-component of projection
     yprj    = 0.000000E+00
Z-component of projection
     zprj    = 0.000000E+00
The functions circent, sp4pt, and project
record their results in the following three parameters:
X-component of a normal
     xnrm    = 0.000000E+00
Y-component of a normal
     ynrm    = 0.000000E+00
Z-component of a normal
     znrm    = 0.000000E+00
```

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```
X-component from cross product crprod function
     xcrprod = 0.000000E+00
Y-component from cross product crprod function
     ycrprod = 0.000000E+00
Z-component from cross product crprod function
     zcrprod = 0.000000E+00
Inner product from inprod function
     inprod  = 0.000000E+00
Distance result from distance, circent, and sp4pt functions
     distance= 0.000000E+00
Subtended angle result from subang function
     subang  = 0.000000E+00
Minimum mesh quality measure
     minmea  = 0.000000E+00
Maximum mesh quality measure
     maxmea  = 0.000000E+00
Mean of mesh quality measure
     mean    = 0.000000E+00
Standard deviation of mesh quality measure
     standev = 0.000000E+00
Irrational constant circle circumference/diameter
     pi      = 3.141593E+00
Minimum x-coordinate of the graphics containing box
     xboxmin = 1.000000E+00
Minimum y-coordinate of the graphics containing box
     yboxmin = 1.000000E+00
Minimum z-coordinate of the graphics containing box
     zboxmin = 1.000000E+00
Maximum x-coordinate of the graphics containing box
     xboxmax = 7.000000E+00
Maximum y-coordinate of the graphics containing box
     yboxmax = 1.300000E+01
Maximum z-coordinate of the graphics containing box
     zboxmax = 3.900000E+01
Next node number (befor merging)
     nextnode=       1
Next linear brick
     nextlbrick=       1
Next linear shell element
     nextlshell=       1
Next linear beam
     nextlbeam =       1
Next quadratic brick element
```

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```
     nextqbrick=       1
Next quadratic shell element
     nextqshell=       1
Next quadratic beam element
     nextqbeam =       1
```

array create a multi-dimensional array of parameters

array *name*(*d*$_1$,*d*$_2$,...) *data* **;**

where

d_i	is the maximum value in the ith index
data	is optional and can be one value to be broadcast or a full list of values

Remarks

The indices start at 1. There are two ways to initially data load an array. If only one value is found in the data section of this command, then it is broadcast initially through out the array. Alternatively, a value can be listed for every parameter in the array. When the full set of values are listed after the array definition, the values are ordered with a left index moving faster, following the FORTRAN convention, not the C convention.

These array parameters can be used in the desk calculator (**dc**), in the assignment of a scalar value to one element in the array at a time in the **para** command, in expressions that are enclosed in the square brackets "**[...]**", in a numbered comment using the comment command, as operands in the **if**, **when**, **for**, and **while** control statements, in the formation of a function using the **def** command, and in the equations in the part phase (**x=**, **y=**, **z=**, **t1=**, **t2=**, and **t3=**).

Three automatic arrays are available in the part phase. They are

idxlist	one dimensional array with the full i-indices of the part
jdxlist	one dimensional array with the full j-indices of the part
kdxlist	one dimensional array with the full k-indices of the part

The same rules apply to these as to other arrays except they are only available in the part phase and they cannot be set using the para command. Related to tese automatic arrays are three scalar automatic parameters:

maxrudi	maximum reduced i-index
maxrudj	maximum reduced j-index
maxrudk	maximum reduced k-index

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Examples

The following shows an array named height with 2 dimensions containing a total number of values. All values are assigned in this statement. For example, the value %height(2,1) is 1.2.

```
array height(2,3) 1 1.2 2 2.3 3 3.4;
```

In this next example, all 64 values of the parameter array called switches is initialized to 0.

```
array switches(2,2,2,2,2,2) 0;
```

This 1 dimensional array has 3 parameters, none of which have been assigned an initial value.

```
array biggest(3)
```

There are various examples of the use of an array in the following. In particular, note that when an assignment to the array called switches can include a parameter as an index into the array.

```
para l 1 h 2;
dc %height(2,2)/sqrt(2)
para switches(1,%l,2,2,1,1) [%height(%l,%h)/10];
pb 1 2 3 1 2 3 x %biggest(1)
if(%biggest(2).lt.2.0)then
   x=x+%height(i,j)
endif
```

In this next example, the array is used to assign values to coef. Then a 3D spline curve is defined using the values in coef. Also notice that the for statement is embedded within the curd command. This is possible because the for statement (s well as the while and when statements) are macros.

```
para mxi 6;
array coef(10);
para coef(1) 1 coef(2) 1;
for idx 3 %mxi 1
   para coef(%idx) [%coef(%idx-2)+%coef(%idx-1)];
endfor
curd 1 csp3 00
for idx 2 %mxi 1
   [%idx-2] [%coef(%idx)/%coef(%idx-1)] 0.0
endfor
;;;
```

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becho echo and beep

becho *text*

Remarks

A string of up to 80 characters is printed in the text window and the computer beeps.

bulc locate butterfly triple point

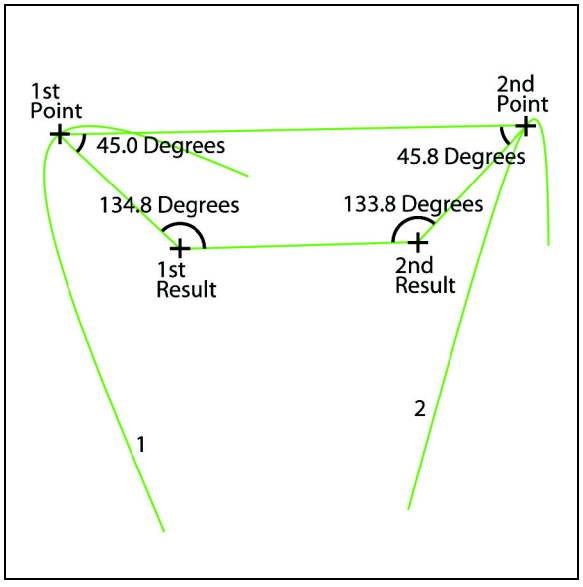
bulc *curve$_1$ curve$_2$ x y z ratio*

Remarks

This command locates an interior vertex of a butterfly structure based only on the shape of two exterior curves. The goal is to produce a pair of points from the two exterior curves and two additional points from this command to form a trapezoid with angles approximately 45, 45, 135, and 135 degrees. This command is needed twice to form the two interior points. The coordinates are saved in the parameters %xprj, %yprj, and %zprj. The point that is input is used to select the closest point on each of the two specified curves. Then a right isosceles triangle is fitted through these points in a plane that is nearly orthogonal to both curves. Actually, there are two such triangles that are formed. The **orpt** command is used to select one from this pair. Then the leg of the triangle that is closest to the initial point is chosen. A point is selected along this leg of the triangle so that the ratio of the leg of the resulting trapezoid and the base of the trapezoid matches the specified ratio.

Example

```
curd 1 csp3 00
   -4.7833413e-01   2.6501161e-01
-8.3939201e-01
   -7.0889658e-01   6.8763685e-01
-1.3136336e-01
   -5.9971255e-01   4.8465800e-01
6.3912874e-01;;;
curd 2 csp3 00
   -6.2983650e-01  -1.9638570e-02
-7.7306795e-01
   -9.9466640e-01   9.3395844e-02
```



Trapezoid using bulc

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```
-3.7190955e-02
 -6.6715145e-01 4.7235709e-02 7.4401599e-01;;;
orpt + 0 0 0
bulc 1 2 -9.9454612e-01 9.5007576e-02 5.5500008e-03 .5
c   -7.752336E-01   1.707245E-01   9.306287E-03
bulc 1 2 -7.1137702e-01 6.9195741e-01 -3.5788484e-02 .5
c   -6.328765E-01   4.716224E-01  -3.093115E-02
```

c beginning of a comment

c *text*

Remarks

The character **c** represents a comment when it appears at the beginning of a line and is followed by a space, or when it appears in the middle of a line and is both preceded and followed by a space. Comments will be preserve in the tsave (session) file.

This feature is especially useful for commenting out commands in a batch input file. Another way to insert comments is to use a dollar sign. If you have a large body of text or a section of commands you wish to turn into comments, encase the text with the curved brackets {...}. This works across multiple lines of text. When pairs are embedded, **True*Grid*®** counts the number of embeddings so that a larger section can be commented out, even when a subsection is already commented out. For example:

```
command 1
command 2
{
command 3
}
command 4
command 5
```

comments out command 3, while

```
command 1
{
command 2
{
command 3
}
command 4
```

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```
}
command 5
```

comments out commands 2, 3, and 4.

cd change directory

cd *new_directory_name*

Remarks

On a WINDOWS system, the default working directory is initially set by the user preferences window (i.e. **tgpref.exe**, also known as **TGControls**). On a UNIX/LINUX system, the default working directory is the window in which **True*Grid*®** is executed. The **cd** changes the working directory. This working directory is the directory for the input/output files for commands such as **iges**, **igesfile**, **saveiges**, **useiges**, **wiges include**, **savepart**, **getbb**, **readmesh**, **vpsd**, **wrsd**, **stlsd**, **edgefile**, **postscript**, trugrdo/**mof** and the **sd** command with options **stl** and **bstl**. However, the tsave and tghist files are not affected by the **cd** command. They will be saved in the default working directory. See the **pwd** command to determine the current working directory.

circent center, radius, and normal of a circle

circent $x_1\ y_1\ z_1\ x_2\ y_2\ z_2\ x_3\ y_3\ z_3$
where (x_i, y_i, z_i) is a point on the circle for i=1,2,3

Remarks

This command finds the center, radius, and normal to a circle that passes through 3 non-colinear points. The center is stored in the automatic parameters %xprj, %yprj, and %zprj, the radius is stored in the automatic parameter %distance, and the normal is stored in the automatic parameters %xnrm, %ynrm, and %znrm.

Example

```
circent 1 1 1 2 -2 2 -3 -3 3
 center =  -8.71622E-01  -1.11486E+00   2.02703E+00
 radius =    3.0050633
 normal =  -1.16248E-01  -3.48743E-01  -9.29981E-01
```

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crprod cross product

Calculates the cross product of two vectors.

crprod x_1 y_1 z_1 x_2 y_2 z_2

Remarks

The predefined paramaters %xcrprod, %ycrprod, and %zcrprod are assigned the result of this function.

Also, add these paramaters names to the index - xcrprod, ycrprod, zcrprod.

Example

```
crprod [cos(70)] [sin(70)] 0 [cos(160)] [sin(160)] 0
c  (x,y,z)=   0.00000E+00   0.00000E+00   1.00000E+00
c  normalized (x,y,z)=   0.00000E+00   0.00000E+00   1.00000E+00
```

curtyp default attach command

curtyp *type*

where *type* can be

cur	equivalence the attach button to the **cur** command
curs	equivalence the attach button to the **curs** command
cure	equivalence the attach button to the **cure** command
curf	equivalence the attach button to the **curf** command

Remarks

This command controls the type of 3D curve attachment used when the ***attach*** button is selected in the environment window while in the part phase. The following summarizes the differences between these options. See the full description of each of these commands for further information.

cur: the selected edge is attached to a curve with the end vertices placed at the closest points on the curve. All interior nodes, including any interior vertices are distributed along the curved, based on the position of the end vertices. Essentially, the interior vertices have no degrees of freedom.

curs: (Default) each component edge of the selected composite edge is placed on the curve, independently. This means that each vertex along the selected component edge is placed on the curve independently. Then the interior nodes of each component edge is distributed along the curve based

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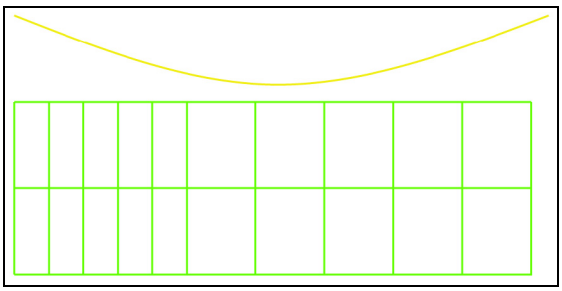
on its end vertices. The same thing can be accomplished by using the **cur** command for each component edge of a composite edge. The **curs** command has the advantage that only one command is needed and if a partition is added with the **insprt**, the **curs** will do the interpolation that is usually expected.

cure: the end vertices of the selected edge are placed at opposite ends of the 3D curve, depending on their initial positions. Then the same interpolation as **cur** is applied. **curf**: the edge is placed onto the 3D curve just as in the **cur** command. However, the nodes are frozen to this 3D curve. Projections to surfaces will have no affect on this edge.

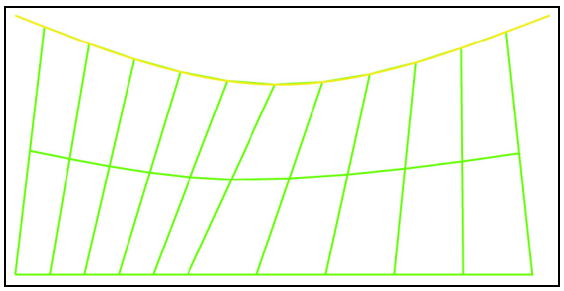
Examples

The following examples all start with the same common input. A simple 2 block part is used with one composite edge attached to a 3D curve. This demonstrates the differences between the **cur**, **curs**, and **cure** commands.

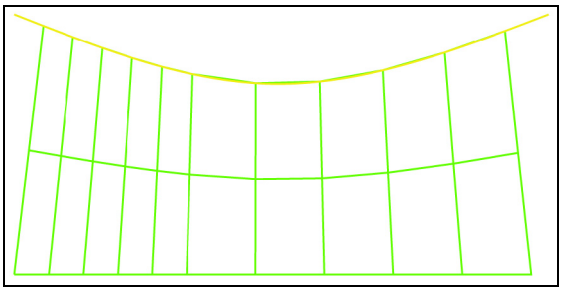
```
block 1 6 11;1 3;-1;.5 1 2 0 .5 0
curd 1 csp3 00 .5 .75 0 1.25 .55 0 2.05 .75 0 ;;;
```



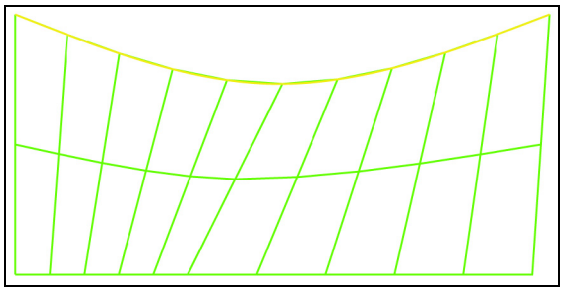
Part & Curve before attachment



Cur attachment



Curs attachment



Cure attachment

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dc desk calculator

Prints the value of a Fortran-like expression.

dc *expression*

Remarks

With **dc**, you can see the value of an expression. For details on the syntax of the expression, see the discussion on expressions at the beginning of this section. The expression must end at the end of the line. Use the **dc** function to learn the value of a parameter.

Example

```
para c2 2 c1 1 c0 -1;
dc (-%c1+sqrt(%c1^2-4*%c2*%c0))/(2*%c2)
 =  5.000000E-01
dc (-%c1-sqrt(%c1^2-4*%c2*%c0))/(2*%c2)
 = -1.000000E+00
```

distance distance between two points

distance x_1 y_1 z_1 x_2 y_2 z_2

Remarks

This command calculates the distance between two points. Use one of the many ways to select a point in the picture using the mouse. Type the **F7** function key to print the coordinates of that point into the **distance** dialogue box or into the text window, if you are typing commands. Select the second point in the picture and type the **F7** function key again. Execute the command to get the distance.

Example

```
distance [cos(45)] [sin(45)] 0 [2*cos(45)] [2*sin(45)] 0
distance =    1.0
```

echo echo a string

echo *text*

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Remarks

A string of up to 80 characters is printed in the text window.

end terminate True*Grid*® with no more output

end *(no arguments)*

errmod set error handler mode

errmod *mode*

where *mode* can be

0	for warning and error messages (default)
1	for avoiding warning messages
2	for error interrupt mode
3	for avoiding warning messages and error interrupt mode

Remarks

These options affect the way warnings and errors are handled.

If the mode 1 or 3 is selected, no warning messages will be issued.

If the mode 2 or 3 is selected, then when an error in the input is encountered while in batch mode, **True*Grid*®** will become interactive as though it had encountered an **interrupt** command in the batch file. Any problems due to the error can then be fixed interactively. Then the **resume** command can be issued to continue the batch command file. This is not easy to do sometimes. Here are three cases that one should be aware of:

1. The command in the batch file is in error and it has many arguments. Any arguments that remain after the error is detected must still be processed once the **resume** command is issued. These arguments will also cause an error and a subsequent interrupt. Warning: do not try to modify the command file at this stage. It will only cause more problems since this command file is being buffered.

2. Some errors are not detected until the **endpart** command is encountered in the batch file. This will cause the part to be completed with errors. Then the interactive mode will be entered in the control phase. There will be little that can be done abut the part at this stage, but you will be able to proceed with a **resume** command if you wish.

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3. When an **insprt** command is encountered, errors in some previously issued commands, such as the **bb** and **trbb** commands, may be detected. This is because the **insprt** command, like the **endpart** command, cause the entire mesh to be re-calculated.

expressions FORTRAN-like expressions

This FORTRAN-like expression interpreter is not a command but is a general feature that can be used anyplace in any command that requires a number. You do this by enclosing the expression in square brackets. For example, you may wish to make the number of elements within a single region of the mesh be a function of a density parameter:

```
block 1 [1+2.7*%d];1 [1+3.2*%d];1 [1+4.9*%d];0 2.7;1 4.2;-1 3.9;
```

You could calculate the distance between two points with

```
[ sqrt(%a*%a + %b*%b) ]
```

For more information on the use of parameters, see the **para** command. Such an expression can be continued across multiple lines. The closing square bracket terminates the expression. There is a limit of 240 symbols in an expression.

Expressions are also used to define a function, a function curve, a function surface, and some loads. The square brackets are not used in this case and to continue a long expression to another line, end the line with a space and ampersand, ("&").

You can enter numbers in the usual FORTRAN formats for integer, floating point, and exponential numbers.

The expressions feature can interpret the arithmetic operators + and - and the binary arithmetic operators: + - * / ** and ^ . All calculations are done in floating point. All trigonometric functions are for angles in degrees.

This feature also interprets the following FORTRAN-like functions:

int(*x***)**	truncates *x* to an integer
nint(*x***)**	rounds *x* to the nearest integer
abs(*x***)**	absolute value of *x*
mod(*a***,***b***)**	*a* modulo *b*
sign(*a***,***b***)**	transfer the sign of *b* to *a*
max(*x*$_1$,*x*$_2$,...,*x*$_n$**)**	maximum number

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min(x_1,x_2,...,x_n) minimum number
sqrt(*x*) square root of *x*; *x* must be positive
exp(*x*) exponential of *x*; *x* must not exceed 85.19
log(*x*) natural logarithm of *x*; *x* must be positive
log10(*x*) common logarithm (base 10) of *x*; *x* must be positive
sin(*x*) sine of *x* (in degrees)
cos(*x*) cosine of *x* (in degrees)
tan(*x*) tangent of *x*; *x* cannot be 90, -90, 180, -180, etc.
asin(*x*) arc sine (in degrees) of *x*; *x* must be between -1 and 1.
acos(*x*) arc cosine (in degrees) of *x*; *x* must be between -1 and 1.
atan(*x*) arc tangent of *x*
atan2(*y*,*x*) arc tangent of *y*/*x*
sinh(*x*) hyperbolic sine of *x*
cosh(*x*) hyperbolic cosine of *x*
rand pseudo-random number from a uniform distribution of unit length, *mean* 0
rand(*seed*) pseudo-random number from a uniform distribution of unit length, *mean* 0, computed from the given *seed*
rand(*seed*,*mean*) pseudo-random number from a uniform distribution between *mean*-½ and *mean*+½, computed from the given *seed*
norm pseudo-random number from a normal distribution with mean 0 and standard deviation 1
norm(*seed*) pseudo-random number from a normal distribution with mean 0 and standard deviation 1, computed from the given *seed*
norm(*seed*,*mean*) pseudo-random number from a normal distribution with mean *mean* and standard deviation 1, computed from the given *seed*
norm(*seed*,*mean*,*sig*) pseudo-random number from a normal distribution with mean *mean* and standard deviation *sig*, computed from the given *seed*; *sig* must be positive

For the random number functions, the default seed is 0.0.

def define a function

def *function_name* **(***a,b,...***)** = *expression*

Remarks

A function can be used in any expression or equation. Following the function name is a pair of parenthesis enclosing a list of dummy arguments that are used in the expression which defines the

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function. The dummy arguments, unlike parameters, do not have a % character at the beginning of the use of a dummy argument in the definition of the function.

Example

```
def len(x1,y1,z1,x2,y2,z2)=sqrt((x1-x2)**2+(y1-y2)**2+(z1-z2)**2)
dc len(1,1,1,2,2,2)
 =  1.732051E+00
dc sqrt(3)
 =  1.732051E+00
```

include execute commands from batch file

include *filename*

Remarks

The specified file will be executed as a batch command file. The commands you have previously issued will remain in effect. Errors will be handled as usually when in batch mode.

The tsave (session) file will incorporate all of the commands that you have issued, whether interactive, from the original batch file, or from an included batch file. Thus, no **include** commands will appear in the tsave file.

The **include** statements can be nested up to 19 statements deep. There is no limit on the number of non-nested **include** commands.

If you forget to select an input file, type or use the include dialogue to initiate the execution of a command file.

A file to be included should be found in the working directory or a path must be included with the file name. The working directory is the directory in which you are running **True*Grid*®**. If you are on WINDOWS, the working directory is specified in the **TGControls** program included with **True*Grid*®**. Alternatively, use the **cd** command to change the working directory.

Example

Suppose there is a file named ring.inc with the following contents:

```
cylinder 1 3;1 91;-1;%r1 %r2;0 360;0;
```

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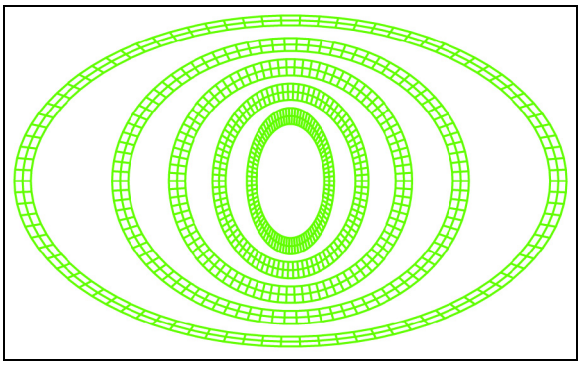
```
lct 1 xsca %xs ysca %ys;lrep 1 ;
endpart
```

Then the following commands will produce the set of rings shown below:

```
para r1 .7 r2 .9 xs .6 ys 1;
include ring.inc
para r1 1 r2 1.2 xs .8 ys 1;
include ring.inc
para r1 1.3 r2 1.5 xs 1 ys 1;
include ring.inc
para r1 2 r2 2.2 xs 1 ys .8;
include ring.inc
para r1 3.2 r2 3.4 xs 1 ys .6;
include ring.inc
merge
```



Concentric rings w/ parameters & include

inprod inner or dot product

This command computes the inner or dot product of two vectors.

inprod $x_1\ y_1\ z_1\ x_2\ y_2\ z_2$

Remarks

The result of this function is stored in the automatic parameter %inprod.

Example

```
inprod [cos(20)] [sin(20)] 0 [cos(30)] [sin(30)] 0
 dot product =   9.84808E-01
dc cos(10)
 =  9.848077E-01
```

interrupt switch from batch commands to interactive mode

When **True*Grid*®** is executing batch commands, this command halts batch execution and begins interactive execution; i.e. with keyboard and mouse.

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interrupt *<no arguments>*

Remarks

Upon reaching this command in a batch file, **True*Grid*®** halts execution of commands in the file and becomes interactive. This gives you an opportunity to interactively modify the model. When you wish to continue execution of the batch file, you can issue the **resume** command.

This feature is useful in making a change to an existing mesh. Using a text editor, enter the interrupt command into the command file at the end of the list of commands for each part that needs modification. Then run the command file. When **True*Grid*®** becomes interactive issue the commands to make the desired modifications and click on the ***Resume*** button (or type the **resume** command into the test window).

Keep in mind that **True*Grid*®** will start executing commands from the command file from where it stopped after encountering the **interrupt** command. If, for example, there was an **endpart** command just after the **interrupt** command in the command file, then it would cause an error if you were to type the **endpart** command into the text window before you resumed execution of the commands from the command file. **True*Grid*®** would see two **endpart** commands in sequence, and the second one would be an error.

After you have completed the execution of all of the commands from the command file and the commands you issued interactively, you can exit **True*Grid*®** and save the new sequence of commands found in the session file (tsave file).

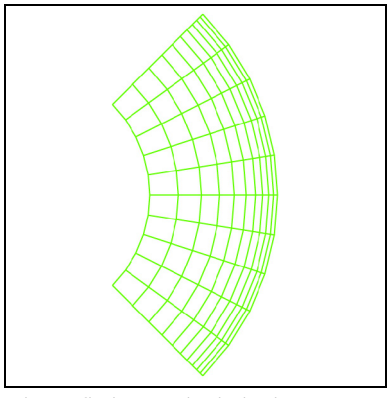
This procedure can also be used to hunt down a bug in your command file. After entering the interactive mode, use the ***History*** window to help locate the command(s) in error. This is also a good way to understand the topology and techniques used to create a model that you are not familiar with.

intyp set default mesh interpolation

intyp *option*

where *option* can be

1	linear interpolation (initial default) (**lin** and **lini** commands)
2	transfinite interpolation (**tf** and **tfi** commands)



Linear interpolation Transfinite interpolation

Remarks

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The linear interpolation is cheap and is good for most problems. The transfinite interpolation is better when there is a lot of curvature and when the nodes along the edges are no distributed evenly.

Example

```
block 1 11;1 11;-1;1 3 -1 1 0
sfi -1;;-1; cy 0 0 0 0 0 1 1
sfi -2;;-1; cy 0 0 0 0 0 1 2
sfi 1 2;-2;-1; plan 0 0 0 1 -1 0
sfi 1 2;-1;-1; plan 0 0 0 1 1 0
res 1 1 1 2 2 1 i .8
```

lcrtgl local coordinate system vector loads

lcrtgl *switch*

where *switch* can be

on	use local coordinate system
off	use global coordinate system

Remarks

When a part is replicated using the **lrep**, **grep**, or **lev** commands, all of the loads are also replicated. The default is to replicate the load in the global coordinate system. If, for example, there is a fixed displacement, the corresponding nodes in each replication have the same displacement as if the displacement were specified in the global coordinate system. If the part is rotated or reflected, the load vector is not.

This command causes the vector loads to be rotated or reflected along with the replicated part. This new command applies to the following commands that define loads in the part phase. It applies to all loads of the type below within that part. It is an internal variable and continues in effect until this command is issue with a different option.

fc, **fci**, **fcc**, **fcci**, **fcs**, **fcsi** - fixed force
fd, **fdi**, **fdc**, **fdci**, **fds**, **fdsi** - fixed displacements
fv, **fvi**, **fvc**, **fvci**, **fvs**, **fvsi** - fixed velocity
fvv, **fvvi**, **fvvc**, **fvvci**, **fvvs**, **fvvsi** - fixed velocity
acc, **acci**, **accc**, **accci**, **accs**, **accsi** - fixed velocity
vacc, **vacci**, **vaccc**, **vaccci**, **vaccs**, **vaccsi** - fixed velocity
ve, **vei** - initial velocities
dis, **disi** - initial displacements

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painfo print information about parameters

This command gives you information on the currently defined parameters.

painfo *<no arguments>*

Remarks

This command helps you remember what you did with the **para** command.

para define parameters

para $name_1$ $value_1$ $name_2$ $value_2$ ... $name_n$ $value_n$;
or
parameter $name_1$ $value_1$ $name_2$ $value_2$ ... $name_n$ $value_n$;

Remarks

This command is used to define parameters. Each is a symbol that stands for a number. They can be used in any other commands that need numbers or expressions. You can define up to 10,000 parameters.

Parameters enable you to set up your model in such a fashion that you can quickly and easily change the design whenever you want.

The arguments to the **para** command are pairs, first the symbolic name and then its value. The value can be a number or an expression. This value will be used until you redefine the parameter with another **para** command.

To reference a parameter, put a **%** before its symbolic name. For example:

```
para  a 1.2  b [ sqrt(%a) ]
sd  1  plan 0 0 %b 0 0 1
```

will define a plane passing through the point

$$(x = 0,\ y = 0,\ z = \sqrt{1.2})$$

A parameter can be used in place of any number. It can be used in expressions and equations. It can

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be used to assign a value to another parameter. In all cases, you must assign a value to the parameter before referencing it. An error message will be issued if you reference a parameter without a value. All parameters are stored internally as floating point values. If a parameter is used for an integer, then the value will be truncated when it is used.

For the name of the parameter, you can use any character string that obeys the following rules:

• characters may be any of the standard (ASCII) printing characters except for:
+ , % * / ^ ( ) [ ] = & $
• the first 16 characters must be unique
• the name must be unique when case is ignored

To see the currently defined parameter values, issue the **painfo** command.

There are some predefined automatic parameters. They are:

distance - result of the **distance** function or
resulting radius of the circle from the **circent** function
inprod - result of the **inprod** function
nextbb - 1 greater than the largest block boundary interface number - see the **bb** command
nextcrv - 1 greater than the largest 3D curve number - see the **curd** command
nextlbeam - 1 greater than the largest linear beam element number from the previous parts
nextlbrick - 1 greater than the largest linear brick element number from the previous parts
nextlshell - 1 greater than the largest linear shell element number from the previous parts
nextlc - 1 greater than the largest load curve number - see the **lcd** command
nextln - 1 greater than the largest 2D curve number - see the **ld** command
nextmat - 1 greater than the largest material number - see the material menu
nextnode - 1 greater than the largest node number of the previous part
nextprt - 1 greater than the last part number
nextqbeam - 1 greater than the largest quadratic beam element number from the previous parts
nextqbrick - 1 greater than the largest quadratic brick element number from the previous parts
nextqshell - 1 greater than the largest quadratic shell element number from the previous parts
nextsrf - 1 greater than the largest surface number - see the **sd** command
node - the node number from the **ajnp** or **sajnp** command or
from the Environment window **Pick>node** selection in the merge phase or
from the **F5** function key to pick a node graphically in the merge phase
subang - result of the **subang** function
xprj, yprj, zprj - coordinates due to the **project** command or
results of the **tricent** function or
resulting center of the circle from the **circent** function or
results of the **trapt** function or

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results of the **bulc** function or
results of the **nodcor** or **snodcor** function or
intersection of 2 2D curves using **ld lpil**

xnrm, ynrm, znrm - normal to the 1st surface from the **project** function
resulting normal to the circle from the **circent** function or

pi - 3.1415926...

mxiridx - maximum i-index of the present part

mxjridx - maximum j-index of the present part

mxkridx - maximum k-index of the present part

xcrprod, ycrprod, zcrprod - coordinates due to the **crprod** function

xboxmin - minimum x-coordinate of the grid in the graphics (**rest** command)

yboxmin - minimum y-coordinate of the grid in the graphics (**rest** command)

zboxmin - minimum z-coordinate of the grid in the graphics (**rest** command)

xboxmax - maximum x-coordinate of the grid in the graphics (**rest** command)

yboxmax - maximum y-coordinate of the grid in the graphics (**rest** command)

zboxmax - maximum z-coordinate of the grid in the graphics (**rest** command)

ptcoor parametric coordinates from a surface or 3D curve

ptcoor *point_symbol*

where the *point_symbol* can be

s.i.j for a labeled point on a surface where
- *s* is the surface number
- *i* is the first index of the point on the surface
- *j* is the second index of the point on the surface

c.i for a labeled point on a 3D curve where
- *c* is the 3D curve number
- *i* is the point index on the curve

Remarks

This command is used primarily to assign the parameters **xprj**, **yprj**, and **zprj** the coordinates of a labeled point on a surface or 3D curve. A full discussion of labeled points on surfaces can be found in the discussion on surface tessellation and labeling (106). Examples of labeled points along a 3D curve can be found under the labels command (?).

pptcd parameterized coordinates of a 3D curve

pptcd *curve_number parameter*

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where

curve_number	an existing 3D curve number
parameter	number from 0.0 to 1,0

Remarks

This command interpolates a point along a 3D curve. The 3D curve is parameterized by relative arc length ranging from 0.0 (start) to1.0 (end). The coordinates are stored in the automatic parameters **xprj**, **yprj**, and **zprj** respectively.

pwd get current working directory

pwd (no arguments)

Remarks

On a WINDOWS system, the default working directory is initially set by the user preferences window (i.e. **tgpref.exe**, also known as **TGControls**). On a UNIX/LINUX system, the default working directory is the window in which **True*Grid*®** is executed. See the **cd** to change the working directory. This working directory is the directory for the input/output files for commands such as **iges**, **igesfile**, **saveiges**, **useiges**, **wiges include**, **savepart**, **getbb**, **readmesh**, **vpsd**, **wrsd**, **stlsd**, **edgefile**, **postscript**, trugrdo/**mof** and the **sd** command with options **stl** and **bstl**. However, the tsave and tghist files are not affected by the **cd** command. They will be saved in the default working directory. The **pwd** command prints the current working directory.

resume resume executing batch commands

resume *<no arguments>*

Remarks

This command resumes execution of a batch file after the **interrupt** command has halted it. See the discussion on the **interrupt** command for the usage of the **interrupt** and **resume** commands.

sp4pt center and radius of a sphere

sp4pt *x1 y1 z1 x2 y2 z2 x3 y3 z3 x4 y4 z4*

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Remarks

The center and radius of a sphere is determined from 4 points. These points cannot be co-linear or co-planer.

title assign title to the problem

title *text*

Remarks

This command assigns a title to the model. The title should be a one-line description of the model. The title is used as the default graphics caption and is usually found somewhere in the output file. You can change the caption without changing the title by issuing the **caption** command.

tpara typed parameters

tpara *type symbol_1 value_1 type symbol_2 value_2 ... ;*

Remarks

This is used to pass parameters into REFLEQS, a fluids simulation code.

tricent find optimal center of a triangular structure

tricent *type* x_1 y_1 z_1 x_2 y_2 z_2 x_3 y_3 z_3

where *type* can be

1	coordinates at the corners
2	coordinates along the midpoint of edges

where

x_1 y_1 z_1	Cartesian coordinates of first point
x_2 y_2 z_2	Cartesian coordinates of second point
x_3 y_3 z_3	Cartesian coordinates of third point

Remarks

The optimal is where the three edges that meet at the center, meet at 120 degrees. There are two options. The first option bases the calculation of the coordinates of the three corners of the triangle.

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It assumes the midpoints of the three edges of the triangle will be connected to the center vertex. In the second option, the intermediate points along the three edges are used to calculate the optimal position for the center vertex. The latter is more versatile since it does not assume midpoints along the edges. The coordinates of the center point are immediately stored in the automatic parameters %xprj, %yprj, and %zprj.

Examples

Both of the following examples start with this input:

```
block 1 6 11;1 6 11;-1;0 1 2 0 1 2
4
sd 1 sp 0 0 0 4
sfi ;; -1; sd 1
dei 2 3; 2 3; -1;
pb 2 2 1 3 3 1 xy 1 1

tricent 1 0 0 4
  1.7888 0 3.5777
  0 1.7888 3.5777
pb 2 2 1 2 2 1 xyz
  %xprj %yprj %zprj
```

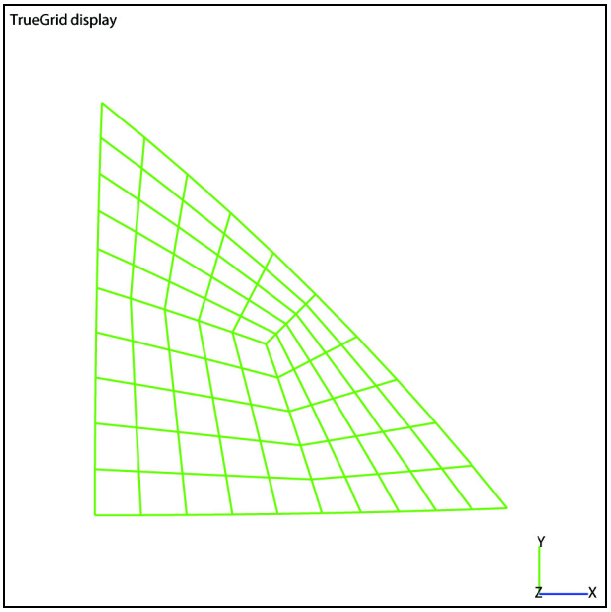

Figure 309 Optimal center based on corners

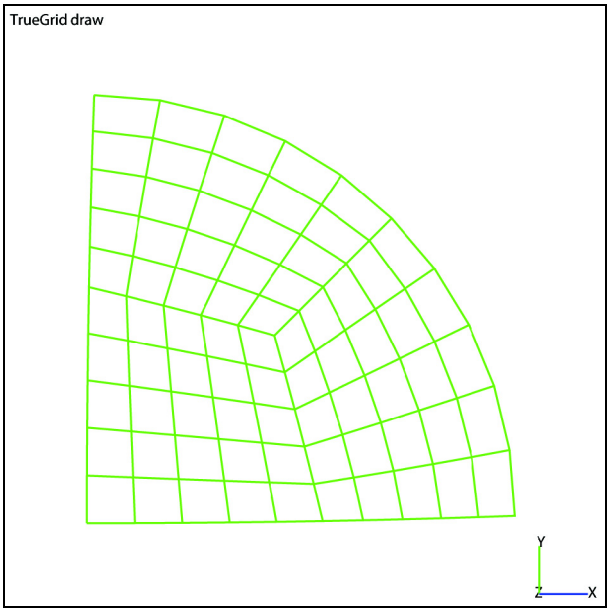

Figure 310 Optimal center based on edges

```
sd 2 cy 0 0 0 0 0 1 2
sd 2 cy 0 0 0 0 0 1 1.789
sfi 1 2; -3; -1; sd 2
sfi -3; 1 2; -1; sd 2
tricent 2 .97 0 3.88
  1.265 1.265 3.577
  0 9.7 3.88
pb 2 2 1 2 2 1 xyz
  %xprj %yprj %zprj
```

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subang angle between two intersecting lines

subang x_1 y_1 z_1 x_2 y_2 z_2 x_3 y_3 z_3

Remarks

The angle subtended by two intersecting lines is printed. The automatic parameter %subang will be assigned the angle in degrees.

Example

```
subang 1 1 [sqrt(2)] 0 0 0 1 1 [-sqrt(2)]
angle =    90.0
```

caption assign a caption to the physical window

caption *option*

where the *option* can be

cap *text*
on
off

Remarks

The title defined with the **title** command is the default caption.

mxp change number of mesh convergence passes

mxp #_*passes*

Remarks

This command is used to handle rare cases where inter-dependencies in the building of a part are not fully resolved. This can occur when there are many interpolations/smoothing commands that intersect and cross each other.

This is the maximum number of passes used in the projection algorithm to resolve dependencies within the **block**, **cylinder**, and **blude** parts. These dependencies occur when a multiple region is interpolated and its boundary is the interior of another multiple region interpolation. The default is 4. The minimum

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is 1. 2 handles almost all cases. If there are dependencies, it will take a longer time to build the part when the maximum number of passes is increased.

trapt transform a point

trapt *x y z type id*

where the *type* can be

lct for a local coordinate transformation

gct for a global coordinate transformation

where *id* is the number of the transformation

Remarks

Trapt will transform a 3D coordinate using a local or global coordinate transformation formed in the **lct** or **gct** commands, respectively. The results are stored in the parameters %xprj, %yprj, and %zprj.

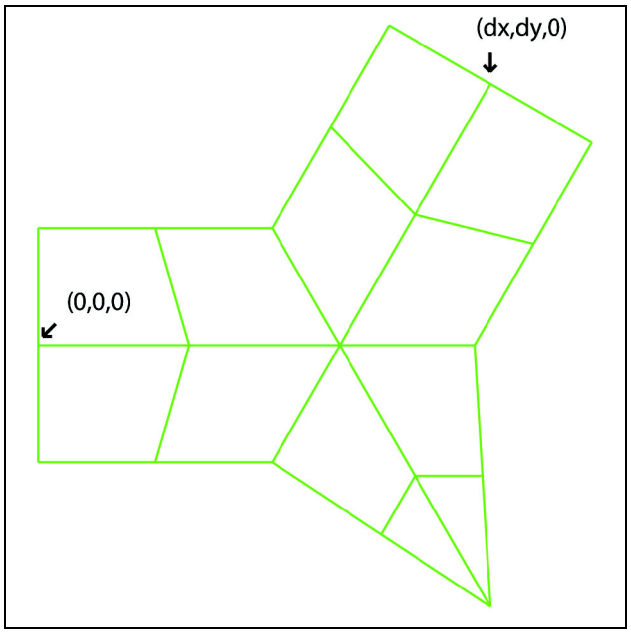
This feature can be used when building a sequence of parts that must start where the previous part ended. Each component can be constructed in a local coordinate and then transformed to the final position using a coordinate transformation. This command can be used to maintain a fiducial point as parts are added.

It is possible to do without this command by using only the replication commands. In some cases, calculating the translation vector can be complicated. This method simplifies the calculation of that translation vector.

Example

This example started with a single component by creating a command file named why.inc with the following commands:

```
block 1 3 0 4 5 6 0 8 10;
 1 2 3 0 5 7;
 1 2;
  [-2-1/sqrt(3)] [-1/sqrt(3)] 0
 -1 0 1 0
 [1/sqrt(3)] [2+1/sqrt(3)];
 -1 0 1 0 [1/sqrt(3)]
```



Component starting at the origin

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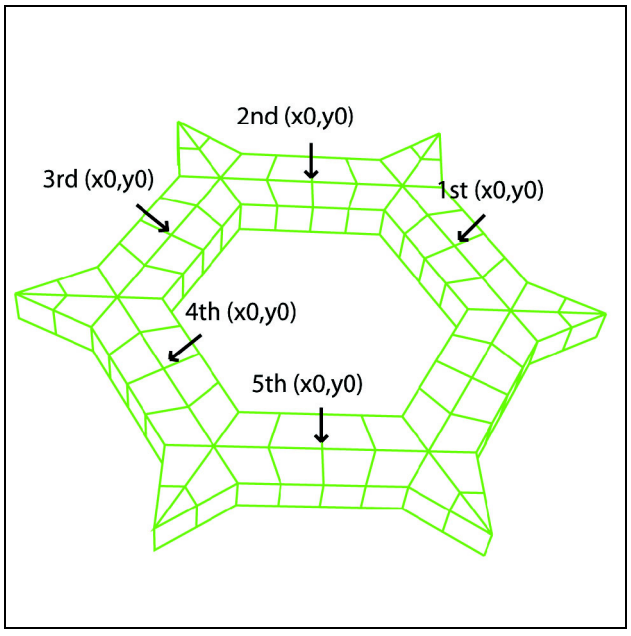
```
 [2+1/sqrt(3)];
 0 1;
dei 4 6; 1 3;;
dei 8 9; 5 6; 1 2;
dei 1 2; 5 6; 1 2;
tr 8 1 1 9 3 2 rz -60;
tr 4 5 1 6 6 2 rz -30;
pb 9 1 1 9 3 2
 xy 1.288675 -2.232051
pb 2 2 1 2 2 2 xy 0 0
pb 5 5 1 5 5 2 xy 0 0
pb 8 2 1 8 2 2 xy 0 0
mb 1 1 1 9 6 2 x [2+1/sqrt(3)]
```

To complete this example, a second command file is formed with the commands below. The component part starts at the origin, to simplify the calculations. Only the x and y-coordinates are maintained, since the part is formed in a plane, again for simplification. The vector (dx,dy,0) is formed between the starting and ending positions of the replicated part. This vector is transformed, just like the component part, to maintain the translation vector for the next replication.

```
para dx [2+1/sqrt(3)+cos(60)*(2+1/sqrt(3))] c differential vector
     dy [sin(60)*(2+1/sqrt(3))];
include why.inc
endpart
para  x0  %dx  y0  %dy;  c  1st
fiducial
include why.inc
gct 1 rz 60 mx %x0 my %y0;
grep 1;
endpart
trapt %dx %dy 0 gct 1
para x0%xprj c 2nd fiducial
     y0%yprj;
include why.inc
gct 1 rz 120 mx %x0 my %y0;
grep 1;
endpart
trapt %dx %dy 0 gct 1
para x0%xprj c 3rd fiducial
     y0%yprj;
include why.inc
gct 1 rz 180 mx %x0 my %y0;
```



Components rotated and translated

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```
grep 1;
endpart
trapt %dx %dy 0 gct 1
para x0%xprj c 4th fiducial
     y0%yprj;
include why.inc
gct 1 rz 240 mx %x0 my %y0;
grep 1;
endpart
trapt %dx %dy 0 gct 1
para x0%xprj c 5th fiducial
     y0%yprj;
include why.inc
gct 1 rz 300 mx %x0 my %y0;
grep 1;
endpart
merge
stp .001
```

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IV. Output

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There are numerous output formats that can be selected for the output file containing the mesh. There are two fundamental file format types. The default is the Finite Element format type. Examples of this are ABAQUAS, ANSYS, LS-DYAN, and NASTRAN. Alternatively, there is the block structured format type such as CFX4, GRIDGEN, or TASCFlow. If you wish to choose the latter, you must do so at the beginning of the input or interactive session.

You must be in the **merge** phase to create the mesh file in one of these formats. Usually you issue a merge nodes command first and then create the file with the **write** command.

1. Output File Format Commands

A full description of the output commands can be found in the **True*Grid*®** Output Manual. The write command is used to **write** the output file containing the specified output file format.

comment pass a comment to the DYNA3D output file

comment *[option [id]] text*

where the *option* can be

no	for the beginning of the output file
mt	for a material
np	for the nodes
he	for the hexahedron elements
be	for the beam elements
se	for the shell elements
ts	for the thick shell elements
sp	for a slide planes
nh	for the node history
hh	for the hexahedron element history
bh	for the beam element history
sh	for the shell element history
th	for the thick shell element history
br	for the brode
lc	for a load curve
nl	for the nodal loads
pr	for the pressure
ve	for the velocity
sw	for a stone wall
nc	for the nodal constraints
in	for the initial conditions
si	for a sliding interface

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ti	for the tied nodes with failure
ex	for extra nodes w/ rigid body
jt	for a joint
xa	for the x-acceleration
ya	for the y-acceleration
za	for the z-acceleration
xv	for the x-velocity
yv	for the y-velocity
zv	for the z-velocity
dt	for a detonator
sd	for the springs and dampers
nr	for the non-reflecting boundaries
md	for momentum deposition in solid elements
1d	for 1D slide line

where *id* can be a number to identify the number of the type object being commented on

Remarks

Up to 100 comments can be saved. Not all types have an id associated with it and if a number is not the first thing encountered after the type, then the text is assumed to be general and not applied to a specific object. The id can be a parameter of an expression in square brackets.

elvotest switch solid element test for zero volume

elvotest *switch*

where *switch* can be

on	turn of the test when the output file is written
off	turn off the test when the output file is written

Remarks

The default is off. This test is applied to the mesh when the output file is written. It applies only to the following output options: ANSYS, ABAQUS, CFX5, EXODUSII, FIDAP, FLUENT, ICFEP, LSDYNA, KIVA, NASTRAN, and the NEUTRAL formats.

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epb element print block for DYNA3D and LSDYNA

epb *set_name*

where

set_name	name of an element set

mof mesh output file name

mof [*path*]*file_name*

where

path	path to the file - optional
file_name	file name including the suffix

Remarks

This function changes the name of the mesh output file to be written. It does not change the name of a mesh output file that has already been written.

The same thing can be accomplished by placing the o= option on the execute line. Under WINDOWS, the o= on the execute line is only possible when you run from a Command Prompt window. If you specify a path, it will be relative to the working directory.

When you start **True*Grid*®** from the desk top in WINDOWS, the working directory is determined from the settings in the **TGControls** window. Then the output file named in this command will be found relative to the working directory selected in **TGControls**.

Be sure to issue the **write** command in the Merge phase to actually create the desired mesh output file.

ndigits number of digits written for coordinates

ndigits *n*

where *n* is the number of digits

Remarks

The default is 7 and the minimum is 5. This applies only to the Abaqus, Ale3d, ICFEP, MPACT, and CFX output options. The maximum is 14, but they may not all be used if the machine dos not supple that much precision. 7 digits are used in a floating point representation other than for the mantissa. Therefore, there are at a maximum 7 digits in the mantissa for a 32 bit version and 13 digits for the 64 bit version.

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npb node print block for DYNA3D and LSDYNA

npb *node_selection*

where *node_selection* can be one of the following:

n *node*	select a node
rt *x y z*	select the node nearest to a point in Cartesian coordinates
cy *rho theta z*	select the node nearest to a point in cylindrical coordinates
sp *rho theta phi*	select the node nearest to a point in spherical coordinates
nset *set_name*	select all nodes in a node set

save dump buffered data to the tsave file

save *(no arguments)*

Remarks

On many systems, data is buffered before it is written to a file on disk. This command forces the data in the buffer to be written to disk. This guarantees that if there is a failure, the data will be saved in the tsave file. This is usually not needed, because if, in the unlikely event, **True*Grid*®** should crash, the tsave file is properly closed first. However, this command is most useful if you want to use the data dumped to the tsave file without terminating **True*Grid*®**. By issuing the save command, you are assured that the tsave file is up to date.

verbatim write verbatim to the output file

verbatim
text
...
endverbatim

Remarks

The primary purpose for this command is to supplement the mesh data written to the ASCII data file produced for one of the output formats in **True*Grid*®**. Although there are many features supporting the various output options, not every feature in a simulation code is supported by **True*Grid*®**. The **verbatim** command makes it possible to include such features in the ASCII data file by typing the

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exact text between the **verbatim** and **endverbatim** keywords. If you find that you are using this command frequently, contact the office of XYZ Scientific Applications and request a new feature that will automate that you are doing.

The text can be any number of lines of text. Anything on the same line and following the verbatim command is ignored by **True*Grid*®**. This text, including the $ character, is saved. When the **write** command is issued, this text is written to the output file if the output file format is keyword driven so that the text can be in any order. This is needed because text is always placed near the beginning of the output file. Some of the formats supported by this feature are **Abaqus**, **Ansys**, **LSdyna**, **Marc**, and **Nastran**. Parameters can be used with this command. The character '$' is not treated as the start of a comment when included in the scope of this command.

The **verbatim/endverbatim** command data that is written to some of the output formats can include the '{' and '}' characters if they are preceded by the backslash character '\'.

Example

verbatim
This is a test of the \{ and \} characters.
endverbatim

Then the output file from **True*Grid*®** would contain the following line:

This is a test of the { and } characters.

write write the output file containing the mesh

write (no arguments)

Remarks

You must be in the merge phase to write the output file. Usually you want to issue one of the merge nodes commands before issuing the write command. There is no default format. You must choose a format, for example ABAQUS, ANSYS, NASTRAN, or LS-DYNA, before writing the file.

The default file name is trugrdo. It will be found relative to the working directory. If you issue the write command a second time, it will be called trugrdo.0001. Each additional write will cause the file name to be incremented. If there exists a trugrdo file before you issue the first write, then the first write will overwrite the trugrdo file. When you issue a write command, a message will appear indicating the name of the file that was written.

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The output file name can be selected if you are running in a Command Prompt Window under WINDOWS or if you are running on UNIX/LINUX/OSX using the o= option on the execute line. You can also switch from the default name by issuing the MOF command.

abaqus ABAQUS output format

abaqus (no arguments)

ale3d ALE3D output format

ale3d *(no arguments)*

ansys ANSYS output format

ansys *(no arguments)*

autodyn AUTODYN-3D output format

autodyn (no arguments)

cf3d Convective Flow output format

cf3d (no arguments)

cfd-ace CFD-ACE output format

cfd-ace (no arguments)

cfx CFX output format

cfx (no arguments)

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dyna3d DYNA3D output format

dyna3d (no arguments)

es3d ES3D output format

es3d (no arguments)

exodusii Exodus II output format

exodusii (no arguments)

fidap FIDAP output format

fidap (no arguments)

fluent FLUENT output format

fluent (no argument)

gemini GEMINI output format

gemini (no arguments)

gridgen3d GRIDGEN output option

gridgen (no arguments)

iri IRI output format

iri (no arguments)

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lsdyna LS-DYNA output format

lsdyna *type*
where
type can be:

fixed	for the fixed format
keyword	for the keyword format

lsnike3d LS-NIKE3D output format

lsnike3d (no arguments)

icfep Imperial College of London finite element program

icfep (no arguments)

Remarks

This should only be used for **quadratic** elements. The **icfepmat** command assigns the element type to a material. For 2 dimensional problems, be sure that all of the z-coordinates have absolute value less than or equal to **ztol**. Be sure to merge the nodes in the merge phase before issuing the **write** command.

kiva3v KIVA3V output option

kiva3v (no arguments)

Remarks

Kiva3v is a block structured code only and the kiva3v command must be issued before any parts are generated.. Use the **kiva3bc** and the **kiva3bci** commands to define boundary conditions in the part phase.

kiva KIVA4 output option

kiva (no arguments)

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Remarks

Use the **kivabc** and the **kivabci** commands to define boundary conditions in the part phase. Use the **kivabc** command to define boundary conditions in the merge phase.

In the merge phase, use the **write** command to write the kiva4grid file. Use material numbers 10 for squish, 11 for bowl, 14 for dome, and 20, 30, 40, and 50 for the cells forming the ports. Material number 1 is the default material in **True*Grid*®** and it will be interpreted as squish. Material 2 will be interpreted as inactive. Use the **kivabc** command to set face types, except for periodic conditions. Use the **bb** and **trbb** with the kiva periodic options to impose periodic boundary conditions and face types. The nodes of a master block boundary interface will appear first in the list of paired periodic vertices. The default face type is fluid.

marc MARC output format

marc (no arguments)

mpact MPACT finite element code by Pedro Marcal

mpact (no arguments)

nastran NASTRAN output format

nastran (no arguments)

nek5000 NEK5000 output format

nek5000 (no arguments)

nike3d NIKE3D output format

nike3d (no arguments)
enike3d (no arguments)
nnike3d (no arguments)
fnike3d (no arguments)

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nekton2d NEKTON2D output format

nekton2d (no arguments)

nekton3d NEKTON3D output format

nekton3d (no arguments)

ne/nastran NE/NASTRAN output format

ne/nastran (no arguments)

neutral Neutral output Format

neutral (no arguments)

refleqs REFLEQS output format

refleqs (no arguments)

sap2000 SAP2000 FEA output file in IGES format

sap2000 (no arguments)

Remarks

The output file name is sapmesh.igs.

starcd STARCD output format

starcd (no arguments)

stl Stereo Lithography (STL) output format

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stl *mode*

where *mode* can be

1	for ASCII 3d Printing
2	for Binary 3D Printing
3	for ASCII exterior mesh surface extraction
4	for Binary exterior mesh surface extraction

Remarks

This command sets the format used when the **write** command is issued in the merge phase. The data file produced is either ASCII or binary. This is the data file sometimes used for StereoLithography (frequently called additive manufacturing). It can also be used to extract the exterior of a mesh and turn it into a surface. For example, an existing mesh can be imported using the **readmesh** command and the exterior faces written to a STL file. In a second run of **True*Grid*®**, the STL file can be read where the surfaces can be used for projections. The visible faces of each material of the original mesh forms a surface in the STL file. To read the STL file into **True*Grid*®**, use the ASCII or binary STL option in the **sd** command or the **stlsd** command in the CAD menu if it is ASCII and there are multiple surfaces (materials) in the STL file.

If the STL is to be used for a 3D printer, choose the option for 3D printers. This is needed because the surface in the STL file must have coordinates in all three directions starting at zero. On the other hand, if the surface is being extracted for other reasons and the part is not to be translated to zero, then choose the option for exterior mesh surface extraction.

These files can get large and the binary format will be considerably shorter than the ASCII format file with the same data. The user is responsible for making the positive normals of shell elements all point outward, if this file is to be used with a 3D printer. All brick faces are automatically pointing outward. It is also the responsibility of the user to make the surface water tight by merging nodes.

Use the **etd** command to select certain element types in the picture. Use the various ways of selecting parts and materials to be drawn in the graphics. Only the faces of the element types, materials, and parts shown in the picture will be recorded in the STL file.

It is possible to write multiple STL files, but each needs a unique name or only the last STL file will survive. Use the **mof** command to change the STL file name between writes. After selecting the parts, materials, element types, and the file name, issue the **write** command. As usual in **True*Grid*®**, you must be in the merge phase to write an outut file.

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plot3d PLOT3D output format

plot3d (no arguments)

poly3d Generic output format

poly3d (no arguments)

stl StereoLithography format

stl *mode*

where *mode* can be

1	for ASCII 3D Printing
2	for Binary 3D Printing
3	for ASCII exterior mesh surface extraction
4	for Binary exterior mesh surface extraction

Remarks

This output command writes a STL file in ASCII or binary format. This is the data file sometimes used for StereoLithography (frequently called additive manufacturing). It can also be used to extract the exterior of a mesh and turn it into a surface. Only the visible faces of the mesh are placed into the STL file.

If the STL is to be used for a 3D printer, choose the option for 3D printers. This is needed because the surface in the STL file must have coordinates in all three directions starting at zero. On the other hand, if the surface is being extracted for other reasons and the part is not to be translated to zero, then choose the option for exterior mesh surface extraction.

These files can get large and the binary format will be considerably shorter than the ASCII format file with the same data. The user is responsible for making the positive normals of shell elements all point outward, if this file is to be used with a 3D printer. All brick faces are automatically pointing outward. It is also the responsibility of the user to make the surface water tight by merging nodes.

Use the **etd** command to select certain element types in the picture. Use the various ways of selecting parts and materials to be drawn in the graphics. Only the faces of the element types, materials, and parts shown in the picture will be be recorded in the STP file.

It is possible to write multiple STL files, but each needs a unique name or only the last STL file will

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survive. Use the **mof** command to change the STL file name between writes. After selecting the parts, materials, element types, and the file name, issue the **write** command. As usual in **True*Grid*®** , you must be in the merge phase to write an output file.

To read the STL file into **True*Grid*®** , use the ASCII or binary STL option in the **sd** command.

tascflow TASCflow output format

tascflow (no arguments)

topaz3d TOPAZ3D output format

topaz3d (no arguments)

topaz3d2 TOPAZ3D version 2000 output format

topaz3d2 (no arguments)

viewpoint VIEWPOINT output format

viewpoint (no arguments)

write WRITE the mesh to a formatted file

write (no arguments)

Remanrks

This command is only effective if an output option has been selected. The format options are listed under the output options. The default name for the output file is "**trugrdo**". There are several exceptions, so check the details for the output format . Use the "o =" option on the execute line (when executing from a command prompt window if you are running a WINDOWS system) and the **mof** command during execution to choose a file name other than the default file name.

2. Analysis Options

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For control parameters and other such input options which are specific to the simulation code, use the appropriate command from this section.

To issue one of these commands, use its dialogue box. For full details, see the **True*Grid*®** Output Manual. But for the sake of completeness, the command names are given below.

XYZ Scientific Applications is actively adding support of more simulation codes. We welcome your requests for more output formats!

abaqstep **ABAQUS analysis step**

ansyopts **ANSYS analysis option**

dynaopts **DYNA3D analysis options**

lsdyopts **LS-DYNA analysis and database options**

marcopts **MARC analysis options**

nastopts **NASTRAN analysis options**

nenstopt **NE/NASTRAN analysis options**

nekopts **NEK5000 analysis options**

nikeopts **NIKE3D analysis options**

lsnkopts **LS-NIKE3D analysis options**

tz3dopts **TOPAZ3D analysis options**

3. Material Definitions

True*Grid*® generates input files for any of a large number of simulation codes. Each code has its own set of material models and its own way to specify them. You can specify your material models in **True*Grid*®** using a **True*Grid*®** dialogue box. For detailed information about material specifications, refer to the manuals of the appropriate simulation code. It is not practical to reproduce those manuals here.

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For the sake of completeness, the currently available material definition commands are listed in the **True*Grid*®** Output Manual. To use these commands, use **True*Grid*®**'s graphical user interface.

XYZ Scientific Applications is actively adding support of more simulation codes. We welcome your requests for even more simulation codes!

abaqmats **ABAQUS materials**

ansymats **ANSYS materials**

dynaeos **DYNA3D equation of state**

dynamats **DYNA3D materials**

fluemats **FLUENT materials**

nastmats **NASTRAN materials**

nenstmats **NE/NASTRAN materials**

lsdymats **LS-DYNA materials**

lsdythmt **LS-DYNA thermal materials**

lsdyeos **LS-DYNA3D equation of state**

nikemats **NIKE3D materials**

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lsnkmats **LS-NIKE3D materials**

marcmats **MARC materials**

patsmats **PATRAN materials**

tz3dmats **TOPAZ3D materials**

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V. Appendix

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Cartesian coordinate system

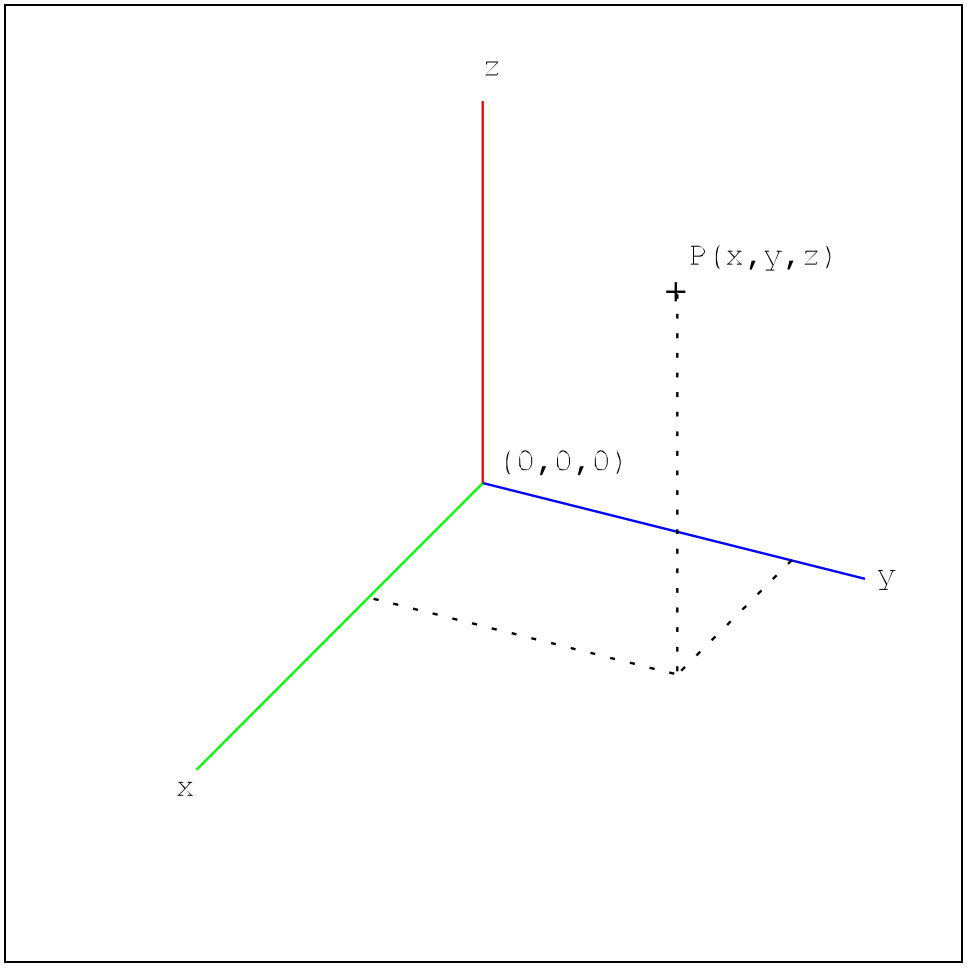

Figure 313 Cartesian coordinate system

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Cylindrical coordinate system

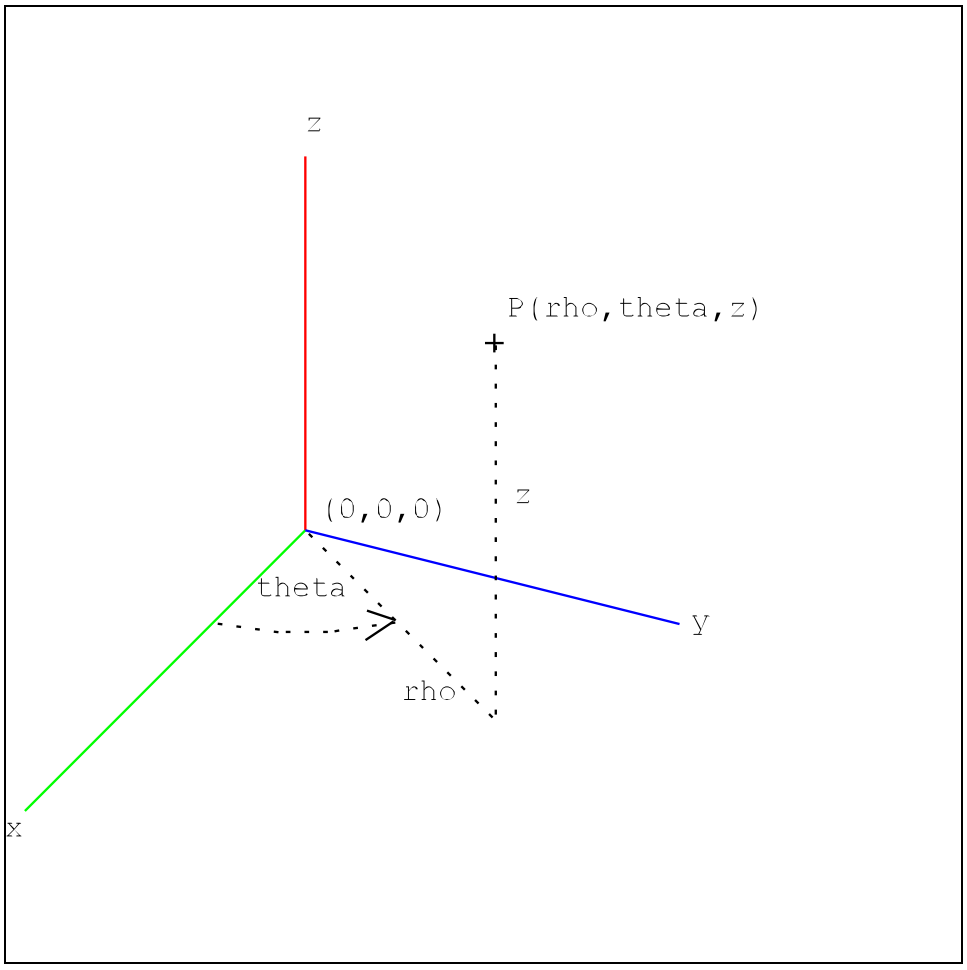

Figure 314 Cylindrical coordinate system

$$x = \cos(\theta)\rho$$

$$y = \sin(\theta)\rho$$

$$z = z$$

where ρ is radius, Θ is angle in x-y plane

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Spherical coordinate system

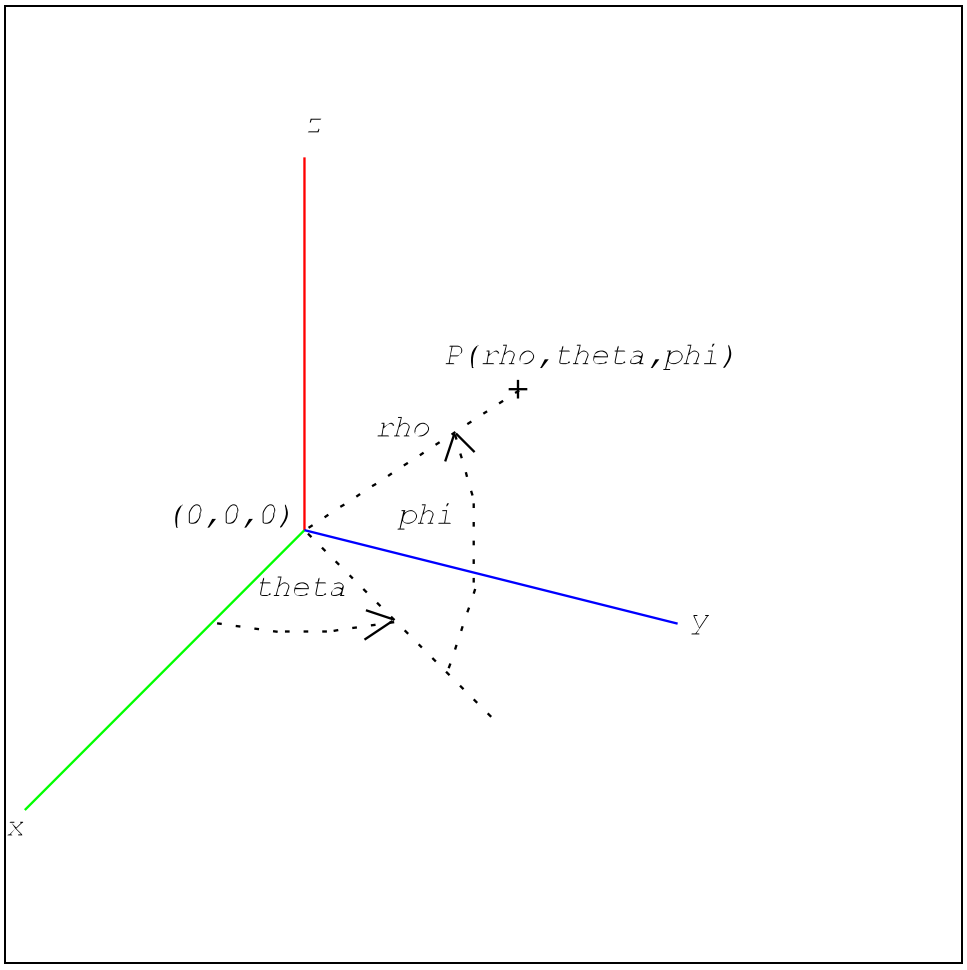

Figure 315 Spherical coordinate system

$$x = \cos(\theta)\cos(\varphi)\rho$$

$$y = \sin(\theta)\cos(\varphi)\rho$$

$$z = \sin(\varphi)\rho$$

where
ρ is radius, Θ is angle in x-y plane, Φ is angle from x-y plane

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