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function out=beam3kustom(mode,b,c,d,e)

% BEAM3 does as listed below. It is an Euler-Bernoulli
% beam/rod/torsion model.
% Beam properties (bprops) are in the order
% bprops=[E G rho A1 A2 A3 J1 J2 J3 Ixx1 Ixx2 Ixx3 Iyy1 Iyy2 Iyy3]
% Third node is in the middle.
% Fourth "node" defines the beam y plane and is actually from the
% points array.
%%
% Defining beam element properties in wfem input file:
% element properties
%   E G rho A1 A2 A3 J1 J2 J3 Izz1 Izz2 Izz3 Iyy1 Iyy2 Iyy3
% Torsional rigidity, $J$, must be less than or equal
% to $Iyy+Izz$ at any given cross section.
%
% Defining beam3 element in wfem input file:
%   node1 node2 node3 pointnumber materialnumber

%
% See wfem.m for more explanation.
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Variables (global):
% -----
% K      :   Global stiffness matrix
% Ks     :   Global stiffness buckling matrix
% M      :   Global mass matrix
% nodes  :   [x y z] nodal locations
global ismatnewer
global K
global Ks
global M
global nodes % Node locations
global elprops
global element
global points
global Fepsn % Initial strain "forces".
global lines
global restart
global reload
global curlineno
global DoverL
global surfs
%
% Variables (local):
% -----
% bnodes :   node/point numbers for actual beam nodes 1-2-3 and point
% k       :   stiffness matrix in local coordiates
% kg      :   stiffness matrix rotated into global coordinates
% m       :   mass matrix in local coordiates
% mg      :   mass matrix rotated into global coordinates
%

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[illegible]

[illegible]

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% Define beam node locations for easy later referencing
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%

x1=nodes(bnodes(1),1);
y1=nodes(bnodes(1),2);
z1=nodes(bnodes(1),3);
x2=nodes(bnodes(2),1);
y2=nodes(bnodes(2),2);
z2=nodes(bnodes(2),3);
x3=points(bnodes(3),1);
y3=points(bnodes(3),2);
z3=points(bnodes(3),3);

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Shape functions for higher order beam.
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
% Shape functions in matrix polynomial form (polyval style) for bending
bn1 = [0.25 0 -0.75 0.5];
bn1d = [0.75 0 -0.75];
bn1dd = [1.5 0];
bn2 = [0.25 -0.25 -0.25 0.25];
bn2d = [0.75 -0.5 -0.25];
bn2dd = [1.5 -0.5];
bn3 = [-0.25 0 0.75 0.5];
bn3d = [-0.75 0 0.75];
bn3dd = [-1.5 0];
bn4 = [0.25 0.25 -0.25 -0.25];
bn4d = [0.75 0.5 -0.25];
bn4dd = [1.5 0.5];

% Shape functions in matrix polynomial form (polyval style) for
% torsion/rod
rn1= [-0.5 0.5];
rn1d= [-0.5];
rn2= [.5 .5];
rn2d= [0.5];
numbeamgauss=5; % Number of Gauss points for integration of beam
element
[bgpts,bgpw]=gauss(numbeamgauss);
kb1=zeros(4,4); % For this beam, 2 nodes, 2DOF each, is a 4 by 4
                % matrix.
kb2=kb1; %Stiffness matrix for the x-z plane beam element.
l=norm([x2 y2 z2]-[x1 y1 z1]);
propertynum=num2str(element(elnum).properties);
% Allowable aspect ratio. I recommend D/l=.1
if isempty(DoverL)==1
    DoverL=.1;
end
%Euler bernoulli beams must be slender. Warn if not.

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if sqrt(A1*4/pi)/l>DoverL|sqrt(A2*4/pi)/l>DoverL
    warndlg({'Dimensions of element ' num2str(elnum) ' using properties
'...
        propertynum ' are more suitable for a Timoshenko beam.'];...
        'radius divided by length is too large'},...
        'Improper application of element.','replace')
end
% This took some work, but provide bounds on other values.
if (Izz1+Iyy1)<(1/2.1*A1^2/pi)|(Izz2+Iyy2)<(1/2.1*A2^2/pi)
    %2.0 would be exact for a circle
    warndlg({'Iyy+Izz for properties number' propertynum ' can't be as
'...
        'low as have been given.'];...
        'Nonphysical properties.'},['Impossible cross sectional' ...
        ' properties'],'replace')
end
slenderness=min([sqrt((Izz1+Iyy1)/A1) sqrt((Izz2+Iyy2)/A2)])/l;
% Check if this is a beam or something so thin that its really a
% string.
if slenderness<.002
    disp([num2str(elnum) ['is a ridiculously thin element. Please' ...
        ' check numbers.']] )
end

Jac=1/2;% Beam Jacobian. valid only if node three is in the
        % middle of the beam. Luck for us, it always is (or the
        % code yells at you)
        % Local Bending in x-y plane
for i=1:numbeamgauss
    beamsfs=[polyval(bn1dd,bgpts(i))/Jac^2;%evaluating second
                                                    %derivatives of shape
                                                    %functions to use in
                                                    %generating stiffness
                                                    %matrix. (at gauss point)

        polyval(bn2dd,bgpts(i))/Jac;
        polyval(bn3dd,bgpts(i))/Jac^2;
        polyval(bn4dd,bgpts(i))/Jac];
    Izz=polyval(rn1*Izz1+rn2*Izz2,bgpts(i));%Find Izz at
                                                    %Gauss point
    kb1=kb1+bgpw(i)*beamsfs*beamsfs'*Izz*E*Jac;%This is the Gauss
                                                    %integration part.
end
% Local Bending in x-z plane
for i=1:numbeamgauss
    beamsfs=[polyval(bn1dd,bgpts(i))/Jac^2;
        -polyval(bn2dd,bgpts(i))/Jac;
        polyval(bn3dd,bgpts(i))/Jac^2;
        -polyval(bn4dd,bgpts(i))/Jac];
    Iyy=polyval(rn1*Iyy1+rn2*Iyy2,bgpts(i));
    kb2=kb2+bgpw(i)*beamsfs*beamsfs'*Iyy*E*Jac;
end

% Local Extension in x, torsion about x
numrodgauss=3;% Number of points to use for gauss point integration

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[rgpts,rgpw]=gauss(numrodgauss);
krod=zeros(2,2);
ktor=zeros(2,2);
for i=1:numrodgauss
    rodsfs=[polyval(rn1d,rgpts(i))/Jac;
            polyval(rn2d,rgpts(i))/Jac];
    if (J1>(Iyy1+Izz1))|(J2>(Iyy2+Izz2))
        if (J1>(Iyy1+Izz1))
            disp('WARNING: J1 must be <= Iyy1+Izz1')%More checks for reality
        end
        if (J2>(Iyy2+Izz2))
            disp('WARNING: J2 must be <= Iyy2+Izz2')%More checks for reality
        end
        disp(['Error in element properties number '...
            num2str(element(elnum).properties) ...
            'used by element ' num2str(elnum) ' on line'...
            num2str(element(elnum).lineno) '.'])
    end
    J=polyval(rn1*J1+rn2*J2,bgpts(i));% J at gauss point.
    A=polyval(rn1*A1+rn2*A2,bgpts(i));% A at gauss point
    krod=krod+rgpw(i)*rodsfs*rodsfs'*A*E*Jac;%Since the shape
                                                %functions and Gauss
                                                %points are the same,
                                                %we are doing the rod
                                                %and torsion rod
                                                %together.

    ktor=ktor+rgpw(i)*rodsfs*rodsfs'*J*G*Jac;
end

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Derivation of Mass matrices
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
numbeamgauss=numbeamgauss+3; %Need more gauss points for the mass
                                %matrix.
[bgpts,bgpw]=gauss(numbeamgauss);
mb1=zeros(4,4); %initialize empty mass matrix
% Local Bending in x-y plane
for i=1:numbeamgauss
    beamsfs=[polyval(bn1,bgpts(i));
            polyval(bn2,bgpts(i))*Jac;
            polyval(bn3,bgpts(i));
            polyval(bn4,bgpts(i))*Jac];
    A=polyval(rn1*A1+rn2*A2,bgpts(i));
    mb1=mb1+bgpw(i)*beamsfs*beamsfs'*rho*A*Jac;%pause, and reflect
                                                %(OK, this was for
debugging)
end

% Local Bending in x-z plane
mb2=zeros(4,4);
for i=1:numbeamgauss
    beamsfs=[polyval(bn1,bgpts(i));

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        -polyval(bn2,bgpts(i))*Jac;
        polyval(bn3,bgpts(i));
        -polyval(bn4,bgpts(i))*Jac];
    A=polyval(rn1*A1+rn2*A2,bgpts(i));
    mb2=mb2+bgpw(i)*beamsfs*beamsfs'*rho*A*Jac;
end

% Local Extension in x, torsion about x
numrodgauss=numrodgauss+1; %Need more gauss points for the mass
                             %matrix.
[rgpts,rgpw]=gauss(numrodgauss);
mrod=zeros(2,2); %initialize empty mass matrix
mtor=zeros(2,2);
for i=1:numrodgauss
    rodsfs=[polyval(rn1,rgpts(i));
            polyval(rn2,rgpts(i))];
    J=polyval(rn1*(Iyy1+Izz1)+rn2*(Iyy2+Izz2),bgpts(i));
    A=polyval(rn1*A1+rn2*A2,bgpts(i));
    mrod=mrod+rgpw(i)*rodsfs*rodsfs'*A*rho*Jac;
    mtor=mtor+rgpw(i)*rodsfs*rodsfs'*J*rho*Jac;
end

% Assembling each stiffness matrix into the complete elemental
% stiffness matrix. We're just telling the sub-elements to be put
% into the correct spots for the total element.
k=zeros(12,12);
k([2 6 8 12],[2 6 8 12])=kb1;
k([3 5 9 11],[3 5 9 11])=kb2;
k([1 7],[1 7])=krod;
k([4 10],[4 10])=ktor;

% Assembling each mass matrix into the complete elemental
% mass matrix
m=zeros(12,12);
m([2 6 8 12],[2 6 8 12])=mb1;
m([3 5 9 11],[3 5 9 11])=mb2;
m([1 7],[1 7])=mrod;
m([4 10],[4 10])=mtor;

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Coordinate rotations
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%

R1=([x2 y2 z2]-[x1 y1 z1]);% Vector along element
lam1=R1/norm(R1);% Unit direction
R2=([x3 y3 z3]-[x1 y1 z1]);% Unit direction to point
R2perp=R2-dot(R2,lam1)*lam1;% Part of R2 perpendicular to lam1
udirec=0;
while norm(R2perp)<10*eps% If R2perp is too small, (point in line
                        % with element, we need to cover the
                        % users a$$ and generate a point that
                        % isn't. We should put out a warning,

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                                % but I commented it out.
    udirec=udirec+1;
    %disp('oops'); %This was my warning.
    %pause
    [minval,minloc]=min(lam1);
    R2perp=zeros(1,3);
    R2perp(udirec)=1;
    R2perp=R2perp-dot(R2perp,lam1)*lam1;
end
%Make the unit direction vectors for rotating and put them in the
%rotation matrix.
lam2=R2perp/norm(R2perp);
lam3=cross(lam1,lam2);
lamloc=[lam1;lam2;lam3];
lam=sparse(12,12);
lam(1:3,1:3)=lamloc;
lam(4:6,4:6)=lamloc;
lam(7:9,7:9)=lamloc;
lam(10:12,10:12)=lamloc;
%   lam(13:15,13:15)=lamloc;
%   lam(16:18,16:18)=lamloc;

% $$$      lam=[lamloc z z z z z;
% $$$          z lamloc z z z z;
% $$$          z z lamloc z z z;
% $$$          z z z lamloc z z;
% $$$          z z z z lamloc z;
% $$$          z z z z z lamloc];
element(elnum).lambda=lam;
element(elnum).m=m;
element(elnum).k=k;

kg=lam'*k*lam;
mg=lam'*m*lam;

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Assembling matrices into global matrices
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%

bn1=bnodes(1);bn2=bnodes(2);
indices=[bn1*6+(-5:0) bn2*6+(-5:0)] ;

K(indices,indices)=K(indices,indices)+kg;
M(indices,indices)=M(indices,indices)+mg;

% At this point we also know how to draw the element (what lines
% and surfaces exist). For the beam3 element, 2 lines are
% appropriate. Just add the pair of node numbers to the lines
% array and that line will always be drawn.
numlines=size(lines,1);
lines(numlines+1,:)=[bn1 bn2];

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%If I have 4 nodes that I want to use to represent a surface, I
%do the following.
panelcolor=[1 0 1];% This picks a color. You can change the
                % numbes between 0 and 1.
%Don't like this color? Use colorui to pick another one. Another
%option is that if we can't see the elements separately we can
%chunk up  $x*y*z$ , divide by  $x*y*x$  of element, see if we get
%integer powers or not to define colors that vary by panel.


% You need to uncomment this line and assign values to node1,
% node2, node3, and node4 in order to draw A SINGLE SURFACE. For
% a brick, you need 6 lines like this.
%surfs=[surfs;node1 node2 node3 node4 panelcolor];


%Each surface can have a different color if you like. Just change
%the last three numbers on the row corresponding to that
%surface.


%diag(M)
elseif strcmp(mode,'istrainforces')
    % You don't need this
    % We need to have the stiffness matrix and the coordinate roation
    matrix.


elseif strcmp(mode,'draw')
elseif strcmp(mode,'buckle')
end

```