

```

function [DEFL,REACT,ELE_FOR,AFLAG] = ud_3d1el(...

nnodes,coord,concen,fixity,nele,ends,A,Izz,Iyy,J,Cw,ISym,Ysc,Zsc,Betay,Betaz,Betaw
,Zzz,Zyy,Ayy,Azz,...
    E,v,Fy,YldSurf,Wt,webdir,beta_ang,w,thermal,truss,anatype);

% Code developed by Mrunmayi Mungekar and Devasmit Dutta
%
% UD_3D1EL performs a user defined three-dimensional
% first-order elastic analysis of a structural system.
%
%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Functions Called
%
% List of all functions called within this file and their purpose:
%
% MD_member_id - returns a matrix that lists the degree of freedom
%                for each element
%
% MD_concen_load_dof - returns a matrix that lists the
concentrated
%                    load for each degree of freedom
%
% MD_computeMemberFEFs - returns a vector of member forces and
%                        moments for a member with no end flexure
%                        releases
%
% MD_computeMemberFEFs_1stnode_MyMz_release - returns a vector of
%                member forces and moments for a member
with
%                a flexure release at the start node
%
% MD_computeMemberFEFs_2ndnode_MyMz_release - returns a vector of
%                member forces and moments for a member
with
%                a flexure release at the finish node
%
% MD_computeMemberFEFs_bothnode_MyMz_release - returns a vector of
%                member forces and moments for a member
with
%                flexure releases at both ends
%
% MD_estiff - returns a member's local stiffness matrix for a
%            member with no end flexure releases
%
% MD_estiff_1stnode_MyMz_release - returns a member's local
stiffness
%            matrix for a member with a flexure release at the
start
%            node
%
% MD_estiff_2ndnode_MyMz_release - returns a member's local
stiffness
%            matrix for a member with a flexure release at the
finish
%            node
%
% MD_estiff_bothnode_MyMz_release - returns a member's local
stiffness
%            matrix for a member with flexure releases at both
ends
%
% MD_etran - returns a member's local to global transformation
matrix
%
% MD_estiff - returns a member's local stiffness matrix for a
%            member with no end flexure releases
%
% MD_estiff_1stnode_MyMz_release - returns a member's local
stiffness

```

```

%               matrix for a member with a flexure release at the
start
%               node
%               MD_estiff_2ndnode_MyMz_release - returns a member's local
stiffness
%               matrix for a member with a flexure release at the
finish
%               node
%               MD_estiff_bothnode_MyMz_release - returns a member's local
stiffness
%               matrix for a member with flexure releases at both
ends
%

% Dictionary of Variables
%   Input Information:
%       nnodes      == total number of nodes
%       coord(i,1:3) == node i's coordinates
%                   coord(i,1) = X coordinate
%                   coord(i,2) = Y coordinate
%                   coord(i,3) = Z coordinate
%       concen(i,1:6) == concentrated loads for node i's 6 d.o.f.
%                   concen(i,1) = force in global X direction
%                   concen(i,2) = force in global Y direction
%                   concen(i,3) = force in global Z direction
%                   concen(i,4) = moment about global X axis
%                   concen(i,5) = moment about global Y axis
%                   concen(i,6) = moment about global Z axis
%       fixity(i,1:6) == prescribed displacements for node i's 6 d.o.f.
%                   Note: A free d.o.f. will have a value of NaN
%                   and hence, you will find the Matlab function
%                   isnan very useful.
%                   Examples: If fixity(15,3) is set to NaN, then node 15's
%                           Z-disp component is free;
%                           If fixity(2,6) is set to 0.0, then node 2's
%                           Z-rotation component is supported;
%                           If fixity(5,2) is set to -2.1, then node 5's
%                           Y-disp component is supported and defined
%                           with a settlement of -2.1 units.
%                   fixity(i,1) = prescribed disp. in global X direction
%                   fixity(i,2) = prescribed disp. in global Y direction
%                   fixity(i,3) = prescribed disp. in global Z direction
%                   fixity(i,4) = prescribed rotation about global X axis
%                   fixity(i,5) = prescribed rotation about global Y axis
%                   fixity(i,6) = prescribed rotation about global Z axis
%       nele        == total number of elements
%       ends(i,1:14) == element i's nodal information
%                   ends(i,1) = start node #
%                   ends(i,2) = finish node #
%                   ends(i,3) = flag to indicate whether or not flexural
%                           moments are released at start node.  ends(i,3)=0 both
not
%                   released (rigid connection); ends(i,3)=1 both flexural
%                   moments are released (pinned connection); ends(i,3)=2
%                   at least one of the flexural moments are partially or
fully
%                   released (see below for connection stiffness
attributes)
%                   ends(i,4) = flag to indicate whether or not flexural

```

```

%           moments are released at finish node.  ends(i,4)=0 both
not
%           released (rigid connection); ends(i,4)=1 both flexural
%           moments are released (pinned connection); ends(i,4)=2
%           at least one of the flexural moments are partially or
fully
%           released (see below for connection stiffness
attributes)
%           ends(i,5) = flag to indicate the degree of warping
%           restraint at start node.  ends(i,5)=0 warping free;
%           ends(i,5)=1 warping fixed; ends(i,5)=2 warping
continuous
%           ends(i,6) = flag to indicate the degree of warping
%           restraint at finish node.  ends(i,6)=0 warping free;
%           ends(i,6)=1 warping fixed; ends(i,6)=2 warping
continuous
%           ends(i,7) = rotational spring stiffness at the start
%           node and about element i's local z-z axis.
%           ends(i,8) = rotational spring stiffness at the start
%           node and about element i's local y-y axis.
%           ends(i,9) = rotational spring stiffness at the finish
%           node and about element i's local z-z axis.
%           ends(i,10) = rotational spring stiffness at the finish
%           node and about element i's local y-y axis.
%           ends(i,11) = connection moment capacity Mpz at the
start
%           node and about element i's local z-z axis.
%           ends(i,12) = connection moment capacity Mpy at the
start
%           node and about element i's local y-y axis.
%           ends(i,13) = connection moment capacity Mpz at the
finish
%           node and about element i's local z-z axis.
%           ends(i,14) = connection moment capacity Mpy at the
finish
%           node and about element i's local y-y axis.
%           A(i) == element i's cross sectional area
%           Izz(i) == element i's moment of inertia about its local z-z axis
%           Iyy(i) == element i's moment of inertia about its local y-y axis
%           J(i) == element i's torsional constant
%           Cw(i) == element i's warping constant
%           Zzz(i) == element i's plastic section modulus about its local z-z
axis
%           Zyy(i) == element i's plastic section modulus about its local y-y
axis
%           Ayy(i) == element i's effective shear area along its local y-y
axis
%           Azz(i) == element i's effective shear area along its local z-z
axis
%           E(i) == element i's material elastic modulus, Young's Modulus
%           v(i) == element i's material Poisson's ratio
%           Fy(i) == element i's material yield strength
%           YldSurf(i) == element i's yield surface maximum values
%                   YldSurf(i,1) = maximum P/Py value
%                   YldSurf(i,2) = maximum Mz/Mpz value
%                   YldSurf(i,3) = maximum My/Mpy value
%           Wt(i) == element i's material weight density
%           (Assume that gravity is directed in the negative global
Y dir)

```

```

%      webdir(i,1:3) == element i's unit web vector. This is a unit vector
%                      that defines the element's local y-y axis with respect
%                      to the global coordinate system. It is based on the
%                      structure's undeformed geometry.
%                      webdir(i,1) = x component of element's unit web
vector
%                      webdir(i,2) = y component of element's unit web
vector
%                      webdir(i,3) = z component of element's unit web
vector
%
%      NOTE: An element's 3x3 rotation matrix, [g], is
constructed
%
%      as follows: First, calculate a unit vector, x_vect, that
%      describes the element's local x-axis. Second, take the
%      cross product of x_vect and webdir(i,:) to obtain
z_vect,
%
%      i.e. z_vect = cross(x_vect,webdir(i,:)). Third, set
z_vect
%
%      to a unit vector, i.e. z_vect = z_vect/norm(z_vect).
%      Finally, the first row of [g] is x_vect, its second row
is
%      webdir(i,:), and its third row is z_vect.
%      beta_ang(i) == element i's web rotation angle. These values are
%                  provided for those students who are required to
calculate
%
%      their own unit web vectors (see above). It is based
%      on the structure's undeformed geometry.
%      Note: MASTAN2 uses the following convention for
%            defining a member's default web orientation:
%            A vector defing the element's local y-axis
%            with respect to the global coordinate system
%            will have a positive component in the global
%            Y direction. If the element's local x-axis,
%            its length axis, is aligned with the global Y
%            axis, then element's local y-axis is aligned
%            with global negative X axis. After this initial
%            orientation, element i may be rotated about
%            its local x-axis by the amount defined by
%            its web rotation angle, beta_ang(i). The
%            angle is in radians and assumes a right-hand
%            convention about the local x-axis which runs from
%            the element's start node to its finish node.
%
%      w(i,1:3) == element i's uniform load which references its
%                  local coordinate system
%                  w(i,1) = x component of uniform load
%                  w(i,2) = y component of uniform load
%                  w(i,3) = z component of uniform load
%
%      thermal(i,1:4) == element i's thermal strain effects which reference its
%                       local coordinate system
%                       thermal(i,1) = coefficient of thermal expansion
%                       thermal(i,2) = change in temperature at centroid
%                       thermal(i,3) = linear temperature gradient in local
y-dir
%
%                       = (T_up_y - T_btm_y) / depth_y
%                       thermal(i,4) = linear temperature gradient in local
z-dir
%
%                       = (T_up_z - T_btm_z) / width_z
%
%      truss == flag to indicate if structure is a truss or not
%            truss = 0 System is not a truss

```

```

%          truss = 1    System is a truss
%          anatype == flag to indicate which type of analysis is requested
%          anatype = 1  First-Order Elastic
%          anatype = 2  Second-Order Elastic
%          anatype = 3  First-Order Inelastic
%          anatype = 4  Second-Order Inelastic
%          anatype = 5  Elastic Buckling (Eigenvalue)
%          anatype = 6  Inelastic Buckling (Eigenvalue)
%
% Local Information:
%      < to be defined by the student >
%
% Output Information:
%      DEFL(i,1:6) == node i's calculated 6 d.o.f. deflections
%                  DEFL(i,1) = displacement in X direction
%                  DEFL(i,2) = displacement in Y direction
%                  DEFL(i,3) = displacement in Z direction
%                  DEFL(i,4) = rotation about X direction
%                  DEFL(i,5) = rotation about Y direction
%                  DEFL(i,6) = rotation about Z direction
%      REACT(i,1:6) == reactions for supported node i's 6 d.o.f.
%                  REACT(i,1) = force in X direction
%                  REACT(i,2) = force in Y direction
%                  REACT(i,3) = force in Z direction
%                  REACT(i,4) = moment about X direction
%                  REACT(i,5) = moment about Y direction
%                  REACT(i,6) = moment about Z direction
%      ELE_FOR(i,1:12) == element i's internal forces and moments
%                      Note: All values reference the element's local
%                           coordinate system.
%                      ELE_FOR(i,1) = x-force at start node
%                      ELE_FOR(i,2) = y-force at start node
%                      ELE_FOR(i,3) = z-force at start node
%                      ELE_FOR(i,4) = x-moment at start node
%                      ELE_FOR(i,5) = y-moment at start node
%                      ELE_FOR(i,6) = z-moment at start node
%                      ELE_FOR(i,7) = x-force at end node
%                      ELE_FOR(i,8) = y-force at end node
%                      ELE_FOR(i,9) = z-force at end node
%                      ELE_FOR(i,10) = x-moment at end node
%                      ELE_FOR(i,11) = y-moment at end node
%                      ELE_FOR(i,12) = z-moment at end node
%                      If you are not programming warping torsion, the
ELE_FOR
%                      array needs to contain only 12 columns, i.e.
ELE_FOR(i,1:12)
%                      For those programming warping torsion, the bimoments
and
%                      rates of twist should be stored as follows.
%                      ELE_FOR(i,13) = bimoment at start node
%                      ELE_FOR(i,14) = bimoment at end node
%                      ELE_FOR(i,15) = rate of twist at start node
%                      ELE_FOR(i,16) = rate of twist at end node
%      AFLAG == logical flag to indicate if a successful
%             analysis has been completed
%             AFLAG = 1    Successful
%             AFLAG = 0    Unstable Structure
%             AFLAG = inf  No analysis code available
%

```

```

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%
% Start by defining all output arrays to be empty
%
DEFL=[]; REACT=[]; ELE_FOR=[];

% Create array that list degree of freedom for each element
memb_id = MD_member_id(nnodes, nele, ends);

disp('member id');
disp(memb_id);

% Create array that list concentrated load for degree of freedom
concen_applied_load_dof = MD_concen_load_dof(concen, nnodes);

% Create array that list member forces for each element
FEF = zeros(6*nnodes,1);

for i =1:nele
    start_node = ends(i,1);
    end_node = ends(i,2);
    start_coord = coord(start_node,:);
    end_coord = coord(end_node,:);
    L = norm(end_coord - start_coord);

    % Check for end flexure release condition and use appropriate function for local
member forces
    if ends(i,3) == 0 && ends(i,4) == 0
        memberlocalFEF = MD_computeMemberFEFs(w(i,:),L);

    elseif ends(i,3) == 1 && ends(i,4) == 0
        memberlocalFEF = MD_computeMemberFEFs_1stnode_MyMz_release(w(i,:),L);

    elseif ends(i,3) == 0 && ends(i,4) == 1
        memberlocalFEF = MD_computeMemberFEFs_2ndnode_MyMz_release(w(i,:),L);

    elseif ends(i,3) == 1 && ends(i,4) == 1
        memberlocalFEF = MD_computeMemberFEFs_bothnode_MyMz_release(w(i,:),L);
    end

    % Transform local member forces to global member forces
    gamma = MD_etran(start_coord,end_coord,webdir(i,:));
    memberglobalFEF = gamma'*memberlocalFEF;
    FEF(memb_id(i,:),1) = memberglobalFEF + FEF(memb_id(i,:),1);

end

disp('concen_applied_load_dof');
disp(concen_applied_load_dof);

disp('FEF');
disp(FEF);

% Initiate deflection vector for free DOFs
D = fixity';
D = D(:);
freeDOF = find(isnan(D));

```

```

supportDOF = D == 0;
displacedDOF = find(D~=0 & ~isnan(D));

% Create global stiffness matrix
kstructureglobal = zeros(6*nnodes,6*nnodes);
for i =1:nele
    start_node = ends(i,1);
    end_node = ends(i,2);
    start_coord = coord(start_node,:);
    end_coord = coord(end_node,:);
    L = norm(end_coord - start_coord);

    % Check for end flexure release condition and use appropriate function for local
    member stiffness, convert each to global
    if ends(i,3) == 0 && ends(i,4) == 0
        kele_local = MD_estiff(A(i), Izz(i), Iyy(i), J(i), Ayy(i), Azz(i), E(i),
v(i), L);

        elseif ends(i,3) == 1 && ends(i,4) == 0
            kele_local = MD_estiff_1stnode_MyMz_release(A(i), Izz(i), Iyy(i), J(i),
Ayy(i), Azz(i), E(i), v(i), L);

        elseif ends(i,3) == 0 && ends(i,4) == 1
            kele_local = MD_estiff_2ndnode_MyMz_release(A(i), Izz(i), Iyy(i), J(i),
Ayy(i), Azz(i), E(i), v(i), L);

        elseif ends(i,3) == 1 && ends(i,4) == 1
            kele_local = MD_estiff_bothnode_MyMz_release(A(i), E(i), L);
        end

    % Transform local member stiffness to global member stiffness
    gamma = MD_etran(start_coord,end_coord,webdir(i,:));
    kele_global = gamma'*kele_local*gamma;
    kstructureglobal(memb_id(i,:),memb_id(i,:)) = kele_global +
kstructureglobal(memb_id(i,:),memb_id(i,:));

end

% Separate global stiffness matrix into free, support and displaced DOF blocks
Kff = kstructureglobal(freeDOF,freeDOF);
Kfn = kstructureglobal(freeDOF,displacedDOF);
Knf = kstructureglobal(displacedDOF,freeDOF);
Ksf = kstructureglobal(supportDOF,freeDOF);
Ksn = kstructureglobal(supportDOF,displacedDOF);
Knn = kstructureglobal(displacedDOF,displacedDOF);

% Solve for free DOF deflections
Pf = concen_applied_load_dof(freeDOF);
FEFf = FEF(freeDOF);
FEFs = FEF(supportDOF);
FEFn = FEF(displacedDOF);
Delta_n = D(displacedDOF);
Delta_f = (Kff)\(Pf - FEFf - Kfn*Delta_n);

disp('Delta_f');
disp(vpa(Delta_f,6));

D_all = zeros(nnodes*6,1); % Initiate with size
D_all(displacedDOF) = Delta_n;

```

```

D_all(freeDOF) = Delta_f;

% Populate start and end internal forces for each element
for i =1:nele
    start_node = ends(i,1);
    end_node = ends(i,2);
    start_coord = coord(start_node,:);
    end_coord = coord(end_node,:);
    L = norm(end_coord - start_coord);

    % Check for end flexure release condition and use appropriate function for local
    member stiffness
    if ends(i,3) == 0 && ends(i,4) == 0
        kele_local = MD_estiff(A(i), Izz(i), Iyy(i), J(i), Ayy(i), Azz(i), E(i),
v(i), L);

        elseif ends(i,3) == 1 && ends(i,4) == 0
            kele_local = MD_estiff_1stnode_MyMz_release(A(i), Izz(i), Iyy(i), J(i),
Ayy(i), Azz(i), E(i), v(i), L);

        elseif ends(i,3) == 0 && ends(i,4) == 1
            kele_local = MD_estiff_2ndnode_MyMz_release(A(i), Izz(i), Iyy(i), J(i),
Ayy(i), Azz(i), E(i), v(i), L);

        elseif ends(i,3) == 1 && ends(i,4) == 1
            kele_local = MD_estiff_bothnode_MyMz_release(A(i), E(i), L);
        end

    % Get transformation matrix
    gamma = MD_etran(start_coord,end_coord,webdir(i,:));

    % Get local DOF displacement
    Dele_global = D_all(memb_id(i,:));
    Dele_local = gamma*Dele_global;

    % Check for end flexure release condition and use appropriate function for local
    member forces
    if ends(i,3) == 0 && ends(i,4) == 0
        memberlocalFEF = MD_computeMemberFEFs(w(i,:),L);

        elseif ends(i,3) == 1 && ends(i,4) == 0
            memberlocalFEF = MD_computeMemberFEFs_1stnode_MyMz_release(w(i,:),L);

        elseif ends(i,3) == 0 && ends(i,4) == 1
            memberlocalFEF = MD_computeMemberFEFs_2ndnode_MyMz_release(w(i,:),L);

        elseif ends(i,3) == 1 && ends(i,4) == 1
            memberlocalFEF = MD_computeMemberFEFs_bothnode_MyMz_release(w(i,:),L);
        end

    localMemberForces = kele_local*Dele_local + memberlocalFEF;

    % Populate start and end internal forces for each element
    ELE_FOR(i,:) = localMemberForces';

end

% Create a matrix of internal forces variation along the length of each member as a
function of local x coordinate

```



```

syms("internal_forces_variation", [nele,6]);
for i =1:nele

    % Compute the force variation based on forces of start node
    syms x;
    internal_forces_variation(i,1:3) = - ELE_FOR(i,1:3) - w(i,1:3)*x;
    internal_forces_variation(i,4) = - ELE_FOR(i,4);
    internal_forces_variation(i,5) = - ELE_FOR(i,5) + ELE_FOR(i,3)*x + w(i,3)*x^2/2;
    internal_forces_variation(i,6) = - ELE_FOR(i,6) + ELE_FOR(i,2)*x + w(i,2)*x^2/2;

    % Plot graphs if necessary

    % figure(i+1);
    % subplot(2,3,1);
    % fplot(internal_forces_variation(i,1),[0,L]);
    % val = max(subs(internal_forces_variation(i,1), linspace(0,L,100)));
    % if( val<1 && val>-1)
    %     ylim([-1,1]);
    % end
    % title('P variation');
    %
    % subplot(2,3,2);
    % fplot(internal_forces_variation(i,2),[0,L]);
    % val = max(subs(internal_forces_variation(i,2), linspace(0,L,100)));
    % if( val<1 && val>-1)
    %     ylim([-1,1]);
    % end
    % title('Vy variation');
    %
    % subplot(2,3,3);
    % fplot(internal_forces_variation(i,3),[0,L]);
    % val = max(subs(internal_forces_variation(i,3), linspace(0,L,100)));
    % if( val<1 && val>-1)
    %     ylim([-1,1]);
    % end
    % title('Vz variation');
    %
    % subplot(2,3,4);
    % fplot(internal_forces_variation(i,4),[0,L]);
    % val = max(subs(internal_forces_variation(i,4), linspace(0,L,100)));
    % if( val<1 && val>-1)
    %     ylim([-1,1]);
    % end
    % title('T variation');
    %
    % subplot(2,3,5);
    % fplot(internal_forces_variation(i,5),[0,L]);
    % val = max(subs(internal_forces_variation(i,5), linspace(0,L,100)));
    % if( val<1 && val>-1)
    %     ylim([-1,1]);
    % end
    % title('My variation');
    %
    % subplot(2,3,6);
    % fplot(internal_forces_variation(i,6),[0,L]);
    % val = max(subs(internal_forces_variation(i,6), linspace(0,L,100)));
    % if( val<1 && val>-1)
    %     ylim([-1,1]);
    % end

```

```

% title('Mz variation');
%
% sgtitle("Member "+i+" internal forces variation");

end

disp('Member forces variation');
disp(vpa(internal_forces_variation,6));

disp('Member forces');
disp(vpa(ELE_FOR,6));

% Compute the reactions at the supports and prescribed displacement DOFs
Rs = FEFs + Ksf*Delta_f + Ksn*Delta_n;
Rn = FEFn + Knf*Delta_f + Knn*Delta_n;

Rall = zeros(nnodes*6,1); % Initiate with size
Rall(supportDOF) = Rs;

REACT = zeros(nnodes,6);
for i = 1:nnodes
    REACT(i,:) = Rall(6*i-5:6*i);
end

disp('Rs');
disp(vpa(Rs,6));

disp('Rn');
disp(Rn);

disp('back-calculating freeDOF load');
disp(FEFF + Kff*Delta_f + Kfn*Delta_n);

AFLAG = 1;
%
% STUDENT NOTE:
%     In order for this routine to become fully active AFLAG
%     must be changed.

% Good luck CE Student!!!
%
```