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function Phi = Kepler_STM_UV(dts, varargin)
%-----
% STM_UNIVERSAL Compute 6x6 state transition matrix using universal variables
%-----
% Inputs are Variation but should be:
%   State - 6x1 initial position [m] and velocity [m/s] in cartesian
%   dt     - time of flight (t - t0) [s]
%   mu    - gravitational parameter [m^3/s^2] (optional; default Earth)
%
% Output:
%   Phi   - 6x6 state transition matrix
%
% Uses universal variable formulation (Goodyear)

% standard error
error(nargchk(3, 9, nargin)); %#ok

% parse input
narg = nargin;
time_unit = 'days';
if (narg >= 8)
    % otherwise
    r1 = [varargin{1:3}];
    v1 = [varargin{4:6}];
    muC = varargin{7};
end
if (narg == 9), time_unit = varargin{8}; end
if (narg <= 4)
    % otherwise
    r1 = varargin{1}(1:3);
    v1 = varargin{1}(4:6);
    muC = varargin{2};
end
if (narg == 4), time_unit = varargin{3}; end

% force everything to be row-matrices
r1 = r1(:).' ; v1 = v1(:).';

% initialize output
zero                  = zeros(numel(dts), 1);
exitflag              = zero;
final_states           = NaN(numel(dts), 6);
output.Kepler_iterations = zero;
output.cont_frac_iterations = zero;
output.time_error       = zero;

% progress only one orbit (to command window, debugging purposes only)
if numel(r1) ~= 3
    error('progress_orbit:only_one_body_allowed',...
          'I can only progress one orbit at a time. Please use ARRAYFUN().');
end

% times are given in days, unless specified otherwise

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if strcmpi(time_unit, 'days')
    dts = dts * 86400;
elseif ~strcmpi(time_unit, 'seconds')
    error('progress_orbit:invalid_timeunit',...
        'Only ''seconds'' and ''days'' are valid timeunits.');
end

% initialize
nu0 = r1*v1.';
r1m = sqrt(r1*r1.');
beta = 2*muC/r1m - v1*v1.';

% period effects
DeltaU = zero;
if (beta > 0)
    P = 2*pi*muC*beta^(-3/2);
    n = floor((dts + P/2 - 2*nu0/beta)/P);
    DeltaU = 2*pi*n*beta^(-5/2);
end

% loop through all requested time steps
for i = 1:numel(dts)

    % extract current time step
    dt = dts(i);

    % quick exit for trivial case
    if (dt == 0)
        final_states(i, :) = [r1,v1];
        exitflag(i) = 1;
        output.Kepler_iterations(i) = 0;
        output.cont_frac_iterations(i) = 0;
        output.time_error(i) = 0;
        continue;
    end

    % loop until convergence of the time step
    u = 0; t = 0; qisbad = false; iter = 0; cont_frac = 0; deltaT = t-dt;
    while abs(deltaT) > 1 % one second accuracy seems fine

        % increase iterations
        iter = iter + 1;

        % compute q
        % NOTE: [q] may not exceed 1/2. In principle, this will never
        % occur, but the iterative nature of the procedure can bring
        % it above 1/2 for some iterations.
        bu = beta*u*u;
        q = bu/(1 + bu);

        % escape clause;
        % The value for [q] will almost always stabilize to a value less
        % than 1/2 after a few iterations, but NOT always. In those
        % cases, just use repeated coordinate transformations
    end
end

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if (iter > 25) || (q >= 1), qisbad = true; break; end

% evaluate continued fraction (when q < 1, always converges)
A = 1; B = 1; G = 1; n = 0;
k = -9; d = 15; l = 3; Gprev = inf;
while abs(G-Gprev) > 1e-14
    k = -k; l = l + 2;
    d = d + 4*l; n = n + (l+k)*l;
    A = d/(d - n*A*q); B = (A-1)*B;
    Gprev = G; G = G + B;
    cont_frac = cont_frac + 1;
end % continued fraction evaluation

% continue kepler loop
U0w2 = 1 - 2*q;
U1w2 = 2*(1-q)*u;
U = 16/15*U1w2^5*G + DeltaU(i);
U0 = 2*U0w2^2-1;
U1 = 2*U0w2*U1w2;
U2 = 2*U1w2^2;
U3 = beta*U + U1*U2/3;
r = r1m*U0 + nu0*U1 + muC*U2;
t = r1m*U1 + nu0*U2 + muC*U3;
deltaT = t - dt;
% Newton-Raphson method works most of the time, but is
% not too stable; the method fails far too often for my
% liking...
% u = u - deltaT/4/(1-q)/r;
% Halley's method is much better in that respect. Working
% out all substitutions and collecting terms gives the
% following simplification:
u = u - deltaT/((1-q)*(4*r + deltaT*beta*u));
end % time loop

% do it the slow way if state transition matrix fails for some q
if qisbad
    % repeated coordinate transformations
    [aa, ee, ii, OO, oo, nu] = ijk2keplerian(r1, v1);

    % Compute mean anomaly
    cosE = (ee + cos(nu)) ./ (1 + ee*cos(nu));
    sinE = sqrt(1 - ee^2) .* sin(nu) ./ (1 + ee*cos(nu));
    E = atan2(sinE, cosE); % robust quadrant
    E = mod(E, 2*pi);
    M = E - e .* sin(E);

    % Update mean anomaly
    M = M + sqrt(muC/abs(aa)^3)*dt;

    % Convert true anomaly
    M = mod(M, 2*pi);
    E = M; % initial guess
    tol = 1e-12;
    diff = 1;

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while abs(diff) > tol
    diff = E - ee*sin(E) - M;
    E = E - diff ./ (1 - ee*cos(E));
end
nu = 2 * atan2( sqrt(1+ee)*sin(E/2), sqrt(1-ee)*cos(E/2) );
nu = mod(nu, 2*pi);

[x, y, z, xd, yd, zd] = keplerian2ijk(aa, ee, ii, oo, oo, M);
final_states(i, :) = [x, y, z, xd, yd, zd];
% this means failure
exitflag(i) = -1;

% use state transition matrix if all went well
else
    % Kepler solution
    f = 1 - muC/r1m*U2;      F = -muC*U1/r/r1m;
    g = r1m*U1 + nu0*U2;      G = 1 - muC/r*U2;
    % create new position and velocity matrices
    final_states(i, :) = [r1*f+v1*g, r1*F+v1*G];
    % all went fine
    exitflag(i) = 1;
end % "q is bad" clause

Phi = [eye(3)*f eye(3)*g; eye(3)*F eye(3)*G];

% process output
output.Kepler_iterations(i) = iter;
output.cont_frac_iterations(i) = cont_frac;
output.time_error(i) = abs(t-dt);

end % loop through [dt]

% generate properly formatted output
narg = nargout;
if (narg <= 3)
    varargout{1} = final_states;      % and output array
    varargout{2} = exitflag;          % exitflag
    varargout{3} = output;           % output
elseif (narg > 3)
    varargout{1} = final_states(:, 1); varargout{4} = final_states(:, 4);
    varargout{2} = final_states(:, 2); varargout{5} = final_states(:, 5);
    varargout{3} = final_states(:, 3); varargout{6} = final_states(:, 6);
    varargout{7} = exitflag;          % exitflag
    varargout{8} = output;           % output
end % process output

end % progress orbit

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