

One-dimensional Conservation Law Solver

Computing Burgers solution using a DG-FEM routine with Polynomial truncation for the nonlinear terms.

for solving the following problem:

$$u_t + f(u)_x = s(u)$$

where $f(u)$ and $S(u)$ can be For linear advection eq.: $f(u) = a*u$ and $s(u) =$ any function of u For non-linear advection: $f(u) = u^2/2$ and $s(u) =$ any function of u

Function residual will be defined as:

$$Residue(u) = -f(u)_x + s(u)$$

Based on ideas of the following papers:

1. TVB Runge-Kutta Local Projection Discontinuous Galerkin Finite Element Method for conservation laws II: General Framework. (1989)
2. Runge-Kutta Discontinuous Galerkin Method Using WENO Limiters. (2005)

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Clear Work Space

```
clear all; close all; clc;
```

Simulation Parameters

```
k          = 4;          % Space order / Number of degrees of freedom: 0 to k
np         = k+1;        % Number of points per Cell/Element
quadn      = 3;          % element grid: {1}sLeg, {2}Lobatto, {3}Leg, {4}Radau
RKs        = 3;          % Time Int. Scheme {1} no-RK, {2}TVD-RK2, {3}TVD-RK3,
```

```

                                % {4}SSP-RK2, {5}SSP-RK3, {6}SSP-RK4S5
flux_type = 3;                % {1}Roe, {2}Global LF, {3}LLF, {4}Upwind (non-conservative)
equation = 2;                 % {1} scalar advection, {2} burgers equation
include_s = 0;                % {1} include source term, {0} do NOT include source term
a = 1.0;                      % for scalar advection speed
cfl = 1/(2*k+1);              % Courant Number
tEnd = 3.10;                  % Final Time for computation
nx = 10;                      % Number of Cells/Elements
MM = 0.01;                    % TVB constant M
IC_case = 3;                  % {1} Gaussian , {2} Square, {3} sine, {4} Riemann.
plot_figs = 1;                % {1}Plot figures, {0}Do NOT plot figures
w_output = 0;                 % Write output: {1} YES please!, {2} NO

```

Define Grid Cell's (Global) nodes

Building nodes for cells/elements:

```

x_left = 0; x_right = 1; dx = (x_right-x_left)/nx;
x_nodes = x_left : dx : x_right; % cells nodes

```

flux function

```

switch equation
case{1} % Scalar advection Eq. flux:
    F = @(w) a * w;
    % and derivate of the flux function
    dF = @(w) a*ones(size(w));
case{2} % Inviscid Burgers Eq. flux:
    F = @(w) w.^2/2;
    % and derivate of the flux function
    dF = @(w) w;
end

```

Source term function

```

switch include_s
case{0} % no source term
    S = @(w) zeros(size(w));
case{1} % with source term
    % example source term
    S = @(w) w.^2;
end

```

SETUP

1. Build Cells/Elements (Local) inner points (quadrature points).
2. Build

Weighting values for our local. 3. Build Vandermonde Matrix for our local quadrature points.

```
[x,xi,w,V] = setup(k,x_nodes,quadn);

% Compute Math Objects:
if quadn == 1; bmath = 1; else bmath = 2; end;
switch bmath
    case{1} % Build Math objects for scaled Legendre polynomials. See Ref.[1]
        % M matrix
        Mcoef = [1 1/12 1/180 1/2800 1/44100 1/698544 1/11099088 1/176679360];
        M = diag(Mcoef(1:k+1));
        % invM matrix
        invM = inv(M);
        % D matrix
        Dcoef = [ ...
            0, 1, 0, (1/10), 0, (1/126), 0, (1/1716); ...
            0, 0, (1/6), 0, (1/70), 0, (1/924), 0; ...
            0, 0, 0, (1/60), 0, (1/756), 0, (1/10296); ...
            0, 0, 0, 0, (1/700), 0, (1/9240), 0; ...
            0, 0, 0, 0, 0, (1/8820), 0, (1/120120); ...
            0, 0, 0, 0, 0, 0, (1/116424), 0; ...
            0, 0, 0, 0, 0, 0, 0, (1/1585584); ...
            0, 0, 0, 0, 0, 0, 0, 0];
        D = Dcoef(1:k+1,1:k+1);
        % Scaled Legendre polynomials of deg 'l' evaluated at x = +1/2
        Ln = zeros(k+1,1); % column vector
        for l = 0:k; % Polynomials degree
            Ln(l+1) = sLegendreP(l,0.5);
        end
        % Scaled Legendre polynomials of deg 'l' evaluated at x = -1/2
        Lp = zeros(k+1,1); % column vector
        for l = 0:k; % Polynomials degree
            Lp(l+1) = sLegendreP(l,-0.5);
        end

    case{2} % Build Math objects for non-scaled Legendre polynomials See Ref.[3]
        l = (0:k)'; % all polynomials degree
        % M matrix
        M = diag(dx./(2*l+1));
        % invM matrix
        invM = inv(M);
        % D matrix
        D = zeros(np,np);
        for ll = 0:k % For all degrees of freedom
```

```

        i = ll+1;          % Dummy index
        for j=1:np         % For all local points
            if j>i && rem(j-i,2)==1
                D(i,j)=2;   % D or differentiated Legendre Matrix
            end
        end
    end
end
% Scaled Legendre polynomials of degree 'l' evaluated at x = +1
Ln = (1).^(1); % LegP @ x_{i+1/2}^(-)
% Scaled Legendre polynomials of degree 'l' evaluated at x = -1
Lp = (-1).^(1); % LegP @ x_{i-1/2}^(+)
end

```

Load Initial Condition, $u(x,0) = u_0$

```

u0 = u_zero(x,IC_case);
f0 = F(u0);
s0 = S(u0);

```

Computing the evolution of the residue ' $L(u)$ ', $du/dt = L(u)$

Load Initial conditions

```

u = u0;

% Transform u(x,t) to degrees of freedom u(t)_{l,i} for each i-Cell/Element
ut = V\u;

% Set Initial time step
t = 0; % time
n = 0; % counter
tic;
while t <= tEnd
    % Time step 'dt'
    u_reshaped = reshape(u,1,nx*np);
    dt = dx*cfl/max(abs(u_reshaped));
    t = t + dt; % iteration time / increment time
    n = n + 1; % update counter

    % Plot solution every time step
    if plot_figs == 1; plot(x,u,'o-');
        title('u_t + f(u)_x = s(u)')
        xlabel('x'); ylabel('u')
        grid on; %axis([0,1,-5,5]);
    end
end

```

```

end;

switch RKs % time integration scheme
case{1} % no integration scheme
    ut_next = ut + dt*AdvecResidue(ut...
        ,F,dF,S,Ln,Lp,V,D,invM,flux_type);

case{2} % TVD-RK2
    ut_1 = ut + dt*AdvecResidue(ut...
        ,F,dF,S,Ln,Lp,V,D,invM,flux_type);
    ut_next = 1/2*ut + 1/2*ut_1 + 1/2*dt*AdvecResidue(ut_1 ...
        ,F,dF,S,Ln,Lp,V,D,invM,flux_type);

case{3} % TVD-RK3
    ut_1 = ut + dt*AdvecResidue(ut ...
        ,F,dF,S,Ln,Lp,V,D,invM,flux_type);
    ut_2 = 3/4*ut + 1/4*ut_1 + 1/4*dt*AdvecResidue(ut_1 ...
        ,F,dF,S,Ln,Lp,V,D,invM,flux_type);
    ut_next = 1/3*ut + 2/3*ut_2 + 2/3*dt*AdvecResidue(ut_2 ...
        ,F,dF,S,Ln,Lp,V,D,invM,flux_type);

case{4} % 2nd Order SSP-RK
    u_1 = u + dt*residual(u,t);
    u_next = 1/2*(u + u_1 + dt*residual(u_1,t+dt));

case{5} % 3rd Order SSP-RK
    u_1 = u + dt*residual(u,t);
    u_2 = 1/4*(3*u + u_1 + dt*residual(u_1,t+dt));
    u_next = 1/3*(u + 2*u_2 + 2*dt*residual(u_2,t+0.5*dt));

case{6} % 5-stages, 4th-order SSP-RK
    % "It is not possible to construc a fourth-order, four-stage
    % SSP-RK schemes where all coefficients are positive." [3]
    % However, one can derive a fourth-order scheme by allowing a
    % fifth stage. The optimal scheme is given as:
    %
    % Low storage Runge-Kutta coefficients
    rk4a = [
        0.0 ...
        -567301805773.0/1357537059087.0 ...
        -2404267990393.0/2016746695238.0 ...
        -3550918686646.0/2091501179385.0 ...
        -1275806237668.0/842570457699.0];
    rk4b = [ 1432997174477.0/9575080441755.0 ...
        5161836677717.0/13612068292357.0 ...
        1720146321549.0/2090206949498.0 ...
        3134564353537.0/4481467310338.0 ...

```

```

                2277821191437.0/14882151754819.0];
rk4c = [
            0.0 ...
            1432997174477.0/9575080441755.0 ...
            2526269341429.0/6820363962896.0 ...
            2006345519317.0/3224310063776.0 ...
            2802321613138.0/2924317926251.0];
resu = 0;
for s = 1:5
    timelocal = t + rk4c(s)*dt;
    [rhsu] = residual(u,timelocal);
    resu = rk4a(s)*resu + dt*rhsu;
    u = u + rk4b(s)*resu;
end
u_next = u;

otherwise
    error ('Scheme not defined')
end

% UPDATE info
ut = ut_next;

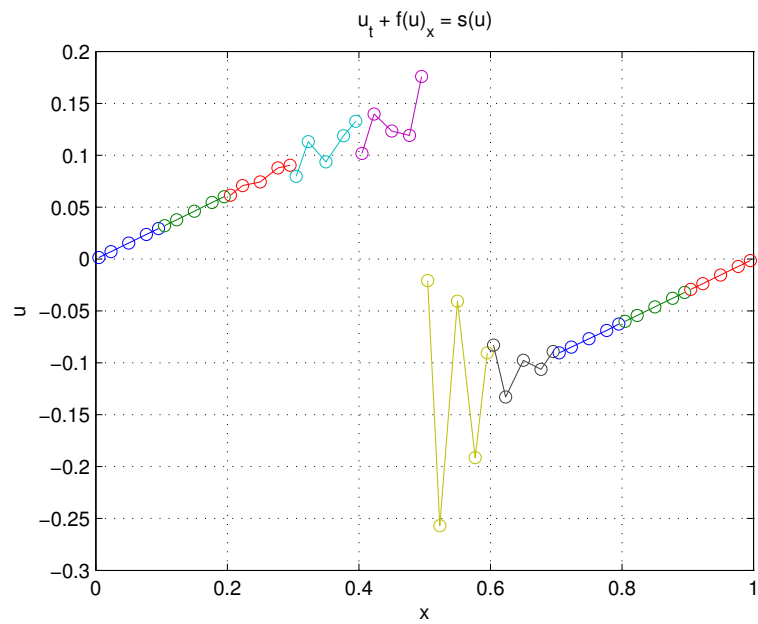
% Update plot
drawnow

% Transform degrees u(t)_{l,i} into values u(x,t)
u = V*ut;

end % time loop
toc;

Elapsed time is 2.390232 seconds.

```



Write Output

Write output to tecplot

```
if w_output == 1;
    % write to tecplot subroutine
end;
```