

## BEARCLAW 2D ADVECTION

### 1 Theory

The conservation law

$$q_t + (uq)_x + (vq)_y = 0,$$

describes advection of the passive scalar  $q(t, x, y)$  by the velocity field  $(u, v)$ .

### 2 Implementation

Define the BEARCLAW problem module.

#### 2.1 Global definitions

A module is constructed with global definitions.

<pre> MODULE problem   USE NodeInfoDef   IMPLICIT NONE   SAVE   PRIVATE   PUBLIC setprob, afterrun, qinit, b4step, setaux, src, physflux, problemBC, &amp;          afterstep, afterfixup, problemIO, problemBadCFL, problemErrFlag   REAL (KIND=qPrec), DIMENSION(2) :: ubar   CONTAINS </pre>	<p>Definition of problem module</p> <ul style="list-style-type: none"> <li>Uses Bearclaw Info structure</li> <li>All variables have to be declared</li> <li>Save variables between calls</li> <li>Internal variables cannot be seen</li> <li>Public entry points</li> </ul> <p><math>(u, v)</math> is global</p>
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#### 2.2 Problem definition: setprob

<pre> SUBROUTINE setprob   CHARACTER*60 :: comment </pre>	Input parameters defining the problem from the setprob.data file.
<pre>   OPEN(unit=7, file='setprob.data', status='old', form='formatted')   READ(7, *) ubar(1), comment ! advection x-velocity   READ(7, *) ubar(2), comment ! advection y-velocity </pre>	Open the setprob.data file and read $(u, v)$ advection velocity.
<pre>   CLOSE(7) END SUBROUTINE setprob </pre>	Close the setprob.data file.

- [setprob.data](#) - [\(update\)](#)

1.0e0 Advection velocity along x direction 1.0e0 Advection velocity along y direction

#### 2.3 Problem definition: afterrun

<pre> SUBROUTINE afterrun END SUBROUTINE afterrun </pre>	Actions after run completion
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#### 2.4 Problem definition: problemBC

<pre> SUBROUTINE problemBC(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE problemBC </pre>	Define problem-specific BCs
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#### 2.5 Problem definition: qinit

<pre> SUBROUTINE qinit(Info)   TYPE (NodeInfo) :: Info   INTEGER i,j   REAL (KIND=xPrec) :: x,y   y=Info%Xlower(2)+Info%dX(2)/2.0   DO j=1,Info%MX(2)     x=Info%Xlower(1)+Info%dX(1)/2.0     DO i=1,Info%MX(1)       IF ( x&gt;8 .AND. x&lt;24 .AND. y&gt;8 .AND. y&lt;24 ) THEN         Info%q(i,j,1,1,1)=1.0       ELSE         Info%q(i,j,1,1,1)=0.0       END IF       IF ( x&gt;10 .AND. x&lt;22 .AND. y&gt;10 .AND. y&lt;22 ) THEN         Info%q(i,j,1,1,1)=0.0       END IF       x=x+Info%dX(1)     END DO     y=y+Info%dX(2)   END DO END SUBROUTINE qinit </pre>	<p>Define initial condition</p> <p>Work on Info grid structure</p> <p>Local variables</p> $q(0, x, y) = 1 \text{ if } (x, y) \in S - S_1$ $\text{else } q(0, x, y) = 0$ $S = [8, 24] \times [8, 24]$ $S_1 = [10, 22] \times [10, 22]$
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## 2.6 Problem definition: b4step

<pre> SUBROUTINE b4step(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE b4step </pre>	Actions before each time step
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## 2.7 Problem definition: afterstep

<pre> SUBROUTINE afterstep(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE afterstep </pre>	Actions after each time step
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## 2.8 Problem definition: afterfixup

<pre> SUBROUTINE afterfixup(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE afterfixup </pre>	Actions after coarse grid update from fine grid values
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## 2.9 Problem definition: problemIO

<pre> SUBROUTINE problemIO(nframe,tnow,IOfrequest,Info,qmax,qmin)   INTEGER :: nframe,IOfrequest; REAL (KIND=qPrec) :: tnow   TYPE (NodeInfo), OPTIONAL :: Info   REAL (KIND=qPrec), OPTIONAL, DIMENSION(:) :: qmax,qmin   INTEGER, PARAMETER :: UserBeforeGridIO=-1, UserAfterGridIO=-2   INTEGER, PARAMETER :: UserIOMinMax=0,UserBeforeOutput=1, &amp;     UserAfterOutput=2 END SUBROUTINE problemIO </pre>	Application-specific I/O
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## 2.10 Problem definition: setaux

<pre> SUBROUTINE setaux(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE setaux </pre>	Define initial auxilliary variables
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## 2.11 Problem definition: `src`

<pre>SUBROUTINE src(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE src</pre>	Define source term
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## 2.12 Problem definition: `problemBadCFL`

<pre>SUBROUTINE problemBadCFL(Info)   TYPE (NodeInfo) :: Info END SUBROUTINE problemBadCFL</pre>	Actions if CFL > 1
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## 2.13 Problem definition: `problemErrFlag`

<pre>SUBROUTINE problemErrFlag(Info,CoarseInfo)   TYPE (NodeInfo) :: Info           ! Current grid   TYPE (NodeInfo) :: CoarseInfo    ! Coarsened version of current grid   Info%ErrorFlags=1 END SUBROUTINE problemErrFlag</pre>	Application-specific refinement criteria
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## 2.14 Problem definition: `physflux`

### 2.14.1 Variable declarations

<pre>SUBROUTINE physflux(ixy,indx,irequest,Info,q,f,s,A)   INTEGER ixy   INTEGER indx(MaxDims)    INTEGER irequest    TYPE (NodeInfo) :: Info ! Data associated with this grid   REAL (KIND=qPrec), POINTER, DIMENSION (:,:) :: q,f,s   REAL (KIND=qPrec), POINTER, DIMENSION (:,:,) :: A   !:physfluxInternalDeclarations:!   INTEGER mbc,mx,nq,tdir,j   INTEGER, POINTER, DIMENSION(:) :: iCell,iEdge,iLft,iRgt   REAL (KIND=qPrec) :: stran,zero=0.   ! Shorter local names for the wave propagation quantities   REAL (KIND=qPrec), POINTER, DIMENSION (:,:) :: Apdq,Amdq,speed   REAL (KIND=qPrec), POINTER, DIMENSION (:,:) :: Asdq,BmAsdq,BpAsdq   REAL (KIND=qPrec), POINTER, DIMENSION (:,:,) :: wave</pre>	Physical fluxes, solve Riemann prob. Direction along which to solve Indices of slice, <code>indx(ixy)</code> has no significance, other indices give position of this 1D slice in index space Request code, definitions in <code>NodeInfoGlobal.f90</code>
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### 2.14.2 Variable definitions

<pre>mx=Info%mxnow mbc=Info%mbc nq=Info%NrVars tdir=Info%tdir ! Index ranges: iEdge=&gt;Info%i1D           ! Interior faces: 2-mbc &lt;= iEdge &lt;= mx+mbc iCell=&gt;Info%i1Dcells     ! Cell centers: 1-mbc &lt;= iCell &lt;= mx+mbc iLft=&gt;Info%i1Dleft      ! Left faces: 1-mbc &lt;= iLft &lt;= mx+mbc-1 iRgt=&gt;Info%i1Dright     ! Right faces: 2-mbc &lt;= iRgt &lt;= mx+mbc Apdq=&gt;Info%Apdq; Amdq=&gt;Info%Amdq speed=&gt;Info%speed; wave=&gt;Info%wave Asdq=&gt;Info%Asdq; BmAsdq=&gt;Info%BmAsdq; BpAsdq=&gt;Info%BpAsdq</pre>	Get information from Info structure
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### 2.14.3 Request responses

SELECT CASE (irequest)	Identify current request
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#### Initialize

CASE (Initialize)	
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#### Finalize

CASE (Finalize)	
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#### RequestNormalWaves

<pre> CASE (RequestNormalWaves)   Apdq=0.; Amdq=0.; speed=0.; wave=0.   wave(iEdge,1,1)=q(iRgt,1)-q(iLft,1)   speed(iEdge,1)=ubar(ixy)   IF (ubar(ixy)&lt;0.) THEN     Amdq(iEdge,1)=speed(iEdge,1)*wave(iEdge,1,1)   ELSE     Apdq(iEdge,1)=speed(iEdge,1)*wave(iEdge,1,1)   END IF </pre>	Solve 1D Riemann problem along normal direction and compute left,right-going fluctuations
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#### RequestTransverseWaves

<pre> CASE (RequestTransverseWaves)   BmAsdq=0.; BpAsdq=0.   stran = ubar(tdir)   IF (stran&lt;0.) THEN     BmAsdq(iEdge,1) = stran * Asdq(iEdge,1)   ELSE     BpAsdq(iEdge,1) = stran * Asdq(iEdge,1)   END IF END SELECT </pre>	Decompose solution of normal Riemann problem into transverse fluctuations
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### 2.15 Close problem module

<pre> END SUBROUTINE physflux END MODULE problem </pre>	
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## 3 Simulations

### 3.1 Data files

- [bear.data](#) - [\(update\)](#)

```

=====
! BEARCLAW bear.data input file. Global parameters valid for all root-level grids.
=====
!:RunFlags:~ | Variable | Description
===== F 0 Restart,
Frame Resume from checkpoint data dump F LevelEqSets Solve different equations on grid levels F
LevelMethods Apply different algorithms on grid levels F SaveAtFixedTimes F=maintain CFL, T=save data
at desired times F MaintainAuxArrays Treat aux similarly to q in MPI runs F InitialAMRonly Generate
initial AMR structure and stop F OutputStyleParams Outputstyle line contains additional formatting
===== !:RunParameters:~
===== 1 nRootGrids Number

```

of root-level grids 3 MaxLevels Maximum number of grid refinement levels 2 2 2 2 2 CoarsenRatio ... of child grid to obtain parent spacing 4 MinimumGridPoints ... along one dimension 0 TimeStepMethod 0 fixed dt, 1 variable dt 0.d0 t0 initial time (if not Restart) 10.00d0 tfinal final time 4 0.5 MaxCFLRetry, rCFL Try reducing CFL by this ratio this many times 3 OutputStyle 1 AMRCLAW, 2 TECPLOT, 3 HDF, 9 GnuPlot, 11 VTK 10 OutputFrames Number of data checkpoints T T T T T T OutputLevel Level output flag  
 !=====

- [grid.data](#) - ([update](#))

```

!=====
! BEARCLAW grid.data input file. Parameters specific to root-level grid.
!===== !:GridParameters:!
Variable Description !=====
2 nDim Grid spatial dimensions 4 MaxLevel Max grid refinement levels for this grid 32 mx Cells in
x direction 32 my Cells in y direction 1 32 mGlobal(1) Global index extents of this grid x 1 32
mGlobal(2) Global index extents of this grid (y-direction) 0.0d0 xlower Left edge of computational
domain 32.0d0 xupper Right edge of computational domain 0.0d0 ylower Bottom edge of computational
domain 32.0d0 yupper Top edge of computational domain 2 mbc Number of ghost cells at each boundary
1 mthbc(1) Left boundary condition code 1 mthbc(2) Right boundary condition code 1 mthbc(3)
Bottom boundary condition code 1 mthbc(4) Top boundary condition code 0.95d0 dtv(1) Initial time
step (constant dt TimeStepMethod=0) 1.0d99 dtv(2) Max allowable time step 1.00d0 cflv(1) Max
allowable Courant number 1.00d0 cflv(2) Desired Courant number 1.0 cflv(3) Time step relaxation
parameter !=====
!:MultiphysicsParameters:! - one value if LevelEqSets==F else (>=MaxLevel) values
!===== ! NrVars = Number
of primary field variables 1 ! nEquationSet = Equation set for these fields 1 ! maux = Number of
auxilliary fields 0 !=====
!:GridRefinementParameters:! - (>=MaxLevel) values for each parameter
!===== ! qTolerance =
Field variable tolerances that trigger refinement 1.0e-3 1.0e-6 1.0e-9 1.0e-3 1.0e-3 1.0e-3 ! xTolerance
= Spatial tolerances that trigger refinement 5.0e-2 5.0e-2 5.0e-2 5.0e-2 5.0e-2 5.0e-2 ! iBuffer
= Size of buffer arround area flagged for refinement 2 2 2 2 2 2 ! DesiredFillRatios= New subgrids
should have this percentage of flagged cells 0.85 0.85 0.85 0.85 0.85 0.85 ! InterpOpt = Interpolation
method used to obtain child data from parent ! (0=minmod, 1=constant, 2=centered, 3=left, 4=right,
5=piecewise) 0 0 0 0 0 0 ! ErrorFlagOpt = Error flag method ! (0=child, 1=parent, 2=apriori 3=user)
0 0 0 0 0 0 !=====
!:NumericalSchemeParameters:! one value if LevelMethods==F else (>=MaxLevel) values
!===== 0 method(1) =
(reserved) 1 method(2) = convergence order 1 method(3) = transverse convergence order 0 method(4) =
verbosity of wavebear output 0 method(5) = source term splitting 0 method(6) = 0 split q differences, 1
split flux differences 0 method(7) = radius of slab around current 1D array of cells
1 mwaves = number of waves in each Riemann solution 4 mthlim(mw) = limiter for each wave (mw=1,
mwaves) !=====
!:UserRootLevelParameters:!
!===== ! (none for this
application) !=====

```

## 3.2 Results

Shell session inside TeXmacs pid = 22857

```
Shell] make clean > /dev/null; make outclean > /dev/null; make distclean > /dev/null
```

```
Shell] make > /dev/null
```

```
Shell] xbear > run.log
```

```
Shell] make anim.gif > /dev/null
```

```
Shell] animate anim.gif & > /dev/null
```

```
[1] 26245
```

```
Shell] make frames
```

```
convert anim.gif frame%05d.png
```

```
[1]+ Done animate anim.gif
```

```
Shell]
```

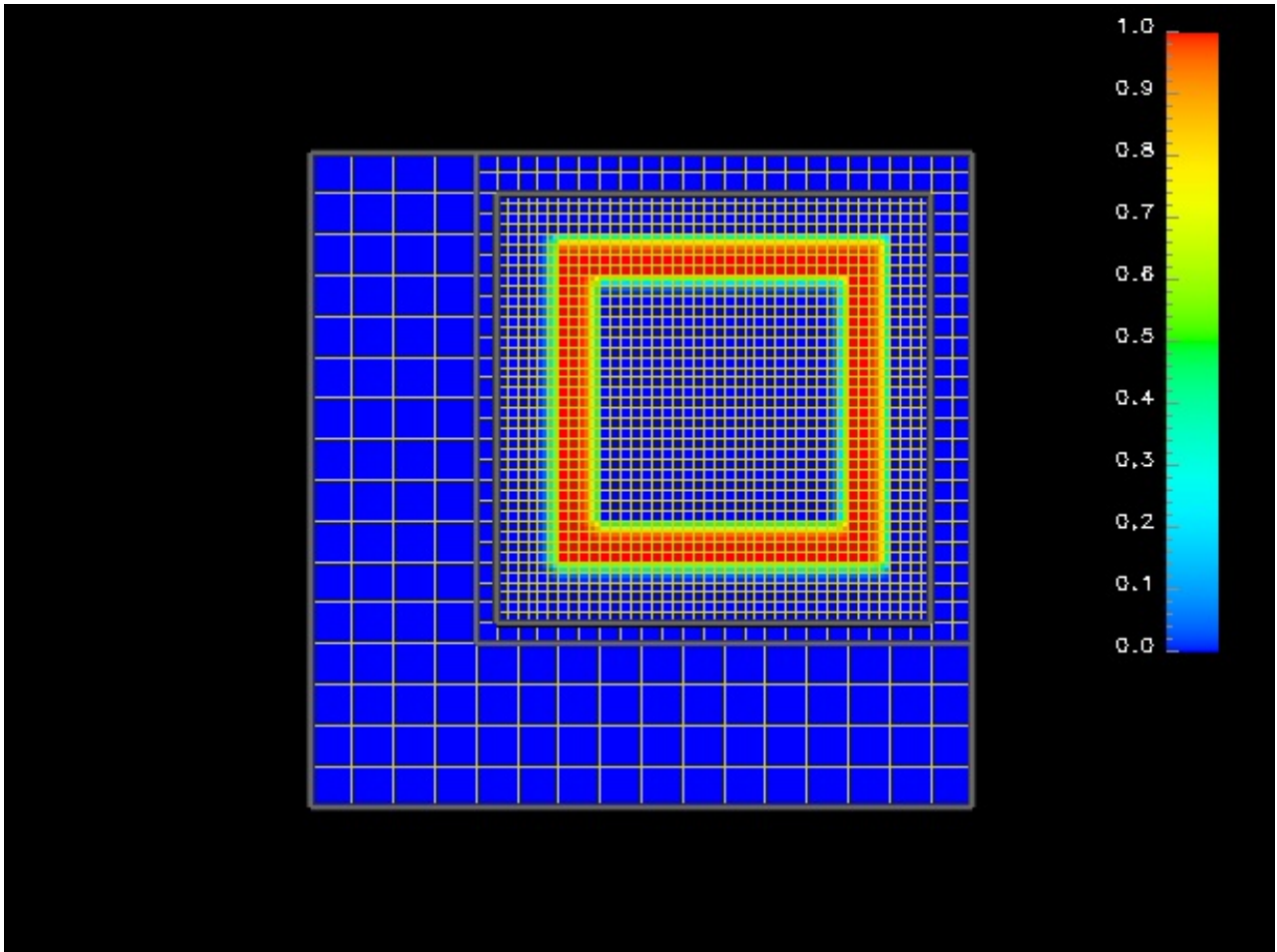
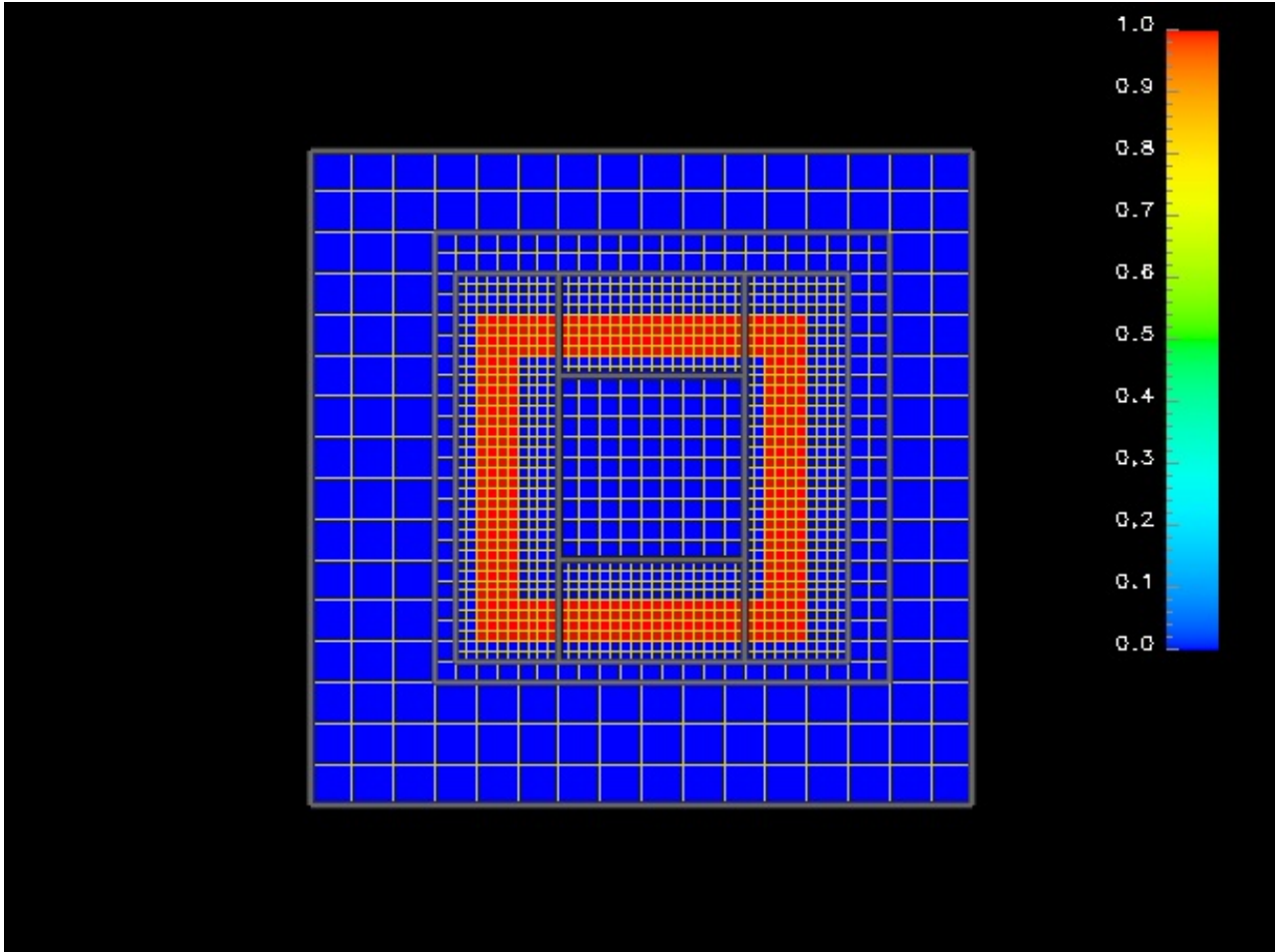


Figure 1. Initial and final frames

