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SOFTWARE FOR THE PARALLEL
CONSERVATIVE SCHEME FOR
THE SHALLOW WATER EQUATIONS

Levi Lustman
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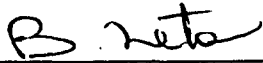
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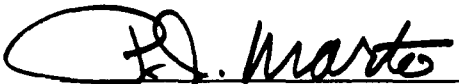
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Software for the Parallel Conservative Scheme for
the Shallow Water Equations

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Abstract

This report contains the host and node programs for the solution of the shallow water equations with topography on an INTEL iPSC/2 hypercube. Finite difference scheme conserving potential enstrophy and energy is employed in each subdomain.

1. Introduction

In this report we supply software for the numerical solution of the shallow water equations on an INTEL iPSC/2 hypercube. The method is based on domain decomposition with overlap. Finite difference scheme is used to solve in each subdomain. The scheme conserves potential enstrophy and energy (Arakawa C grid [1].) See also Neta and Lustman [2]. The efficiency of the algorithm is 81% when using 8 processors.

2. Host program

```
PROGRAM HOST
c
parameter(lparm0=10)

c cube parameters
parameter(nprocs=8,node9=nprocs-1)
parameter(lbuf=10*nprocs+30+lparm0,leninit=lbuf*4)
c                                     4 bytes per float
parameter(inityp=914)
parameter(nodes=-1,idhost=2,nodepid=3)
c domain constants
parameter(maxx=6000,maxy=2000,meshy=50,meshx=50)
parameter(j9=maxy/meshy,i9global=maxx/meshx-1)
parameter(i9dim=4+(1+i9global)/nprocs)
c
c notice that the domain size and mesh define the
c dimensions of all the arrays
c
parameter( myslv=1 , ibev=2 , iav=3 , i0gv=4)
parameter( iminv=5 , imaxv=6 , i9v=7)

common/dtcom/dt,trepo,ttot
common/physcom/coriof,corica,coriob,g
common/uvh0com/u0,v0,top,width,h0,parm0(lparm0)

c
parameter(lenuvh=4*(i9dim+1)*3*(j9+1))
parameter(ku=1,kv=2,kh=3)

logical done(0:node9),alldone
dimension uvh(0:j9,3,0:i9dim)
dimension buf(lbuf)
dimension b00kp(10,0:node9)
equivalence (b00kp,buf)

call getcube('arakawa',' ',' ',1)
call setpid(idhost)
nproc=numnodes()
print*, ' got the maximal cube, ',nproc, ' nodes'
call load('node',nodes,nodepid)

print*, ' domain parameters xmax,ymax,meshx,meshy='
, ,maxx,maxy,meshx,meshy
print*, ' this generates a grid (0:',i9global
, ,',0:',j9,')'
do 13 i=0,node9
13 done(i)=.false.
call bokeph(buf)
```

```

        l=10*nprocs+1
        call getdata
        buf(1)=dt
        l=l+1
        buf(1)=trepo
        l=l+1
        buf(1)=ttot
        nstep9=ttot/dt
c          the only parameter host must know is the maximal
c          number of steps, to decide whether the nodes
c          are done
c
        l=l+1
        buf(1)=coriof
        l=l+1
        buf(1)=corioa
        l=l+1
        buf(1)=coriob
        l=l+1
        buf(1)=g
        l=l+1
        buf(1)=u0
        l=l+1
        buf(1)=v0
        l=l+1
        buf(1)=top
        l=l+1
        buf(1)=width
        l=l+1
        buf(1)=h0
        l=l+1
        do 976 j=1,lparm0
        buf(1)=parm0(j)
        l=l+1
976      continue

        call csend(inityp,buf,leninit,nodes,nodepid)

c
c now the host will wait for results
c all come in uvh, but must be unscrambled using
c the message type to be printed
c
        n0=0
        open(UNIT=11,FORM='FORMATTED',FILE='ARA')
1132      continue
        call crecv(-1,uvh,leuvh)
c          any message at all
        ityp=infotype()
        n=ityp/100

```

```

c               number of steps
      if(n.ge.n0) then
        write(UNIT=11,FMT=7500)n*dt,n*dt/3600,n*dt/24/3600
7500      format(' Report after ',f10.0,' sec =',f10.2,' hours ='
        ,f10.2,' days')
        n0=n+1
c
c to print the title just once, not for each processor
c
      endif
      node=mod(ityp,100)
      imin=b00kp(iminv,node)
      imax=b00kp(imaxv,node)
      i0g=b00kp(i0gv,node)
c               to convert i to global i
      do 76 i=imin,imax
        iglo=mod(i0g+i , i9global+1)
        do 76 j=0,j9
          write(UNIT=11,FMT=7600)n,j+1,iglo+1
          ,uvh(j,ku,i),uvh(j,kv,i),uvh(j,kh,i)
76      continue
7600      format(i7,2i4,3g20.5)
          done(node)= (n.ge.nstep9)
452      continue
          alldone=done(0)
          do 23 i=1,node9
23      alldone=alldone.and.done(i)
          if(.not.all done) goto 1132
c               more messages coming
c
c the test above replaces waitall, which cannot be used
c
      call relcube('arakawa')
      stop
      end
      subroutine bokeph (b00kp)
c
c each node sees its data as an array (0:j9,0:i9)
c
c the column (,i) in the node data is the same as (,i+i0global) in
c the full matrix, which is (0:j9,0:i9global)
c
      parameter(lparm0=10)

c cube parameters
      parameter(nprocs=8,node9=nprocs-1)
c domain constants
      parameter(maxx=6000,maxy=2000,meshy=50,meshx=50)
      parameter(j9=maxy/meshy,i9global=maxx/meshx-1)
      parameter(i9dim=4+(1+i9global)/nprocs)

```

```

c
c notice that the domain size and mesh define the
c dimensions of all the arrays
c
      parameter( myslv=1 , ibev=2 , iav=3 , i0gv=4)
      parameter( iminv=5 , imaxv=6 , i9v=7)

      dimension b00kp(10,0:node9)
      integer gray,ginv
      integer i0w(0:node9)
      integer i9w(0:node9)

      linnod=(i9global+1)/nprocs
      do 1 i=0,node9
      i0w(i)=i*linnod
      i9w(i)=i0w(i)+linnod-1
1      continue
      iad=0
15     continue
      if(i9w(node9).ne.i9global) then
      i9w(iad)=i9w(iad)+1
      do 2 i=iad+1,node9
      i0w(i)= i0w(i)+1
      i9w(i)= i9w(i)+1
2      continue
      iad=iad+1
      goto 15
      endif
      print*,i0w
      print*,i9w

c
c at this stage, i0w(slice) to i9w(slice) are the columns that
c belong to slice, and will be advanced in time by
c the processor that treats slice
c
c but it may need four additional columns, to compute
c neighborhood averages
c
      do 333 iam=0,node9
      myslice=ginv(iam)
      i0wm=i0w(myslice)-2
      i0m=i0wm
      if (i0wm.lt.0) i0wm=i0wm+i9global+1
      i0w(myslice)=i0wm
      i9wm=i9w(myslice)+2
      i9m=i9wm
      if (i9wm.gt.i9global) i9wm=i9wm-i9global-1
      i9w(myslice)=i9wm
      i9=i9m-i0m
      ibefore=-1
      iafter=-1

```

```

        myp=mod(3*nprocs+myslice+1,nprocs)
        mym=mod(3*nprocs+myslice-1,nprocs)
        iafter=gray(myp)
        ibefore=gray(mym)
C=====
C
C      also define imin, imax
C these are the lines which this node should advance in time,
C the other lines are for information only
C
        imin=2
        imax=i9 -2
C=====
        b00kp(i0gv,iam)=i0w(myslice)
        b00kp(myslv,iam)=myslice
        b00kp(ibev,iam)=ibefore
        b00kp(iav,iam)=iafter
        b00kp(i9v,iam)=i9
        b00kp(imaxv,iam)=imax
        b00kp(iminv,iam)=imin
        print*, ' ..... I am ',iam
        print*, ' my slice is ', myslice
        print*, ' procs. before,after me=',ibefore,iafter
        print*, ' my data have dim (,0:',i9,')'
        ,imin, ' to ',imax,' meaningful'
        iii=b00kp(i0gv,iam)
        print*, ' my column 0 is global column ',iii
333      continue
        return
        end
        subroutine getdata

        parameter(lparm0=10)
C domain constants
        parameter(maxx=6000,maxy=2000,meshy=50,meshx=50)

        common/dtcom/dt,trepo,ttot
        common/physcom/coriof,corioa,coriob,g
        common/uvh0com/u0,v0,top,width,h0,parm0(lparm0)
        character*3 name

        minmsh=min(meshx,meshy)
        dt=150.*minmsh/125
        ttot=24*3600*5

C              5 days
        trepo= -ttot

C
C      silly programming trick, so ttot,trepo
C      may be entered in any order
C
        coriof=1.e-4
        corioa=0

```

```

coriob=0
g=9.8e-3
c          notice! length in km !
u0=20e-3
v0=0
top=2
width=1000
h0=5
do 1 i=1,lparm0)
1   parm0(i)=0
c
c all these are defaults. then more data may be read
c
1132      continue
      read100,name
100      format(a3)
      if(name.eq.'par') then
1131      read*,i,x
      if(i.gt.0.and.i.le.lparm0) then
          parm0(i)=x
          goto 1131
      else
          goto 1132
      endif
      else
          if(name.eq.'end') goto 9999
          if(name.eq.'sto') goto 9999
          if(name.eq.'dt ') then
              read*,dt
              goto 1132
          endif
          if(name.eq.'tre') then
              read*,trepo
              goto 1132
          endif
          if(name.eq.'tto') then
              read*,ttot
              if(trepo.lt.0) trepo= -ttot
              goto 1132
          endif
          if(name.eq.'cor') then
              read*,coriof
              goto 1132
          endif
          if(name.eq.'coa') then
              read*,corioa
              goto 1132
          endif
          if(name.eq.'cob') then
              read*,coriob
              goto 1132
          endif
      endif
  
```

```

        if(name.eq.'g ') then
            read*,g
            goto 1132
        endif
        if(name.eq.'u0 ') then
            read*,u0
            goto 1132
        endif
        if(name.eq.'v0 ') then
            read*,v0
            goto 1132
        endif
        if(name.eq.'h0 ') then
            read*,h0
            goto 1132
        endif
        if(name.eq.'top') then
            read*,top
            goto 1132
        endif
        if(name.eq.'wid') then
            read*,width
            goto 1132
        endif
        stop
    endif
9999    continue
        trepo=abs(trepo)
        print*, 'dt,treport,ttotal='
        , ,dt,trepo,ttot
        print*, 'coriolis f,corioa,coriob,g='
        , ,coriof,corioa,coriob,g
        print*, 'u0,v0,top,width,h0='
        , ,u0,v0,top,width,h0
        do 2 i=1,lparm0
2        print*,parm0(i)
        return
    end

```

3. Node program

```
PROGRAM NODE
c
  parameter(lparm0=10)

c cube parameters
  parameter(nprocs=8,node9=nprocs-1)
  parameter(nodes=-1,idhost=2,nodepid=3)

c domain constants
  parameter(maxx=6000,maxy=2000,meshy=50,meshx=50)
  parameter(j9=maxy/meshy,i9global=maxx/meshx - 1)
  parameter(i9dim=4+(1+i9global)/nprocs)

c
c notice that the domain size and mesh define the
c dimensions of all the arrays
c

c
c The array UVH contains the data u,v,h, in a format that allows
c fast message passing.
c At a fixed i, just send a buffer beginning at UVH(0,1,i)
c with length 2(columns)*3(variables)*(j9+1) words
c
  parameter(lenUVH =6*(j9+1)*4 )
c
  parameter(4 bytes per real
  parameter(inddt0=9140,inddt9=9149)
  parameter(ku=1,kv=2,kh=3)
  common/allcom/ UVH(0:j9,3,0:i9dim)
  _,zeta(0:i9dim,0:j9),hq(0:i9dim,0:j9),q(0:i9dim,0:j9)
  _,f(0:i9dim,0:j9),alfa(0:i9dim,0:j9),beta(0:i9dim,0:j9)
  _,gama(0:i9dim,0:j9),delta(0:i9dim,0:j9)
  _,eps(0:i9dim,0:j9),fi(0:i9dim,0:j9)
  _,cay(0:i9dim,0:j9),ustar(0:i9dim,0:j9),vstar(0:i9dim,0:j9)
  _,dudt(0:i9dim,0:j9),dvd(0:i9dim,0:j9),dhdt(0:i9dim,0:j9)
  _,hs(0:i9dim,0:j9),hu(0:i9dim,0:j9),hv(0:i9dim,0:j9)

  common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
  _,i0global,i9,i8,imin,imax

  common/dtcom/dt,trepo,ttot

  common/physcom/coriof,corioa,coriob,g

  common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

  common/seqcom/lmnpr
```

```

c
c during debug only
c
    dimension uold(0:i9dim,0:j9)
    _,vold(0:i9dim,0:j9),hold(0:i9dim,0:j9)

c
c during debug only
c
    if(mynode().ge.nprocs) stop
    iam=mynode()
    lmnpr=1000 +100*iam

    m0dul0=i9global+1
    call init
    nreport=trepo/dt
    nstep9=ttot/dt
    dthalf=dt/2
    twodt=2*dt

    nstep=0
    call report(nstep,nreport,nstep9)

c=====
c sending data to the neighbors THE VERY FIRST TIME
c the data get to different buffers, so the CALLs are synchronous
c
    msgbefo=isend(inddt9,UVH(0,1,imin),lenUVH,ibefore,nodepid )
    msgaftr=isend(inddt0,UVH(0,1,imax-1),lenUVH,iafter,nodepid )
c=====

10      continue
c
c Heun step
c
    call defddt

c
c until now, UVH has not changed. before it changes,
c ensure the sending worked
c
    call msgwait(msgbefo)
    call msgwait(msgaftr)

    do 11 i=0,i9
    do 11 j=0,j9
    uold(i,j)=UVH(j,ku,i)
    UVH(j,ku,i) =uold(i,j)+dudt(i,j)*dthalf
    vold(i,j)=UVH(j,kv,i)
    UVH(j,kv,i) =vold(i,j)+dvd(i,j)*dthalf
    hold(i,j)=UVH(j,kh,i)

```

```

        UVH(j,kh,i) =hold(i,j)+dhdt(i,j)*dthalf
11      continue
c=====
c
c the step ends with sending data to the neighbors
c the data get to different buffers, so the CALLs are synchronous
c
        msgbefo=isend(inddt9,UVH(0,1,imin),lenUVH,ibefore,nodepid )
        msgaftr=isend(inddt0,UVH(0,1,imax-1),lenUVH,iafter,nodepid )
        call defddt
c
c until now, UVH has not changed. before it changes,
c ensure the sending worked
c
        call msgwait(msgbefo)
        call msgwait(msgaftr)
c=====
        do 12 i=0,i9
        do 12 j=0,j9
        UVH(j,ku,i) =uold(i,j)+dudt(i,j)*dt
        UVH(j,kv,i) =vold(i,j)+dvdt(i,j)*dt
        UVH(j,kh,i) =hold(i,j)+dhdt(i,j)*dt
12      continue
c=====
c
c the step ends with sending data to the neighbors
c the data get to different buffers, so the CALLs are synchronous
c
        msgbefo=isend(inddt9,UVH(0,1,imin),lenUVH,ibefore,nodepid )
        msgaftr=isend(inddt0,UVH(0,1,imax-1),lenUVH,iafter,nodepid )
        nstep=nstep+1
        call report(nstep,nreport,nstep9)
125     continue
c
c following leapfrog steps
c
        call defddt
c
c until now, UVH has not changed. before it changes,
c ensure the sending worked
c
        call msgwait(msgbefo)
        call msgwait(msgaftr)
c=====
        do 13 i=0,i9
        do 13 j=0,j9
        temp=UVH(j,ku,i)
        UVH(j,ku,i) =uold(i,j)+dudt(i,j)*twodt
        uold(i,j)=temp
        temp=UVH(j,kv,i)
        UVH(j,kv,i) =vold(i,j)+dvdt(i,j)*twodt
        vold(i,j)=temp

```

```

        temp=UVH(j,kh,i)
        UVH(j,kh,i) =hold(i,j)+dhdt(i,j)*twodt
        hold(i,j)=temp
13      continue
c=====
c
c the step ends with sending data to the neighbors
c the data get to different buffers, so the CALLs are synchronous
c
        msgbefo=isend(inddt9,UVH(0,1,imin),lenUVH,ibefore,nodepid )
        msgaftr=isend(inddt0,UVH(0,1,imax-1),lenUVH,iafter,nodepid )
c=====
        nstep=nstep+1
        call report(nstep,nreport,nstep9)
        if( mod(nstep+1,300).eq.1) then
            goto 10
        else
            goto 125
        endif
        end
        subroutine UVHpr(UVH,t,kUVH, leadim,i9,j9 )
        parameter(ku=1,kv=2,kh=3)
        character*(*) t
        common/seqcom/lmnpr
        dimension UVH(0:j9,3,0:leadim)
        common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
        ,i0global,i90,i8,imin,imax
c
c comment out the return during debug
c
        return
        print2000,10000*lmnpr,t,leadim,i9,j9,imin,imax,iam
2000      format(i10,2x,a20,' leadim,i9,j9=',3i4,' imin,imax=',3i4)
        do 1 i=0,i9
            ii=mod(i+i0global ,m0dul0)
            if(i.ge.imin.and.i.le.imax) then
                iamy=iam
            else
                iamy=iam+10
            endif
            do 1 j=0,j9
                print1000,ii+100*(j+100*lmnpr),t,j,ii
                ,UVH(j,kUVH, i),iamy
1          continue
1000      format(i10,2x,a10,2x,2i4,f20.7,i3)
            lmnpr=lmnpr+1
            return
        end

```

```

      subroutine bokepn(b00kp)
c
c the node sees its data as an array (0:i9dim,0:j9)
c
c the column (,i) in the node data is the same as (,i+i0global) in
c the full matrix, which is (0:i9dimglobal,0:j9)
c
c most of the work was done by the host, and
c data is just rearranged here
c
      parameter(lparm0=10)

c cube parameters
      parameter(nprocs=8,node9=nprocs-1)
c connection parameters
      parameter( myslv=1,ibev=2,iav=3,i0gv=4)
      parameter( iminv=5,imaxv=6,i9v=7)

      dimension b00kp(10,0:node9)
      common/seqcom/lmnpr
      common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
      _,i0global,i9,i8,imin,imax

      iam=mynode()
      myslice=b00kp(myslv,iam)
      ibefore=b00kp(ibev,iam)
      iafter=b00kp(iav,iam)
      i0global=b00kp(i0gv,iam)
      i9=b00kp(i9v,iam)
      imin=b00kp(iminv,iam)
      imax=b00kp(imaxv,iam)
      i8=i9-1

c
c print during debug only
c
cdebug      key=lmnpr*10000+100*iam
cdebug      print*,key,' ... I am ',iam
cdebug      key=key+1
cdebug      print*,key,' slice=', myslice
cdebug      key=key+1
cdebug      print*,key,' nodes bef,aft=',ibefore,iafter
cdebug      key=key+1
cdebug      print*,key,' data dim (0:' ,i9,',)'
cdebug      key=key+1
cdebug      print*,key,'      ',imin, ' to ',imax,' meaningful'
cdebug      key=key+1
cdebug      print*,key,' my column 0=global ',i0global
cdebug      lmnpr=lmnpr+1
      return
      end

```

```

function coriofun(i,j,meshx,meshy)
parameter(lparm0=10)
common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
_,i0global,i9,i8,imin,imax
common/dtcom/dt,trepo,ttot
common/physcom/coriof,corioa,coriob,g
common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

x=meshx*mod(i0global+i,m0dul0)
y=meshy*(j)
c                               x,y for f --like zeta-- from C-mesh
coriofun=coriof

c
c for the simplest case, coriolis is constant, but in general
c it might be computed using x,y and the parm0 data
c or corioa, coriob -- tangent plane, if needed
c
return
end
subroutine defddt

c
c the results always in dudt, dvdt, dhdt in allcom
c

parameter(lparm0=10)

c cube parameters
parameter(nprocs=8,node9=nprocs-1)
c domain constants
parameter(maxx=6000,maxy=2000,meshy=50,meshx=50)
parameter(j9=maxy/meshy,i9global=maxx/meshx-1)
parameter(i9dim=4+(1+i9global)/nprocs)

c
c notice that the domain size and mesh define the
c dimensions of all the arrays
c

c
c The array UVH contains the data u,v,h, in a format that allows
c fast message passing.
c At a fixed j, just send a buffer beginning at UVH(0,1,j)
c with length 6*(i9+1) words = 2 rows of 3 vars each
parameter(lenUVH =6*(j9+1)*4 )
c                               4 bytes per real
parameter(inddt0=9140,inddt9=9149)

c
parameter(ku=1,kv=2,kh=3)

```

```

common/allcom/ UVH(0:j9,3,0:i9dim)
_,zeta(0:i9dim,0:j9),hq(0:i9dim,0:j9),q(0:i9dim,0:j9)
_,f(0:i9dim,0:j9),alfa(0:i9dim,0:j9),beta(0:i9dim,0:j9)
_,gama(0:i9dim,0:j9),delta(0:i9dim,0:j9)
_,eps(0:i9dim,0:j9),fi(0:i9dim,0:j9)
_,cay(0:i9dim,0:j9),ustar(0:i9dim,0:j9),vstar(0:i9dim,0:j9)
_,dudt(0:i9dim,0:j9),dvdt(0:i9dim,0:j9),dhdt(0:i9dim,0:j9)
_,hs(0:i9dim,0:j9),hu(0:i9dim,0:j9),hv(0:i9dim,0:j9)

```

```

common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
_,i0global,i9,i8,imin,imax

```

```

common/dtcom/dt,trepo,ttot

```

```

common/physcom/coriof,corioa,coriob,g

```

```

common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

```

```

i7=i8-1
dx=meshx
dy=meshy
j8=j9-1
j7=j9-2

```

```

C=====

```

```

C
C any such step begins with receiving data from the neighbors
C the data get to different buffers, so the CALLs are synchronous
C

```

```

in0=irecv(inddt0,UVH(0,1,0),lenUVH )
in9=irecv(inddt9,UVH(0,1,i8),lenUVH )
call msgwait(in0)
call msgwait(in9)
CALL UVHpr(UVH,'u after irecv',ku,i9dim,i9,j9)
CALL UVHpr(UVH,'v after irecv',kv,i9dim,i9,j9)
CALL UVHpr(UVH,'h after irecv',kh,i9dim,i9,j9)

```

```

C
C The big mess is that u,v,h are not defined everywhere!
C      u defined for 0<=i<=i9, 0<=j < j9
C      v defined for 0<=i < i9, 0<=j<=j9
C      h defined for 0<=i < i9, 0<=j < j9
C
C But because of periodicity in i , which is maintained by the
C send+receive :
C      v defined for 0<=i<=i9, 0<=j<=j9
C      h defined for 0<=i<=i9, 0<=j < j9
C
C So, finally, only the following are MISSING:
C      u(,j9) , h(,j9)

```

```

C
C NOW WE DEFINE ALL VARIABLES WHEREVER POSSIBLE
C THEN,
C SEE IF d/dt IS DEFINED WHERE IT SHOULD:
C     imin<=i<=imax, jmin<=j<=jmax
C
C     imin,imax from bookkeeping,
C     jmin,jmax depend on the variable u,v,h
C
C IT TURNS OUT THAT EVERYTHING IS SAFE PROVIDED: 2<=imin , imax < i8
C
C     N.B. imax less than i8 !
C
C
C 3.12 in the paper follows
C
C
C     do 3120 i=1,i9
C     do 3121 j=1,j8
C
C both i,j safe
C
C     zeta(i,j)=(UVH(j-1,ku,i) - UVH(j,ku,i))/dy
C     _+(UVH(j,kv,i) -UVH(j,kv,i-1))/dx
3121 continue
C
C THE PAPER HAS THE COMPUTATIONAL BD.CD. ZETA=0
C
C     zeta(i,0)=0
C     zeta(i,j9)=0
3120 continue
312 continue
cdebug CALL matpr('zeta',zeta,i9dim,i9,j9)
C
C
C 3.15 in the paper follows
C
C
C     do 3150 j=1,j8
C     do 3151 i=1,i9
C
C both i,j safe, periodic bd.cd in i automatic
C
C     hq(i,j)=(UVH(j,kh,i)+UVH(j,kh,i-1)+UVH(j-1,kh,i-1)
C     ++UVH(j-1,kh,i))/4
3151 continue
3150 continue

```

```

c
c this still leaves hq undefined at j=0
c
c do some extrapolation, maybe zeta=0 will take care of
c instability
c
      do 3156 i=1,i9
        hq(i,0)=3*hq(i,1)-3*hq(i,2)+hq(i,3)
3156      continue
c
c this still leaves hq undefined at j=j9
c
c do some extrapolation, maybe zeta=0 will take care of
c instability
c
      do 3157 i=1,i9
        hq(i,j9)=3*hq(i,j9-1)-3*hq(i,j9-2)+hq(i,j9-3)
3157      continue
315      continue
cdebug          CALL matpr('hq',hq,i9dim,i9,j9)


c
c 3.11 in the paper follows
c
      do 311 i=1,i9
        do 311 j=0,j9
          q(i,j)=(f(i,j)+ zeta(i,j)) / hq(i,j)
311      continue
cdebug          CALL matpr('q',q,i9dim,i9,j9)
c
c 3.34 in the paper follows
c
      do 334 j=0,j8

c the same j loop all over

      do 3340 i=1,i8
        eps(i,j)=(q(i+1,j+1)+ q(i,j+1)-q(i,j)-q(i+1,j))/24
        fi(i,j)=(-q(i+1,j+1)+ q(i,j+1)+q(i,j)- q(i+1,j))/24
c
c these have indices i+1/2, j+1/2 only, like h
c
3340      continue

      do 3341 i=2,i8
        alfa(i,j)=(2*q(i+1,j+1)+q(i,j+1)+ 2*q(i,j)+ q(i+1,j))/24
        beta(i,j)=(q(i,j+1)+ 2*q(i-1,j+1)+ q(i-1,j)+ 2*q(i,j))/24
        gama(i,j)=(2*q(i,j+1)+ q(i-1,j+1)+ 2*q(i-1,j)+ q(i,j))/24
        delta(i,j)=(q(i+1,j+1)+2*q(i,j+1)+ q(i,j)+2*q(i+1,j))/24
3341      continue

```

c closing the j loop

334 continue

cdebug CALL matpr('eps',eps,i9dim,i9,j9)

cdebug CALL matpr('fi',fi,i9dim,i9,j9)

cdebug CALL matpr('alfa',alfa,i9dim,i9,j9)

cdebug CALL matpr('beta',beta,i9dim,i9,j9)

cdebug CALL matpr('gama',gama,i9dim,i9,j9)

cdebug CALL matpr('delta',delta,i9dim,i9,j9)

c

c 3.36-3.37 in the paper follow

c

do 336 j=0,j8

do 336 i=1,i9

hu(i,j)=(UVH(j,kh,i-1) +UVH(j,kh,i))/2

336 continue

do 337 i=0,i9

do 337 j=1,j9

hv(i,j)=(UVH(j-1,kh,i) + UVH(j,kh,i))/2

c

c at j=0 and j=j9, hv may be anything, because v=0

c

337 continue

cdebug CALL matpr('hu',hu,i9dim,i9,j9)

cdebug CALL matpr('hv',hv,i9dim,i9,j9)

c

c 3.41 in the paper follows

c

do 341 i=0,i8

do 341 j=0,j8

cay(i,j)=(UVH(j,ku,i) **2 +UVH(j,ku,i+1) **2

+UVH(j,kv,i) **2 +UVH(j+1,kv,i) **2)/4

341 continue

cdebug CALL matpr('cay',cay,i9dim,i9,j9)

c

c 3.3-3.4 in the paper follow

c

do 33 j=0,j8

do 33 i=1,i9

ustar(i,j)= hu(i,j)* UVH(j,ku,i)

33 continue

```

        do 34 i=0,i9
          vstar(i,0)=0
c
c bd.cd          v=0
c
        do 34 j=1,j9
          vstar(i,j)= hv(i,j)* UVH(j,kv,i)
34      continue
cdebug      CALL matpr('ustar',ustar,i9dim,i9,j9)
cdebug      CALL matpr('vstar',vstar,i9dim,i9,j9)

c
c 3.1-3.2 in the paper follow
c
        do 32 i=1,i8
          do 32 j=0,j8

c
c this is where dhdt values are defined. they are needed for:
c   imin <= i <= imax
c   0      <= j <= j8
c           NOT needed for j=j9
c
          dhdt(i,j)=(ustar(i,j)-ustar(i+1,j))/dx
          _+(vstar(i,j)-vstar(i,j+1))/dy
31      continue
32      continue

cdebug      CALL matpr('dhdt',dhdt,i9dim,i9,j9)

c
c 3.5 in the paper follows to compute du/dt
c
c
        do 35 j=0,j8
          do 35 i=2,i8

c
c this is where dudt is defined . it is needed for:
c   imin <= i <= imax
c   0      <= j <= j8
c           NOT needed for j=j9
c
c 2.5 in the paper follows, defining capital phi in terms of h,hs
c
          caphi1=g*(UVH(j,kh,i-1) +hs(i-1,j))
          caphi2=g*(UVH(j,kh,i) +hs(i,j))

          dudt(i,j)= alfa(i,j)* vstar(i,j+1)+beta(i,j)* vstar(i-1,j+1)
          _+gama(i,j)* vstar(i-1,j)
          _+delta(i,j)* vstar(i,j)-eps(i,j)*ustar(i+1,j)
          _+eps(i-1,j)* ustar(i-1,j)
          _+(cay(i-1,j)+ caphi1 -cay(i,j)- caphi2)/dx
35      continue

```

```

cdebug          CALL matpr('dudt',dudt,i9dim,i9,j9)

c
c 3.6 in the paper follows to compute dv/dt
c
      do 36 i=2,i8-1
      do 36 j=1,j8
c
c this is where dvdt is defined
c it is needed for
c      imin <= i <= imax
c      1      <= j <= j8
c
c dvdt NOT needed at j=0 or j=j9, because there v=0 always
c
c 2.5 in the paper follows, defining capital phi in terms of h,hs
c
      caphi3=g*(UVH(j-1,kh,i) +hs(i,j-1))
      caphi4=g*(UVH(j,kh,i) +hs(i,j))

      dvdt(i,j)= -gama(i+1,j)* ustar(i+1,j)- delta(i,j)* ustar(i,j)
      - alfa(i,j-1)*ustar(i,j-1)- beta(i+1,j-1)* ustar(i+1,j-1)
      +fi(i,j-1)*vstar(i,j-1)+fi(i,j)* vstar(i,j+1)
      +(caphi3+cay(i,j-1)-caphi4 - cay(i,j))/dy
36      continue
cdebug          CALL matpr('dvdt',dvdt,i9dim,i9,j9)
      return
      end
      function hsfun(i,j,meshx,meshy)
      parameter(lparm0=10)
      parameter(maxx=6000,maxy=2000)
      common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
      -,i0global,i9,i8,imin,imax
      common/dtcom/dt,trepo,ttot
      common/physcom/coriof,corioa,coriob,g
      common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

      x=meshx*(mod(i0global+i,m0dul0)+0.5)
      y=meshy*(j+0.5)
c
c          x,y for h from C-mesh
c
      xx=abs(x-0.5*maxx)
      if(xx.lt.width) then
      hsfun=top*(1-xx/width)
      else
      hsfun=0
      endif
c
c this is the triangular ridge. more generally,
c hsfun might be computed using x,y and the parm0 data
c
      return
      end

```



```

        l=l+1
        coriob=buf(l)
        l=l+1
        g=buf(l)
        l=l+1
        u0=buf(l)
        l=l+1
        v0=buf(l)
        l=l+1
        top=buf(l)
        l=l+1
        width=buf(l)
        l=l+1
        h0=buf(l)
        l=l+1
        do 976 j=1,lparm0
            parm0(j)=buf(l)
            l=l+1
976      continue
c
c the common allcom is filled with trash, to check
c glitches in index manipulation
c
c but some variables get meaningful values: f,vstar,...
c
        do 1 i=0,i9
            do 1 j=0,j9
                UVH(j,ku,i) =1.e6+j+100*i
                zeta(i,j)=3.e6+j+100*i
                UVH(j,kh,i) =4.e6+j+100*i
                hq(i,j)=5.e6+j+100*i
                q(i,j)=6.e6+j+100*i
                f(i,j)=coriofun(i,j,meshx,meshy)
                alfa(i,j)=8.e6+j+100*i
                beta(i,j)=9.e6+j+100*i
                gama(i,j)=-1.e6-j-100*i
                delta(i,j)=-2.e6-j-100*i
                eps(i,j)=-3.e6-j-100*i
                fi(i,j)=-4.e6-j-100*i
                hu(i,j)=-5.e6-j-100*i
                hv(i,j)=-6.e6-j-100*i
                cay(i,j)=-7.e6-j-100*i
                ustar(i,j)=-8.e6-j-100*i
                UVH(j,kv,i) =0
                vstar(i,j)=0
c
c implementation of WALL bd.cd
c
        dudt(i,j)=0
        dvdt(i,j)=0
        dhdt(i,j)=0
1      continue

```

```

c
c initial data, defined on PART OF the arrays
c

      j8=j9-1
      do 170 i=0,i9
      do 170 j=0,j8
      UVH(j,ku,i) =u0fun(i,j,meshx,meshy )
170      continue
      do 171 i=0,i9
      do 171 j=1,j8

c
c 1<=j<=j8  <=>  implementation of WALL bd.cd
c
      UVH(j,kv,i) =v0fun(i,j,meshx,meshy )
171      continue
      do 172 i=0,i9
      do 172 j=0,j9
      hs(i,j)=hsfun(i,j,meshx,meshy)

c
c hs is defined everywhere, it is the geography
c
172      continue
      do 271 i=0,i9
      do 271 j=0,j8
      UVH(j,kh,i) =h0-hs(i,j )
271      continue

      CALL UVHpr(UVH,'u init',ku,i9dim,i9,j9)
      CALL UVHpr(UVH,'v init',kv,i9dim,i9,j9)
      CALL UVHpr(UVH,'h init',kh,i9dim,i9,j9)

      return
      end
      subroutine matpr(t,a,leadim, i9 ,j9)
      common/seqcom/lmnpr
      common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
      ,i0global,i90,i8,imin,imax
      character*(*) t
      dimension a(0:leadim,0:j9)

c
c comment out the return during debug
c

      return
      print2000,10000*lmnpr,t,leadim,i9,j9,imin,imax,iam
2000      format(i10,2x,a10,' leadim,i9,j9=',3i4,' imin,imax=',3i4)
      do 1 i=0,i9
      ii=mod(i+i0global ,m0dul0)
      if(i.ge.imin.and.i.le.imax) then
      iamy=iam
      else

```

```

        iamy=iam+10
    endif
    do 1 j=0,j9
        print1000,ii+100*(j+100*lmnpr),t,j,ii,a(i,j),iamy
1      continue
1000    format(i10,2x,a10,2x,2i4,f20.7,i3)
        lmnpr=lmnpr+1
        return
    end
    subroutine report(n,nrep,nmax)
c
c n is the number of steps
c
        parameter(lparm0=10)

c cube parameters
        parameter(nprocs=8,node9=nprocs-1)
        parameter(nodes=-1,idhost=2,nodepid=3)
c domain constants
        parameter(maxx=6000,maxy=2000,meshy=50,meshx=50)
        parameter(j9=maxy/meshy,i9global=maxx/meshx - 1)
        parameter(i9dim=4+(1+i9global)/nprocs)
c
c notice that the domain size and mesh define the
c dimensions of all the arrays
c
c
c The array UVH contains the data u,v,h, in a format that allows
c fast message passing.
c
        parameter(len=12*(i9dim+1)*(j9+1))
        parameter(ku=1,kv=2,kh=3)
        common/allcom/ UVH(0:j9,3,0:i9dim)
        _ ,zeta(0:i9dim,0:j9),hq(0:i9dim,0:j9),q(0:i9dim,0:j9)
        _ ,f(0:i9dim,0:j9),alfa(0:i9dim,0:j9),beta(0:i9dim,0:j9)
        _ ,gama(0:i9dim,0:j9),delta(0:i9dim,0:j9)
        _ ,eps(0:i9dim,0:j9),fi(0:i9dim,0:j9)
        _ ,cay(0:i9dim,0:j9),ustar(0:i9dim,0:j9),vstar(0:i9dim,0:j9)
        _ ,dudt(0:i9dim,0:j9),dvdt(0:i9dim,0:j9),dhdt(0:i9dim,0:j9)
        _ ,hs(0:i9dim,0:j9),hu(0:i9dim,0:j9),hv(0:i9dim,0:j9)

        common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
        _ ,i9global,i9,i8,imin,imax

        common/dtcom/dt,trepo,ttot

        common/physcom/coriof,corioa,coriob,g

        common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

```

```

        if(mod(n,nrep).ne.0) return
        itype=n*100+iam
452      continue
        myh=myhost()
        call csend(itype,UVH,len,myh,idhost )
        if(n.ge.nmax) stop 5144
        return
      end
      function u0fun(i,j,meshx,meshy)
      parameter(lparm0=10)
      common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
      _,i0global,i9,i8,imin,imax
      common/dtcom/dt,trepo,ttot
      common/physcom/coriof,corioa,coriob,g
      common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

      x=meshx*mod(i0global+i,m0dul0)
      y=meshy*(j+0.5)
c
c      x,y for u from C-mesh
      u0fun=u0
c
c for the simplest case, u is constant, but in general
c it might be computed using x,y and the parm0 data
c
      return
      end
      function v0fun(i,j,meshx,meshy)
      parameter(lparm0=10)
      common/bkkpcom/m0dul0,iam,myslice,ibefore,iafter
      _,i0global,i9,i8,imin,imax
      common/dtcom/dt,trepo,ttot
      common/physcom/coriof,corioa,coriob,g
      common/UVH0com/u0,v0,top,width,h0,parm0(lparm0)

      x=meshx*(mod(i0global+i,m0dul0)+0.5)
      y=meshy*(j)
c
c      x,y for v from C-mesh
      v0fun=v0
c
c for the simplest case, v is constant, but in general
c it might be computed using x,y and the parm0 data
c
      return
      end

```

4. Makefile

```
# This file is used to compile and link the host.f, node.f
#
# The command "make all" causes compilation and linking.
```

```
all :   host node
```

```
host.o: host.f
```

```
node.o: node.f
```

```
host:   host.o
        f77 -o host host.o -host
```

```
node:   node.o
        f77 -o node node.o -node
```

5. Input File

dt
60
ttotal
79800
top
2
width
1000
u0
20.e-3
end

Acknowledgements

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