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! Skeleton 2-1/2D Electromagnetic OpenMP PIC code
! written by Viktor K. Decyk, UCLA
    program mbpic2
    use mbpush2_h
    use omplib_h
    implicit none
    integer, parameter :: indx = 9, indy = 9
    integer, parameter :: npx = 3072, npy = 3072
    integer, parameter :: ndim = 3
    real, parameter :: tend = 10.0, dt = 0.04, qme = -1.0
    real, parameter :: vtx = 1.0, vty = 1.0, vx0 = 0.0, vy0 = 0.0
    real, parameter :: vtz = 1.0, vz0 = 0.0
    real :: ax = .912871, ay = .912871, ci = 0.1
! idimp = dimension of phase space = 5
! relativity = (no,yes) = (0,1) = relativity is used
    integer :: idimp = 5, ipbc = 1, relativity = 1
    real :: wke = 0.0, we = 0.0, wf = 0.0, wm = 0.0, wt = 0.0
! sorting tiles, should be less than or equal to 32
    integer :: mx = 16, my = 16
! fraction of extra particles needed for particle management
    real :: xtras = 0.2
! declare scalars for standard code
    integer :: np, nx, ny, nxh, nyh, nxe, nye, nxeh, nxyh, nxhy
    integer :: mx1, my1, mxy1, ntime, nloop, isign
    real :: qbme, affp, dth
!
! declare scalars for OpenMP code
    integer :: nppmx, nppmx0, ntmax, npbm, irc
    integer :: nvp
!
! declare arrays for standard code
    real, dimension(:,:), pointer :: part
    real, dimension(:,:), pointer :: qe
    real, dimension(:,:,:), pointer :: cue, fxyze, bxyze
    complex, dimension(:,:,:), pointer :: exyz, bxyz
    complex, dimension(:,:), pointer :: ffc
    integer, dimension(:), pointer :: mixup
    complex, dimension(:), pointer :: sct
!
! declare arrays for OpenMP (tiled) code
    real, dimension(:,:,:), pointer :: ppart, ppbuff
    integer, dimension(:), pointer :: kp
    integer, dimension(:,:), pointer :: ncl
    integer, dimension(:,:,:), pointer :: ihole
!
! declare and initialize timing data
    real :: time
    integer, dimension(4) :: itime
    real :: tdpost = 0.0, tguard = 0.0, tfft = 0.0, tfield = 0.0
    real :: tdjpost = 0.0, tpush = 0.0, tsort = 0.0
    double precision :: dtime
!
    irc = 0

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! nvp = number of shared memory nodes (0=default)
  nvp = 0
!   write (*,*) 'enter number of nodes:'
!   read (5,*) nvp
! initialize for shared memory parallel processing
  call INIT_OMP(nvp)
!
! initialize scalars for standard code
  np = npx*ncpy; nx = 2**indx; ny = 2**indy; nxh = nx/2; nyh = ny/2
  nxe = nx + 2; nye = ny + 1; nxeh = nxe/2
  nxyh = max(nx,ny)/2; nxhy = max(nxh,ny)
  mx1 = (nx - 1)/mx + 1; my1 = (ny - 1)/my + 1; mxy1 = mx1*my1
  nloop = tend/dt + .0001; ntime = 0
  qbme = qme
  affp = real(nx*ny)/real(np)
  dth = 0.0
!
! allocate and initialize data for standard code
  allocate(part(idimp,np))
  allocate(qe(nxe,nye),fxyz(ndim,nxe,nye))
  allocate(cue(ndim,nxe,nye),bxyz(ndim,nxe,nye))
  allocate(exyz(ndim,nxeh,nye),bxyz(ndim,nxeh,nye))
  allocate(ffc(nxh,nyh),mixup(nxhy),sct(nxyh))
  allocate(kpic(mxy1))
!
! prepare fft tables
  call WFFT2RINIT(mixup,sct,indx,indy,nxhy,nxyh)
! calculate form factors
  isign = 0
  call MPOIS23(qe,fxyz,isign,ffc,ax,ay,affp,we,nx,ny,nxeh,nye,nxh, &
    &nyh)
! initialize electrons
  call DISTR2H(part,vtx,vty,vtz,vx0,vy0,vz0,npx,ncpy,idimp,np,nx,ny, &
    &ipbc)
!
! initialize transverse electromagnetic fields
  exyz = cmplx(0.0,0.0)
  bxyz = cmplx(0.0,0.0)
!
! find number of particles in each of mx, my tiles: updates kpic, nppmx
  call DBLKP2L(part,kpic,nppmx,idimp,np,mx,my,mx1,mxy1,irc)
  if (irc /= 0) then
    write (*,*) 'DBLKP2L error, irc=', irc
    stop
  endif
! allocate vector particle data
  nppmx0 = (1.0 + xtras)*nppmx
  ntmax = xtras*nppmx
  npbm = xtras*nppmx
  allocate(ppart(idimp,nppmx0,mxy1))
  allocate(ppbuff(idimp,npbm,mxy1))
  allocate(ncl(8,mxy1))
  allocate(ihole(2,ntmax+1,mxy1))
! copy ordered particle data for OpenMP

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        call PPMOVIN2L(part,ppart,kpic,nppmx0,idimp,np,mx,my,mx1,my1,irc)
        if (irc /= 0) then
            write (*,*) 'PPMOVIN2L overflow error, irc=', irc
            stop
        endif
! sanity check
        call PPCHECK2L(ppart,kpic,idimp,nppmx0,nx,ny,mx,my,mx1,my1,irc)
        if (irc /= 0) then
            write (*,*) 'PPCHECK2L error: irc=', irc
            stop
        endif
!
        if (dt > 0.45*ci) then
            write (*,*) 'Warning: Courant condition may be exceeded!'
        endif
!
! * * * start main iteration loop * * *
!
500 if (nloop <= ntime) go to 2000
!     write (*,*) 'ntime = ', ntime
!
! deposit current with OpenMP: updates ppart, cue, ncl, ihole, irc
        call dtimer(dtime,itime,-1)
        cue = 0.0
        if (relativity==1) then
            call GRJPPOSTF2L(ppart,cue,kpic,ncl,ihole,qme,dth,ci,nppmx0, &
&idimp,nx,ny,mx,my,nxe,nye,mx1,my1,ntmax,irc)
        else
            call GJPPOSTF2L(ppart,cue,kpic,ncl,ihole,qme,dth,nppmx0,idimp, &
&nx,ny,mx,my,nxe,nye,mx1,my1,ntmax,irc)
        endif
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tdjpost = tdjpost + time
        if (irc /= 0) then
            if (relativity==1) then
                write (*,*) 'GRJPPOSTF2L error: irc=', irc
            else
                write (*,*) 'GJPPOSTF2L error: irc=', irc
            endif
            stop
        endif
!
! reorder particles by cell with OpenMP:
! updates ppart, ppbuff, kpic, ncl, and irc
        call dtimer(dtime,itime,-1)
        call PORDERF2L(ppart,ppbuff,kpic,ncl,ihole,idimp,nppmx0,mx1,my1, &
&nxbmx,ntmax,irc)
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tsort = tsort + time
        if (irc /= 0) then
            write (*,*) 'current PORDERF2L error: ntmax, irc=', ntmax, irc
            stop
        endif

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        endif
!
! deposit charge with OpenMP: updates qe
    call dtimer(dtime,itime,-1)
    qe = 0.0
    call GPPOST2L(ppart,qe,kpic,qme,nppmx0,idimp,mx,my,nxe,nye,mx1,    &
&mxy1)
    call dtimer(dtime,itime,1)
    time = real(dtime)
    tdpost = tdpost + time
!
! add guard cells with OpenMP: updates cue, qe
    call dtimer(dtime,itime,-1)
    call ACGUARD2L(cue,nx,ny,nxe,nye)
    call AGUARD2L(qe,nx,ny,nxe,nye)
    call dtimer(dtime,itime,1)
    time = real(dtime)
    tguard = tguard + time
!
! transform charge to fourier space with OpenMP: updates qe
    call dtimer(dtime,itime,-1)
    isign = -1
    call WFFT2RMX(qe,isign,mixup,sct,indx,indy,nxeh,nye,nxhy,nxyh)
    call dtimer(dtime,itime,1)
    time = real(dtime)
    tfft = tfft + time
!
! transform current to fourier space with OpenMP: updates cue
    call dtimer(dtime,itime,-1)
    isign = -1
    call WFFT2RM3(cue,isign,mixup,sct,indx,indy,nxeh,nye,nxhy,nxyh)
    call dtimer(dtime,itime,1)
    time = real(dtime)
    tfft = tfft + time
!
! take transverse part of current with OpenMP: updates cue
    call dtimer(dtime,itime,-1)
    call MCUPERP2(cue,nx,ny,nxeh,nye)
    call dtimer(dtime,itime,1)
    time = real(dtime)
    tfield = tfield + time
!
! calculate electromagnetic fields in fourier space with OpenMP:
! updates exyz, bxyz
    call dtimer(dtime,itime,-1)
    if (ntime==0) then
        call MIBPOIS23(cue,bxyz,ffc,ci,wm,nx,ny,nxeh,nye,nxh,nyh)
        wf = 0.0
        dth = 0.5*dt
    else
        call MMAXWEL2(exyz,bxyz,cue,ffc,ci,dt,wf,wm,nx,ny,nxeh,nye,nxh,&
&nyh)
    endif
    call dtimer(dtime,itime,1)

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        time = real(dtime)
        tfield = tfield + time
!
! calculate force/charge in fourier space OpenMP: updates fxyz
        call dtimer(dtime,itime,-1)
        isign = -1
        call MPOIS23(qe,fxyz,isign,ffc,ax,ay,affp,we,nx,ny,nxeh,nye,nxh, &
&nyh)
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tfield = tfield + time
!
! add longitudinal and transverse electric fields with OpenMP:
! updates fxyz
        call dtimer(dtime,itime,-1)
        isign = 1
        call MEMFIELD2(fxyz,exyz,ffc,isign,nx,ny,nxeh,nye,nxh,nyh)
! copy magnetic field with OpenMP: updates bxyz
        isign = -1
        call MEMFIELD2(bxyz,bxyz,ffc,isign,nx,ny,nxeh,nye,nxh,nyh)
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tfield = tfield + time
!
! transform electric force to real space with OpenMP: updates fxyz
        call dtimer(dtime,itime,-1)
        isign = 1
        call WFFT2RM3(fxyz,isign,mixup,sct,indx,indy,nxeh,nye,nxhy,nxyh)
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tfft = tfft + time
!
! transform magnetic force to real space with OpenMP: updates bxyz
        call dtimer(dtime,itime,-1)
        isign = 1
        call WFFT2RM3(bxyz,isign,mixup,sct,indx,indy,nxeh,nye,nxhy,nxyh)
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tfft = tfft + time
!
! copy guard cells with OpenMP: updates fxyz, bxyz
        call dtimer(dtime,itime,-1)
        call BGUARD2L(fxyz,nx,ny,nxe,nye)
        call BGUARD2L(bxyz,nx,ny,nxe,nye)
        call dtimer(dtime,itime,1)
        time = real(dtime)
        tguard = tguard + time
!
! push particles with OpenMP: updates ppart, ncl, ihole, wke, irc
        wke = 0.0
        call dtimer(dtime,itime,-1)
        if (relativity==1) then
            call GRBPPUSHF23L(ppart,fxyz,bxyz,kpic,ncl,ihole,qbme,dt,dth,&
&ci,wke,idimp,nppmx0,nx,ny,mx,my,nxe,nye,mx1,mxy1,ntmax,irc)

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else
    call GBPPUSHF23L(ppart,fxyze,bxyze,kpic,ncl,ihole,qbme,dt,dth, &
&wke,idimp,nppmx0,nx,ny,mx,my,nxe,nye,mx1,my1,ntmax,irc)
endif
call dtimer(dtime,itime,1)
time = real(dtime)
tpush = tpush + time
if (irc /= 0) then
    if (relativity==1) then
        write (*,*) 'GRBPPUSHF23L error: irc=', irc
    else
        write (*,*) 'GBPPUSHF23L error: irc=', irc
    endif
    stop
endif
endif
!
! reorder particles by cell with OpenMP:
! updates ppart, ppbuff, kpic, ncl, and irc
    call dtimer(dtime,itime,-1)
    call PORDERF2L(ppart,ppbuff,kpic,ncl,ihole,idimp,nppmx0,mx1,my1, &
&npxbm,ntmax,irc)
    call dtimer(dtime,itime,1)
    time = real(dtime)
    tsort = tsort + time
    if (irc /= 0) then
        write (*,*) 'push PORDERF2L error: ntmax, irc=', ntmax, irc
        stop
    endif
endif
!
if (ntime==0) then
    wt = we + wf + wm
    write (*,*) 'Initial Total Field, Kinetic and Total Energies:'
    write (*, '(3e14.7)') wt, wke, wke + wt
    write (*,*) 'Initial Electrostatic, Transverse Electric and Mag&
&netic Field Energies:'
    write (*, '(3e14.7)') we, wf, wm
endif
ntime = ntime + 1
go to 500
2000 continue
!
! * * * end main iteration loop * * *
!
    write (*,*) 'ntime, relativity = ', ntime, relativity
    wt = we + wf + wm
    write (*,*) 'Final Total Field, Kinetic and Total Energies:'
    write (*, '(3e14.7)') wt, wke, wke + wt
    write (*,*) 'Final Electrostatic, Transverse Electric and Magnetic&
& Field Energies:'
    write (*, '(3e14.7)') we, wf, wm
!
    write (*,*)
    write (*,*) 'deposit time = ', tdpost
    write (*,*) 'current deposit time = ', tdjpost

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    tdpost = tdpost + tdjpost
    write (*,*) 'total deposit time = ', tdpost
    write (*,*) 'guard time = ', tguard
    write (*,*) 'solver time = ', tfield
    write (*,*) 'fft time = ', tfft
    write (*,*) 'push time = ', tpush
    write (*,*) 'sort time = ', tsort
    tfield = tfield + tguard + tfft
    write (*,*) 'total solver time = ', tfield
    time = tdpost + tpush + tsort
    write (*,*) 'total particle time = ', time
    wt = time + tfield
    write (*,*) 'total time = ', wt
    write (*,*)

!
    wt = 1.0e+09/(real(nloop)*real(np))
    write (*,*) 'Push Time (nsec) = ', tpush*wt
    write (*,*) 'Deposit Time (nsec) = ', tdpost*wt
    write (*,*) 'Sort Time (nsec) = ', tsort*wt
    write (*,*) 'Total Particle Time (nsec) = ', time*wt

!
    stop
end

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