

On the Role of Directives in High-Performance Computing

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CCDSC, October 2016

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OpenACC KERNELS as a construct ...

```
% pgfortran -fast -acc -Minfo -c PdV_kernel.f90
pdv_kernel:
...
77, Loop is parallelizable
79, Loop is parallelizable
Accelerator kernel generated
Generating Tesla code
77, !$acc loop gang, vector(4) ! blockidx%y
! threadidx%x
79, !$acc loop gang, vector(32)! blockidx%x
! threadidx%x
...
```

```
% pgfortran -fast -ta=multicore ... PdV_kernel.f90
pdv_kernel:
77, Loop is parallelizable
Generating Multicore code
77, !$acc loop gang
79, Loop is parallelizable
3 loop-carried redundant expressions removed
with 9 operations and 9 arrays
Innermost loop distributed: 2 new loops
Generated vector sse code for the loop
Generated 2 prefetch instructions for the loop
Generated 12 prefetch instructions for the loop
...
```

```
75 !$ACC KERNELS
76 !$ACC LOOP INDEPENDENT
77 DO k=y_min,y_max
78 !$ACC LOOP INDEPENDENT PRIVATE(right_flux,left_flux,top_flux,bottom_flux,total_flux,
min_cell_volume,energy_change,recip_volume)
79 DO j=x_min,x_max
80
81 left_flux= (xarea(j ,k )*(xvel0(j ,k )+xvel0(j ,k+1) &
82 +xvel0(j ,k )+xvel0(j ,k+1)))*0.25_8*dt*0.5 &
83 right_flux= (xarea(j+1,k )*(xvel0(j+1,k )+xvel0(j+1,k+1) &
84 +xvel0(j+1,k )+xvel0(j+1,k+1)))*0.25_8*dt*0.5 &
85 bottom_flux=(yarea(j ,k )*(yvel0(j ,k )+yvel0(j+1,k ) &
86 +yvel0(j ,k )+yvel0(j+1,k )))*0.25_8*dt*0.5 &
87 top_flux= (yarea(j ,k+1)*(yvel0(j ,k+1)+yvel0(j+1,k+1) &
88 +yvel0(j ,k+1)+yvel0(j+1,k+1)))*0.25_8*dt*0.5 &
89 total_flux=right_flux-left_flux+top_flux-bottom_flux
90
91 volume_change(j,k)=volume(j,k)/(volume(j,k)+total_flux)
92
93 min_cell_volume=MIN(volume(j,k)+right_flux-left_flux+top_flux-bottom_flux &
94 ,volume(j,k)+right_flux-left_flux &
95 ,volume(j,k)+top_flux-bottom_flux)
96
97 recip_volume=1.0/volume(j,k)
98
99 energy_change=(pressure(j,k)/density0(j,k)+viscosity(j,k)/density0(j,k))*...
100
101 energy1(j,k)=energy0(j,k)-energy_change
102
103 density1(j,k)=density0(j,k)*volume_change(j,k)
104
105 ENDDO
106 ENDDO
107 !$ACC END KERNELS
```

... and a path to standard parallel languages

Fortran 2008/2015 DO CONCURRENT

- + True Parallel Loops
- + Loop-scope shared/private data
- No support for reductions
- No support for data management

```
78 DO CONCURRENT (k=y_min:y_max)
79 DO CONCURRENT (j=x_min:x_max) LOCAL(right_flux,left_flux,top_flux,bottom_flux,
total_flux,min_cell_volume,energy_change,
recip_volume)
80
81 left_flux= (xarea(j ,k )*(xvel0(j ,k )+xvel0(j ,k+1) &
82 +xvel0(j ,k )+xvel0(j ,k+1)))*0.25_8*dt*0.5 &
83 right_flux= (xarea(j+1,k )*(xvel0(j+1,k )+xvel0(j+1,k+1) &
84 +xvel0(j+1,k )+xvel0(j+1,k+1)))*0.25_8*dt*0.5 &
85 bottom_flux=(yarea(j ,k )*(yvel0(j ,k )+yvel0(j+1,k ) &
86 +yvel0(j ,k )+yvel0(j+1,k )))*0.25_8*dt*0.5 &
87 top_flux= (yarea(j ,k+1)*(yvel0(j ,k+1)+yvel0(j+1,k+1) &
88 +yvel0(j ,k+1)+yvel0(j+1,k+1)))*0.25_8*dt*0.5 &
89 total_flux=right_flux-left_flux+top_flux-bottom_flux
90
91 volume_change(j,k)=volume(j,k)/(volume(j,k)+total_flux)
92
93 min_cell_volume=MIN(volume(j,k)+right_flux-left_flux+top_flux-bottom_flux &
94 ,volume(j,k)+right_flux-left_flux &
95 ,volume(j,k)+top_flux-bottom_flux)
97 recip_volume=1.0/volume(j,k)
99 energy_change=(pressure(j,k)/density0(j,k)+viscosity(j,k)/density0(j,k))*...
101 energy1(j,k)=energy0(j,k)-energy_change
103 density1(j,k)=density0(j,k)*volume_change(j,k)
105 ENDDO
106 ENDDO
107
```