

# OpenMP

## ■ Introducción

## ■ Directivas

- ✓ Regiones paralelas
- ✓ Worksharing
- ✓ sincronizaciones
- ✓ Visibilidad datos

## ■ Implementación

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# OpenMP: introduction

## ■ Standard promoted by main manufacturers

- <http://www.openmp.org>
- Fortran
  - ✓ V1.0: Oct. 1997
  - ✓ V2.0: Nov. 2000
- C
  - ✓ V1.0: Oct. 1998
  - ✓ V2.0: March 2002
- Structure: Directives, clauses and run time calls

## ■ Managed by the ARB (Architectural Review Board)

- Companies: Intel, SGI, KAI, IBM, Compaq, HP, Fujitsu, ...
- Committees Fortran, C, Futures,...
- Teleconferences

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## Basic example

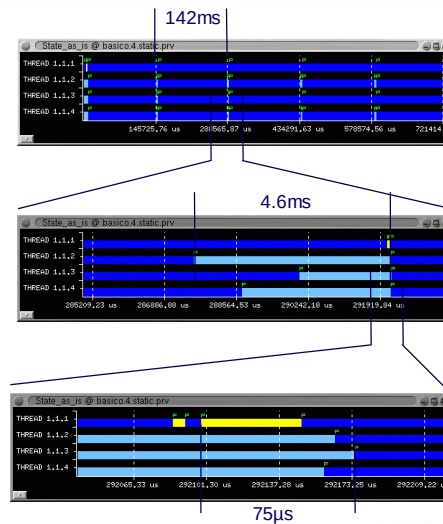
```

PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001
time = 10

DO 150 iter=1,5
C$OMP PARALLELDO SCHEDULE(STATIC)
C$OMP SHARED(dummy) PRIVATE(I,J)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    call delay(time)
200 CONTINUE
150 CONTINUE

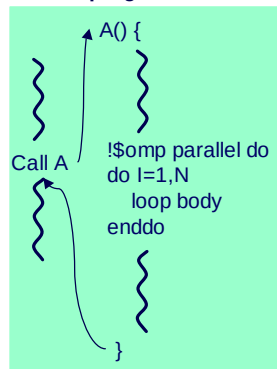
END
    
```



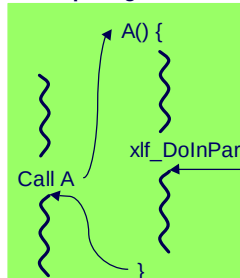
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## OpenMP compilation and Run Time

### Source program



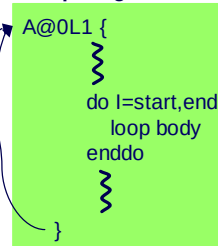
### Compiler generated



### libomp

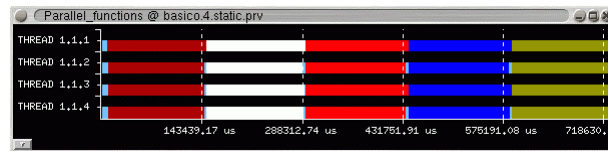


### Compiler generated



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## OpenMP outlined routines



- Display of the identifier (color encoded) of the outline routine being executed
- The compiler unrolls the iter loop !!

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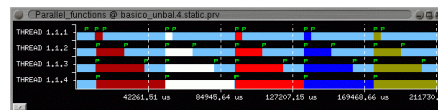
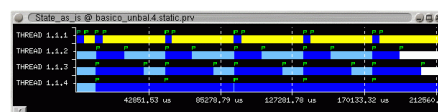
## Unbalanced loop

```
PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001

DO 150 iter=1,5
C$OMP PARALLELDO SCHEDULE(STATIC)
C$OMP SHARED(dummy) PRIVATE(I,J,time)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    time = i/100
    call delay(time)
200 CONTINUE
150 CONTINUE

END
```

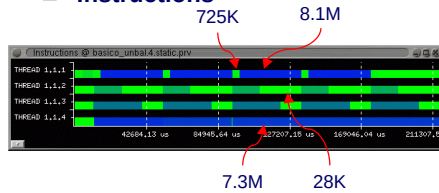


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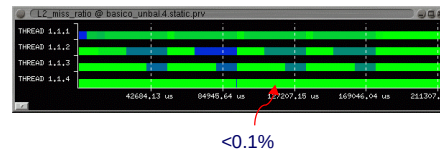


## Performance indices

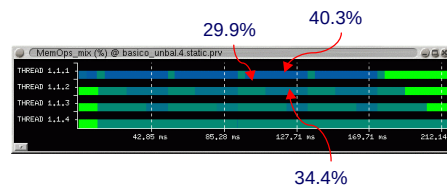
### Instructions



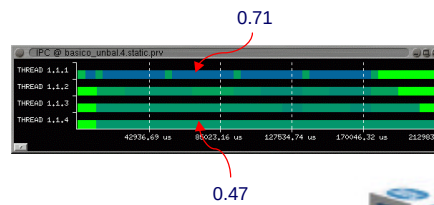
### L2 miss ratio



### Mem Ops mix



### IPC

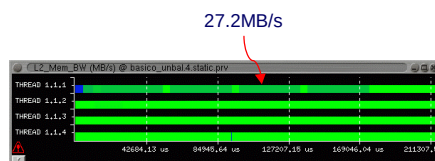


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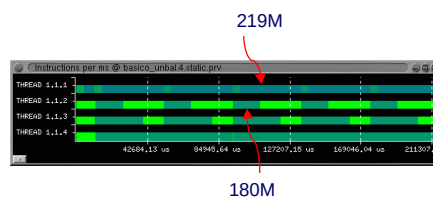


## Performance indices

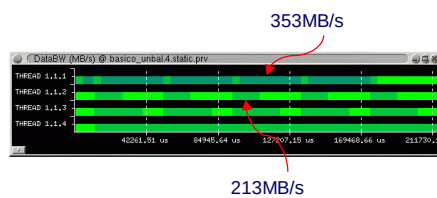
### Memory Bandwidth



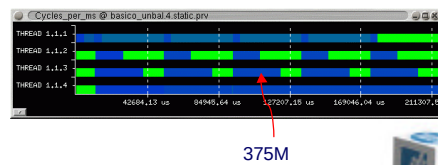
### MIPS



### CPU Bandwidth



### Cycles/s



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## Dynamic scheduling

```

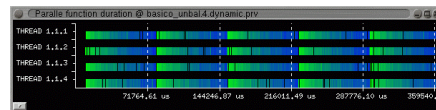
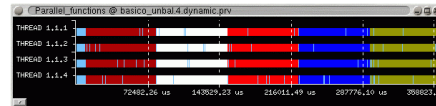
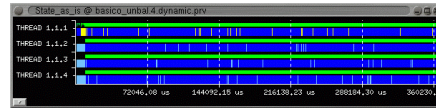
PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001

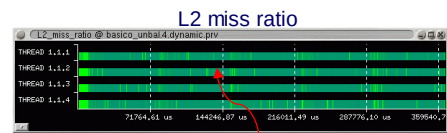
DO 150 iter=1,5
C$OMP PARALLELDO SCHEDULE(DYNAMIC)
C$OMP SHARED(dummy) PRIVATE(I,J,time)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    time = i/100
    call delay(time)
200 CONTINUE
150 CONTINUE

END
    
```

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Outline routine duration



Avg. 2.3%

## Dynamic scheduling

```

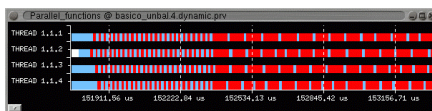
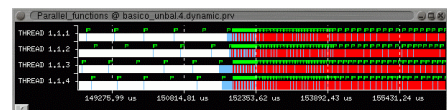
PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001

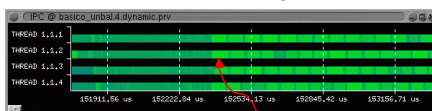
DO 150 iter=1,5
C$OMP PARALLELDO SCHEDULE(DYNAMIC)
C$OMP SHARED(dummy) PRIVATE(I,J,time)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    time = i/100
    call delay(time)
200 CONTINUE
150 CONTINUE

END
    
```

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IPC



0.25

## Coarser Grain Dynamic

```

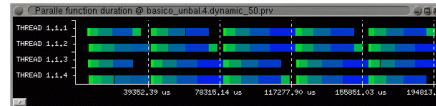
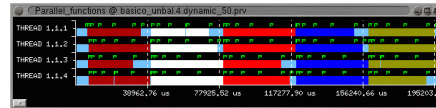
PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001

DO 150 iter=1,5
C$OMP PARALLEL DO SCHEDULE(DYNAMIC,50)
C$OMP SHARED(dummy) PRIVATE(I,J,time)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    time = i/100
    call delay(time)
200 CONTINUE
150 CONTINUE

END
    
```

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Parallel function duration

- Less overhead
- Some unbalance:
  - Heavy chunks towards the end



## Guided Scheduling

```

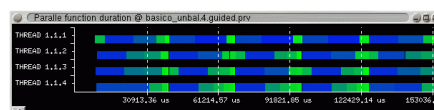
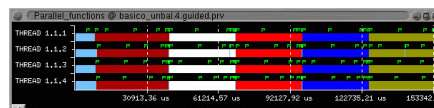
PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001

DO 150 iter=1,5
C$OMP PARALLEL DO SCHEDULE(GUIDED)
C$OMP SHARED(dummy) PRIVATE(I,J,time)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    time = i/100
    call delay(time)
200 CONTINUE
150 CONTINUE

END
    
```

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Parallel function duration

- Less overhead
- Good load balance:
  - Heavy chunks towards the beginning
- Dynamic:
  - Non repetitive pattern





## Always possible to fool a schedule

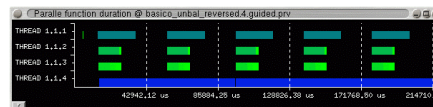
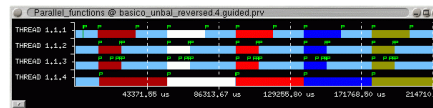
```

PROGRAM test
PARAMETER (N=1024)
REAL dummy(N), factor, var
REAL a(64000)
INTEGER i,j,time
common /varios/a

factor=1/1.0000001

DO 150 iter=1,5
C$OMP PARALLEL DO SCHEDULE(GUIDED)
C$OMP SHARED(dummy) PRIVATE(I,J,time)
DO 200 i=0,N
    dummy(i)= dummy(i)*factor
    time = (N-i)/100
    call delay(time)
200 CONTINUE
150 CONTINUE

END
    
```



Parallel function duration

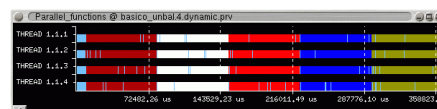
- **Dynamic:**
  - Repetitive pattern (not enforced)

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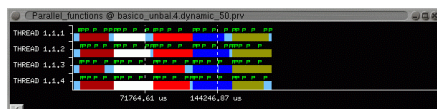


## Comparison

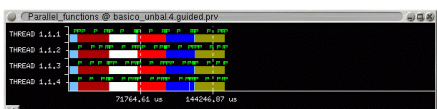
- **Dynamic**



- **Dynamic,50**



- **Guided**



Same scale

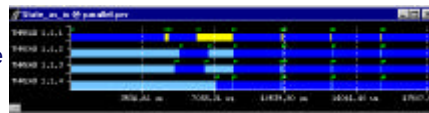
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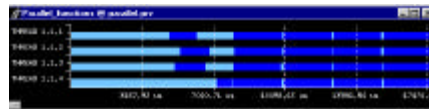
## Back to basics: Parallel directive

```
DO iter=1,5
C$OMP PARALLEL PRIVATE(time)
  time = 10
  call delay(time)
C$OMP ENDPARALLEL
ENDDO
```

State



Parallel functions



### ■ Team

- Threads to work on a parallel

### ■ Parallel

- All threads execute the body

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## Worksharings

### ■ Directives within parallel

### ■ Switch from replicated work to partitioned work

### ■ Work sharing constructs

#### • loops

- ✓ C\$OMP [END] DO [clause[[,] clause]...]
- ✓ C\$OMP [END] PARALLEL DO ...

#### • sections

- ✓ C\$OMP [END] SECTIONS [clause[[,] clause]...]
- ✓ C\$OMP SECTION
- ✓ C\$OMP [END] PARALLEL SECTIONS ...

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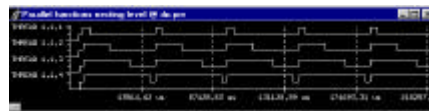
## Do

```
DO iter=1,5
C$OMP PARALLEL PRIVATE(time)
  time = 50
  call delay(time)
C$OMP DO SCHEDULE (STATIC)
  DO I=1,N
    time = i/100
    call delay(time)
  ENDDO
C$OMP ENDPARALLEL

ENDDO
```

Parallel  
function

Parallel  
function  
nesting  
level



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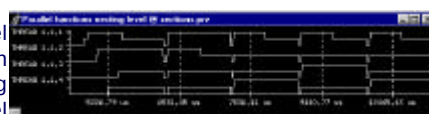
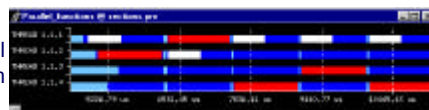
## Sections

```
DO iter=1,5
C$OMP PARALLEL PRIVATE(time)
C$OMP SECTIONS
C$OMP SECTION
  time = 30
  call delay(time)
C$OMP SECTION
  time = 60
  call delay(time)
C$OMP END SECTIONS
C$OMP ENDPARALLEL

ENDDO
```

Parallel  
function

Parallel  
function  
nesting  
level



### ■ Dynamic:

- Sections randomly taken by threads

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## Shortcuts

### ■ Parallel do

- Shortcut for do within parallel
- Allows more efficient implementation

### ■ Parallel sections

- Shortcut for sections within parallel
- Allows more efficient implementation

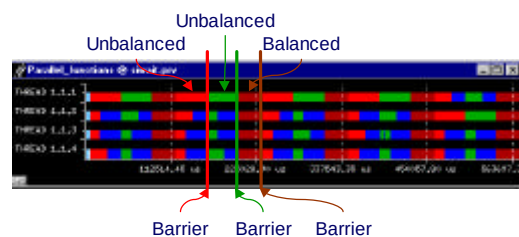
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## Relaxing synchronizations

```

DO iter=1,5
  CSOMP PARALLEL SHARED(dummy) PRIVATE(I,J, time)
  CSOMP DO SCHEDULE(GUIDED)
    DO i=0,N
      time = (N-i)/100
      call delay(time)
    ENDDO
  CSOMP DO SCHEDULE(GUIDED)
    DO i=0,N
      time = (N-i)/100
      call delay(time)
    ENDDO
  CSOMP ENDDO
  CSOMP DO SCHEDULE(GUIDED)
    DO i=0,N
      time = i/100
      call delay(time)
    ENDDO
  CSOMP END PARALLEL
ENDDO
  
```



### ■ Globally

- 2 unbalanced loops

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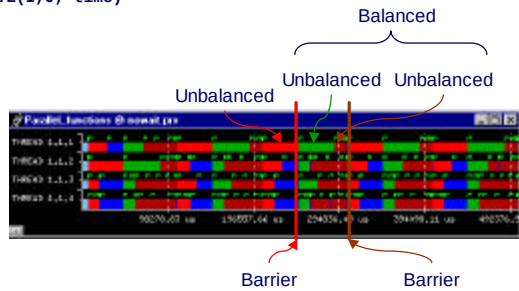


## Relaxing synchronizations: Nowait

```

DO iter=1,5
  CSOMP PARALLEL SHARED(dummy) PRIVATE(I,J, time)
  CSOMP DO SCHEDULE(GUIDED)
  DO i=0,N
    time = (N-i)/100
    call delay(time)
  ENDDO
  CSOMP DO SCHEDULE(GUIDED)
  DO i=0,N
    time = (N-i)/100
    call delay(time)
  ENDDO
  CSOMP ENDDO NOWAIT
  CSOMP DO SCHEDULE(GUIDED)
  DO i=0,N
    time = i/100
    call delay(time)
  ENDDO
  CSOMP END PARALLEL
ENDDO

```



### ■ Globally

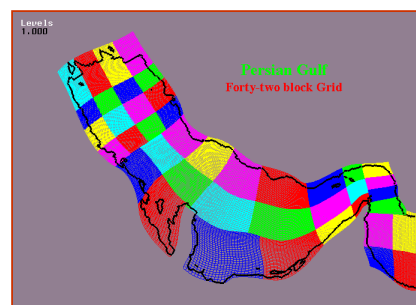
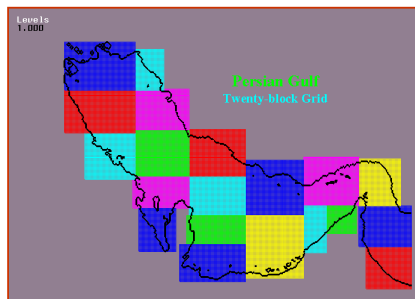
- Achieved global balance

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## SPMD parallelization

- Very coarse grain
  - Similarity to MPI
- Need for Barrier



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# Malleability

Scheduling decisions: Once for all

Explicit code only related to parallelism

```
C$OMP PARALLEL
WhoAmI=RunTimeCall()
myBlock=f(WhoAmI)
...
Call Compute1(myBlock)
...
DO iters=1, #iters
    Call Compute2(myBlock)
END DO
...
C$OMP END PARALLEL
```

```
C$OMP PARALLEL DO
DO Block=1, #blocks
    Call Compute1(Block)
END DO
C$OMP END PARALLEL
...
DO iter=1, #iters
    C$OMP PARALLEL DO
    DO Block=1, #blocks
        Call Compute1(Block)
    END DO
    C$OMP END PARALLEL
END DO
...
C$OMP END PARALLEL
```

myBlock ∈ [1, #processors] ~ f(resources)  
Block ∈ [1, #blocks] ~ f(algorithm)

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# Other computation patterns

## ■ Master

- C\$OMP [END] MASTER
- Only master threads executes. All other wait

## ■ Single

- C\$OMP [END] SINGLE [clause[[],] clause]...
- One thread executes
- Barrier at end

## ■ Ordered

- C\$OMP [END] ORDERED
- Ensures execution in sequential order

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## Other computation patterns

### ■ Critical

- `C$OMP [END] CRITICAL [(name)]`
- Mutual exclusion

### ■ Atomic

- `C$OMP ATOMIC`
- Atomicity of single assignment statement
  - ✓ Optimizable implementation

### ■ Barrier

- `C$OMP BARRIER`
- All threads must reach it before continuing

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## Other computation patterns

### ■ Flush

- `C$OMP FLUSH [(list)]`
- Enforce consistency

### ■ Lock routines:

- `OMP_INIT_LOCK (var)`
- `OMP_DESTROY_LOCK (var)`
- `OMP_SET_LOCK (var)`
- `OMP_UNSET_LOCK (var)`
- `OMP_TEST_LOCK (var)`

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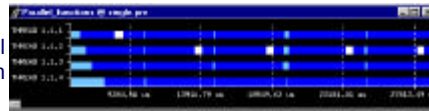
# Single

```

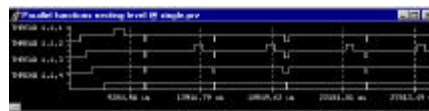
DO iter=1,5
  C$OMP PARALLEL PRIVATE(time)
    time = 50
    call delay(time)
  C$OMP SINGLE
    time = 25
    call delay(time)
  C$OMP END SINGLE
    time = 10
    call delay(time)
  C$OMP ENDPARALLEL
ENDDO

```

Parallel  
function



Parallel  
function  
nesting  
level



## ■ Implementaion

- outlined call
- Other alternatives possible

## ■ Duration

- Not proportional to time !!!
- ???



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# Who/how many?

## ■ Run time calls to find out

- `OMP_GET_NUM_THREADS`
  - ✓ How many are we?
- `OMP_GET_THREAD_NUM`
  - ✓ Who am I

## ■ ... or set

- `OMP_SET_NUM_THREADS (expr)`

## ■ Advice: minimize their use if possible

- Avoid the MPI/SPMD approach
- Write malleable codes



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## Data scope

- **PRIVATE(list)**
  - ~ allocate local space on entry to outlined routine
  - Scalars / vectors
- **FIRSTPRIVATE(list)**
  - Allocate and initialize
- **LASTPRIVATE(list)**
  - Copy value of last iteration to global
- **THREADPRIVATE(list)**
  - Privatization of common blocks

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## Data scope

- **REDUCTION(list)**
  - Perform reduction operation

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## Run time library

### ■ Compute schedules

### ■ Wait modes

- Overhead
- Behavior under overcommitted multiprogrammed workload
- Busy wait, yield, block
  - ✓ When to change mode?

### ■ Locks

### ■ Synchronizations

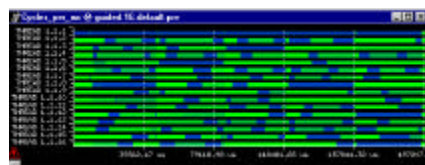
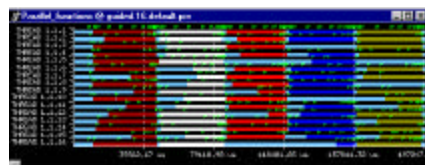
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## Run time library

### ■ Multiprogramming effects

- OMP\_NUM\_THREADS=16
- Guided
- Machine
  - ✓ 16 way SMP
  - ✓ load  $\cong$  9



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## Other topics

### ■ Array worksharings

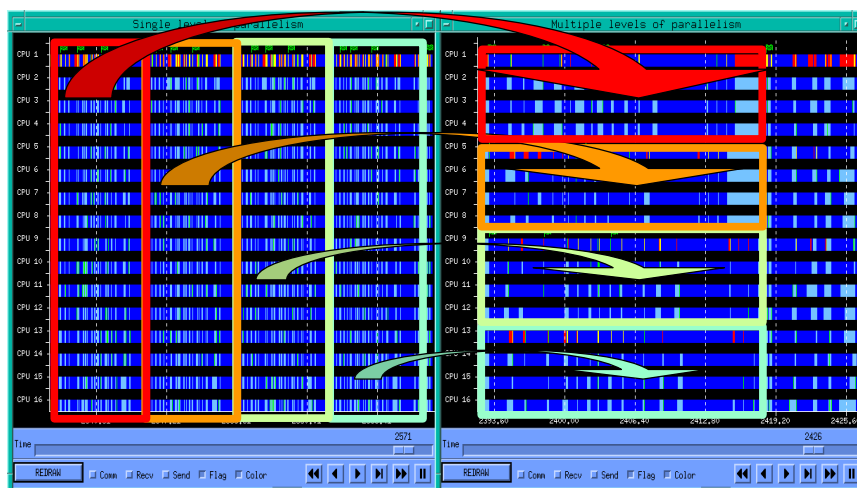
### ■ Future topics

- Nested parallelism
- Task queues
- Precedences
- Further schedules
- Multiprogrammed workload management

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## Nested parallelism



SPEC95 hydro2d



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## Single level precedences

### ■ Sections

- PRED/SUCC at any point

!\$omp parallel sections

!\$omp section **name** (S1) **succ** (S3)

!\$omp section **name** (S2) **succ** (S4)

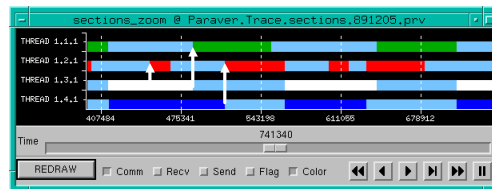
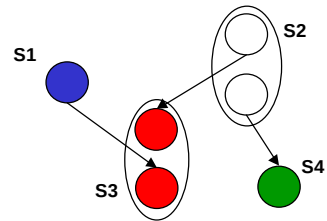
!\$omp **succ** (S3)

!\$omp section **name** (S3) **pred** (S2)

!\$omp **pred** (S1)

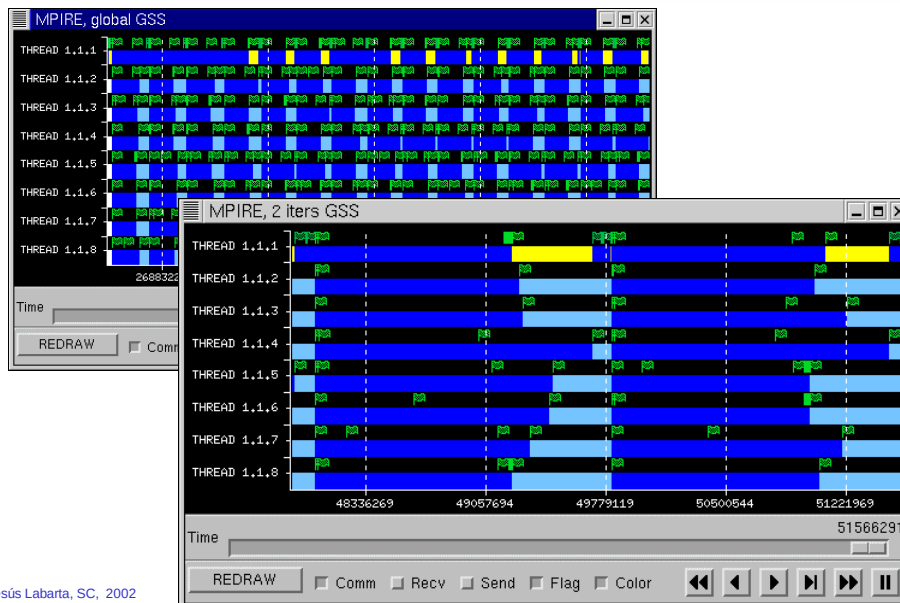
!\$omp section **name** (S4) **pred** (S2)

!\$omp end parallel sections



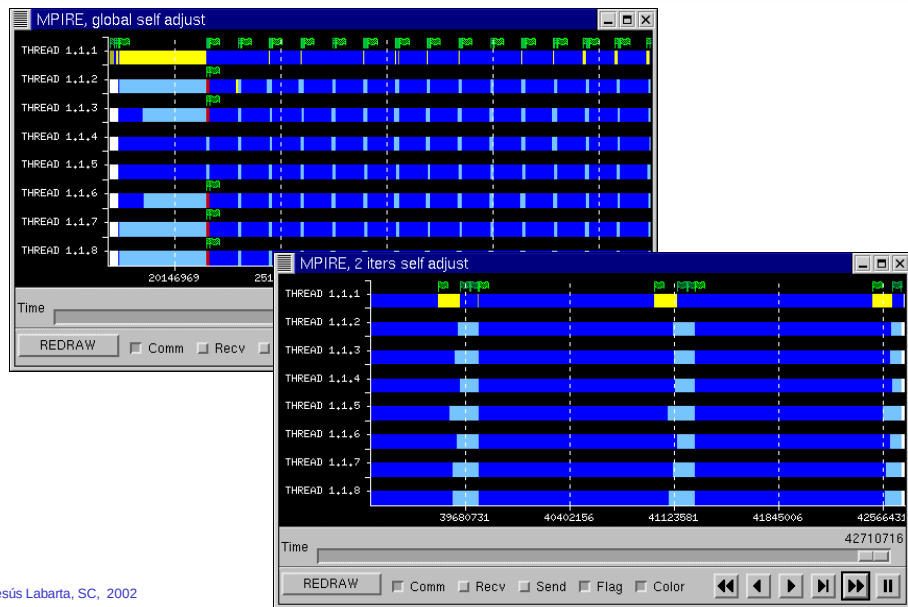
Jesús Labarta, SC, 2002

## Loop scheduling: MPIRE



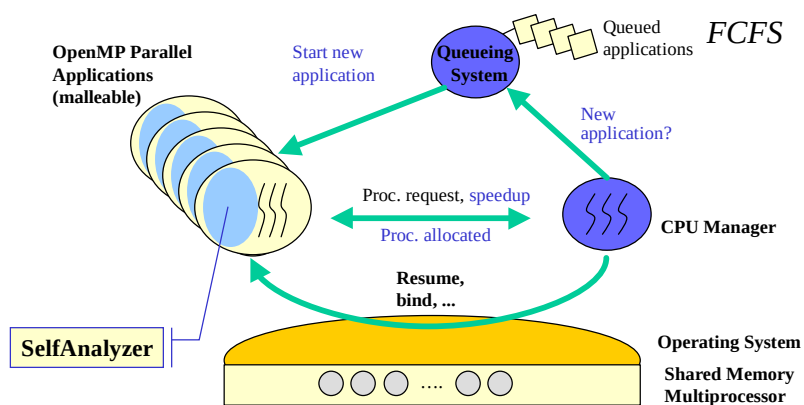
Jesús Labarta, SC, 2002

## Loop scheduling: MPIRE



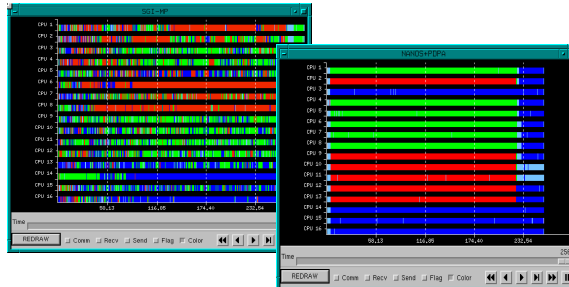
Jesús Labarta, SC, 2002

## NANOS OS scheduling environment



Jesús Labarta, SC, 2002

## CPU scheduling & memory



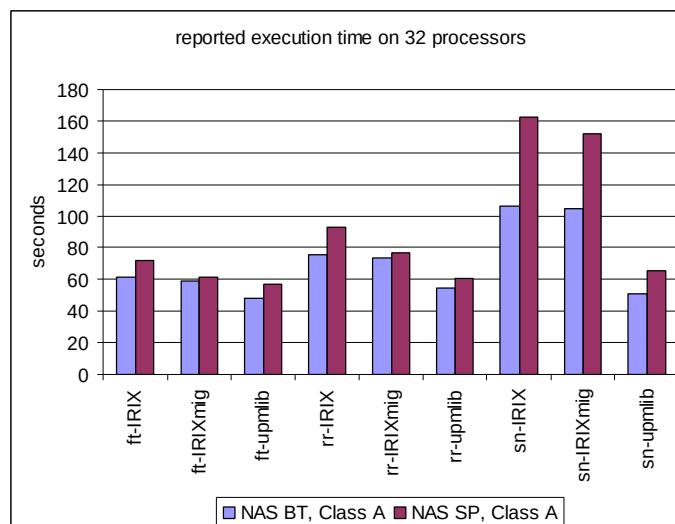
### ■ What is the real difference ?

- Time
- Context switch overhead
- Migrations

Jesús Labarta, SC, 2002



## Page migration



LHP

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