

# Using OpenMP



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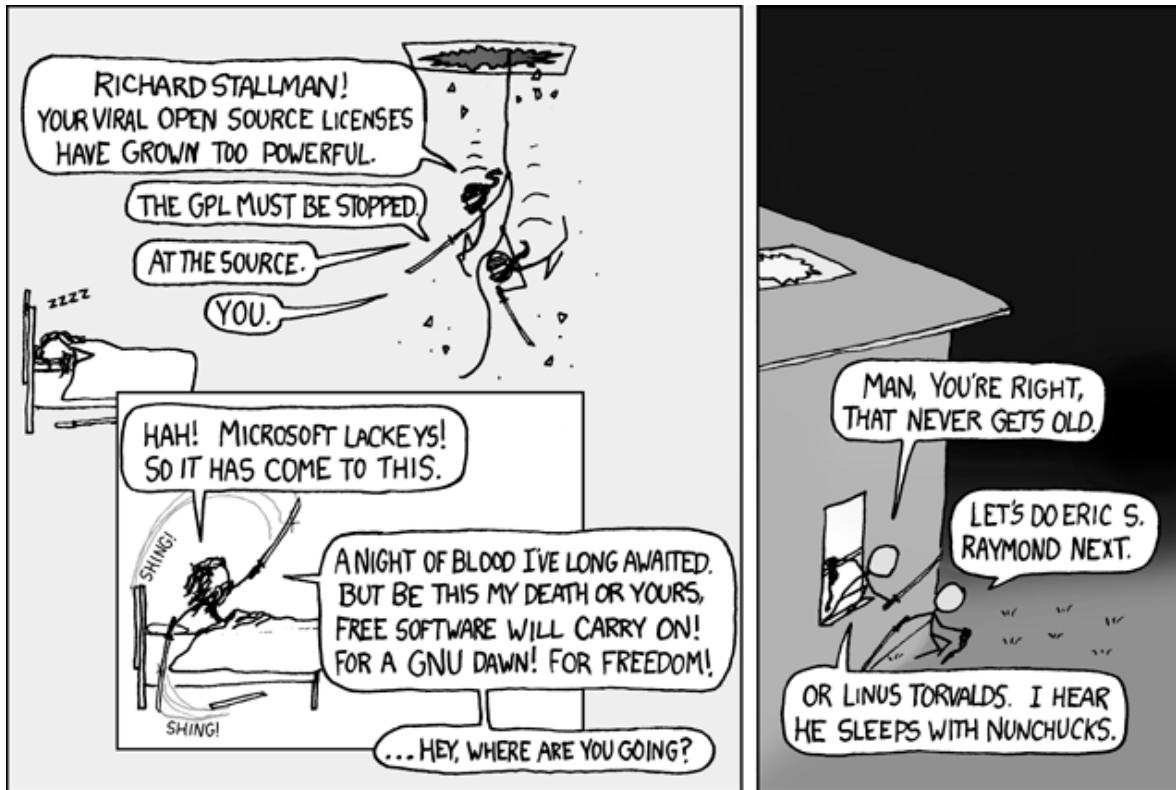


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# Outline

- I. About OpenMP
- II. OpenMP Directives
- III. Data Scope
- IV. Runtime Library Routines and Environment Variables
- V. Using OpenMP
- VI. Project: Computing Pi



# I. ABOUT OPENMP

Source: <http://xkcd.com/225/>

# About OpenMP

- Industry-standard shared memory programming model
- Developed in 1997
- OpenMP Architecture Review Board (ARB) determines additions and updates to standard

# Advantages to OpenMP

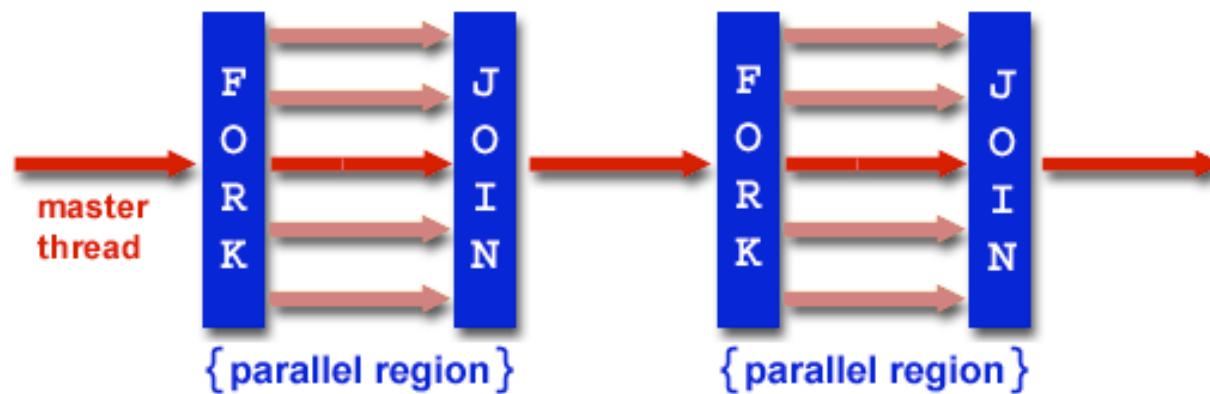
- Parallelize small parts of application, one at a time (beginning with most time-critical parts)
- Can express simple or complex algorithms
- Code size grows only modestly
- Expression of parallelism flows clearly, so code is easy to read
- Single source code for OpenMP and non-OpenMP – non-OpenMP compilers simply ignore OMP directives

# OpenMP Programming Model

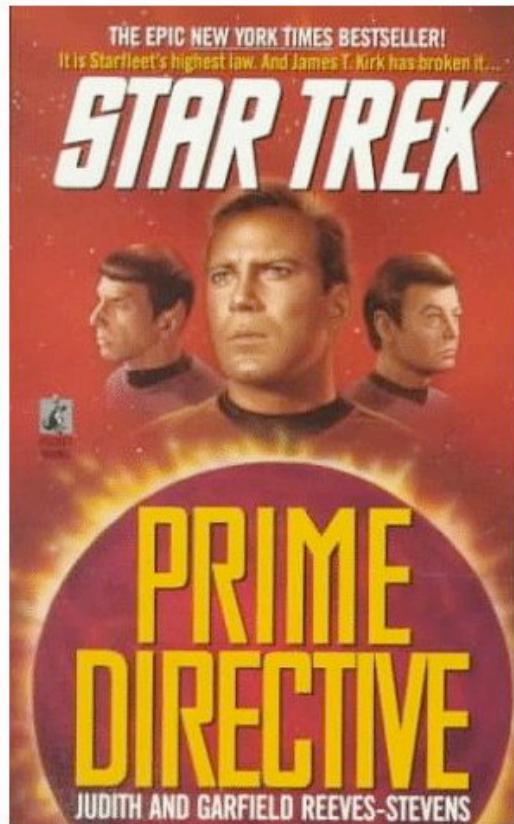
- Application Programmer Interface (API) is combination of
  - Directives
  - Runtime library routines
  - Environment variables
- API falls into three categories
  - Expression of parallelism (flow control)
  - Data sharing among threads (communication)
  - Synchronization (coordination or interaction)

# Parallelism

- Shared memory, thread-based parallelism
- Explicit parallelism (parallel regions)
- Fork/join model



Source: <https://computing.llnl.gov/tutorials/openMP/>



## II. OPENMP DIRECTIVES

*Star Trek: Prime Directive* by Judith and Garfield Reeves-Stevens, ISBN 0671744666

## II. OpenMP Directives

- **Syntax overview**
- **Parallel**
- **Loop**
- **Sections**
- **Synchronization**
- **Reduction**

# Syntax Overview: C/C++

- Basic format

```
#pragma omp directive-name [clause] newline
```

- All directives followed by newline
- Uses pragma construct (pragma = Greek for “thing”)
- Case sensitive
- Directives follow standard rules for C/C++ compiler directives
- Long directive lines can be continued by escaping newline character with \

# Syntax Overview: Fortran

- Basic format:

*sentinel directive-name [clause]*

- Three accepted sentinels: !\$omp \*\$omp c\$omp

- Some directives paired with **end** clause

- Fixed-form code:

- Any of three sentinels beginning at column 1
  - Initial directive line has space/zero in column 6
  - Continuation directive line has non-space/zero in column 6
  - Standard rules for fixed-form line length, spaces, etc. apply

- Free-form code:

- !\$omp only accepted sentinel
  - Sentinel can be in any column, but must be preceded by only white space and followed by a space
  - Line to be continued must end in & and following line begins with sentinel
  - Standard rules for free-form line length, spaces, etc. apply

# OpenMP Directives: Parallel

- A block of code executed by multiple threads
- Syntax:

```
#pragma omp parallel private(list) \
    shared(list)
{
    /* parallel section */
}

 !$omp parallel private(list) &
 !$omp shared(list)
 ! Parallel section
 !$omp end parallel
```

# Simple Example (C/C++)

```
#include <stdio.h>
#include <omp.h>

int main (int argc, char *argv[]) {
    int tid;
    printf("Hello world from threads:\n");
    #pragma omp parallel private(tid)
    {
        tid = omp_get_thread_num();
        printf("<%d>\n", tid);
    }
    printf("I am sequential now\n");
    return 0;
}
```

# Simple Example (Fortran)

```
program hello

integer tid, omp_get_thread_num

write(*,*) 'Hello world from threads:'

!$OMP parallel private(tid)
tid = omp_get_thread_num()
write(*,*) '<', tid, '>'

!$omp end parallel

write(*,*) 'I am sequential now'

end
```

# Output (Simple Example)

## Output 1

Hello world from  
threads:

<0>  
<1>  
<2>  
<3>  
<4>

I am sequential now

## Output 2

Hello world from  
threads:

<1>  
<2>  
<0>  
<4>  
<3>

I am sequential now

***Order of execution is scheduled by OS!!!!!!***

# OpenMP Directives: Loop

- Iterations of the loop following the directive are executed in parallel
- Syntax:

```
#pragma omp for schedule(type [,chunk]) \
private(list) shared(list) nowait
```

```
{  
    /* for loop */  
}
```

```
!$OMP do schedule(type [,chunk]) &  
!$OMP private(list) shared(list)
```

C do loop goes here

```
!$OMP end do nowait
```

- *type* = {static, dynamic, guided, runtime}

- If nowait specified, threads do not synchronize at end of loop

# Which Loops Are Parallelizable?

## Parallelizable

- Number of iterations known upon entry, and does not change
- Each iteration independent of all others
- No data dependence

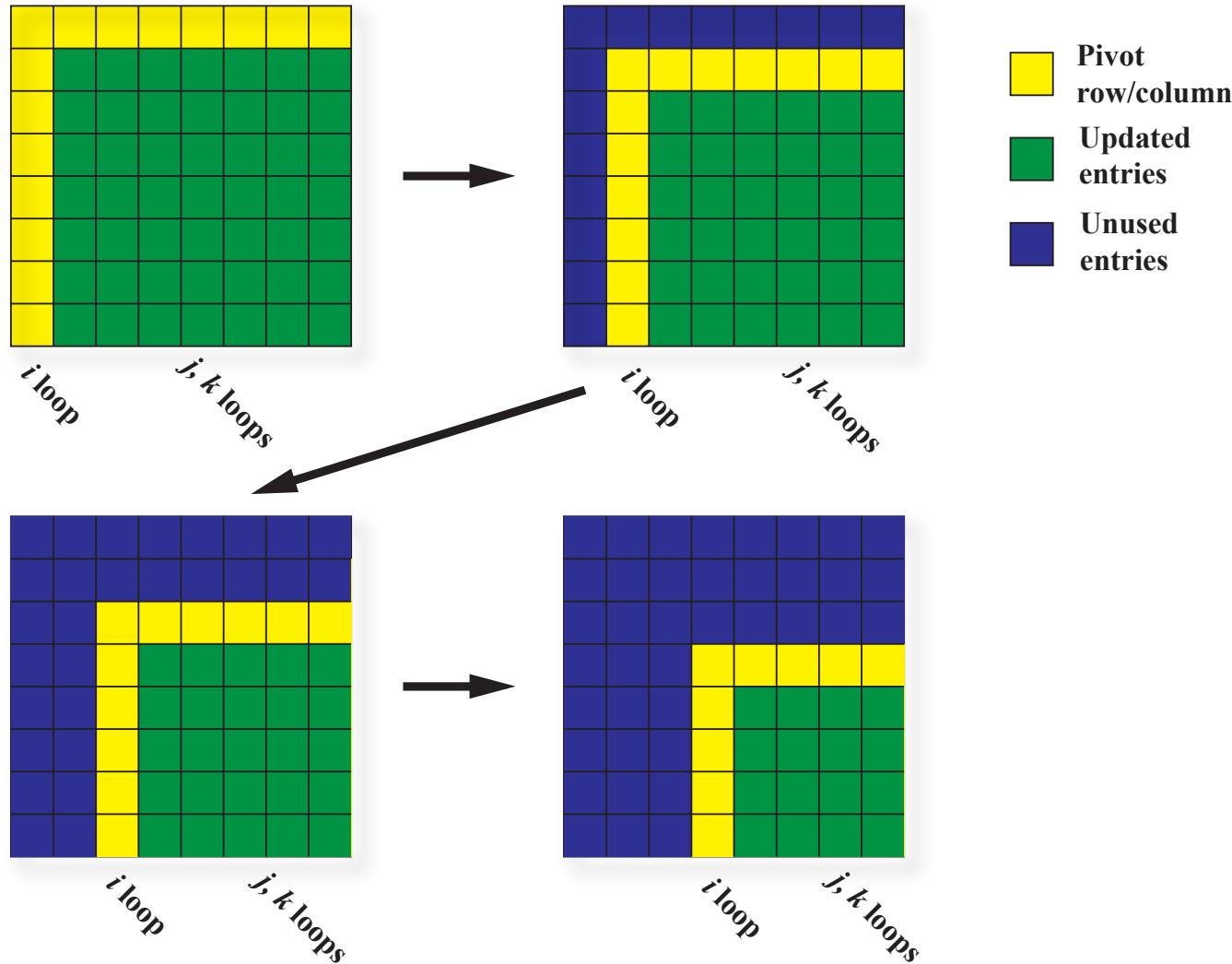
## Not Parallelizable

- Conditional loops (many while loops)
- Iterator loops (e.g., iterating over a `std::list<...>` in C++)
- Iterations dependent upon each other
- Data dependence

# Example: Parallelizable?

```
/* Gaussian Elimination (no pivoting) :  
   x = A\b                                         */  
  
for (int i = 0; i < N-1; i++) {  
    for (int j = i; j < N; j++) {  
        double ratio = A[j][i]/A[i][i];  
        for (int k = i; k < N; k++) {  
            A[j][k] -= (ratio*A[i][k]);  
            b[j] -= (ratio*b[i]);  
        }  
    }  
}
```

# Example: Parallelizable?



# Example: Parallelizable?

- Outermost Loop ( $i$ ):
  - $N-1$  iterations
  - Iterations depend upon each other (values computed at step  $i-1$  used in step  $i$ )
- Inner loop ( $j$ ):
  - $N-i$  iterations (constant for given  $i$ )
  - Iterations can be performed in any order
- Innermost loop ( $k$ ):
  - $N-i$  iterations (constant for given  $i$ )
  - Iterations can be performed in any order

# Example: Parallelizable?

```
/* Gaussian Elimination (no pivoting) :
   x = A\b                                         */

for (int i = 0; i < N-1; i++) {
#pragma omp parallel for
    for (int j = i; j < N; j++) {
        double ratio = A[j][i]/A[i][i];
        for (int k = i; k < N; k++) {
            A[j][k] -= (ratio*A[i][k]);
            b[j] -= (ratio*b[i]);
        }
    }
}
```

Note: can combine parallel and for into single pragma line

# OpenMP Directives: Loop Scheduling

- Default scheduling determined by implementation
- Static
  - ID of thread performing particular iteration is function of iteration number and number of threads
  - Statically assigned at beginning of loop
  - Load imbalance may be issue if iterations have different amounts of work
- Dynamic
  - Assignment of threads determined at runtime (round robin)
  - Each thread gets more work after completing current work
  - Load balance is possible

# Loop: Simple Example

```
#include <omp.h>
#define CHUNKSIZE 100
#define N      1000
int main ()  {
    int i, chunk;
    float a[N], b[N], c[N];
    /* Some initializations */
    for (i=0; i < N; i++)
        a[i] = b[i] = i * 1.0;
    chunk = CHUNKSIZE;
#pragma omp parallel shared(a,b,c,chunk) private(i)
{
    #pragma omp for schedule(dynamic,chunk) nowait
    for (i=0; i < N; i++)
        c[i] = a[i] + b[i];
} /* end of parallel section */
return 0;
}
```

# OpenMP Directives: Sections

- Non-iterative work-sharing construct
- Divide enclosed sections of code among threads
- Section directives nested within sections directive
- Syntax: C/C++

```
#pragma omp sections
{
    #pragma omp section
    /* first section */
    #pragma omp section
    /* next section */
}
```

Fortran

```
!$OMP sections
!$OMP section
C First section
!$OMP section
C Second section
!$OMP end sections
```

# Sections: Simple Example

```
#include <omp.h>
#define N      1000
int main () {
    int i;
    double a[N] , b[N] , c
        [N] , d[N];
    /* Some initializations
     */
    for (i=0; i < N; i++) {
        a[i] = i * 1.5;
        b[i] = i + 22.35;
    }
}
```

```
#pragma omp parallel \
    shared(a,b,c,d) private(i)
{
    #pragma omp sections nowait
    {
        #pragma omp section
            for (i=0; i < N; i++)
                c[i] = a[i] + b[i];
        #pragma omp section
            for (i=0; i < N; i++)
                d[i] = a[i] * b[i];
    } /* end of sections */
} /* end of parallel section */
return 0;
}
```

# OpenMP Directives: Synchronization

- Sometimes, need to make sure threads execute regions of code in proper order
  - Maybe one part depends on another part being completed
  - Maybe only one thread need execute a section of code
- Synchronization directives
  - Critical
  - Barrier
  - Single

# OpenMP Directives: Synchronization

- **Critical**

- Specifies section of code that must be executed by only one thread at a time

- Syntax: C/C++

```
#pragma omp critical [name]
```

Fortran

```
!$OMP critical [name]  
!$OMP end critical
```

- Names are global identifiers – critical regions with same name are treated as same region

- **Single**

- Enclosed code is to be executed by only one thread
  - Useful for thread-unsafe sections of code (e.g., I/O)

- Syntax: C/C++

```
#pragma omp single
```

Fortran

```
!$OMP single  
!$OMP end single
```

# OpenMP Directives: Synchronization

- **Barrier**
  - Synchronizes all threads: thread reaches barrier and waits until all other threads have reached barrier, then resumes executing code following barrier
  - Syntax: C/C++  
`#pragma omp barrier`
  - Fortran  
`!$OMP barrier`
  - Sequence of work-sharing and barrier regions encountered must be the same for every thread

# OpenMP Directives: Reduction

- Reduces list of variables into one, using operator (e.g., max, sum, product, etc.)
- Syntax

```
#pragma omp reduction(op : list)
```

```
!$OMP reduction(op : list)
```

where *list* is list of variables and *op* is one of following:

- C/C++: +, -, \*, &, ^, |, &&, or ||
- Fortran: +, -, \*, .and., .or., .eqv., .neqv., or max, min, iand, ior, ieor



### III. VARIABLE SCOPE

Angled spotting scope. Source: <http://www.spottingscopes.us/angled-scope-328.jpg>

# Variable Scope

- By default, all variables shared except
  - Certain loop index values – private by default
  - Local variables and value parameters within subroutines called within parallel region – private
  - Variables declared within lexical extent of parallel region – private

# Default Scope Example

```
void caller(int *a, int n) {  
    int i,j,m=3;  
    #pragma omp parallel for  
    for (i=0; i<n; i++) {  
        int k=m;  
        for (j=1; j<=5; j++) {  
            callee(&a[i], &k, j);  
        }  
    }  
  
    void callee(int *x, int *y, int  
               z) {  
        int ii;  
        static int cnt;  
        cnt++;  
        for (ii=1; ii<z; ii++) {  
            *x = *y + z;  
        }  
    }  
}
```

Var	Scope	Comment
a	shared	Declared outside parallel construct
n	shared	same
i	private	Parallel loop index
j	shared	Sequential loop index
m	shared	Declared outside parallel construct
k	private	Automatic variable/parallel region
x	private	Passed by value
*x	shared	(actually a)
y	private	Passed by value
*y	private	(actually k)
z	private	(actually j)
ii	private	Local stack variable in called function
cnt	shared	Declared static (like global)

# Variable Scope

- **Good programming practice: explicitly declare scope of all variables**
- **This helps you as programmer understand how variables are used in program**
- **Reduces chances of data race conditions or unexplained behavior**

# Variable Scope: Shared

- Syntax
  - `shared(list)`
- One instance of shared variable, and each thread can read or modify it
- **WARNING:** watch out for multiple threads simultaneously updating same variable, or one reading while another writes
- Example

```
#pragma omp parallel for shared(a)
for (i = 0; i < N; i++) {
    a[i] += i;
}
```

# Variable Scope: Shared – Bad Example

```
#pragma omp parallel for shared(n_eq)
for (i = 0; i < N; i++) {
    if (a[i] == b[i]) {
        n_eq++;
    }
}
```

- **n\_eq will not be correctly updated**
- Instead, put **n\_eq++ ;** in critical block (slow); introduce private variable **my\_n\_eq**, then update **n\_eq** in critical block after loop (faster); or use **reduction pragma (best)**

# Variable Scope: Private

- Syntax
  - `private(list)`
- Gives each thread its own copy of variable
- Example

```
#pragma omp parallel private(i, my_n_eq)
{
    #pragma omp for
    for (i = 0; i < N; i++) {
        if (a[i] == b[i]) my_n_eq++;
    }
    #pragma omp critical (update_sum)
    {
        n_eq+=my_n_eq;
    }
}
```

# Best Solution for Sum

```
#pragma parallel for reduction
(+:_eq)
for (i = 0; i < N; i++) {
    if (a[i] == b[i]) {
        _eq = _eq+1;
    }
}
```



## IV. RUNTIME LIBRARY ROUTINES AND ENVIRONMENT VARIABLES

Mt. McKinley National Monument, July, 1966. Source: National Park Service Historic Photograph Collection,  
[http://home.nps.gov/applications/hafe/hfc/npsphoto4h.cfm?Catalog\\_No=hpc-001845](http://home.nps.gov/applications/hafe/hfc/npsphoto4h.cfm?Catalog_No=hpc-001845)

# OpenMP Runtime Library Routines

- **void omp\_set\_num\_threads(int num\_threads)**  
**subroutine omp\_set\_num\_threads(scalar\_integer\_expression)**
  - Sets number of threads used in next parallel region
  - Must be called from serial portion of code

# OpenMP Runtime Library Routines

- **int omp\_get\_num\_threads()**  
**integer function omp\_get\_num\_threads()**
  - Returns number of threads currently in team executing parallel region from which it is called
- **int omp\_get\_thread\_num()**  
**integer function omp\_get\_thread\_num()**
  - Returns rank of thread
  - $0 \leq \text{omp\_get\_thread\_num}() < \text{omp\_get\_num\_threads}()$

# OpenMP Environment Variables

- Set environment variables to control execution of parallel code
- **OMP\_SCHEDULE**
  - Determines how iterations of loops are scheduled
  - E.g., `setenv OMP_SCHEDULE "guided, 4"`
- **OMP\_NUM\_THREADS**
  - Sets maximum number of threads
  - E.g., `setenv OMP_NUM_THREADS 4`



## V. USING OPENMP

# Conditional Compilation

- Can write single source code for use with or without OpenMP
- Pragmas/sentinels are ignored
- What about OpenMP runtime library routines?
  - `_OPENMP` macro is defined if OpenMP available: can use this to conditionally include `omp.h` header file, else redefine runtime library routines

# Conditional Compilation

```
#ifdef _OPENMP
    #include <omp.h>
#else
    #define omp_get_thread_num() 0
#endif

...
int me = omp_get_thread_num();
...
```



## VI. PROJECT: COMPUTING PI

Source: [http://www.ehow.com/how\\_2141082\\_best-berry-pie-ever.html](http://www.ehow.com/how_2141082_best-berry-pie-ever.html)

# Project Description

- We want to compute  $\pi$
- One method: method of darts\*
- Ratio of area of square to area of inscribed circle proportional to  $\pi$

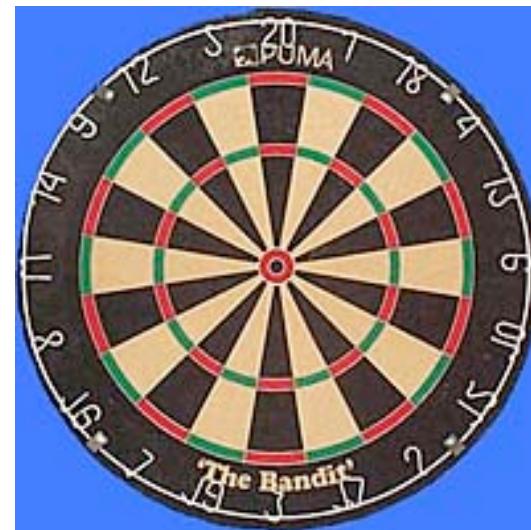


\*Disclaimer: this is a **TERRIBLE** way to compute  $\pi$ . Don't even think about doing it this way except for the purposes of this project!

# Method of Darts

- Imagine dartboard with circle of radius  $R$  inscribed in square

- Area of circle  $= \pi R^2$
- Area of square  $= (2R)^2 = 4R^2$
- $\frac{\text{Area of circle}}{\text{Area of square}} = \frac{\pi R^2}{4R^2} = \frac{\pi}{4}$



# Method of Darts

- So, ratio of areas proportional to  $\pi$
- How to find areas?
  - Suppose we threw darts (completely randomly) at dartboard
  - Could count number of darts landing in circle and total number of darts landing in square
  - Ratio of these numbers gives approximation to ratio of areas
  - Quality of approximation increases with number of darts
- $\pi = \frac{4 \times \# \text{ darts inside circle}}{\# \text{ darts thrown}}$



# Method of Darts

- Okay, Rebecca, but how in the world do we simulate this experiment on computer?
  - Decide on length  $R$
  - Generate pairs of random numbers  $(x, y)$  s.t.  
 $-R \leq x, y \leq R$
  - If  $(x, y)$  within circle (i.e. if  $(x^2+y^2) \leq R^2$ ), add one to tally for inside circle
  - Lastly, find ratio

# The Code (darts.c)\*

```
#include <omp.h>
#include "random.h"
static long num_trials = 10000;

int main() {
    long i;
    long Ncirc = 0;
    double pi, x, y;
    double r = 1.0; // radius of circle
    double r2 = r*r;
    for (i = 0; i < num_trials; i++) {
        x = random();
        y = random();
        if ((x*x + y*y) <= r2)
            Ncirc++;
    }
    pi = 4.0*(((double)Ncirc)/(double)num_trials);
    printf("\n For %d trials, pi = %f\n" ,num_trials, pi);
}
```

# The Code (random.h)\*

```
#include <omp.h>
/* Random number generator -- and not a very good one,
   either */
static long MULTIPLIER = 1366;
static long ADDEND = 150889;
static long PMOD = 714025;
long random_last = 0;

/* This is not a thread-safe random number generator */
double random() {
    long random_next;
    random_next = (MULTIPLIER * random_last + ADDEND) %
PMOD;
    random_last = random_next;
    return ((double)random_next/(double)PMOD);
}
```

# Your Mission (should you choose to accept it)

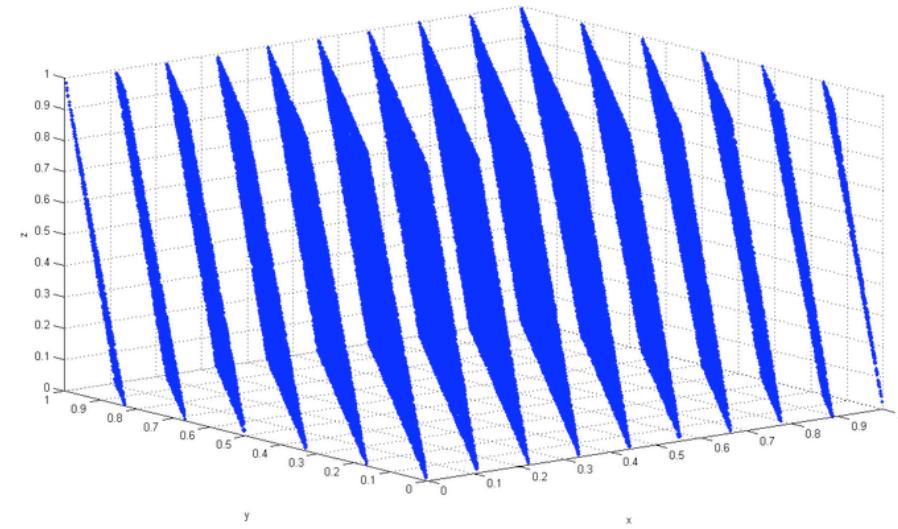
- Take provided code (darts.c, darts.cc, or darts.f) and parallelize with OpenMP
- Run with different numbers of threads and track performance and accuracy of solution
- Oops! Random number generator is not thread-safe. How can we fix this? (Discussion)

# Random Number Generator

- No such thing as random number generation – proper term is pseudorandom number generator (PRNG)
- Generate long sequence of numbers that seems “random”
- Properties of a good PRNG:
  - Very long period
  - Uniformly distributed
  - Reproducible
  - Quick and easy to compute

# Pseudorandom Number Generator

- Generator from `random.h` is Linear Congruential Generator (LCG)
  - Short period (= PMOD, 714025)
  - Not uniformly distributed
    - known to have correlations
  - Reproducible
  - Quick and easy to compute
  - Poor quality (don't do this at home)



Correlation of RANDU LCG (source:  
<http://en.wikipedia.org/wiki/File:Randu.png>)

# Pseudorandom Number Generator

- Generator is not thread-safe – how to fix it?
- Problem: all threads have access to random\_last
  - Second thread grabs random\_last before first thread updates it, resulting in duplicate results
  - Makes reproducible sequence irreproducible – will not happen the same way every time
  - How can we make generator thread-safe?
- Bonus fun: Try different solutions and profile their performance
  - Use `omp_get_wtime()` for timings (elapsed time = end - start)

# Bibliography/Resources: OpenMP

- Chapman, Barbara, Gabrielle Jost, and Ruud van der Pas. (2008) *Using OpenMP*, Cambridge, MA: MIT Press.
- Kendall, Ricky A. (2007) *Threads R Us*,  
[http://www.nccs.gov/wp-content/training/  
scaling\\_workshop\\_pdfs/threadsRus.pdf](http://www.nccs.gov/wp-content/training/scaling_workshop_pdfs/threadsRus.pdf)
- Mattson, Tim, and Larry Meadows (2008) SC08 OpenMP “Hands-On” Tutorial,  
<http://openmp.org/mp-documents/omp-hands-on-SC08.pdf>
- LLNL OpenMP Tutorial,  
<https://computing.llnl.gov/tutorials/openMP/>
- OpenMP.org, <http://openmp.org/>
- OpenMP 3.0 API Summary Cards:
  - Fortran: <http://openmp.org/mp-documents/OpenMP3.0-FortranCard.pdf>
  - C/C++:  
<http://www.openmp.org/mp-documents/OpenMP3.0-SummarySpec.pdf>

## Appendix: Better Ways to Compute $\pi$

- Look it up on the internet, e.g.  
<http://oldweb.cecm.sfu.ca/projects/ISC/data/pi.html>
- Compute using the BBP (Bailey-Borwein-Plouffe) formula

$$\pi = \sum_{n=0}^{\infty} \left( \frac{4}{8n+1} - \frac{2}{8n+4} - \frac{1}{8n+5} - \frac{1}{8n+6} \right) \left( \frac{1}{16} \right)^n$$

- For less accurate computations, try your programming language's constant, or quadrature or power series expansions

# Appendix: Better Ways to Generate Pseudorandom Numbers

- For serial codes
  - Mersenne twister
  - GSL (Gnu Scientific Library), many generators available (including Mersenne twister)  
<http://www.gnu.org/software/gsl/>
- For parallel codes
  - SPRNG, regarded as leading parallel pseudorandom number generator <http://sprng.cs.fsu.edu/>
  - PPRNG, Bill Cochran's new parallel pseudorandom number generator, supposedly superior to SPRNG  
<http://runge.cse.uiuc.edu/~wkcochra/pprng/>