

# Chapter 19 OpenMP

Speaker: Lung-Sheng Chien

Reference: [1] OpenMP C and C++ Application Program Interface v2.0  
[2] OpenMP C and C++ Application Program Interface v3.0  
[3] OpenMP forum, <http://www.openmp.org/forum/>  
[4] OpenMP tutorial: <https://computing.llnl.gov/tutorials/openMP/>  
[5] Getting Started with OpenMP:  
[http://rac.uits.iu.edu/hpc/openmp\\_tutorial/C/](http://rac.uits.iu.edu/hpc/openmp_tutorial/C/)

# OutLine

- OpenMP introduction
  - shared memory architecture
  - multi-thread
- Example 1: hello world
- Example 2: vector addition
- enable openmp in vc2005
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

# What is OpenMP

<http://en.wikipedia.org/wiki/OpenMP>

- The **OpenMP** (Open Multi-Processing) is an **application programming interface** (API) that supports multi-platform **shared memory multiprocessing** programming in **C/C++** and **Fortran** on many architectures, including **Unix** and **Microsoft Windows** platforms. It consists of a set of **compiler directives**, library routines, and **environment variables** that influence run-time behavior.
- OpenMP is a **portable**, scalable model that gives **programmers** a simple and flexible interface for developing parallel applications for platforms ranging from the desktop to the **supercomputer**.
- An application built with the hybrid model of **parallel programming** can run on a **computer cluster** using both OpenMP and **Message Passing Interface** (MPI).  
OpenMP: shared memory  
MPI: distributed memory

# History of OpenMP

- The OpenMP Architecture Review Board (ARB) published its first API specifications, OpenMP for Fortran 1.0, in October 1997. October the following year they released the C/C++ standard.
- 2000 saw version 2.0 of the Fortran specifications with version 2.0 of the C/C++ specifications being released in 2002.
- Version 2.5 is a combined C/C++/Fortran specification that was released in 2005.
- Version 3.0, released in May, 2008, is the current version of the API specifications. Included in the new features in 3.0 is the concept of *tasks* and the **task** construct. These new features are summarized in Appendix F of the [OpenMP 3.0 specifications](#).


# Goals of OpenMP

- **Standardization:**  
Provide a standard among a variety of **shared memory** architectures/platforms.
- **Lean and Mean:**  
establish a simple and limited set of directives for programming shared memory machines. Significant parallelism can be implemented by using just 3 or 4 directives.
- **Ease of Use:**
  - Provide capability to incrementally parallelize a serial program, unlike message-passing libraries which typically require an all or nothing approach
  - Provide the capability to implement both coarse-grain and fine-grain parallelism
- **Portability:**
  - Supports Fortran (77, 90, and 95), C, and C++
  - Public forum for API and membership

Website: <http://openmp.org/wp/>

# OpenMP™

THE OPENMP API SPECIFICATION FOR PARALLEL PROGRAMMING



RSS

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- » [OpenMP Specs](#)
- » [About OpenMP.org](#)
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**Events**

*The 5th International Workshop on OpenMP - Evolving OpenMP in an Age of Extreme Parallelism - will take place in Dresden (Germany) from 3rd June until 5th June 2009.*

**Input Register**

Alert the OpenMP.org

## OpenMP News

» SC08 OpenMP "Hands-On" Tutorial Available

Tim Mattson and Larry Meadows, both of Intel, presented a day-long tutorial introducing parallel programming with OpenMP at SC08 last week in Austin, TX.

The slides and class exercises from that tutorial are now available:

- [Hands-On Introduction to OpenMP](#), Mattson and Meadows, from SC08 (Austin) (PDF)
  - [Code Exercises](#) (zip)

Posted on November 24, 2008

» OpenMP 3.0 Status

» Christian Terboven reports:

SC08 brought us some pretty good news regarding availability of (full) support for OpenMP 3.0:

- Intel 11.0: Linux (x86), Windows (x86) and MacOS (x86)
- Sun Studio Express 11/08: Linux (x86) and Solaris (SPARC + x86)
- PGI 8.0: Linux (x86) and Windows (x86)
- IBM 10.1: Linux (POWER) and AIX (POWER)

GCC 4.4 will have support for OpenMP 3.0 as well, it is currently in *regression fixes and docs only*

**OpenMP.org**

The OpenMP Application Program Interface (API) supports multi-platform shared-memory parallel programming in C/C++ and Fortran. OpenMP is a portable, scalable model with a simple and flexible interface for developing parallel applications on platforms from the desktop to the supercomputer.

» [Read about OpenMP](#)

**Get It**

» [OpenMP specs](#)

**Use It**

» [OpenMP Compilers](#)

**Learn It**

Vendor/Source	Compiler	Information
» GNU	gcc (4.3.2)	Free and open source - Linux, Solaris, AIX, MacOSX, Windows  Compile with <code>-fopenmp</code>  » <a href="#">More information</a>
» Intel	C/C++ / Fortran (10.1)	Windows, Linux, and MacOSX.  Compile with <code>-Qopenmp</code> on Windows, or just <code>-openmp</code> on Linux or Mac OSX  » <a href="#">More information</a>

OpenMP forum: <http://www.openmp.org/forum/>

**OpenMP Forum**  
Discussion on the OpenMP specification run by the OpenMP ARB

[Board index](#)

It is currently

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FORUM	TOPICS	POSTS	LAST POST
 <b>General</b> General OpenMP discussion	321	1222	by Ernst0  on Thu Dec 18, 2008 1:51 pm
 <b>OpenMP 3.0 API Specifications</b> Discuss the OpenMP 3.0 API Specifications document in this forum.	4	14	by DeLoghi  on Fri Dec 05, 2008 11:36 pm
 <b>Draft 3.0 Public Comment</b> The public comment period closed January 31, 2008. This forum is now locked (read only).	19	59	by jakub  on Mon May 19, 2008 8:39 am

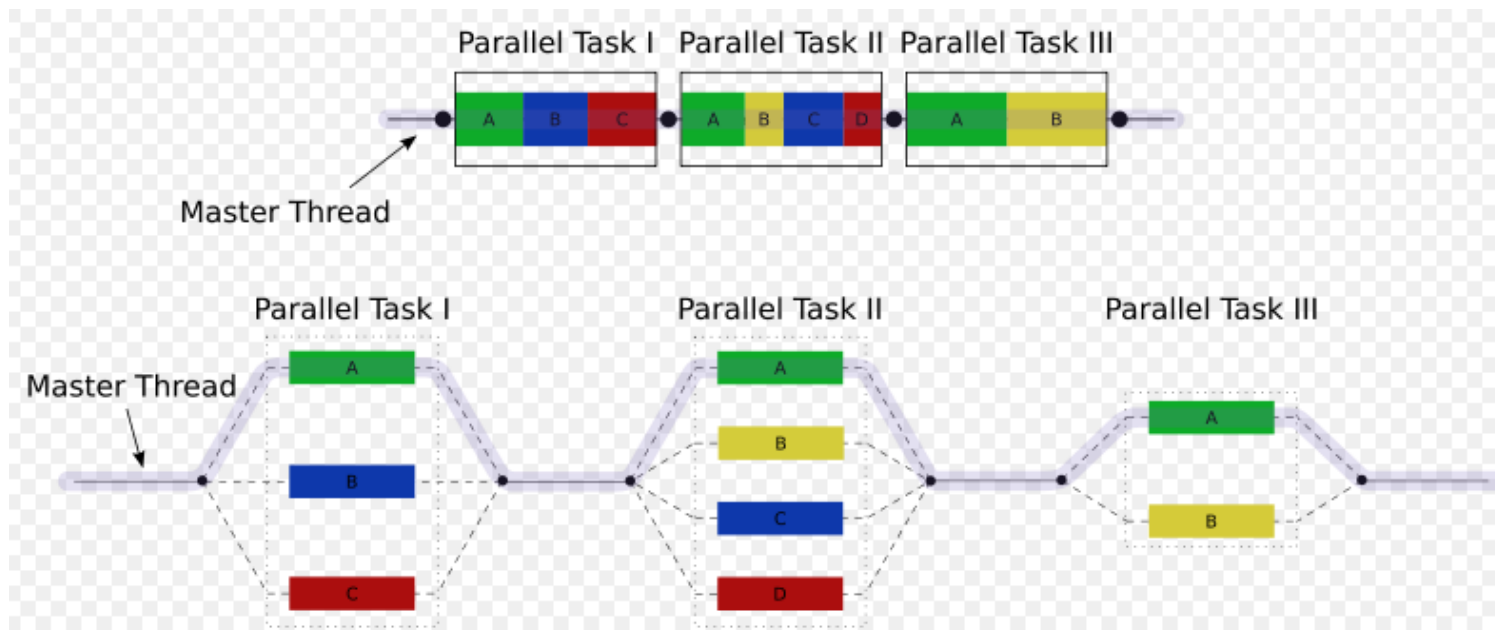
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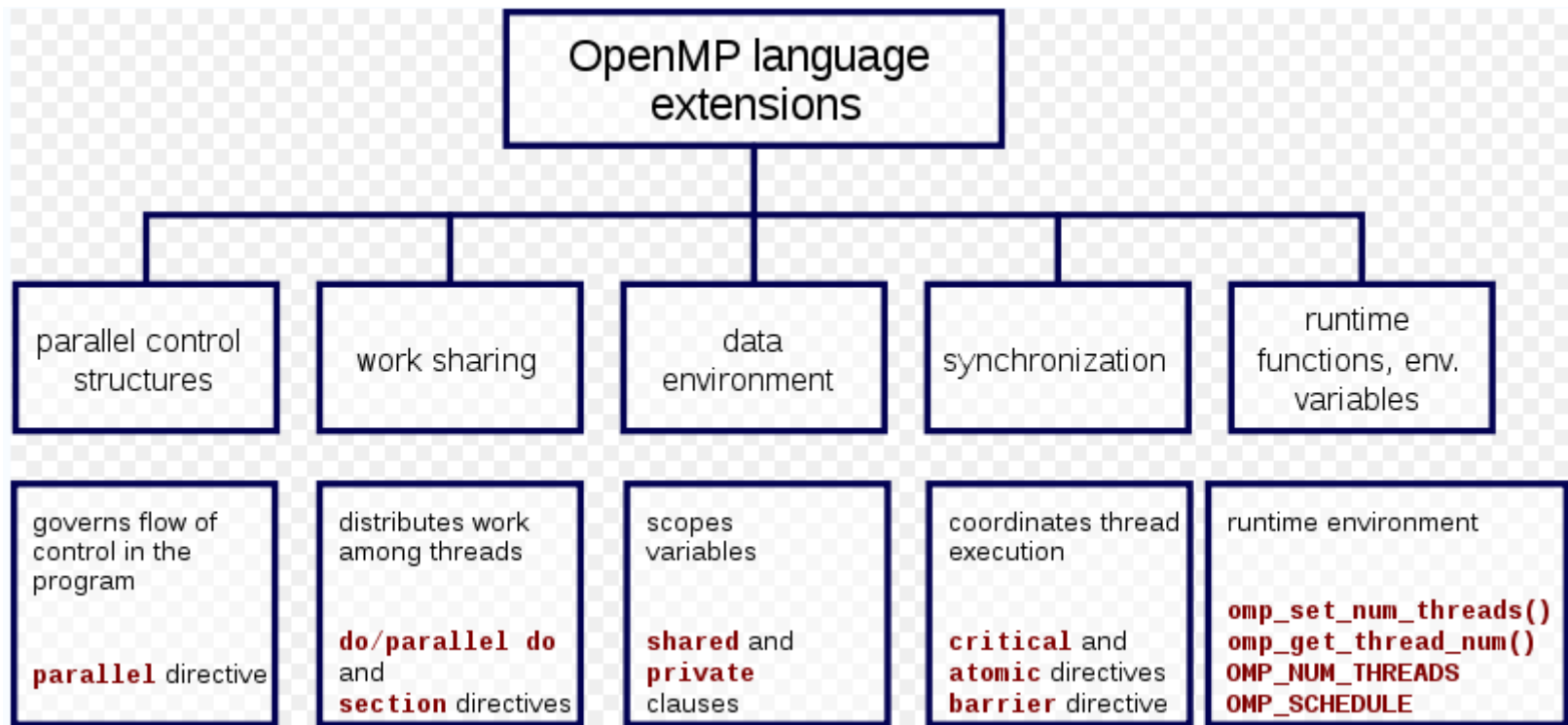
# Multithread (多執行緒)

- OpenMP is an implementation of **multithreading**, a method of parallelization whereby the master "thread" (a series of instructions executed consecutively) "forks" a specified number of slave "threads" and a task is divided among them. The threads then run concurrently, with the **runtime environment** allocating threads to different processors.
- The runtime environment allocates threads to processors depending on usage, machine load and other factors. The number of threads can be assigned by the runtime environment based on environment variables or in code using functions. The OpenMP functions are included in a header file labelled "omp.h" in C/C++





# Core elements

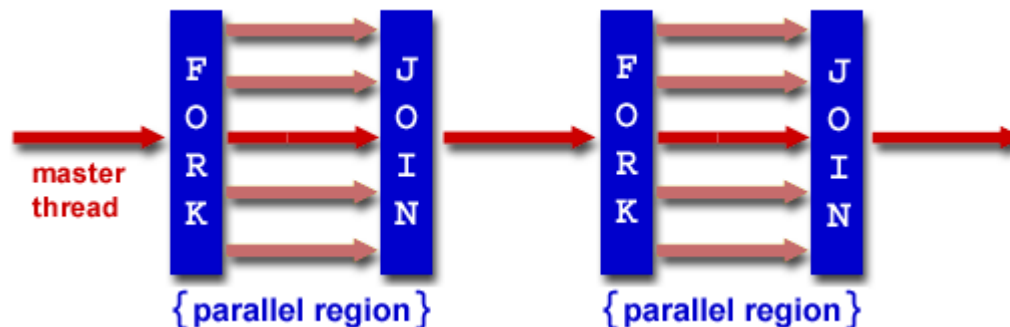


A compiler directive in C/C++ is called a *pragma* (pragmatic information). It is a preprocessor directive, thus it is declared with a hash (#). Compiler directives specific to OpenMP in C/C++ are written in codes as follows:

```
#pragma omp <rest of pragma>
```

# OpenMP programming model [1]

- **Shared Memory, Thread Based Parallelism:**  
OpenMP is based upon the existence of multiple threads in the shared memory programming paradigm. A shared memory process consists of multiple threads.
- **Explicit Parallelism:**  
OpenMP is an explicit (not automatic) programming model, offering the programmer full control over parallelization.
- **Fork - Join Model:**
  - OpenMP uses the fork-join model of parallel execution
  - All OpenMP programs begin as a single process: the **master thread**. The master thread executes sequentially until the first **parallel region** construct is encountered
  - **FORK**: the master thread then creates a **team of parallel threads**
  - The statements in the program that are enclosed by the parallel region construct are then executed in parallel among the various team threads
  - **JOIN**: When the team threads complete the statements in the parallel region construct, they synchronize and terminate, leaving only the master thread



## OpenMP programming model [2]

- **Compiler Directive Based:**  
OpenMP parallelism is specified through the use of compiler directives.
- **Nested Parallelism Support:**
  - The API provides for the placement of parallel constructs inside of other parallel constructs
  - Implementations may or may not support this feature.
- **Dynamic Threads:**
  - The API provides for dynamically altering the number of threads which may used to execute different parallel regions
  - Implementations may or may not support this feature.
- **I/O:**
  - OpenMP specifies nothing about parallel I/O. This is particularly important if multiple threads attempt to write/read from the same file.
  - If every thread conducts I/O to a different file, the issues are not as significant.
  - It is entirely up to the programmer to insure that I/O is conducted correctly within the context of a multi-threaded program.
- **FLUSH Often?:**
  - OpenMP provides a "relaxed-consistency" and "temporary" view of thread memory (in their words). In other words, threads can "cache" their data and are not required to maintain exact consistency with real memory all of the time.
  - When it is critical that all threads view a shared variable identically, the programmer is responsible for insuring that the variable is FLUSHed by all threads as needed.

# OutLine

- OpenMP introduction
- Example 1: hello world
  - parallel construct
- Example 2: vector addition
- enable openmp in vc2005
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

## Example 1: hello world [1]

### hello.c

```
1
2 #include <omp.h>
3 #include <stdio.h>
4
5 int main (int argc, char *argv[])
6 {
7     #pragma omp parallel
8     {
9         printf("Hello World\n");
10    }
11    return 0;
12 }
13
```

### Makefile

```
1
2 hello: hello.c
3     icpc -openmp -mp -c hello.c
4     icpc -openmp -o hello hello.o
5
```

header file "omp.h" is necessary for OpenMP programming

### MSDN library 2005

The **#pragma** directives offer a way for each compiler to offer machine- and operating system-specific features while retaining overall compatibility with the C and C++ languages. Pragas are machine- or operating system-specific by definition, and are usually different for every compiler.

If the compiler finds a pragma it does not recognize, it issues a warning, but compilation continues.

man icpc

**-openmp**

Enable the parallelizer to generate multi-threaded code based on the OpenMP\* directives. The code can be executed in parallel on both uniprocessor and multiprocessor systems. The **-openmp** option works with both -O0 (no optimization) and any optimization level of -O1, -O2 (default) and -O3. Specifying -O0 with **-openmp** helps to debug OpenMP applications.

## Example 1: hello world [2]

```
[macrold@quartet2 hello_word1]$ ls
Makefile hello.c
[macrold@quartet2 hello_word1]$ make hello
icpc -openmp -mp -c hello.c
hello.c(7): (col. 3) remark: OpenMP DEFINED REGION WAS PARALLELIZED.
icpc -openmp -o hello hello.o
[macrold@quartet2 hello_word1]$ ls
Makefile hello hello.c hello.o
[macrold@quartet2 hello_word1]$ ./hello
Hello World
Hello World
Hello World
Hello World
[macrold@quartet2 hello_word1]$
```

hello.c

```
1
2 #include <omp.h>
3 #include <stdio.h>
4
5 int main (int argc, char *argv[])
6 {
7     #pragma omp parallel
8     {
9         printf("Hello World\n");
10    }
11    return 0;
12 }
13
```

Machine quartet2 has 4 cores

```
top - 11:17:15 up 14 days, 22:54, 2 users, load average: 0.00, 0.00, 0.00
Tasks: 138 total, 1 running, 137 sleeping, 0 stopped, 0 zombie
Cpu0  :  0.0%us,  0.0%sy,  0.0%ni,100.0%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu1  :  0.0%us,  0.0%sy,  0.0%ni,100.0%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu2  :  0.0%us,  0.0%sy,  0.0%ni,100.0%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu3  :  0.0%us,  0.0%sy,  0.0%ni,100.0%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Mem:   8201628k total, 3524316k used, 4677312k free, 219744k buffers
Swap:  8193140k total,      0k used, 8193140k free, 2849960k cached
```

PID	USER	PR	NI	VIRT	RES	SHR	S	%CPU	%MEM	TIME+	COMMAND
14400	macrold	20	0	18936	1208	900	R	0	0.0	0:00.01	top
1	root	20	0	10328	688	580	S	0	0.0	0:05.36	init
2	root	15	-5	0	0	0	S	0	0.0	0:00.00	kthreadd

```
[macrold@quartet2 hello_word1]$ cat /proc/cpuinfo
```

```
model name      : Intel(R) Core(TM)2 Quad CPU    Q6600  @ 2.40GHz
stepping        : 11
cpu MHz         : 1596.000
cache size      : 4096 KB
```

## Example 1: hello world [3]

octet1

```
[macrold@octet1 hello_word1]$ ls
Makefile hello hello.c hello.o
[macrold@octet1 hello_word1]$ ./hello
```

```
Hello World
Hello World
Hello World
Hello World
Hello World
Hello World
Hello World
-
```

Machine octet1 has 8 cores (two quad-core)

```
top - 09:58:07 up 80 days, 18:39, 1 user, load average: 0.00, 0.00, 0.00
Tasks: 194 total, 1 running, 193 sleeping, 0 stopped, 0 zombie
Cpu0  :  1.5%us,  0.1%sy,  0.0%ni, 98.4%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu1  :  1.1%us,  0.1%sy,  0.0%ni, 98.8%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu2  :  0.9%us,  0.1%sy,  0.0%ni, 98.8%id,  0.1%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu3  :  0.5%us,  0.0%sy,  0.0%ni, 99.4%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu4  :  1.5%us,  0.1%sy,  0.0%ni, 98.4%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu5  :  0.8%us,  0.0%sy,  0.0%ni, 99.1%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu6  :  1.1%us,  0.2%sy,  0.0%ni, 98.6%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Cpu7  :  0.8%us,  0.0%sy,  0.0%ni, 99.2%id,  0.0%wa,  0.0%hi,  0.0%si,  0.0%st
Mem:  65342468k total, 11726988k used, 53615480k free,  416284k buffers
Swap: 67103496k total,  30464k used, 67073032k free, 10182848k cached
```

PID	USER	PR	NI	VIRT	RES	SHR	S	%CPU	%MEM	TIME+	COMMAND
1	root	20	0	10328	280	252	S	0	0.0	0:13.96	init
2	root	15	-5	0	0	0	S	0	0.0	0:00.03	kthreadd

Question 1: How to impose number of threads in code?

environment variable `OMP_NUM_THREADS`

```
[macrold@quartet2 hello_word1]$
[macrold@quartet2 hello_word1]$ set | grep OMP_NUM
OMP_NUM_THREADS=4
[macrold@quartet2 hello_word1]$
```

```
[macrold@octet1 hello_word1]$
[macrold@octet1 hello_word1]$ set | grep OMP_NUM
OMP_NUM_THREADS=8
[macrold@octet1 hello_word1]$
```

hello.c

```
1
2 #include <omp.h>
3 #include <stdio.h>
4
5 int main (int argc, char *argv[])
6 {
7     #pragma omp parallel
8     {
9         printf("Hello World\n");
10    }
11    return 0;
12 }
13
```

## Example 1: hello world [4]

Question 2: How can we run the same code in sequential mode?

### hello.c

```
1
2 #include <omp.h>
3 #include <stdio.h>
4
5 int main (int argc, char *argv[])
6 {
7     #pragma omp parallel
8     {
9         printf("Hello World\n");
10    }
11    return 0;
12 }
13
```

### Makefile

```
1
2 hello: hello.c
3     icpc -openmp -mp -c hello.c
4     icpc -openmp -o hello hello.o
5
6
7 hello_seq: hello.c
8     icpc -mp -c hello.c
9     icpc -o hello_seq hello.o
10
```

sequential version

### quartet2

```
[macrold@quartet2 hello_word1]$
[macrold@quartet2 hello_word1]$ make hello_seq
icpc -mp -c hello.c
hello.c(7): warning #161: unrecognized #pragma
        #pragma omp parallel
        ^
```

```
icpc -o hello_seq hello.o
[macrold@quartet2 hello_word1]$ ls
Makefile hello hello.c hello.o hello_seq
[macrold@quartet2 hello_word1]$ ./hello_seq
Hello World
[macrold@quartet2 hello_word1]$
```

### octet1

```
[macrold@octet1 hello_word1]$
[macrold@octet1 hello_word1]$ ./hello_seq
Hello World
[macrold@octet1 hello_word1]$
```

only one core executes



## Example 1: hello world [5]

Question 3: How can we issue number of threads explicitly in code?  
hello.c

```
1
2 #include <omp.h>
3 #include <stdio.h>
4
5 int main (int argc, char *argv[])
6 {
7     int th_id, nthreads;
8
9     #pragma omp parallel private(th_id) num_threads(5)
10 {
11     th_id = omp_get_thread_num();
12     printf("Hello World from thread %d\n", th_id);
13
14     #pragma omp barrier
15
16     if ( th_id == 0 ) {
17         nthreads = omp_get_num_threads();
18         printf("There are %d threads\n",nthreads);
19     }
20 }
21 return 0;
22 }
```

every thread has its own copy

use 5 threads (explicit) to execute concurrently

*synchronization*

wait until all 5 threads execute "printf" statement.

The **barrier** directive synchronizes all the threads in a team. When encountered, each thread in the team waits until all of the others have reached this point. The syntax of the **barrier** directive is as follows:

```
#pragma omp barrier new-line
```

After all threads in the team have encountered the barrier, each thread in the team begins executing the statements after the barrier directive in parallel.

## Example 1: hello world [6]

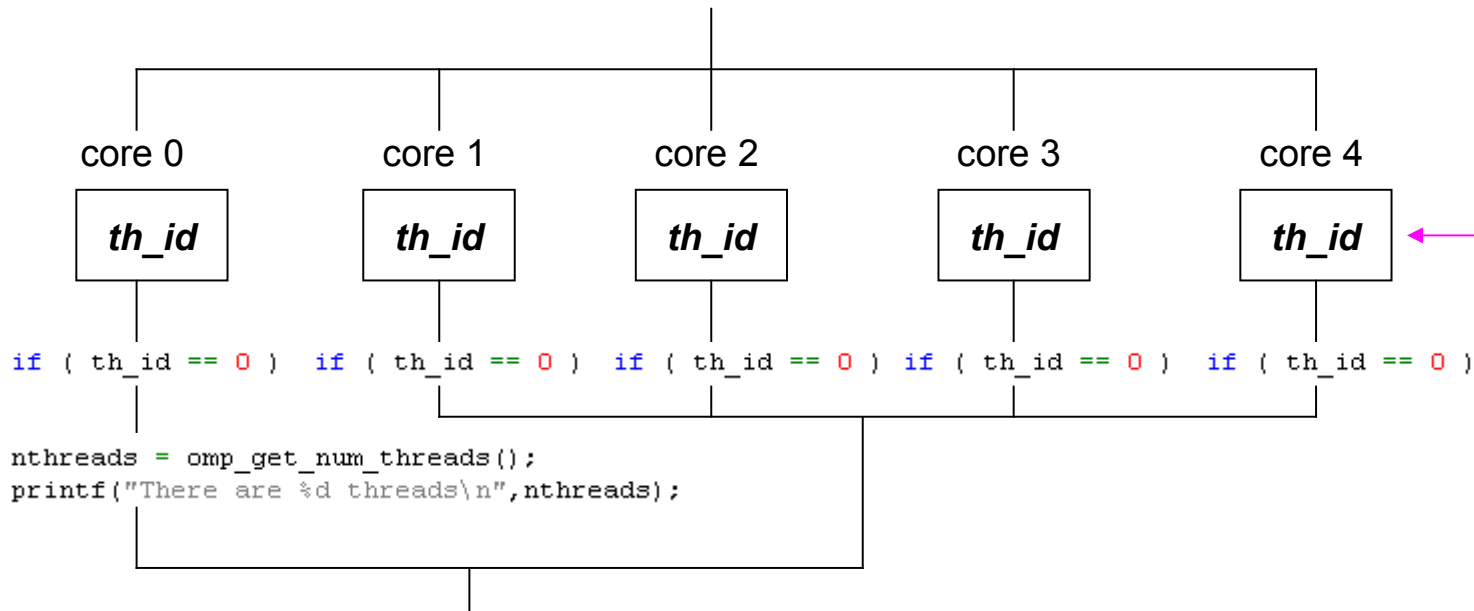
### quartet2

```
[macroid@quartet2 hello_word2]$  
[macroid@quartet2 hello_word2]$ make hello  
icpc -openmp -mp -c hello.c  
hello.c(8): (col. 3) remark: OpenMP DEFINED REGION WAS PARALLELIZED.  
icpc -openmp -o hello hello.o  
[macroid@quartet2 hello_word2]$ ls  
Makefile  hello  hello.c  hello.o  
[macroid@quartet2 hello_word2]$ ./hello  
Hello World from thread 1  
Hello World from thread 2  
Hello World from thread 3  
Hello World from thread 0  
Hello World from thread 4  
There are 5 threads  
[macroid@quartet2 hello_word2]$
```

### octet1

```
[macroid@octet1 hello_word2]$  
[macroid@octet1 hello_word2]$ ./hello  
Hello World from thread 3  
Hello World from thread 0  
Hello World from thread 4  
Hello World from thread 2  
Hello World from thread 1  
There are 5 threads  
[macroid@octet1 hello_word2]$
```

```
th_id = omp_get_thread_num();  
printf("Hello World from thread %d\n", th_id);
```

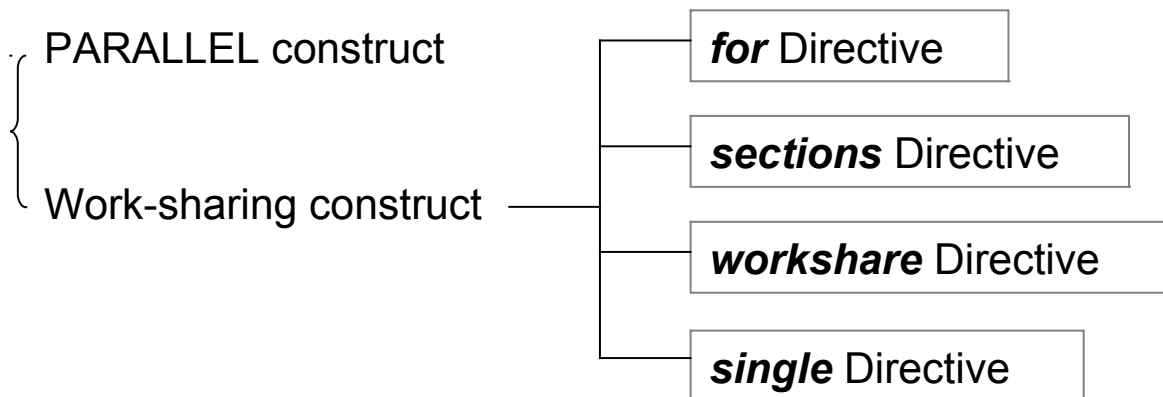


## Directive Format

The syntax of an OpenMP directive is formally specified by the grammar

```
#pragma omp directive-name [clause[ [,] clause]...] new-line
```

Each directive starts with **#pragma omp**, to reduce the potential for conflict with other (non-OpenMP or vendor extensions to OpenMP) *pragma* directives with the same names. White space can be used before and after the **#**, and sometimes white space must be used to separate the words in a directive. Preprocessing tokens following the **#pragma omp** are subject to macro replacement.



Conditional compilation

```
#ifdef _OPENMP  
iam = omp_get_thread_num() + index;  
#endif
```

# Parallel construct

```
#pragma omp parallel [clause ...] newline
    if (scalar_expression)
    private (list)
    shared (list)
    default (shared | none)
    firstprivate (list)
    reduction (operator: list)
    copyin (list)
    num_threads (integer-expression)
```

```
#pragma omp parallel private(th_id) num_threads(5)
{
    th_id = omp_get_thread_num();
    printf("Hello World from thread %d\n", th_id);

    #pragma omp barrier

    if ( th_id == 0 ) {
        nthreads = omp_get_num_threads();
        printf("There are %d threads\n", nthreads);
    }
}
```

*structured\_block*

- The number of physical processors hosting the threads is implementation-defined. Once created, the number of threads in the team remains constant for the duration of that parallel region.
- When a thread reaches a PARALLEL directive, it creates a team of threads and becomes the master of the team. The master is a member of that team and has thread number 0 within that team.
- Starting from the beginning of this parallel region, the code is duplicated and all threads will execute that code.
- There is an implied barrier at the end of a parallel region. Only the master thread of the team continues execution at the end of a parallel region.

## How many threads

- The number of threads in a parallel region is determined by the following factors, in order of precedence:
  - evaluation of the *IF* clause
  - setting of the *NUM\_THREADS* clause
  - use of the *omp\_set\_num\_threads()* library function
  - setting of the *OMP\_NUM\_THREADS* environment variable
  - implementation default - usually the number of CPUs on a node, though it could be dynamic.
- Threads are numbered from 0 (master thread) to N-1.
- Master thread is numbered as 0.

**Question 4:** How to write parallel code such that it is independent of number of cores of host machine?

**Question 5:** What happens if number of threads is larger than number of cores of host machine?

## Private clause

The PRIVATE clause declares variables in its list to be private to each thread.

“private variable” means each thread has its own copy and cannot interchange information.

```
#pragma omp parallel private(th_id) num_threads(5)
{
    th_id = omp_get_thread_num();
    printf("Hello World from thread %d\n", th_id);

    #pragma omp barrier

    if ( th_id == 0 ) {
        nthreads = omp_get_num_threads();
        printf("There are %d threads\n", nthreads);
    }
}
```

- PRIVATE variables behave as follows:
  - a new object of the same type is declared once for each thread in the team
  - all references to the original object are replaced with references to the new object
  - variables declared PRIVATE are **uninitialized** for each thread

**Exercise 1:** modify code of hello.c to show “every thread has its own private variable ***th\_id***”, that is, shows ***th\_id*** has 5 copies.

**Exercise 2:** modify code of hello.c, remove clause “private (th\_id)” in #pragma directive, what happens? Can you explain?

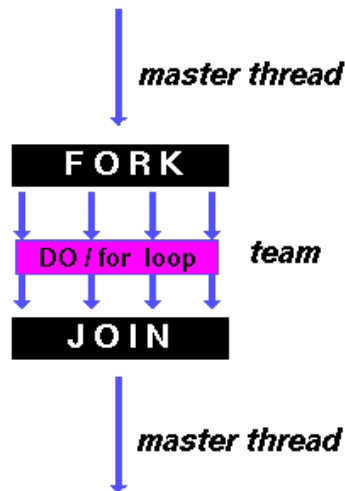
# OutLine

- OpenMP introduction
- Example 1: hello world
- Example 2: vector addition
  - work-sharing construct: for Directive
- enable openmp in vc2005
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

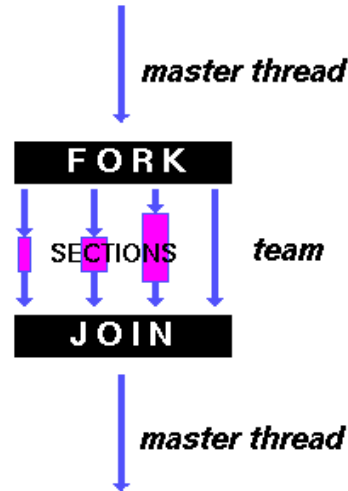
# Work-sharing construct

- A work-sharing construct divides the execution of the enclosed code region among the members of the team that encounter it
- A work-sharing construct must be enclosed dynamically within a parallel region in order for the directive to execute in parallel
- Work-sharing constructs do not launch new threads
- There is no implied barrier upon entry to a work-sharing construct, however there is an implied barrier at the end of a work sharing construct

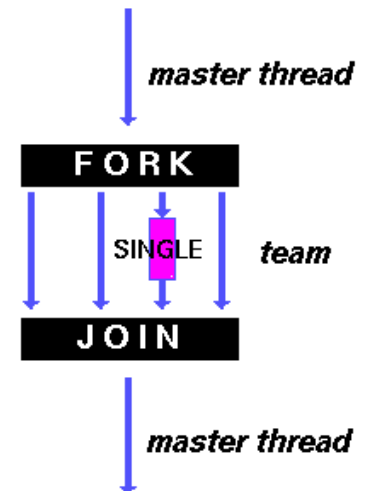
**for:** shares iterations of a loop across the team.  
A type of data parallelism



**sections:** breaks work into separate, discrete sections. Each section is executed by a thread.  
A type of functional parallelism



**single:** serializes a section of code.





## Example 2: vector addition [1]

### vecadd.c

```
1  #include <omp.h>
2  #include <stdio.h>
3  #include <stdlib.h>
4  #include <assert.h>
5
6  double walltime( double *t0 ) ;
7  void randomInit( float* data, int size) ;
8
9  int main(int argc, char *argv[] )
10 {
11     long int N = 200000000 ;
12     int thread_num = 4 ;
13     long int i;
14     float *a, *b, *c;
15     double startTime, elapsedTime; /* for timing */
16     double clockZero = 0.0;
17
18     a = (float*) malloc( sizeof(float) *N ) ; assert(a) ;
19     b = (float*) malloc( sizeof(float) *N ) ; assert(b) ;
20     c = (float*) malloc( sizeof(float) *N ) ; assert(c) ;
21
22     startTime = walltime( &clockZero );
23     randomInit(a, N);
24     randomInit(b, N);
25     elapsedTime = walltime( &startTime );
26     printf("Time to randomize a, b = %6.4f (s)\n", elapsedTime);
```

parameter

### walltime.c

```
4  #include <sys/time.h>
5  // return current_time - t0 in seconds
6  double walltime( double *t0 )
7  {
8      double mic, time;
9      double mega = 0.000001;
10     struct timeval tp;
11     struct timezone tzp;
12     static long base_sec = 0;
13     static long base_usec = 0;
14
15     (void) gettimeofday(&tp, &tzp);
16     if (base_sec == 0)
17     {
18         base_sec = tp.tv_sec;
19         base_usec = tp.tv_usec;
20     }
21
22     time = (double) (tp.tv_sec - base_sec);
23     mic = (double) (tp.tv_usec - base_usec);
24     time = (time + mic * mega) - *t0;
25     return(time);
26 }
```

Tool for measuring time

only valid in Linux system

### vecadd.c

```
49 // Allocates a matrix with random float entries.
50 void randomInit(float* data, int size)
51 {
52     for (int i = 0; i < size; ++i){
53         data[i] = rand() / (float)RAND_MAX;
54     }
55 }
```

## Example 2: vector addition [2]

### vecadd.c

```
27
28     startTime = walltime( &clockZero );
29
30 → #pragma omp parallel default(none) num_threads(thread_num) \
31     shared(a,b,c,N) private(i)
32 {
33 → #pragma omp for schedule( static ) nowait
34     for (i=0; i < N; i++){
35         c[i] = a[i] + b[i];
36     }
37
38 } /* end of parallel section */
39 elapsedTime = walltime( &startTime );
40 double size = ((double)N)*sizeof(float)/1.E6 ;
41 printf("size = %6.2f (MB) \n", size );
42 printf("thread_num = %d, time for vecadd = %6.4f (s)\n",
43     thread_num, elapsedTime) ;
44
45 free(a) ; free(b) ; free(c) ;
46 return 0 ;
47 }
```

### Makefile

```
1
2 vecadd: vecadd.c walltime.c
3     icpc -openmp -mp -O0 -c vecadd.c
4     icpc -c walltime.c
5     icpc -openmp -o vecadd walltime.o vecadd.o
6
7 clean:
8     rm -f *.o
```

“O0” means no optimization

```
[macrold@quartet2 vecadd]$ ls
Makefile vecadd.c walltime.c
[macrold@quartet2 vecadd]$ make vecadd
icpc -openmp -mp -O0 -c vecadd.c
vecadd.c(33): (col. 5) remark: OpenMP DEFINED LOOP WAS PARALLELIZED.
vecadd.c(30): (col. 1) remark: OpenMP DEFINED REGION WAS PARALLELIZED.
icpc -c walltime.c
icpc -openmp -o vecadd walltime.o vecadd.o
[macrold@quartet2 vecadd]$ ./vecadd
Time to randomize a, b = 6.4568 (s)
size = 800.00 (MB)
thread_num = 4, time for vecadd = 0.6257 (s)
[macrold@quartet2 vecadd]$
```

## shared clause and default clause

The SHARED clause declares variables in its list to be shared among all threads in the team

```
#pragma omp parallel default(none) num_threads(thread_num) \
    shared(a,b,c,N) private(i)
{
    #pragma omp for schedule( static ) nowait
    for (i=0; i < N; i++){
        c[i] = a[i] + b[i];
    }

} /* end of parallel section */
```

- A shared variable exists in only one memory location and all threads can read or write to that address (every thread can “**see**” the shared variable)
- It is the programmer's responsibility to ensure that multiple threads properly access SHARED variables (such as via CRITICAL sections)

**Question 6:** Why index *i* must be private variable and *a,b,c,N* can be shared variable? What happens if we change *i* to shared variable? What happens if we change *a,b,c,N* to private variable?

The DEFAULT clause allows the user to specify a default PRIVATE, SHARED, or NONE scope for all variables in the lexical extent of any parallel region.

```
default (shared | none)
```

## Work-Sharing construct: for Directive

```
#pragma omp for [clause ...] newline
    schedule (type [,chunk])
    ordered
    private (list)
    firstprivate (list)
    lastprivate (list)
    shared (list)
    reduction (operator: list)
    nowait
```

*for\_loop*

```
#pragma omp parallel default(none) num_threads(thread_num) \
    shared(a,b,c,N) private(i)
{
    #pragma omp for schedule( static ) nowait
    for (i=0; i < N; i++){
        c[i] = a[i] + b[i];
    }
} /* end of parallel section */
```

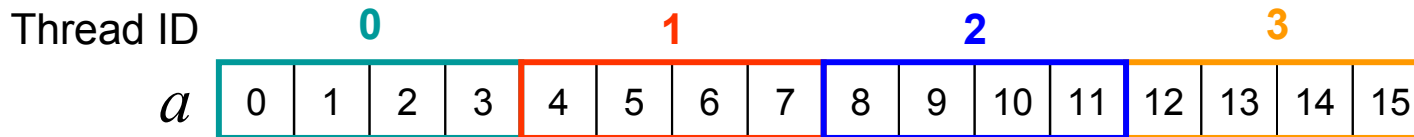
- **SCHEDULE:** Describes how iterations of the loop are divided among the threads in the team
  - **static:** loop iterations are divided into pieces of size *chunk* and then statically assigned to threads. If chunk is not specified, the iterations are evenly (if possible) divided contiguously among the threads
  - **dynamic:** loop iterations are divided into pieces of size *chunk*, and dynamically scheduled among the threads; when a thread finishes one chunk, it is dynamically assigned another.  
The default chunk size is 1.
- **nowait:** If specified, then threads do not synchronize at the end of the parallel loop.

## Example of static schedule

Assume we have 16 array elements, say  $a[16]$ ,  $b[16]$  and  $c[16]$  and use 4 threads

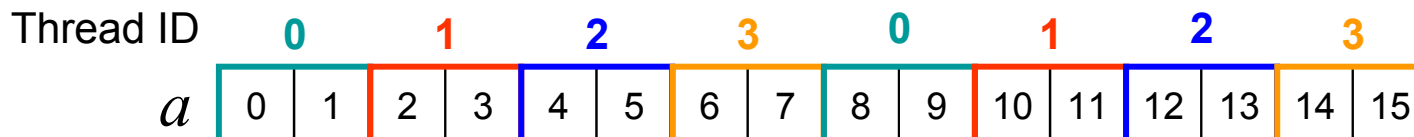
- 1 no chunk is specified, compiler would divide 16 elements into 4 threads

```
#pragma omp parallel default(none) num_threads(thread_num) \
    shared(a,b,c,N) private(i)
{
    #pragma omp for schedule( static ) nowait
    for (i=0; i < N; i++){
        c[i] = a[i] + b[i];
    }
} /* end of parallel section */
```



- 2 chunk = 2

```
#pragma omp parallel default(none) num_threads(thread_num) \
    shared(a,b,c,N) private(i)
{
    #pragma omp for schedule( static, 2 ) nowait
    for (i=0; i < N; i++){
        c[i] = a[i] + b[i];
    }
} /* end of parallel section */
```



## Results of example 2

$$N = 2 \times 10^8$$

compiler: Intel C compiler icpc 10.0

Compiler option: -O0

```
#pragma omp parallel default(none) num_threads(thread_num) \
    shared(a,b,c,N) private(i)
{
    #pragma omp for schedule( static ) nowait
    for (i=0; i < N; i++){
        c[i] = a[i] + b[i];
    }
}

/* end of parallel section */
```

$$\text{Octet1: } \frac{T(\text{single})}{T(8\text{-core})} = \frac{1.5451}{0.483} = 3.199$$

$$\text{quartet2: } \frac{T(\text{single})}{T(4\text{-core})} = \frac{1.6571}{0.5433} = 3.05$$

**Question 7:** the limitation of performance improvement is 3, why? Can you use different configuration of schedule clause to improve this number?

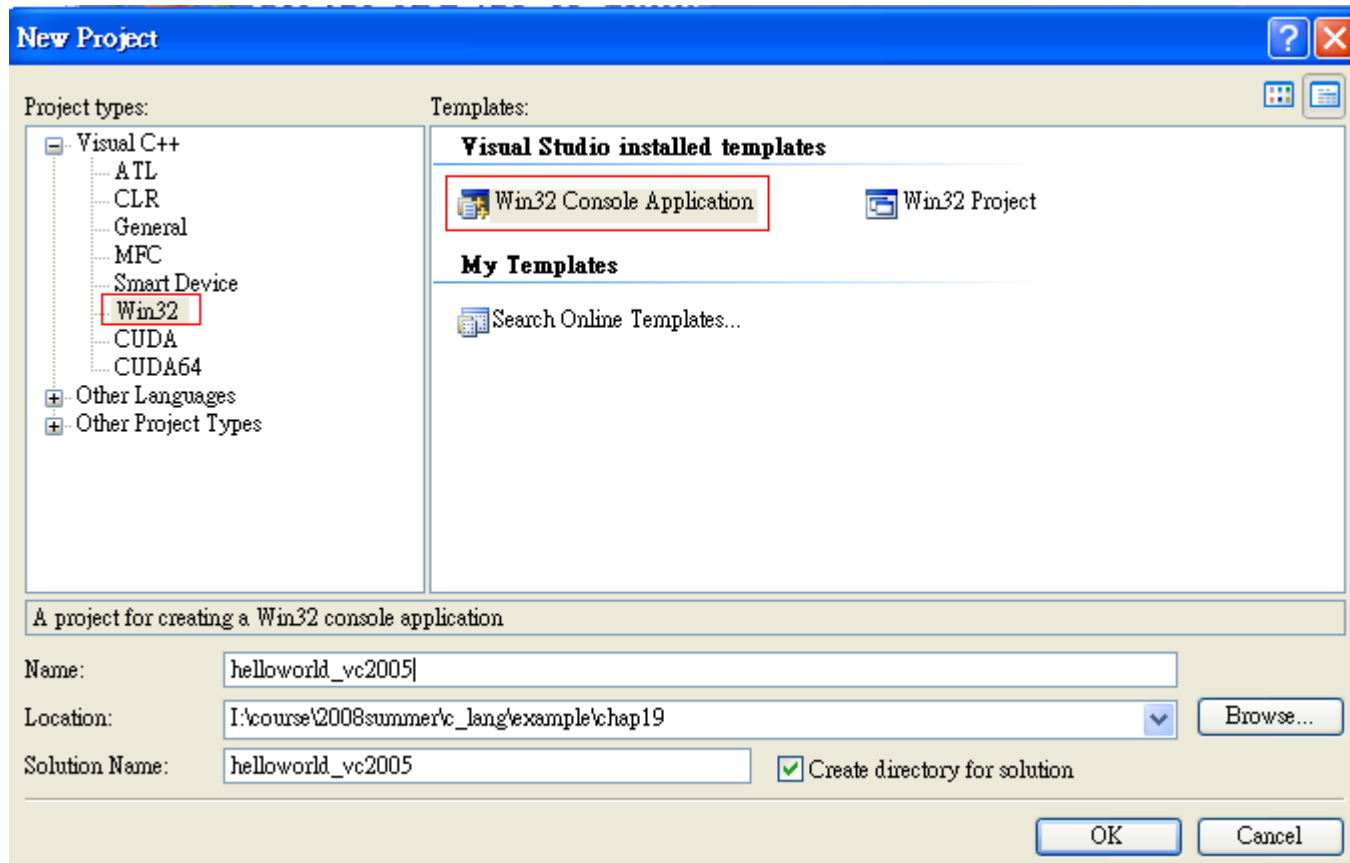
Number of thread	quartet2	Octet1
1	1.6571 (s)	1.5451 (s)
2	0.9064 (s)	0.9007 (s)
4	0.5433 (s)	0.5165 (s)
8	0.6908 (s)	0.4830 (s)
16	0.7694 (s)	0.5957 (s)
32	0.9263 (s)	0.7098 (s)
64	0.9625 (s)	0.7836 (s)

# OutLine

- OpenMP introduction
- Example 1: hello world
- Example 2: vector addition
- enable openmp in vc2005
  - vc2005 supports OpenMP 2.0
  - vc 6.0 does not support OpenMP
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

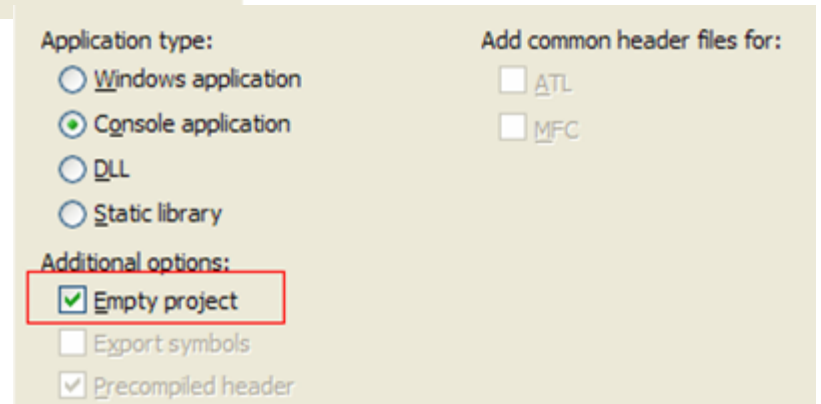
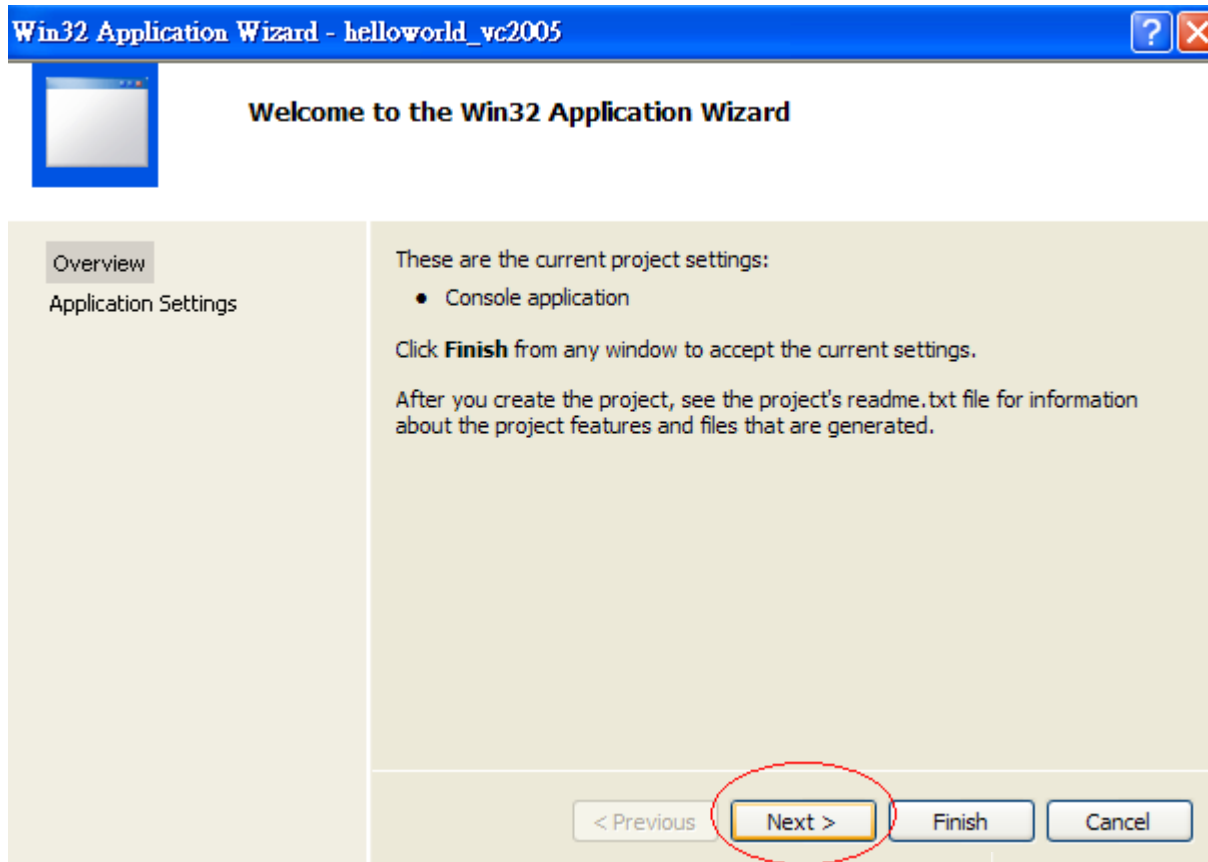
## Example 1 (hello world) in vc2005 [1]

Step 1: create a empty consol application

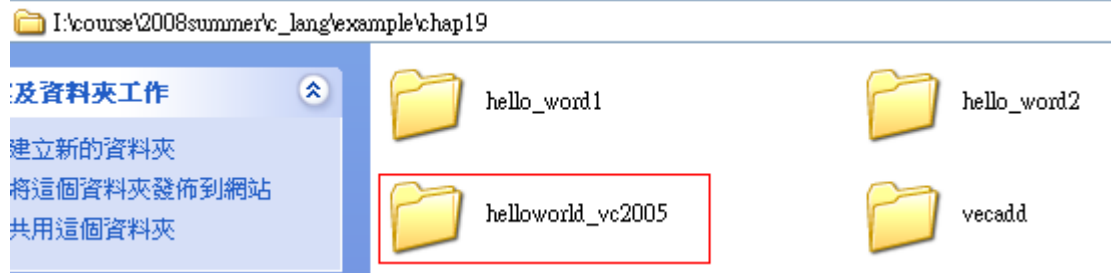




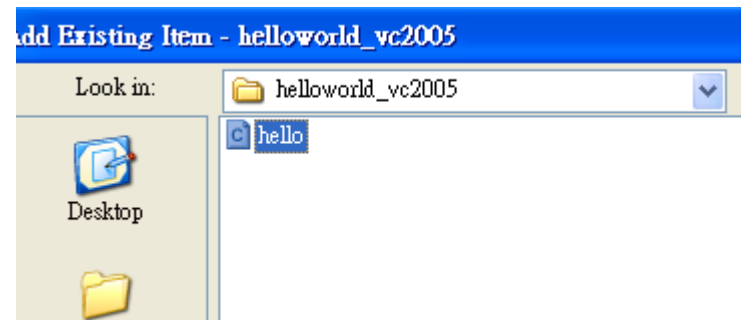
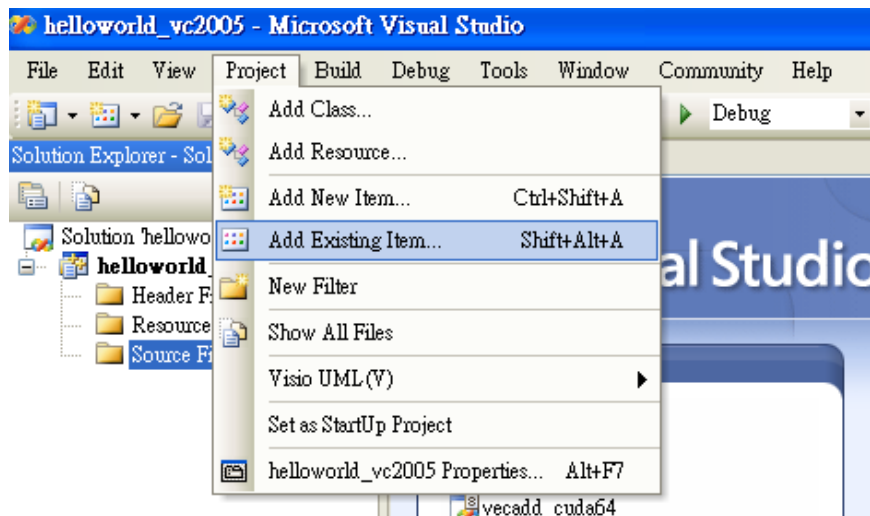
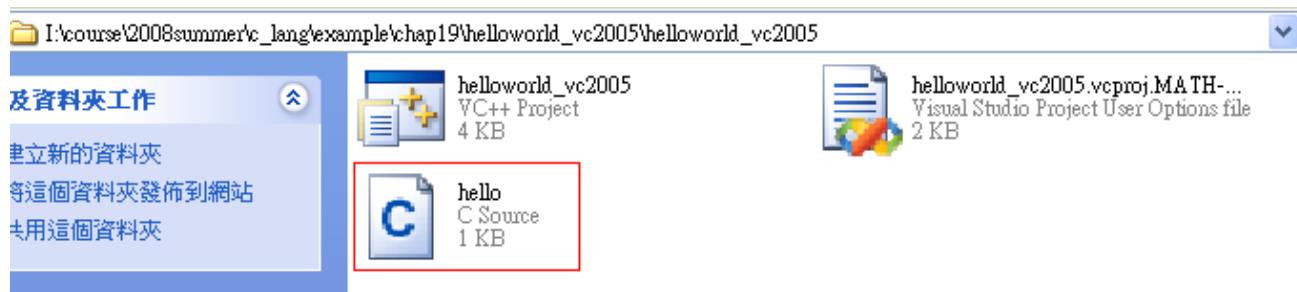
## Example 1 (hello world) in vc2005 [2]



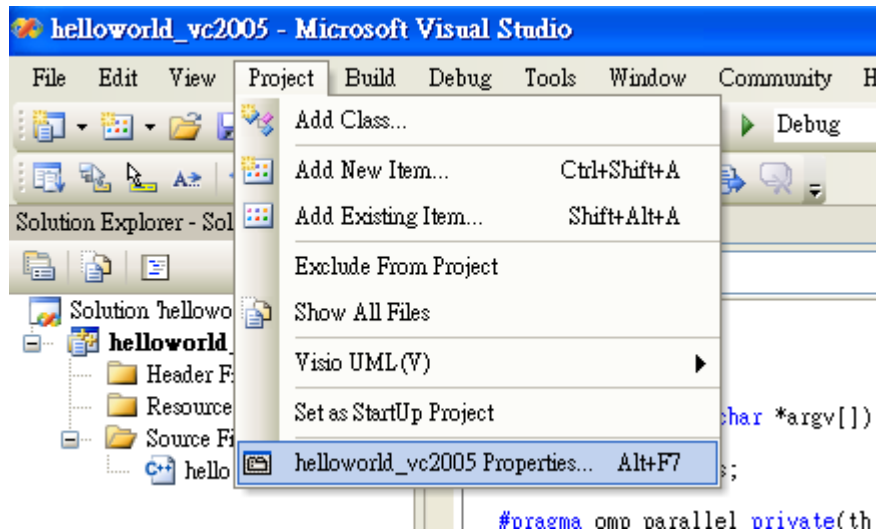
## Example 1 (hello world) in vc2005 [3]



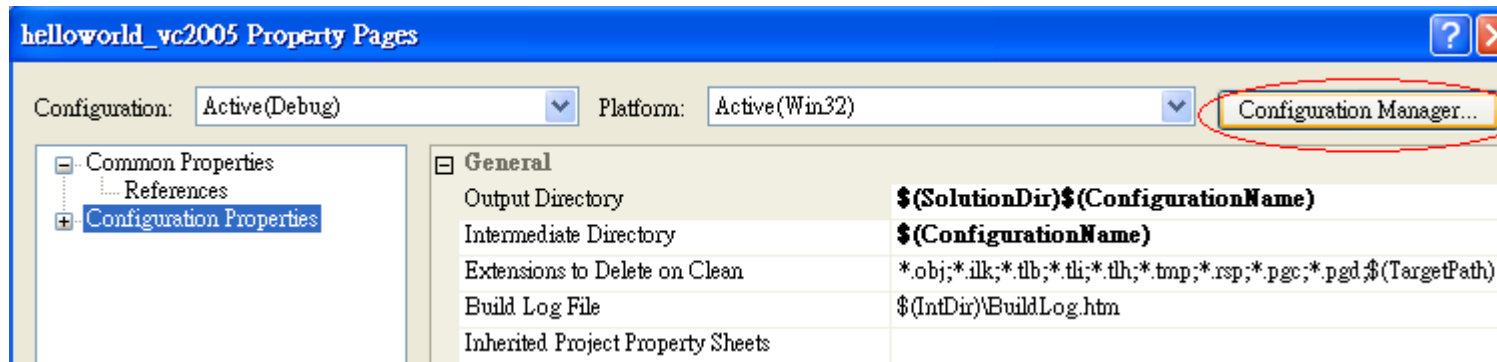
Step 2: copy `hello.c` to this project and add `hello.c` to project manager



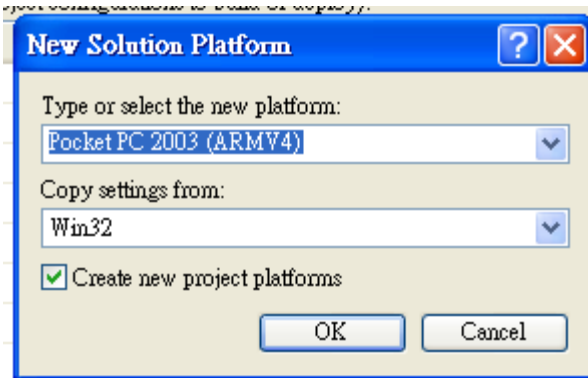
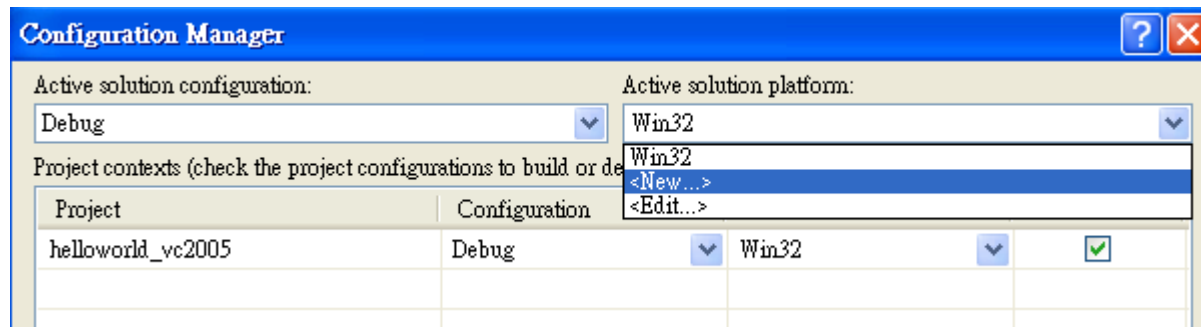
## Example 1 (hello world) in vc2005 [4]



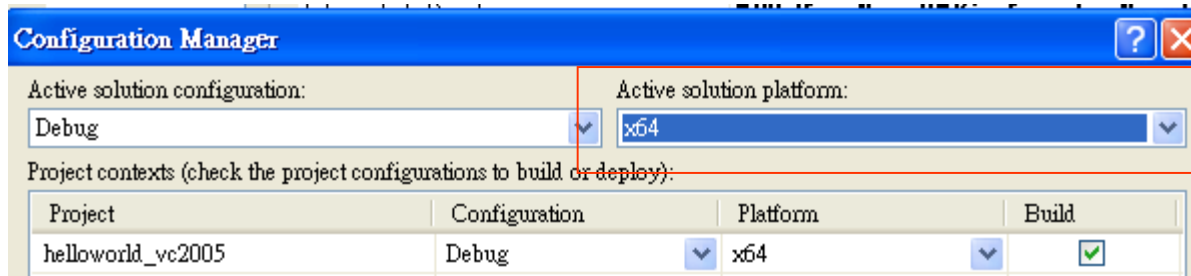
Step 3: change platform to x64



## Example 1 (hello world) in vc2005 [5]



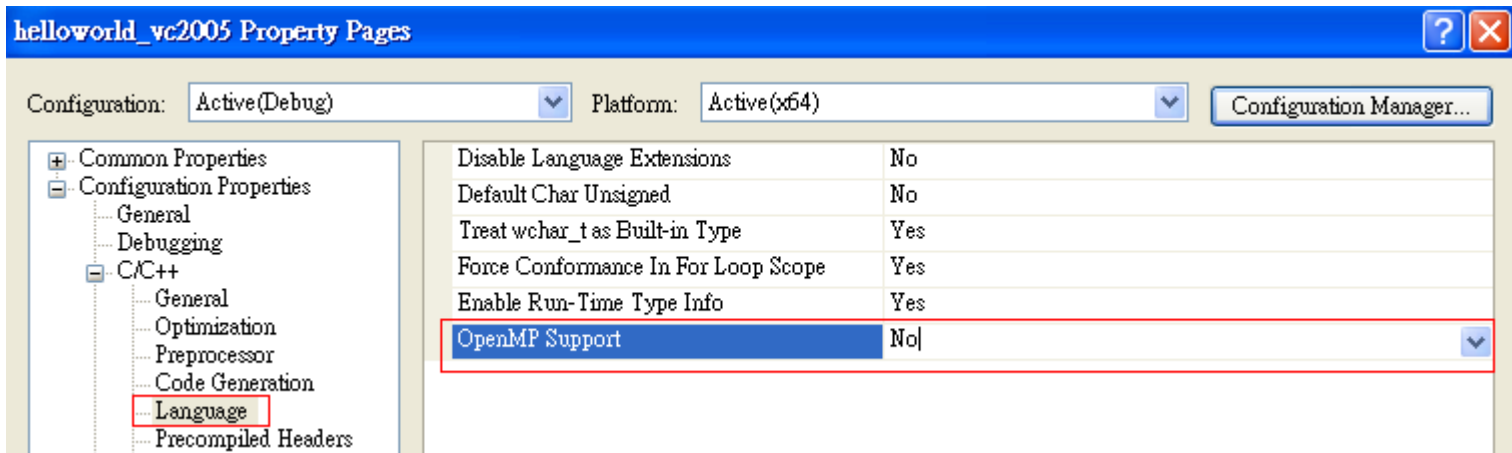
choose option "x64"



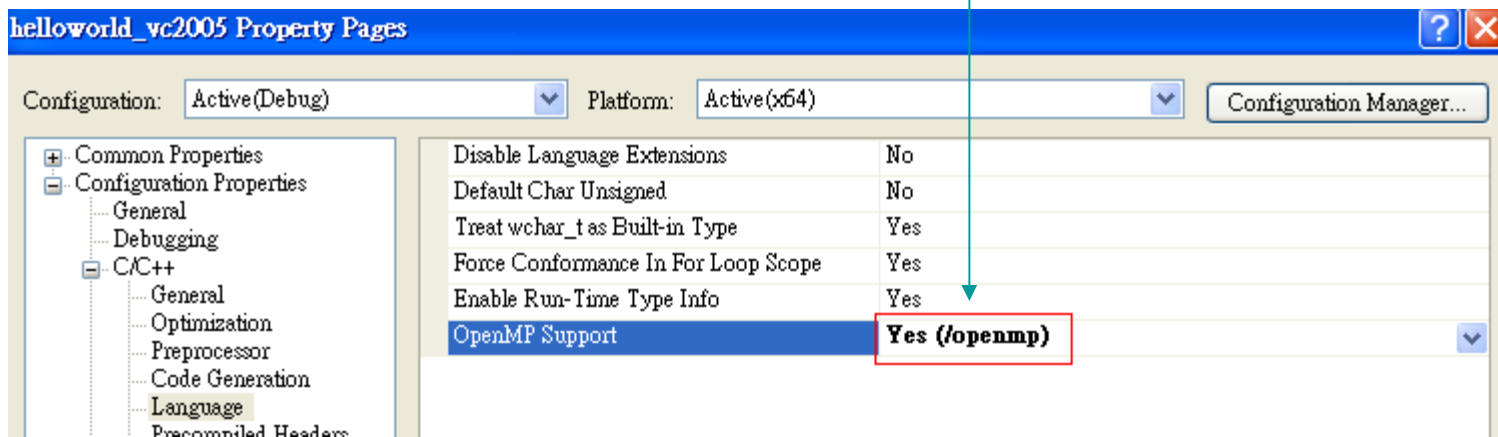
update platform as "x64"

## Example 1 (hello world) in vc2005 [6]

Step 4: enable “openmp” support

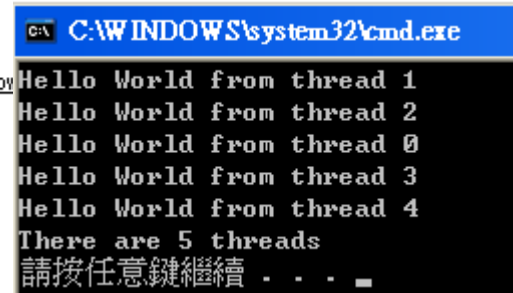
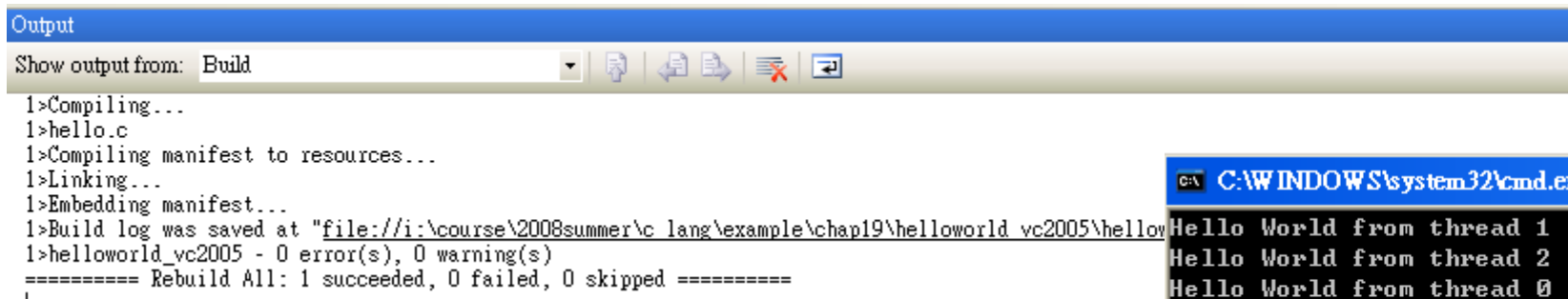
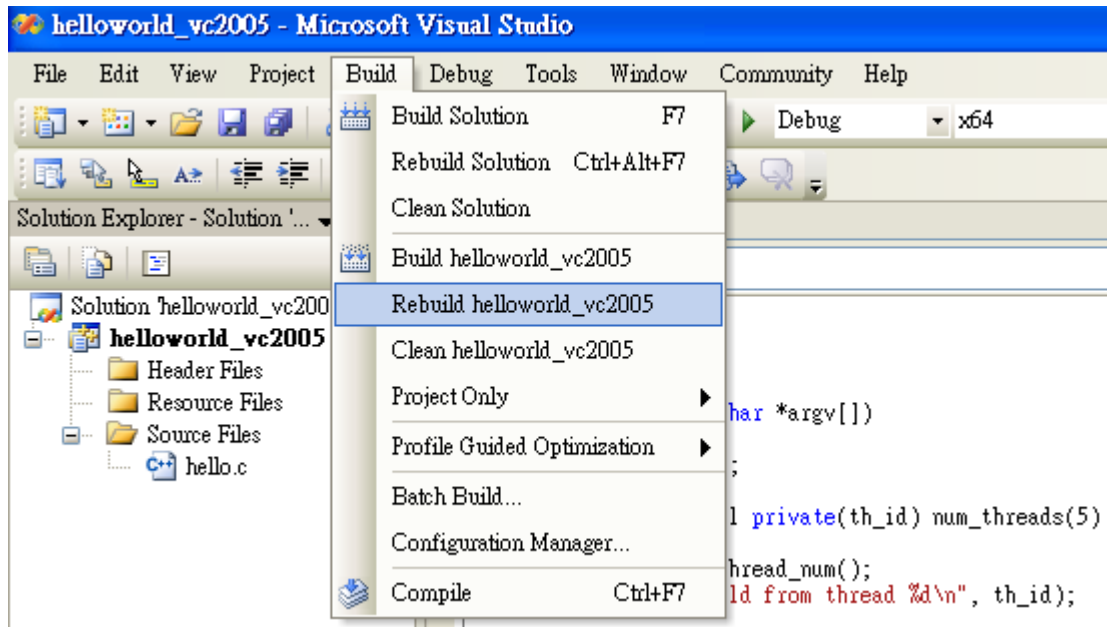


vc 2005 support OpenMP 2.0



## Example 1 (hello world) in vc2005 [7]

Step 5: compile and execute



## Example 2 (vector addition) in vc2005 [1]

walltime.c only works in Linux machine since no "sys/time.h" in windows

In time.h of ANCI C, no function "gettimeofday", hence we give up walltime.c

```
(Global Scope) walltime(double *t0)

/* wall clock */
/* see "Software Optimization for High Performance Computing" p. 135 */

#include <sys/time.h>
// return current time - t0 in seconds
double walltime( double *t0 )
{
    double mic, time;
    double mega = 0.000001;
    struct timeval tp;
    struct timezone tzp;
    static long base_sec = 0;
    static long base_usec = 0;

    (void) gettimeofday(&tp,&tzp);
    if (base_sec == 0)
    {
        base_sec = tp.tv_sec;
        base_usec = tp.tv_usec;
    }

    time = (double) (tp.tv_sec - base_sec);
    mic = (double) (tp.tv_usec - base_usec);
    time = (time + mic * mega) - *t0;
    return(time);
}

Output
Show output from: Build
1>----- Build started: Project: vecadd_vc2005, Configuration: Debug x64 -----
1>Compiling...
1>walltime.c
1>.\walltime.c(4) : fatal error C1083: Cannot open include file: 'sys/time.h': No such file or directory
1>Build log was saved at "file:///i:/course/2008summer/c_lang/example/chap19/vecadd_vc2005/vecadd_vc2005/x64/Debug/BuildLog.htm"
1>vecadd_vc2005 - 1 error(s), 0 warning(s)
===== Build: 0 succeeded, 1 failed, 0 up-to-date, 0 skipped =====
```

## Example 2 (vector addition) in vc2005 [2]

`time_t time( time_t *tp)`

returns the current calendar time or -1 if the time is not available. If tp is not NULL, the return value is also assigned to \*tp.

`double difftime( time_t time_2, time_t time_1)`

returns time\_2 – time\_1 expressed in seconds

### vecadd.cpp

```
16 // double startTime, elapsedTime; /* for timing */
17 time_t startTime, endTime ;
18 double elapsedTime;

24 // startTime = walltime( &clockZero );
25 time( &startTime ) ;
26 randomInit(a, N);
27 randomInit(b, N);
28 // elapsedTime = walltime( &startTime );
29 time( &endTime ) ;
30 elapsedTime = difftime(endTime, startTime) ;
31 printf("Time to randomize a, b = %6.4f (s)\n", elapsedTime);

35 // startTime = walltime( &clockZero );
36 time( &startTime ) ;
37 #pragma omp parallel default(none) num_threads(thread_num) \
38   shared(a,b,c,N) private(i)
39 {
40     #pragma omp for schedule( static ) nowait
41     for (i=0; i < N; i++){
42         c[i] = a[i] + b[i];
43     }
44 }
45 /* end of parallel section */
46 // elapsedTime = walltime( &startTime );
47 time( &endTime ) ;
48 elapsedTime = difftime(endTime, startTime) ;
```



# OutLine

- OpenMP introduction
- Example 1: hello world
- Example 2: vector addition
- enable openmp in vc2005
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

## Example 3: vector addition (Qtime) [1]

### vecadd.cpp

```
1  #include <omp.h>
2  #include <stdio.h>
3  #include <stdlib.h>
4  #include <assert.h>
5
6  #include <qdatetime.h>
7
8  void randomInit( float* data, int size) ;
9
10 int main(int argc, char *argv[] )
11 {
12     long int N = 200*1024*1024 ;
13     int thread_num = 4 ;
14     long int i;
15     float *a, *b, *c ;
16
17     QTime t; // QT timer
18
19     #ifdef _OPENMP
20         printf("OpenMP-compliant implementation\n");
21     #endif
22
23     a = (float*) malloc( sizeof(float) *N ) ; assert(a) ;
24     b = (float*) malloc( sizeof(float) *N ) ; assert(b) ;
25     c = (float*) malloc( sizeof(float) *N ) ; assert(c) ;
```

constructs the time 0 hours, minutes, seconds and milliseconds, i.e. 00:00:00.000 (midnight).

This is a valid time.

- A QTime object contains a clock time, i.e. the number of hours, minutes, seconds, and milliseconds since midnight
- QTime uses the 24-hour clock format; it has no concept of AM/PM. It operates in local time; it knows nothing about time zones or daylight savings time.
- QTime can be used to measure a span of elapsed time using the start(), restart(), and elapsed() functions

## Example 3: vector addition (Qtime) [2]

### vecadd.cpp

```
26     t.start() ;
27
28
29     randomInit(a, N);
30     randomInit(b, N);
31     printf("Time to randomize a, b = %d (ms)\n", t.elapsed());
32
33     t.start() ;
34
35     #pragma omp parallel default(none) num_threads(thread_num) \
36         shared(a,b,c,N) private(i)
37 {
38     #pragma omp for schedule(static) nowait
39     for (i=0; i < N; i++){
40         c[i] = a[i] + b[i];
41     }
42 } /* end of parallel section */
43
44 printf("thread_num = %d, time for vecadd = %d (ms)\n",
45        thread_num, t.elapsed()); ;
46
```

### void QTime::start ()

Sets this time to the current time. This is practical for timing.

### int QTime::elapsed () const

Returns the number of milliseconds that have elapsed since the last time `start()` or `restart()` was called.

Note that the counter wraps to zero 24 hours after the last call to `start()` or `restart()`.

Note that the accuracy depends on the accuracy of the underlying operating system; not all systems provide 1-millisecond accuracy.

### Example 3: vector addition (Qtime) [3]

```
[macrold@octet1 vecadd_gt]$  
[macrold@octet1 vecadd_gt]$ ls  
Makefile vecadd.cpp  
[macrold@octet1 vecadd_gt]$ qmake -project  
[macrold@octet1 vecadd_gt]$ ls  
Makefile vecadd.cpp vecadd_gt.pro  
[macrold@octet1 vecadd_gt]$ qmake -spec linux-icc-openmp vecadd_gt.pro  
[macrold@octet1 vecadd_gt]$ ls  
Makefile vecadd.cpp vecadd_gt.pro  
[macrold@octet1 vecadd_gt]$ make  
icpc -c -w -O2 -openmp -mp -DQT_NO_DEBUG -DQT_SHARED -DQT_THREAD_SUPPORT -I/opt/qt/mkspecs/linux-icc-openmp -I. -I. -I/opt/qt/include -o vecadd.o vecadd.cpp  
vecadd.cpp(36): (col. 5) remark: OpenMP DEFINED LOOP WAS PARALLELIZED.  
vecadd.cpp(33): (col. 1) remark: OpenMP DEFINED REGION WAS PARALLELIZED.  
vecadd.cpp(37): (col. 5) remark: LOOP WAS VECTORIZED.  
icpc -openmp -Qoption,ld,-rpath,/opt/qt/lib -o vecadd_gt vecadd.o -L/opt/qt/lib -L/usr/X11R6/lib -lqt-mt -lXext -lX11 -lm  
[macrold@octet1 vecadd_gt]$ ls  
Makefile vecadd.cpp vecadd.o vecadd_gt vecadd_gt.pro  
[macrold@octet1 vecadd_gt]$ ./vecadd_gt  
OpenMP-compliant implementation  
Time to randomize a, b = 5515 (ms)  
thread_num = 4, time for vecadd = 522 (ms)  
size = 800.00 (MB)  
[macrold@octet1 vecadd_gt]$
```

generate project file **vecadd\_gt.pro**

generate Makefile

#### Makefile

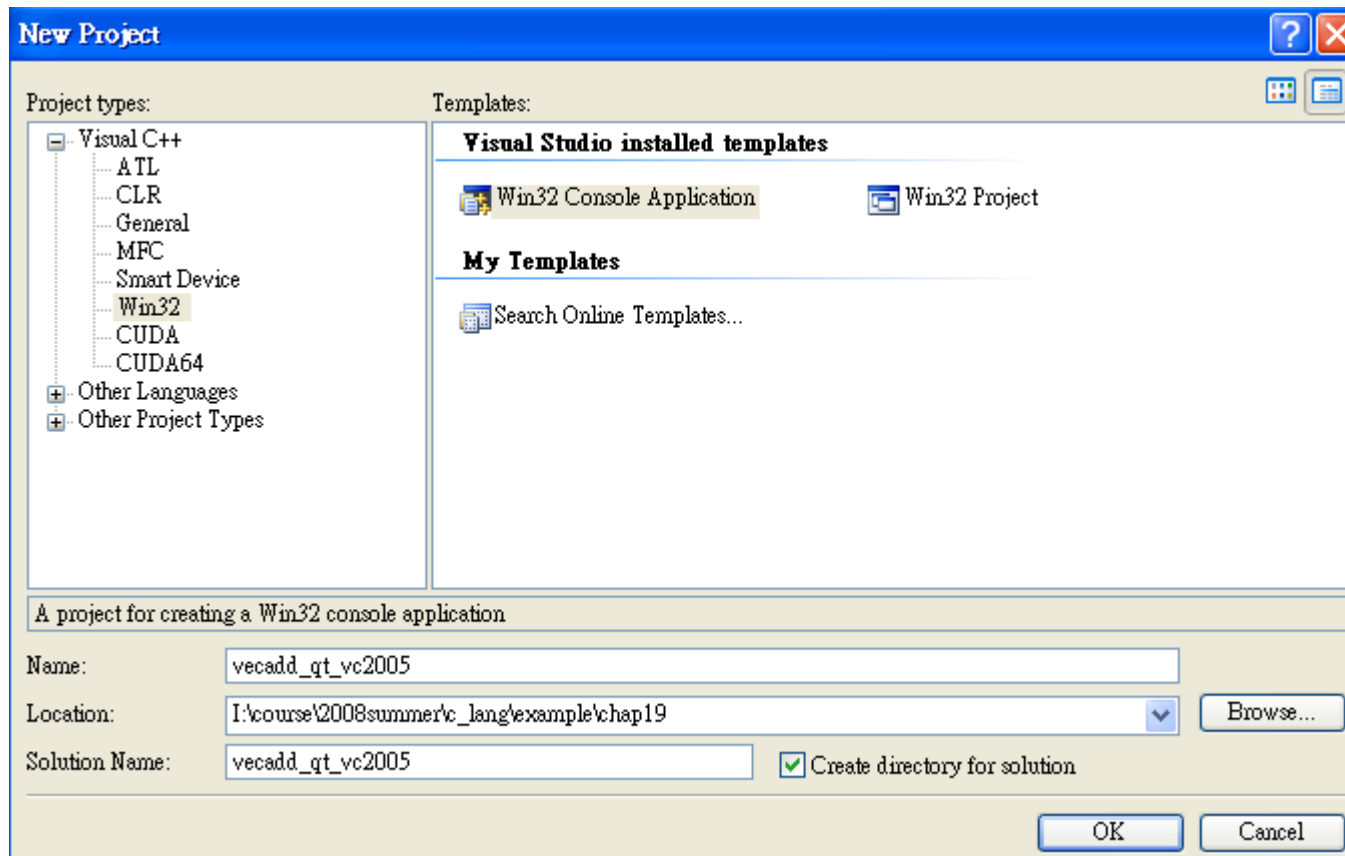
##### Compiler, tools and options

```
CC      = icc  
CXX     = icpc  
LEX     = flex  
YACC    = yacc  
CFLAGS  = -w -O2 -openmp -mp -DQT_NO_DEBUG -DQT_SHARED -DQT_THREAD_SUPPORT  
CXXFLAGS = -w -O2 -openmp -mp -DQT_NO_DEBUG -DQT_SHARED -DQT_THREAD_SUPPORT  
LEXFLAGS =  
YACCFLAGS = -d  
INCPATH = -I/opt/qt/mkspecs/linux-icc-openmp -I. -I. -I$(QTDIR)/include  
LINK    = icpc  
LFLAGS  = -openmp -Qoption,ld,-rpath,$(QTDIR)/lib  
LIBS    = $(SUBLIBS) -L$(QTDIR)/lib -L/usr/X11R6/lib -lqt-mt -lXext -lX11 -lm
```

### Example 3: vector addition (Qtime) [4]

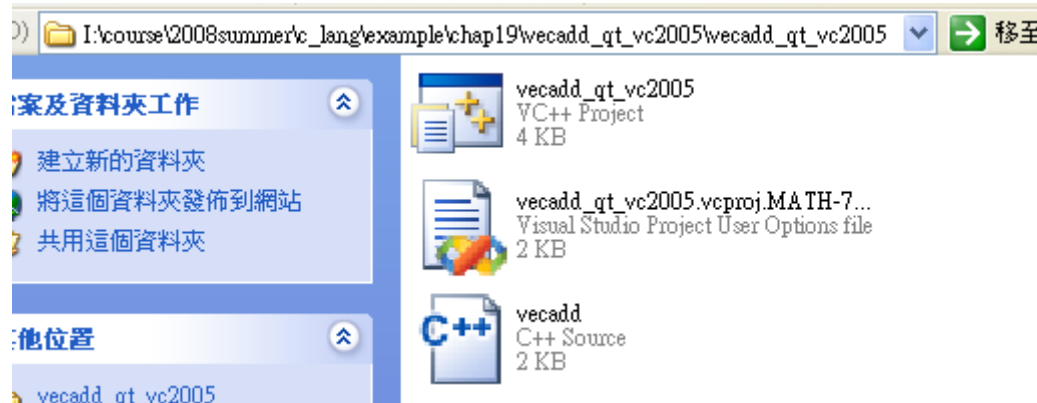
*Embed Qt 3.2.1 non-commercial version into vc 2005*

Step 1: setup an empty project

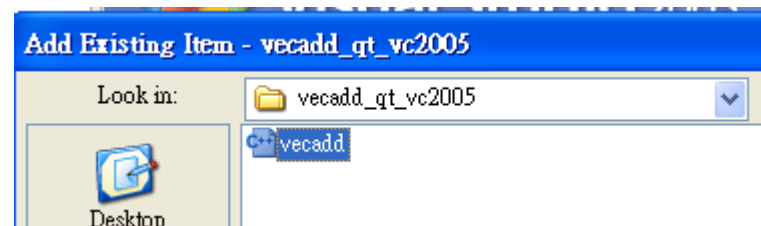
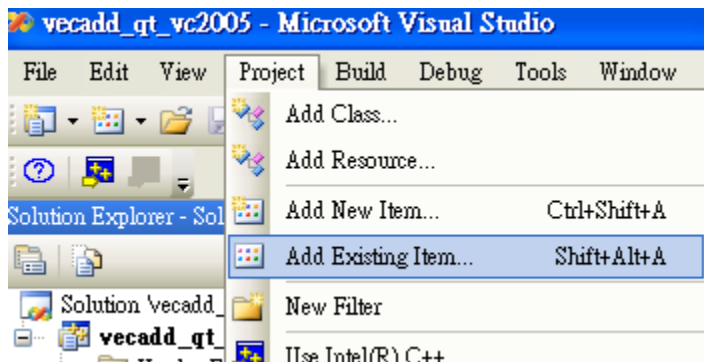


### Example 3: vector addition (Qtime) [5]

Step 2: copy vecadd.cpp into this project



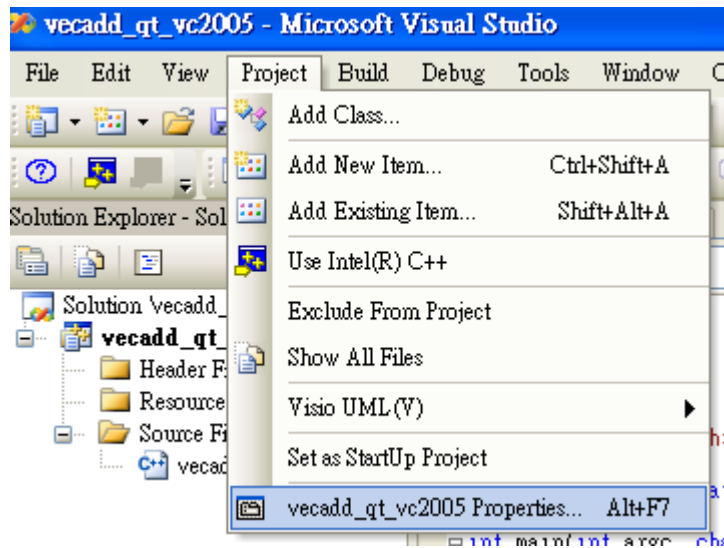
Step 3: add item “vecadd.cpp” in project manager



### Example 3: vector addition (Qtime) [6]

Step 4: project → properties → C/C++ → General → Additional include Directories

.;\$(QTDIR)\include;C:\Qt\3.2.1NonCommercial\mkspecs\win32-msvc

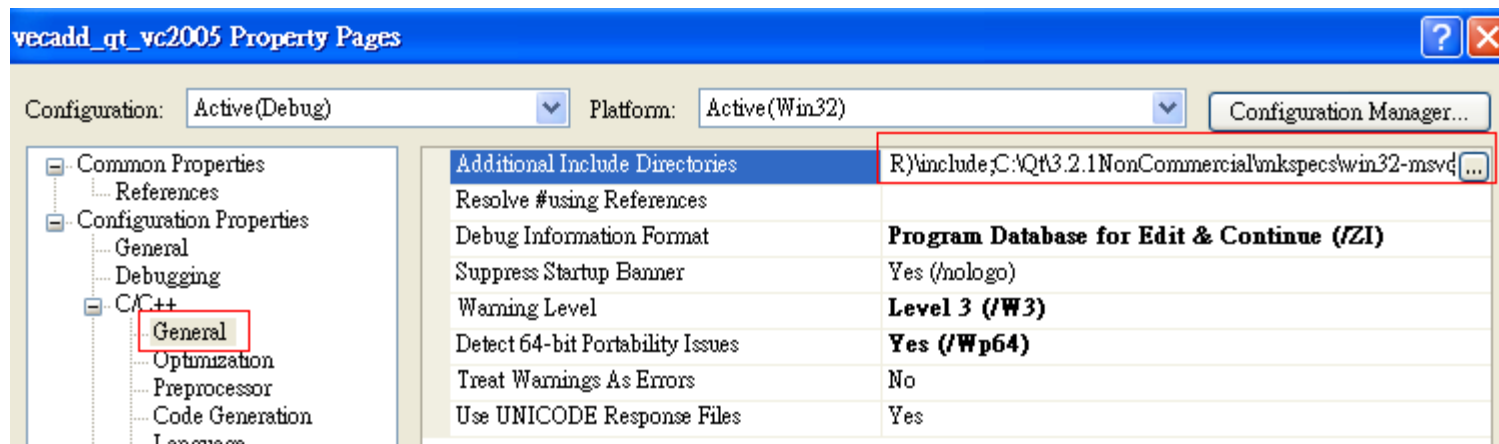


```
C:\Documents and Settings\LungShengChien>set  
ALLUSERSPROFILE=C:\Documents and Settings\All Users  
APPDATA=C:\Documents and Settings\LungShengChien\Application Data  
CC=c1  
CLIENTNAME=Console
```

∩

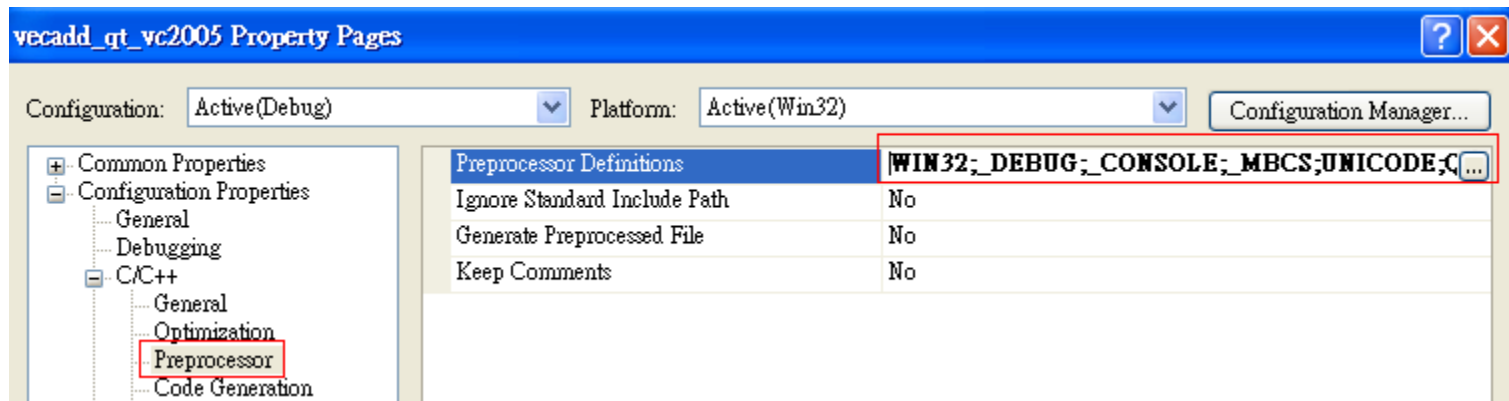
```
QMAKESPEC=win32-msvc  
QTDIR=C:\Qt\3.2.1NonCommercial
```

∩

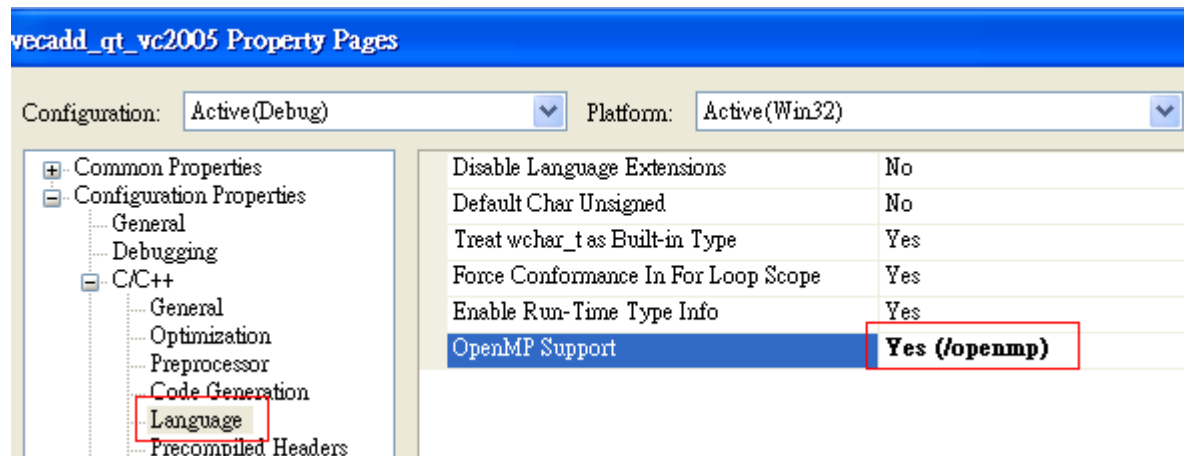


### Example 3: vector addition (Qtime) [7]

Step 5: project → properties → C/C++ → Preprocessor → Preprocessor Definitions  
WIN32;\_DEBUG;\_CONSOLE;\_MBCS;UNICODE;QT\_DLL;QT\_THREAD\_SUPPORT



Step 6: project → properties → C/C++ → Language → OpenMP Support

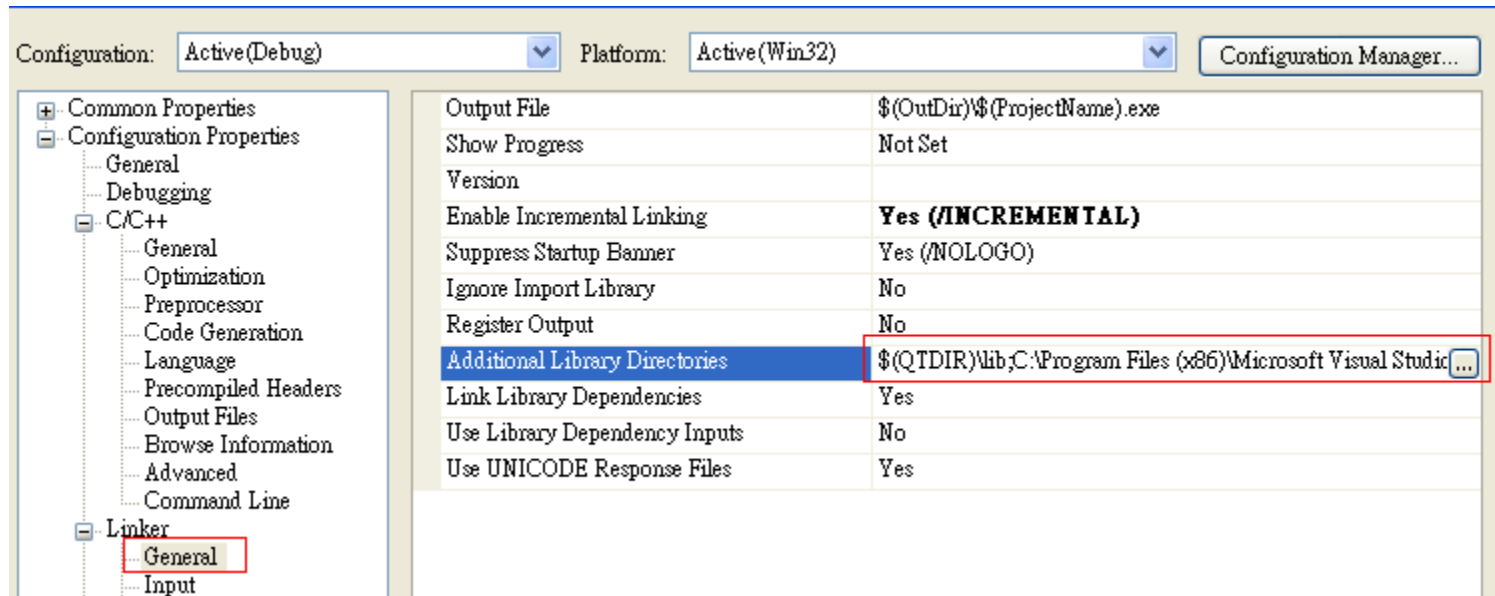




### Example 3: vector addition (Qtime) [8]

Step 7: project → properties → Linker → General → Additional Library Directories

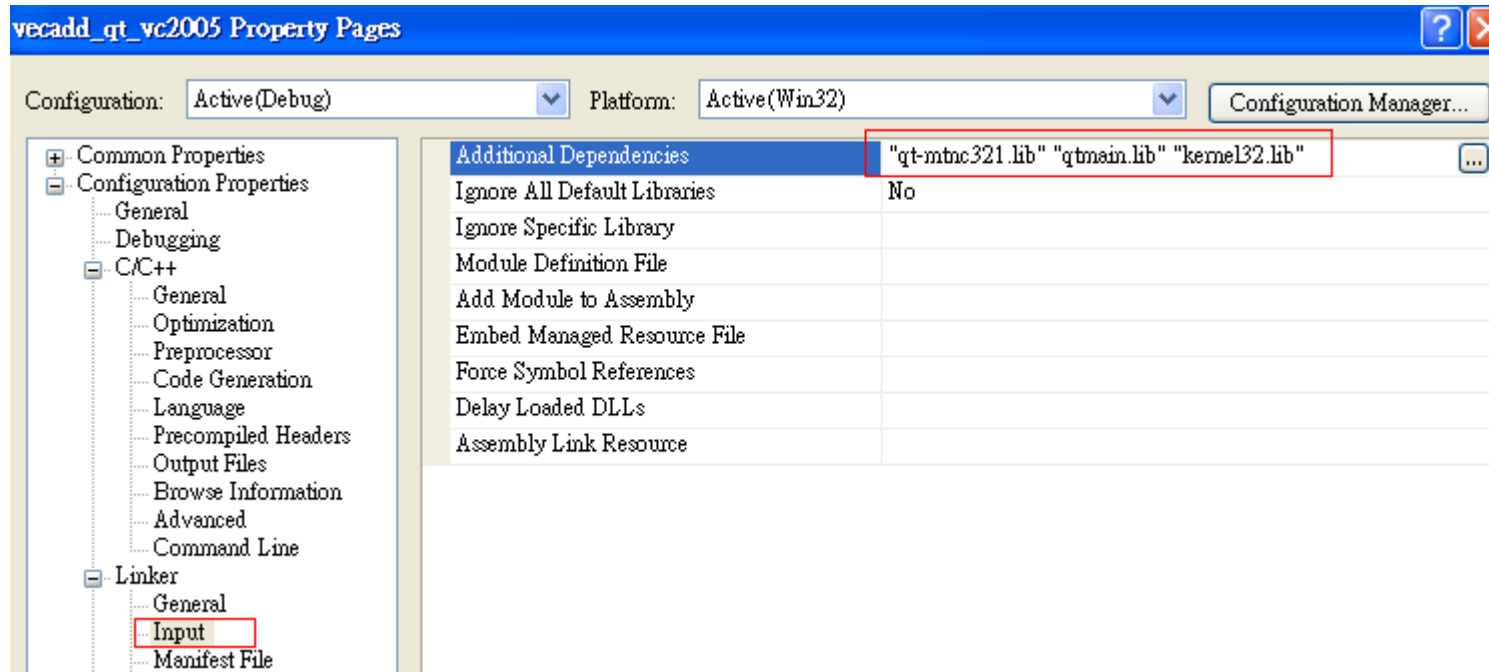
`$(QTDIR)\lib;C:\Program Files (x86)\Microsoft Visual Studio 8\VC\lib`



### Example 3: vector addition (Qtime) [9]

Step 8: project → properties → Linker → Input → Additional Dependence

"qt-mtnc321.lib" "qtmain.lib" "kernel32.lib"



Step 9: compile and execute

**Restriction:** QT3 in windows only support 32-bit application, we must choose platform as "Win32", we will solve this problem after installing QT4

# OutLine

- OpenMP introduction
- Example 1: hello world
- Example 2: vector addition
- enable openmp in vc2005
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

## Example 4: matrix multiplication [1]

matrixMul.h

```

36 #ifndef _MATRIXMUL_H_
37 #define _MATRIXMUL_H_
38
39 // Thread block size
40 #define BLOCK_SIZE 16
41
42 // Matrix dimensions
43 #define WA (25 * BLOCK_SIZE) // Matrix A width
44 #define HA (25 * BLOCK_SIZE) // Matrix A height
45 #define WB (25 * BLOCK_SIZE) // Matrix B width
46 #define HB WA // Matrix B height
47 #define WC WB // Matrix C width
48 #define HC HA // Matrix C height
49
50 #endif // _MATRIXMUL_H_

```

matrixMul.cpp

```

2  #include "matrixMul.h"
3
4  void matrixMul_seq(float* C, const float* A, const float* B,
5      unsigned int hA, unsigned int wA, unsigned int wB ) ;
6
7  void matrixMul_parallel(float* C, const float* A, const float* B,
8      unsigned int hA, unsigned int wA, unsigned int wB, int nthreads ) ;
9
10 // c = a * B
11 void matrixMul_seq(float* C, const float* A, const float* B,
12     unsigned int hA, unsigned int wA, unsigned int wB )
13 {
14     double sum ;
15     unsigned int i, j, k ;
16     double a, b ;
17     for ( i = 0; i < hA; ++i){
18         for (j = 0; j < wB; ++j) {
19             sum = 0;
20             for (k = 0; k < wA; ++k) {
21                 a = A[i * wA + k];
22                 b = B[k * wB + j];
23                 sum += a * b;
24             } // for k
25             C[i * wB + j] = (float)sum;
26         } // for j
27     } // for i
28 }

```

sequential version

$$c_{ij} = \sum_{k=1}^{wA} a_{ik} b_{kj}$$

row-major index

$$a_{ik} = A[i \times wA + k]$$

$$b_{kj} = B[k \times wB + j]$$

$$c_{ij} = C[i \times wC + j]$$

## Example 4: matrix multiplication [2]

### matrixMul.cpp

```
30 void matrixMul_parallel(float* C, const float* A, const float* B,
31     unsigned int hA, unsigned int wA, unsigned int wB, int nthreads )
32 {
33     double sum ;
34     int i, j, k ;
35     double a, b ;
36     #pragma omp parallel default(none) num_threads(nthreads) \
37         shared(A,B,C, hA, wB, wA) private(i,j,k,sum,a,b)
38     {
39         #pragma omp for schedule( static) nowait
40         for (i = 0; i < hA; ++i){
41             for (j = 0; j < wB; ++j) {
42                 sum = 0;
43                 for (k = 0; k < wA; ++k) {
44                     a = A[i * wA + k];
45                     b = B[k * wB + j];
46                     sum += a * b;
47                 } // for k
48                 C[i * wB + j] = (float)sum;
49             } // for j
50         } // for i
51     } // end of parallel section
52 }
53
```

parallel version

**Question 8:** we have three for-loop, one is for “*i*”, one is for “*j*” and last one is for “*k*”, which one is parallelized by OpenMP directive?

**Question 9:** explain why variable *i*, *j*, *k*, *sum*, *a*, *b* are declared as *private*? Can we move some of them to *shared* clause?

## Example 4: matrix multiplication [3]

main.cpp

```
2  #include <stdlib.h>
3  #include <stdio.h>
4  #include <string.h>
5  #include <math.h>
6  #include <assert.h>
7  #include <iostream>
8
9  #include <qdatetime.h> ← use QT timer
10
11 #include "matrixMul.h"
12
13 void runTest(int argc, char** argv) ;
14 void randomInit(float*, int);
15
16 void matrixMul_seq(float* C, const float* A, const float* B,
17     unsigned int hA, unsigned int wA, unsigned int wB ) ;
18
19 void matrixMul_parallel(float* C, const float* A, const float* B,
20     unsigned int hA, unsigned int wA, unsigned int wB, int nthreads ) ;
21
22 int main(int argc, char** argv)
23 {
24     runTest(argc, argv);
25     return 0 ;
26 }
27
28 void runTest(int argc, char** argv)
29 {
30     int nthreads = 2 ;
31     unsigned int total_size = 0 ;
32     QTime t; // QT timer
33
34     // set seed for rand()
35     srand(2006);
```

## main.cpp

```

37 // allocate host memory for matrices A and B
38 unsigned int size_A = WA * HA;
39 unsigned int mem_size_A = sizeof(float) * size_A;
40 float* h_A = (float*) malloc(mem_size_A);
41 assert( h_A );
42
43 total_size += mem_size_A ;
44
45 unsigned int size_B = WB * HB;
46 unsigned int mem_size_B = sizeof(float) * size_B;
47 float* h_B = (float*) malloc(mem_size_B);
48 assert( h_B );
49 total_size += mem_size_B ;
50
51 // initialize host memory
52 randomInit(h_A, size_A);
53 randomInit(h_B, size_B);
54
55 unsigned int size_C = WC * HC;
56 unsigned int mem_size_C = sizeof(float) * size_C;
57 float* h_C = (float*) malloc(mem_size_C);
58 assert( h_C );
59 total_size += mem_size_C ;
60
61 t.start() ;
62 if ( 1 == nthreads ){
63     matrixMul_seq( h_C, h_A, h_B, HA, WA, WB ) ;
64 }else{
65     matrixMul_parallel( h_C, h_A, h_B,
66         HA, WA, WB, nthreads ) ;
67 }
68 printf("threads = %d, matrixMul cost = %d (ms)\n",
69     nthreads, t.elapsed() ) ;
70 printf( "size(A) = (%d,%d)\n", HA, WA );
71 printf( "size(B) = (%d,%d)\n", HB, WB );
72 printf("total memory size = %6.4f (MB)\n",
73     total_size/1048576.0 );
74 // clean up memory
75 free(h_A); free(h_B); free(h_C);
76 }

```

## use qmake to generate Makefile

```

[macrold@quartet2 matrixMul]$ ls
main.cpp  matrixMul.cpp  matrixMul.h
[macrold@quartet2 matrixMul]$ qmake -project
[macrold@quartet2 matrixMul]$ ls
main.cpp  matrixMul.cpp  matrixMul.h  matrixMul.pro
[macrold@quartet2 matrixMul]$ qmake -spec linux-icc-openmp matrixMul.pro
[macrold@quartet2 matrixMul]$ ls
Makefile  main.cpp  matrixMul.cpp  matrixMul.h  matrixMul.pro
[macrold@quartet2 matrixMul]$ make
icpc -c -w -O2 -openmp -mp -DQT_NO_DEBUG -DQT_SHARED -DQT_THREAD_SUPPORT
/qt/mkspecs/linux-icc-openmp -I. -I. -I/opt/qt/include -o main.o main.cpp
icpc -c -w -O2 -openmp -mp -DQT_NO_DEBUG -DQT_SHARED -DQT_THREAD_SUPPORT
/qt/mkspecs/linux-icc-openmp -I. -I. -I/opt/qt/include -o matrixMul.o matr
cpp
matrixMul.cpp(39): (col. 5) remark: OpenMP DEFINED LOOP WAS PARALLELIZED.
matrixMul.cpp(36): (col. 1) remark: OpenMP DEFINED REGION WAS PARALLELIZED
icpc -openmp -Qoption,ld,-rpath,/opt/qt/lib -o matrixMul main.o matrixMul
L/opt/qt/lib -L/usr/X11R6/lib -lqt-mt -lXext -lX11 -lm
[macrold@quartet2 matrixMul]$ ls
Makefile  main.o  matrixMul.cpp  matrixMul.o
main.cpp  matrixMul  matrixMul.h  matrixMul.pro
[macrold@quartet2 matrixMul]$
[macrold@quartet2 matrixMul]$ ./matrixMul
threads = 2, matrixMul cost = 125 (ms)
size(A) = (400,400)
size(B) = (400,400)
total memory size = 1.8311 (MB)
[macrold@quartet2 matrixMul]$

```

## Example 4: matrix multiplication [5]

Let  $BLOCK\_SIZE = 16$  and  $size(A) = size(B) = size(C) = (N \times BLOCK\_SIZE)^2$

total memory usage =  $size(A) + size(B) + size(C)$  float

Platform: oectet1, with compiler icpc 10.0, -O2

$N$	Total size	Thread 1	Thread 2	Thread 4	Thread 8
16	0.75 MB	53 ms	31 ms	21 ms	24ms
32	3 MB	434 ms	237 ms	121 ms	90 ms
64	12 MB	17,448 ms	8,964 ms	6,057 ms	2,997 ms
128	48 MB	421,854 ms	312,983 ms	184,695 ms	92,862 ms
256	192 MB	4,203,536 ms	2,040,448 ms	1,158,156 ms	784,623 ms

Large performance gap amogn  $N = 32$ ,  $N = 64$  and  $N = 128$ , so this algorithm is **NOT** good. Besides improvement of multi-thread is not significant.



## Example 4: matrix multiplication [6]

```
[macrold@octet1 matrixMul]$  
[macrold@octet1 matrixMul]$ ./matrixMul
```

■ running

Use command “top” to see resource usage

```
[macrold@octet1 ~]$  
[macrold@octet1 ~]$ top
```

top - 12:57:59 up 84 days, 21:39, 2 users, load average: 1.38, 0.30, 0.10  
Tasks: 198 total, 4 running, 194 sleeping, 0 stopped, 0 zombie  
Cpu0 : 99.7%us, 0.3%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu1 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu2 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu3 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu4 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu5 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu6 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Cpu7 : 100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st  
Mem: 65342468k total, 11876088k used, 53466380k free, 416924k buffers  
Swap: 67103496k total, 30464k used, 67073032k free, 10193384k cached

PID	USER	PR	NI	VIRT	RES	SHR	S	%CPU	%MEM	TIME+	COMMAND
12290	macrold	20	0	291m	135m	4988	R	799	0.2	1:16.45	matrixMul
12287	macrold	20	0	14668	1144	812	R	0	0.0	0:00.05	top
13120	gdm	20	0	98.8m	3428	2740	S	0	0.0	34:43.42	at-spi-registry
1	root	20	0	10328	280	252	S	0	0.0	0:14.56	init
2	root	15	5	0	0	0	S	0	0.0	0:00.00	bash

CPU usage is 800 %, 8 cores are busy

### Exercise 3: verify subroutine *matrixMul\_parallel*

#### matrixMul.cpp

```
30 void matrixMul_parallel(float* C, const float* A, const float* B,
31     unsigned int hA, unsigned int wA, unsigned int wB, int nthreads )
32 {
33     double sum ;
34     int i, j, k ;
35     double a, b ;
36     #pragma omp parallel default(none) num_threads(nthreads) \
37         shared(A,B,C, hA, wB, wA) private(i,j,k,sum,a,b)
38     {
39         #pragma omp for schedule( static) nowait
40         for (i = 0; i < hA; ++i){
41             for (j = 0; j < wB; ++j) {
42                 sum = 0;
43                 for (k = 0; k < wA; ++k) {
44                     a = A[i * wA + k];
45                     b = B[k * wB + j];
46                     sum += a * b;
47                 } // for k
48                 C[i * wB + j] = (float)sum;
49             } // for j
50         } // for i
51     } // end of parallel section
52 }
53
```

# Combine Parallel Work-sharing constructs

## parallel for Construct

The **parallel for** directive is a shortcut for a **parallel** region that contains only a single **for** directive. The syntax of the **parallel for** directive is as follows:

```
#pragma omp parallel for [clause[,] clause] ...] new-line  
for-loop
```

This directive allows all the clauses of the **parallel** directive and the **for** directive, except the **nowait** clause, with identical meanings and restrictions. The semantics are identical to explicitly specifying a **parallel** directive immediately followed by a **for** directive.

### matrixMul.cpp

```
30 void matrixMul_parallel(float* C, const float* A, const float* B,  
31 unsigned int hA, unsigned int wA, unsigned int wB, int nthreads )  
32 {  
33     double sum ;  
34     int i, j, k ;  
35     double a, b ;  
36     #pragma omp parallel for default(none) num_threads(nthreads) \  
37     shared(A,B,C, hA, wB, wA) private(i,j,k,sum,a,b) \  
38     schedule( static)  
39     for (i = 0; i < hA; ++i){  
40         for (j = 0; j < wB; ++j) {  
41             sum = 0;  
42             for (k = 0; k < wA; ++k) {  
43                 a = A[i * wA + k];  
44                 b = B[k * wB + j];  
45                 sum += a * b;  
46             } // for k  
47             C[i * wB + j] = (float)sum;  
48         } // for j  
49     } // for i  
50 }
```

**Exercise 4:** verify following subroutine *matrix\_parallel*, which parallelizes loop-*j* , not loop-*i*.

1. Performance between loop-*i* and loop-*j*
2. why do we declare index *i* as shared variable? What happens if we declare index *i* as private variable?

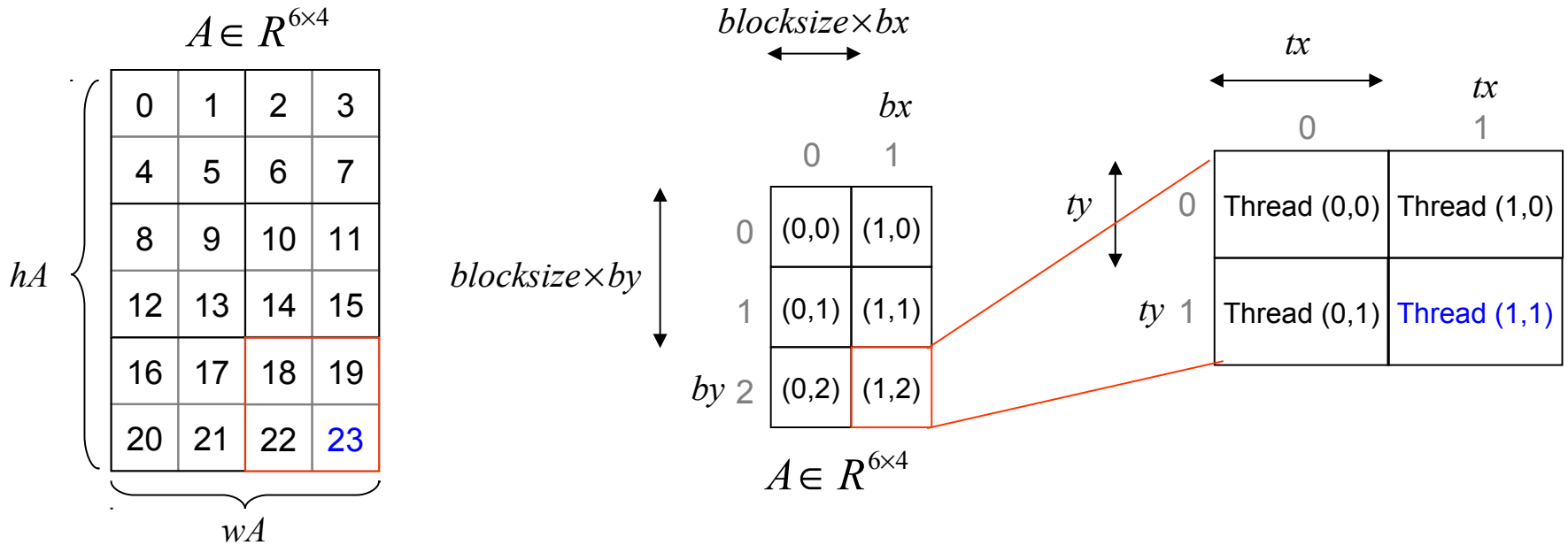
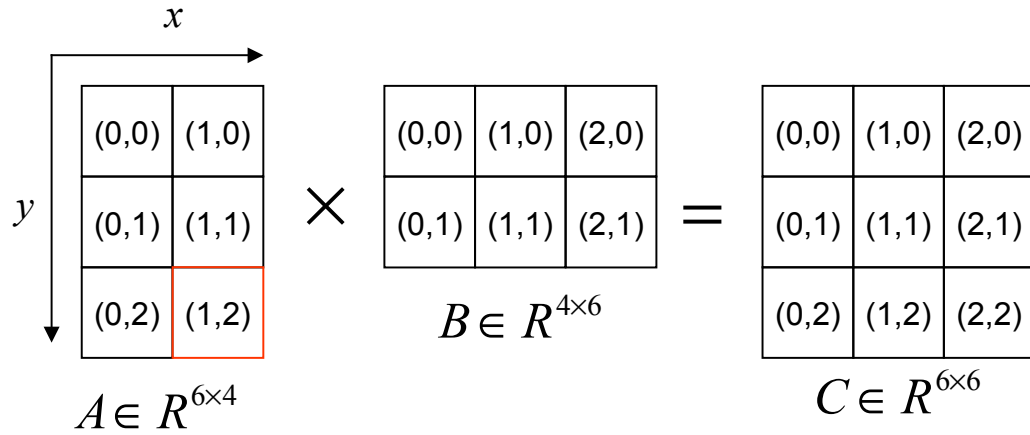
matrixMul.cpp

```
30 void matrixMul_parallel(float* C, const float* A, const float* B,
31   unsigned int hA, unsigned int wA, unsigned int wB, int nthreads )
32 {
33     double sum ;
34     int i, j, k ;
35     double a, b ;
36
37     for (i = 0; i < hA; ++i){
38         #pragma omp parallel for default(none) num_threads(nthreads) \
39           shared(A,B,C, hA, wB, wA, i) private(j,k,sum,a,b) \
40           schedule( static)
41         for (j = 0; j < wB; ++j) {
42             sum = 0;
43             for (k = 0; k < wA; ++k) {
44                 a = A[i * wA + k];
45                 b = B[k * wB + j];
46                 sum += a * b;
47             } // for k
48             C[i * wB + j] = (float)sum;
49         } // for j
50     } // for i
51 }
```

# OutLine

- OpenMP introduction
- Example 1: hello world
- Example 2: vector addition
- enable openmp in vc2005
- Example 3: vector addition + Qtime
- Example 4: matrix multiplication
- Example 5: matrix multiplication (block version)

# Example 5: matrix multiplication (block version) [1]



global index

$$((bx, by), (tx, ty)) \longrightarrow (blocksize \times bx + tx, blocksize \times by + ty) \longrightarrow \text{row-major}$$

## Example 5: matrix multiplication (block version) [2]

matrixMul\_block.cpp

```

7  // store the sub-matrix of A
8  doublereal As[BLOCK_SIZE][BLOCK_SIZE];
9  // store the sub-matrix of B
10 doublereal Bs[BLOCK_SIZE][BLOCK_SIZE];
11
12 // Matrix multiplication on the device: C = A * B
13 void matrixMul_block_seq( doublereal* C, doublereal* A, doublereal* B,
14     const int hA, const int wA, const int wB,
15     const int hA_grid, const int wA_grid, const int wB_grid )
16 {
17     int bx, by ; // Block index
18     int tx, ty ;
19     int aBegin, aEnd, aStep, bBegin, bStep ;
20     int a, b, c, k ;
21 #ifdef HIGH_PRECISION_PACKAGE
22     doublereal Asub, Bsub, Csub ;
23 #else
24     double Asub, Bsub, Csub ;
25 #endif
26
27     unsigned long int i ;
28     unsigned long int size = hA*wB ; // size of matrix C
29 #ifdef HIGH_PRECISION_PACKAGE
30     doublereal* C_ptr = C ;
31     for (i = 0 ; i < size ; i++) {
32         *C_ptr++ = 0.0 ;
33     }
34 #else
35     memset( C, 0, size*sizeof(doublereal) ) ;
36 #endif

```

} Shared memory in GPU

(0,0)	(1,0)
(0,1)	(1,1)
(0,2)	(1,2)

$A \in R^{6 \times 4}$

×

(0,0)	(1,0)	(2,0)
(0,1)	(1,1)	(2,1)

$B \in R^{4 \times 6}$

$hA\_grid = 3 \quad wA\_grid = 2 \quad wB\_grid = 3$

## Example 5: matrix multiplication (block version) [3]

matrixMul\_block.cpp

```

38  for ( by = 0 ; by < hA_grid ; by++ ){
39      for ( bx = 0 ; bx < wB_grid ; bx++ ){
40          // Index of the first sub-matrix of A processed by the block
41          aBegin = wA * BLOCK_SIZE * by;
42          // Index of the last sub-matrix of A processed by the block
43          aEnd   = aBegin + wA - 1;
44          // Step size used to iterate through the sub-matrices of A
45          aStep  = BLOCK_SIZE;
46          // Index of the first sub-matrix of B processed by the block
47          bBegin = BLOCK_SIZE * bx;
48          // Step size used to iterate through the sub-matrices of B
49          bStep  = BLOCK_SIZE * wB;
50
51          // Loop over all the sub-matrices of A and B
52          // required to compute the block sub-matrix
53          for (a = aBegin, b = bBegin;
54              a <= aEnd;
55              a += aStep, b += bStep) {
56              // Load the matrices from main memory
57              for ( ty = 0 ; ty < BLOCK_SIZE ; ty++ ){
58                  for ( tx = 0 ; tx < BLOCK_SIZE ; tx++ ){
59                      As[ty][tx] = A[a + wA * ty + tx];
60                      Bs[ty][tx] = B[b + wB * ty + tx];
61                  } // for tx ;
62              } // for ty

```

} copy global data to small block, why?

(0,0)	(1,0)
(0,1)	(1,1)
(0,2)	(1,2)

×

(0,0)	(1,0)	(2,0)
(0,1)	(1,1)	(2,1)

$B \in R^{4 \times 6}$

$aBegin$  = physical index of first entry in block  $A$   $\begin{pmatrix} (0,1) \end{pmatrix}$

$bBegin$  = physical index of first entry in block  $B$   $\begin{pmatrix} (1,0) \end{pmatrix}$

$A \in R^{6 \times 4}$



## Example 5: matrix multiplication (block version) [4]

matrixMul\_block.cpp

```

64      // Multiply the two matrices together
65      for ( ty = 0 ; ty < BLOCK_SIZE ; ty++ ){
66      for ( tx = 0 ; tx < BLOCK_SIZE ; tx++ ){
67          Csub = 0.0 ;
68      for (k = 0; k < BLOCK_SIZE; ++k ){
69          Asub = As[ty][k ] ;
70          Bsub = Bs[k ][tx] ;
71          Csub += Asub * Bsub ;
72      }
73      c = wB * BLOCK_SIZE * by + BLOCK_SIZE * bx;
74      C[c + wB * ty + tx] += (double) Csub;
75      }// for tx ;
76      }// for ty
77      }// for each submatrix A and B
78
79      }// for bx
80      }// for by
81  }
    
```

Compute submatrix of C sequentially

$$C(i, j) = \sum_{k=1}^{w_A} A(i, k) B(k, j)$$

for all  $(i, j) \in \text{block}(1, 1)$

(0,0)	(1,0)
(0,1)	(1,1)
(0,2)	(1,2)

$$A \in R^{6 \times 4}$$

$\times$

(0,0)	(1,0)	(2,0)
(0,1)	(1,1)	(2,1)

$$B \in R^{4 \times 6}$$

$=$

(0,0)	(1,0)	(2,0)
(0,1)	(1,1)	(2,1)
(0,2)	(1,2)	(2,2)

$$C \in R^{6 \times 6}$$

or equivalently

$$A \begin{bmatrix} (0,1) \end{bmatrix} B \begin{bmatrix} (1,0) \end{bmatrix} + A \begin{bmatrix} (1,1) \end{bmatrix} B \begin{bmatrix} (1,1) \end{bmatrix} = C \begin{bmatrix} (1,1) \end{bmatrix}$$

## Example 5: matrix multiplication (block version) [5]

### Parallel version

```

124 // Loop over all the sub-matrices of A and B
125 // required to compute the block sub-matrix
126 for (a = aBegin, b = bBegin;
127      a <= aEnd;
128      a += aStep, b += bStep) {
129     // Load the matrices from main memory
130     #pragma omp parallel for default(none) num_threads(nthreads) \
131     shared(A,B,As,Bs,a,b,wA,wB) private(ty,tx) \
132     schedule(static)
133     for (ty = 0 ; ty < BLOCK_SIZE ; ty++){
134     for (tx = 0 ; tx < BLOCK_SIZE ; tx++){
135         As[ty][tx] = A[a + wA * ty + tx];
136         Bs[ty][tx] = B[b + wB * ty + tx];
137     } // for tx ;
138 } // for ty
139
140 // Multiply the two matrices together
141 #pragma omp parallel for default(none) num_threads(nthreads) \
142 shared(As,Bs,C,bx,by,wB) private(ty,tx,k,c,Asub,Bsub,Csub) \
143 schedule(static)
144 for (ty = 0 ; ty < BLOCK_SIZE ; ty++){
145 for (tx = 0 ; tx < BLOCK_SIZE ; tx++){
146     Csub = 0.0 ;
147     for (k = 0; k < BLOCK_SIZE; ++k){
148         Asub = As[ty][k] ;
149         Bsub = Bs[k][tx] ;
150         Csub += Asub * Bsub ;
151     }
152     c = wB * BLOCK_SIZE * by + BLOCK_SIZE * bx;
153     C[c + wB * ty + tx] += (double) Csub;
154 } // for tx ;
155 } // for ty
156 } // for each submatrix A and B
157
158 } // for bx
159 } // for by
160 }

```

### GPU code

```

for (int a = aBegin, b = bBegin;
     a <= aEnd;
     a += aStep, b += bStep) {

    // Declaration of the shared memory array As
    // used to store the sub-matrix of A
    __shared__ float As[BLOCK_SIZE][BLOCK_SIZE];

    // Declaration of the shared memory array Bs
    // used to store the sub-matrix of B
    __shared__ float Bs[BLOCK_SIZE][BLOCK_SIZE];

    // Load the matrices from device memory
    // to shared memory; each thread loads
    // one element of each matrix
    AS(ty, tx) = A[a + wA * ty + tx];
    BS(ty, tx) = B[b + wB * ty + tx];

    // Synchronize to make sure the matrices are
    __syncthreads();

    // Multiply the two matrices together;
    // each thread computes one element
    // of the block sub-matrix
    for (int k = 0; k < BLOCK_SIZE; ++k)
        Csub += AS(ty, k) * BS(k, tx);

    // Synchronize to make sure that the precedi
    // computation is done before loading two ne
    // sub-matrices of A and B in the next itera
    __syncthreads();

}

// Write the block sub-matrix to device memory;
// each thread writes one element
int c = wB * BLOCK_SIZE * by + BLOCK_SIZE * bx;
C[c + wB * ty + tx] = Csub;

```

## Example 5: matrix multiplication (block version) [6]

Let  $BLOCK\_SIZE = 16$  and  $size(A) = size(B) = size(C) = (N \times BLOCK\_SIZE)^2$

total memory usage =  $size(A) + size(B) + size(C)$  float

Platform: oectet1, with compiler icpc 10.0, -O2

$N$	Total size	Thread 1	Thread 2	Thread 4	Thread 8
16	0.75 MB	40 ms	34 ms	34 ms	44 ms
32	3 MB	301 ms	309 ms	240 ms	219 ms
64	12 MB	2,702 ms	2,310 ms	1,830 ms	1,712 ms
128	48 MB	24,548 ms	19,019 ms	15,296 ms	13,920 ms
256	192 MB	198,362 ms	151,760 ms	129,754 ms	110,540 ms

## Non-block version

$N$	Total size	Thread 1	Thread 2	Thread 4	Thread 8
16	0.75 MB	53 ms	31 ms	21 ms	24 ms
32	3 MB	434 ms	237 ms	121 ms	90 ms
64	12 MB	17,448 ms	8,964 ms	6,057 ms	2,997 ms
128	48 MB	421,854 ms	312,983 ms	184,695 ms	92,862 ms
256	192 MB	4,203,536 ms	2,040,448 ms	1,158,156 ms	784,623 ms

**Question 10:** non-block version is much slower than block version, why?

## Example 5: matrix multiplication (block version) [7]

Block version, BLOCK\_SIZE = 512

$N$	Total size	Thread 1	Thread 2	Thread 4	Thread 8
2	12 MB	3,584 ms	1,843 ms	961 ms	453 ms
4	48 MB	27,582 ms	14,092 ms	7,040 ms	3,533 ms
8	192 MB	222,501 ms	110,975 ms	55,894 ms	28,232 ms

Block version, BLOCK\_SIZE = 16

$N$	Total size	Thread 1	Thread 2	Thread 4	Thread 8
64	12 MB	2,702 ms	2,310 ms	1,830 ms	1,712 ms
128	48 MB	24,548 ms	19,019 ms	15,296 ms	13,920 ms
256	192 MB	198,362 ms	151,760 ms	129,754 ms	110,540 ms

**Question 11:** larger BLOCK\_SIZE implies better performance when using multi-thread, why?

**Question 12:** small BLOCK\_SIZE is better in single thread, why?

**Question 13:** matrix-matrix multiplication is of complexity  $O(N^3)$ , which algorithm is “good” to achieve this property?

## Example 5: matrix multiplication (block version) [8]

```
[macrold@octet1 matrixMul_block2]$  
[macrold@octet1 matrixMul_block2]$ cat /proc/cpuinfo
```

```
processor       : 0  
vendor_id      : GenuineIntel  
cpu family     : 6  
model          : 15  
model name     : Intel(R) Xeon(R) CPU           X5365  @ 3.00GHz  
stepping       : 11  
cpu MHz        : 2000.000  
cache size     : 4096 KB  
physical id    : 0  
siblings       : 4  
core id        : 0  
cpu cores      : 4  
fpu            : yes  
fpu_exception  : yes  
cpuid level    : 10  
wp             : yes  
flags           : fpu vme de pse tsc msr pae mce cx8 apic sep mtrr pge mca  
cmov pat pse36 clflush dts acpi mmx fxsr sse sse2 ss ht tm pbe syscall nx l  
m constant_tsc arch_perfmon pebs bts rep_good pni monitor ds_cpl vmx est tm  
2 sse3 cx16 xtpr dca lahf_lm  
bogomips       : 5987.36  
clflush size   : 64  
cache_alignme  : 64  
address sizes   : 38 bits physical, 48 bits virtual  
power manage    :
```

Cache has 4 MB, we can have large BLOCK\_SIZE

cache line is 64 byte (16 float)

In CPU

BLOCK\_SIZE = 512  $\longrightarrow size(Bs) = size(As) = 512^2 \text{ float} = 1024^2 \text{ Byte} = 1\text{MB}$

In GPU

BLOCK\_SIZE = 16  $\longrightarrow size(Bs) = size(As) = 16^2 \text{ float} = 1\text{kB}$

**Exercise 5:** verify subroutine *matrixMul\_block\_seq* with non-block version, you can use high precision package.

### Non-block version

```
12 // C = A * B
13 void matrixMul_seq(doublereal* C, const doublereal* A, const doublereal* B,
14     unsigned int hA, unsigned int wA, unsigned int wB )
15 {
16     unsigned int i, j, k ;
17     #ifdef HIGH_PRECISION_PACKAGE
18         doublereal sum ;
19         doublereal a, b ;
20     #else
21         double sum ;
22         double a, b ;
23     #endif
24     for ( i = 0; i < hA; ++i){
25         for (j = 0; j < wB; ++j) {
26             sum = 0.0 ;
27             for (k = 0; k < wA; ++k) {
28                 a = A[i * wA + k];
29                 b = B[k * wB + j];
30                 sum += a * b;
31             } // for k
32             C[i * wB + j] = (doublereal)sum;
33         } // for j
34     } // for i
35 }
```

**Exercise 6:** if we use “double”, how to choose value of BLOCK\_SIZE, show your experimental result.

**Exercise 7:** Can you modify subroutine *matrixMul\_block\_parallel* to improve its performance?

**Exercise 8:** compare parallel computation between CPU and GPU in your host machine