Optimization techniques: an introduction

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Agenda:

- Introduction Philosophy & Performance Issues
- memory hierarchy
  - Cache & other stuff..
- Performance and evaluating process (Profiling and timing your code)
- Optimization techniques
- General Performance techniques
  - Use of Libraries: see next lecture..
The Philosophy...

- Discuss how to measure performances of your cluster/system
- Discuss performance tuning techniques common to most modern architecture (mainly 32/64 bit commodity processor)
- Using optimization techniques users have control over
  - Code modification
  - Compiler options
- Optimization is a dirty work (and dangerous one for your code...)
- Compiler is your best friend..
Memory Hierarchy

- In modern computer system same data is stored in several storage devices during processing
- The storage devices can be described & ranked by their speed and “distance” from the CPU
- There is thus a hierarchy of memory objects
- Programming a machine with memory hierarchy requires optimization for that memory structure.
Memory Hierarchy

• Most programs have a high degree of **locality** in their accesses
• Memory hierarchy tries to exploit locality

---

**Diagram:**
- Processor
  - Control
  - Datapath
  - Registers
  - On-chip cache
- Second level cache (SRAM)
- Main memory (DRAM)
- Secondary storage (Disk)
- Tertiary storage (Disk/Tape)

**Table:**

<table>
<thead>
<tr>
<th>Speed</th>
<th>1ns</th>
<th>10ns</th>
<th>100ns</th>
<th>10ms</th>
<th>10sec</th>
</tr>
</thead>
<tbody>
<tr>
<td>Size</td>
<td>B</td>
<td>KB</td>
<td>MB</td>
<td>GB</td>
<td>TB</td>
</tr>
</tbody>
</table>
Locality of Reference

- **Temporal locality:** Recently referenced items (instr or data) are likely to be referenced again in the near future:
  - iterative loops, subroutines, local variables
  - working set concept
- **Spatial locality:** programs access data which is near to each other:
  - operations on tables/arrays
  - cache line size is determined by spatial locality
- **Sequential locality:** processor executes instructions in program order:
  - branches/in-sequence ratio is typically 1 to 5
Performance Evaluation process

- Monitoring System:
  - Use monitoring tools to better understand your machine’s limits and usage
    • Is the system limit well suited to run my application?
  - Observe both overall system performance and single-program execution characteristics. Monitoring your own code
    • Is the system doing well? Is my program running in a pathological situation?

- Monitoring your own code:
  - Timing the code:
    • timing a whole program (time command :/usr/bin/time)
    • timing a portion (all portions) of the program
  - Profiling the program
What to measure?

- Scire est Mensurare - J.Keppler
- Example measurements to help draw conclusions
- Elapsed execution time of (part of) a program
  - Which algorithm runs faster in practice?
- Number of instructions executed, absolute total or rate per second
  - Compare to theoretical peak performance e.g. instructions per cycle (IPC)
- Rate of floating point operations per second (Mflops)
  - Compare to theoretical or sustained Mflops peak performance
- System events, e.g. cache hits, system calls
  - Determine how much and when/where overhead occurs
- Communication latency
  - Reorganize communications and computation (improve latency hiding)
- Parallel speedup
  - How effective is parallelism?
Useful Monitoring Commands (Linux)

- **Uptime(1)** returns information about system usage and user load
- **ps(1)** lets you see a “snapshot” of the process table
- **top** process table dynamic display
- **free** memory usage
- **vmstat** memory usage monitor

```
top - 08:48:53 up 6 days, 18:35, 5 users, load average: 0.98, 0.55, 0.22
Tasks: 91 total, 2 running, 89 sleeping, 0 stopped, 0 zombie
Cpu(s): 99.3% us, 0.7% sy, 0.0% ni, 0.0% id, 0.0% wa, 0.0% hi, 0.0% si
Mem: 507492k total, 440336k used, 67156k free, 57928k buffers
Swap: 2048248k total, 108456k used, 1939792k free, 99908k cached

   PID USER      PR NI VIRT  RES   SHR S %CPU %MEM    TIME+  COMMAND
18806 cozzini  25  0  2164  464  396 R  90.8  0.1   3:52.85  a.out
  9874 moses    15  0  211m  94m  18m S  7.6  19.0  54:25.92  firefox-bin
  9578 root     15  0  194m  38m  4836 S  7.7  0.0   16:00.59  X
18807 cozzini  16  0  2320  956  756 R  0.3  0.2   0:00.53  top
```
Monitoring your own code (time)

NAME
time - time a simple command or give resource usage

SYNOPSIS
time [options] command [arguments...]

DESCRIPTION
The `time` command runs the specified program command with
the given arguments. When command finishes, time writes a
message to standard output giving timing statistics about
this program..

-------------> time ./a.out
[program output]

real  0m1.361s
user  0m0.770s
sys   0m0.590s

user time: Cpu-time dedicated to your program
sys time: time used by your program to execute
system calls
real time: total time aka walltime
User/System/Walltime

- **Real time (or wall clock time)** is the total elapsed time from start to end of a timed task

- **CPU user time** is the time spent executing in user space
  - Does not include time spent in system (OS calls) and time spent executing other processes

- **CPU system time** is the time spent executing system calls (kernel code)
  - System calls for I/O, devices, synchronization and locking, threading, memory allocation
  - Typically does not include process waiting time for non-ready devices such as disks

- CPU user time + CPU system time < real time
  - CPU percentage spent on process = 100% * (user+system) / real
a top disaster: swapping..

• virtual or swap memory:
  • This memory is actually space on the hard drive. The operating system reserves a space on the hard drive for “swap space”.
  •

• time to access virtual memory VERY large:
• this time is done by the system not by your program!
• sometimes the system assumes a killer to kill your program.
  (oom killer)

```
top - 08:57:02 up 6 days, 19:35, 7 users, load average: 2.77, 0.73, 0.25
Tasks: 86 total, 2 running, 84 sleeping, 0 stopped, 0 zombie
Cpu(s): 0.3% us, 4.8% sy, 0.0% ni, 0.0% id, 94.2% wa, 0.6% hi, 0.0% si
Mem: 507492k total, 506572k used, 920k free, 196k buffers
Swap: 2048248k total, 941984k used, 1106264k free, 4740k cached

   PID USER   PR NI    VIRT    RES   SHR   %CPU  %MEM    TIME+ COMMAND
11656 cozzini   18   0  2172m  408m  260 D   4.3  82.4  0:03.75  a.out
  33 root      15   0    0     0     0 D   0.7   0.0  0:00.54  kswapd0
 3195 root      15   0  20696 1432 1140 D   0.3   0.3  0:06.81  clock-applet
11656 cozzini   17   0  3540  375  700 D   0.2   0.2  0:00.05  http
```
A top disaster example (1) situation is illustrated with code output.

```
[cozzini@stroligo optimization]$ /usr/bin/time ./a.out
Provide an integer (suggested range 100-250)
larger values can be very memory and time-consuming
300
  initialisation time= 11.787208
10.86user 0.98system 0:14.22elapsed 83%CPU (0avgtext+0avgdata
0maxresident)k
0inputs+0outputs (5major+106090minor)pagefaults 0swaps

[cozzini@stroligo optimization]$ /usr/bin/time ./a.out
Provide an integer (suggested range 100-250)
larger values can be very memory and time-consuming
320
Command terminated by signal 2
0.18user 1.81system 0:29.27elapsed 6%CPU (0avgtext+0avgdata
0maxresident)k
0inputs+0outputs (5846major+170788minor)pagefaults 0swaps
```
top disaster example (2)

[cozzini@stroligo optimization]$ /usr/bin/time ./a.out <300 &
[cozzini@stroligo optimization]$ free

```
total used free shared buffers
cached
Mem: 507492 484916 22576 0 1156 10172
-/+ buffers/cache: 473588 33904
Swap: 2048248 78108 1970140
```

[cozzini@stroligo optimization]$ /usr/bin/time ./a.out <320 &
[cozzini@stroligo optimization]$ free

```
total used free shared buffers
cached
Mem: 507492 506412 1080 0 252 3936
-/+ buffers/cache: 502224 5268
Swap: 2048248 546348 1501900
```
Timing portion of the code

- Record the time before portion A
- execute portion A
- record the time after portion A
- print/save the difference in time for subsequent analysis

- C function to compute time:
  - clock
- Fortran90 function to compute time:
  - cpu_time routine (f95)

```c
void clock_t c0, c1;
c0 = clock();
    // section to code..
c1= clock();
cputime = (c1 - c0)/(CLOCKS_PER_SEC );
```
Well written codes have their own timing report.

!!Specific TIMING for section: MD INTEGRATION !!!!!!!!!!!

```
c !Serial subroutines:

!section times  avg-time    max(PE)     min(PE)
!vscale      2    0.0600    0.0600( 0)    0.0600( 0)
!scanpairs_prot  100  72.5500  72.5500( 0)  72.5500( 0)
!vertest_prot   100  2.2600   2.2600( 0)    2.2600( 0)
!link_list      7    70.2900  70.2900( 0)  70.2900( 0)
!spme_prot      100  727.8700 727.8700( 0) 727.8700( 0)
!fill_charge_gri 100  214.5000 214.5000( 0) 214.5000( 0)
!fft_back       100  79.2700  79.2700( 0)  79.2700( 0)
!scalar_sum     100  43.2400  43.2400( 0)    43.2400( 0)
!fft_forw       100  78.8400  78.8400( 0)  78.8400( 0)
!grad_sum       100 303.6600  303.6600( 0) 303.6600( 0)
!ewcorr_prot    100  24.4000  24.4000( 0)  24.4000( 0)
!ewald3_prot    5870100 15.4300  15.4300( 0)  15.4300( 0)
!pair_force_prot 100  0.0000   0.0000( 0)    0.0000( 0)
!srfew2_prot    5855979 817.5300 817.5300( 0) 817.5300( 0)
!dihfrc_prot    100  3.0800   3.0800( 0)    3.0800( 0)
```
Processes/programs

- A *CPU bound* process is compute intensive
  - Very few I/O operations
  - Execution speed of algorithm is determined by CPU
  - Use performance profiling

- A *Memory bound* process is memory intensive
  - Limited I/O operations and large memory footprint
  - Execution speed is determined by Memory performances
  - Use performance profiling with RAM checkers

- An *I/O bound* process is I/O intensive
  - Process includes I/O operations such as file access, message passing over network
  - Execution speed is limited by system’s I/O latencies
  - Performance analysis method depends on I/O operations: operations on file on disk, messages over pipes and sockets,
Analysis Techniques

• Aside from simple timings, there are three generally available techniques for analyzing code performance:
  – Compiler reports and listings
  – Profiling
  – Hardware performance counters
Compiler Reports and Listings

• Compilers on most modern high performance computers are capable of doing a wide range of optimizations,

• By default, compilers generally do not describe in much detail what kinds of optimizations they were able to perform on a given piece of code.

• However, many compilers will optionally generate optimization reports and/or listing files.
  - Optimization reports are typically sent to stderr at compile time and contain messages describing what optimizations could or could not be applied at various points in the source code.
  - Listing files usually consist of a listing of the source code with messages about optimizations interspersed through the listing.
# Reporting and Listing Compiler Options

## GNU compilers

None

## PGI compilers

- **Minfo=option[,option,...]**
  - Prints information to `stderr` on `option`; `option` can be one or more of `time, loop, inline, sym, or all`

- **Mneginfo=option[,option]**
  - Prints information to `stderr` on why optimizations of type `option` were not performed; `option` can be `concur` or `loop`

- **Mlist**
  - Generates a listing file

## Intel compilers

- **-opt_report**
  - Generates an optimization report on `stderr`

- **-opt_report_file filename**
  - Generates an optimization report to `filename`
Profiling

- Profiling is an approach to performance analysis in which the amount of time spent in sections of code is measured (using either a sampling technique or on entry/exit of a code block) and presented as a histogram.
- This allows a developer to identify the routines which are taking the most execution time, as these are typically the best candidates for optimization.
- Profiling can be done at varying levels of granularity:
  - Subroutine
  - Basic block
  - Source code line
- Profiling usually requires special compilation.
  - The specially compiled executable will generate a file containing execution profile data as it runs.
  - This data file can be analyzed after the code is run
  - A profiling analysis program should be employed
Profiling Compiler Options

FORTRAN INTEL:
-prof-dir <d> specify directory for profiling output files (*.dyn and *.dpi)
-prof-file <f> specify file name for profiling summary file
-prof-gen instrument program for profiling
-prof-use enable use of profiling information during optimization
-qp compile and link for function profiling with UNIX gprof tool
-p same as -qp

FORTRAN PGF90
-Mprof[=dwarf|func|hwcts|lines|mpi|time] Generate additional code for profiling
  dwarf Add limited DWARF info for third party profilers
  func Function-level profiling
  hwcts PAPI-based profiling using hardware counters, 64-bit only
  lines Line-level profiling
  mpi Use profiled MPI communication library; implies -Mmpi
  time Sample-based instruction-level profiling

GNU:
-p Generate extra code to write profile information suitable for the analysis program prof.
-pg Generate extra code to write profile information suitable for the analysis program gprof.
program profiled
    integer,parameter :: N=1000,ntimes=100
    real,dimension(N,N) :: b,c,d
    real,dimension(N) :: a
    real :: begin,end
    real,dimension(2) :: rtime

    call random_number(b)
call random_number(c)
call random_number(d)
begin=dtime(rtime)
do it=1,ntimes
    do j=1,N
        a(j)=myvsum(b,j)+myvprod(c,j)*myvsum(d,j)
    end do
    if (mod(it,100).eq.0) write (*,*) a(1),b(1,1),c(1,1),d(1,1)
end do
end=dtime(rtime)
write (*,*) "loop time = ",end," seconds"
flops=(ntimes*N*(3*N+2))/end*1.0e-6
write (*,*) "loop ran at ",flops," MFLOPS"
Demo: Profiling (g95, con’t)

contains

real function myvsum(x,j)
  real,dimension(N,N),intent(IN) :: x
  integer,intent(IN) :: j
  myvsum=x(1,j)
  do i=2,N
    myvsum=myvsum+x(i,j)
  end do
end function myvsum

real function myvprod(x,j)
  real,dimension(N,N),intent(IN) :: x
  integer,intent(IN) :: j
  myvprod=x(1,j)
  do i=2,N
    myvprod=myvprod*x(i,j)
  end do
end function myvprod

end program profiled
**Demo: Profiling (g95 con’t)**

```bash
[cozzini@stroligo optimization]$ g95 -pg profiled.f90
[cozzini@stroligo optimization]$ ./a.out
[cozzini@stroligo optimization]$ gprof a.out gmon.out
```

Flat profile:

<table>
<thead>
<tr>
<th>Flat profile:</th>
</tr>
</thead>
<tbody>
<tr>
<td>Each sample counts as 0.01 seconds.</td>
</tr>
<tr>
<td>%   cumulative   self              self     total</td>
</tr>
<tr>
<td>seconds   seconds    calls   s/call   s/call  name</td>
</tr>
<tr>
<td>time      seconds   seconds</td>
</tr>
<tr>
<td>-----------</td>
</tr>
<tr>
<td>95.63</td>
</tr>
<tr>
<td>2.40</td>
</tr>
<tr>
<td>1.96</td>
</tr>
<tr>
<td>0.01</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>%</th>
</tr>
</thead>
<tbody>
<tr>
<td>the percentage of the total running time of the</td>
</tr>
<tr>
<td>program used by this function.</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>cumulative</th>
</tr>
</thead>
<tbody>
<tr>
<td>a running sum of the number of seconds accounted</td>
</tr>
<tr>
<td>seconds</td>
</tr>
<tr>
<td>for by this function and those listed above it.</td>
</tr>
</tbody>
</table>
Hardware Performance Counters

- Most modern microprocessors have one or more event counters which can be used to count low level processor events such as floating point operations, cache line misses, and total instructions executed.
- The output from these counters can be used to infer how well a program is utilizing the processor.
- In many cases, there are utilities for accessing these hardware counters, through either a library or a command line timing interface.
How to optimize...

- Iterative optimization
  - Check for correct answers (program must be correct!)
  - Profile to find the hotspots, e.g. most time-consuming routines
  - Optimize these routines using compiler options, compiler directives (pragmas), and source code modifications
  - Repeat 1-3

- Optimizing the hotspots of a program improves overall performance

- Programs with “flat profiles” (flat timing histogram)
  - Programs with lots of routines that each take a small amount of time are difficult to optimize
Optimization Methodology

- Optimize one loop/routine at a time
- Start with the most time consuming routines
- Then the second most....
- Then the third most....
- Parallelise your program..
Optimization Techniques

• There are basically three different categories:
  
  – Improve memory performance (The most important)
    
      • Better memory access pattern
      
      • Optimal usage of cache lines (improve spatial locality)
      
      • Re-usage of cached data (improve temporal locality)
  
  – Improve CPU performance
    
      • Create more opportunities to go superscalar (high level)
      
      • Better instruction scheduling (low level)
  
  – Use already highly optimized libraries/subroutines
## Where to optimize?

<table>
<thead>
<tr>
<th></th>
<th>Annual increase</th>
<th>Typical value in 2006</th>
</tr>
</thead>
<tbody>
<tr>
<td>Single-chip floating-point performance</td>
<td>59%</td>
<td>4 GFLOP/s</td>
</tr>
<tr>
<td>Front-side bus bandwidth</td>
<td>23%</td>
<td>1 GWord/s</td>
</tr>
<tr>
<td></td>
<td></td>
<td>= 0.25 word/flop</td>
</tr>
<tr>
<td>DRAM latency</td>
<td>(5.5%)</td>
<td>70 ns</td>
</tr>
<tr>
<td></td>
<td></td>
<td>= 280 FP ops</td>
</tr>
<tr>
<td></td>
<td></td>
<td>= 70 loads</td>
</tr>
</tbody>
</table>
Optimization Techniques for memory

- Loop Interchanges
- Effective Reuse of Data Cache
- Loop Unrolling
- Loop Fusion/Fission
- Prefetching
- Floating Point Division
Storage in Memory

The storage order is language dependent:

Fortran stores “column-wise”

C stores “row-wise”

Accessing elements in storage order greatly enhances the performance for problem sizes that do not fit in the cache(s)

(spatial locality: stride 1 access)
Array Indexing

There are several ways to index arrays:

- **Direct**

  ```
  Do j=1,M
    Do i=1,N
      ..A(i, j)
    END DO
  END DO
  ```

  **Loop carried**

  ```
  Do j=1,M
    Do i=1,N
      k=k+1
      ..A(k)
    END DO
  END DO
  ```

  **Explicit**

  ```
  Do j=1,M
    Do i=1,N
      ..A(i+(j-1)*N)
    END DO
  END DO
  ```

  **Indirect**

  ```
  Do j=1,M
    Do i=1,N
      ..A(index(i,j))..
    END DO
  END DO
  ```

The addressing scheme can (and will) have an impact on the performance.
Data Dependencies

- Independent instructions can be scheduled at the same time on the multiple execution units in superscalar CPU.
- Independent operations can be (software) pipelined on the CPU.

Loop-carried dependencies

\[
\text{do } i=1,n
\]
\[
\quad a(\text{index}(1,i)) = b(i)
\]
\[
\quad a(\text{index}(2,i)) = c(i)
\]
\[
\text{end do}
\]

Standard prog language (F77/F90/C/C++) do not provide explicit information on data dependencies.

Compilers assume worse case for the data dependencies:
- problem for indirectly addressed arrays in Fortran
- problem for all pointers C
Loop Interchange

Basic idea: In a nested loop, examine and possibly change the order of the loop

Advantages:
Better memory access patterns (leading to improved cache and memory usage)
Elimination of data dependencies (to increase the opportunities for CPU optimization and parallelization)

Disadvantage:
May make a short loop innermost (which is not good for optimal performances)
Loop Interchange - Example 1

**Original**

```plaintext
DO i=1,N
   DO j=1,M
      C(i,j)=A(i,j)+B(i,j)
   END DO
END O
```

**Interchanged loops**

```plaintext
DO j=1,M
   DO i=1,N
      C(i,j)=A(i,j)+B(i,j)
   END DO
END DO
```
Loop Interchange in C

In C, the situation is exactly the opposite

- The performance benefit is the same in this case
- In many practical situations, loop interchange is much easier to achieve than index reversal

```c
for (j=0; j<M; j++)
    for (i=0; i<N; i++)
        C[j][i] = A[j][i] + B[j][i];

for (i=0; i<N; i++)
    for (j=0; j<N; j++)
        C[i][j] = A[i][j] + B[i][j];

for (j=0; j<M; j++)
    for (i=0; i<N; i++)
        C[i][j] = A[i][j] + B[i][j];
```
Loop Interchange – Mnemonic rule

- With row-major, the column or "rightmost" index varies most quickly (C/C+)
- With column-major, the row of "leftmost" index varies most quickly. (Fortran/F90)
**Loop Interchange - Example 2**

Timings are in seconds

```fortran
DO i=1,300
    DO j=1,300
        DO k=1,300
            A (i,j,k) = A (i,j,k)+ B (i,j,k)* C (i,j,k)
        END DO
    END DO
END DO
```

<table>
<thead>
<tr>
<th>Loop order</th>
<th>x335 (P4 2.4Ghz)</th>
<th>x330 (P3 1.4Ghz)</th>
</tr>
</thead>
<tbody>
<tr>
<td>i j k</td>
<td>8.77</td>
<td>9.06</td>
</tr>
<tr>
<td>i k j</td>
<td>7.61</td>
<td>6.82</td>
</tr>
<tr>
<td>j i k</td>
<td>2.00</td>
<td>2.66</td>
</tr>
<tr>
<td>j k i</td>
<td>0.57</td>
<td>1.32</td>
</tr>
<tr>
<td>k i j</td>
<td>0.90</td>
<td>1.95</td>
</tr>
<tr>
<td>k j i</td>
<td>0.44</td>
<td>1.25</td>
</tr>
</tbody>
</table>
Loop Interchange Compiler Options

GNU compilers:

None

PGI compilers:

-Mvect

Enable vectorization, including loop interchange

Intel compilers:

-03

Enable aggressive optimization, including loop transformations

TEST IF WHAT THEY CLAIM TO DO IS WHAT THEY ACTUALLY DO
Cache Thrashing

- Typical problem in code performance is *cache thrashing*.
- Cache thrashing happens when data in cache are rewritten and fully reused.
- In the previous case, the cache thrashing was minimized by *loop interchange*.
- Another optimization technique aimed at minimizing cache thrashing is *COMMON block padding*.
- Consider: $a(i) + b(i) = c(i)$

![Diagram showing CPU, L1 cache, L2 cache, and RAM with data distribution and references between them.](image)
**Demo: Cache Thrashing**

program thrash
    integer,parameter :: N=4*1024*1024
    real,dimension(N) :: c,b,a
    real:: begin,end
    real,dimension(2):: rtime
    common/saver/a,b,c
    !DIR$ CACHE_ALIGN /saver/
    call random_number(b)
call random_number(c)
    begin=dtime(rtime)
do i=1,N
    a(i)=b(i)+c(i)
end do
end=dtime(rtime)
print *,"my loop time (s) is ",end
flop=N/end*1.0e-6
print *,"loop runs at ",flop,' MFLOP'
print *,a(1),b(1),c(1)
end program thrash

~56 Mflops

Considerable cache thrashing occurs as c(1) knocks out the line b(1) is in, then b(2) knocks out the line c(1) was in and so on
demo on cache trashing

In the next two programs arrays \( b \) and \( c \) both fill the cache so when they are used in the addition loop considerable cache thrashing occurs as \( c(1) \) knocks out the line \( b(1) \) is in, then \( b(2) \) knocks out the line \( c(1) \) was in and so on.

The second program illustrates the use of common-block padding. By putting a fake array - \( \text{space}(4) \) - between \( b \) and \( c \) in the COMMON block we shift the memory addresses of \( b \) and \( c \) and thus shift the cache lines they map to. Arrays \( b \) and \( c \) no longer collide in the cache.
Solution: COMMON Block Padding

```fortran
program pad
  integer,parameter :: N=4*1024*1024
  real,dimension(N) :: c,b,a
  real:: begin,end
  real,dimension(2):: rtime
  common/saver/a,b,space(4),c
!DIR$ CACHE_ALIGN /saver/
call random_number(b)
call random_number(c)
begin=dtime(rtime)
do i=1,N
  a(i)=b(i)+c(i)
end do
end=dtime(rtime)
print *,'my loop time (s) is ',end
flop=N/end*1.0e-6
print *,'loop runs at ',flop,' MFLOP'
print *,a(1),b(1),c(1)
end program pad
```

By putting the fake array, `space(4)` , in the COMMON block, arrays `b` and `c` no longer collide in the cache.

~83 MFlops
COMMON Block Padding Compiler Options

GNU compilers:
- \texttt{-malign-double}
  Align double precision variables on 64-bit boundaries, including in COMMON blocks

PGI compilers:
- \texttt{-Mdalign}
  Align doubles [i.e. REAL*8] in COMMON blocks and structures on 8-byte boundaries

Intel compilers:
- \texttt{-zp8}
  Specify alignment constraint for structures on 8-byte boundaries, including in COMMON blocks; default
- \texttt{-pad}
  Enable changing of variable and array memory layout
- \texttt{-nopad}
  Disable changing of variable and array memory layout; default

\textbf{TEST IF WHAT THEY CLAIM TO DO IS WHAT THEY ACTUALLY DO}
Prefetching

- *Prefetching* is the retrieval of data from memory to cache before it is needed in an upcoming calculation. This is an example of general optimization technique called *latency hiding* in which communications and calculations are overlapped and occur simultaneously.

- The actual mechanism for prefetching varies from one machine to another.
Prefetching Compiler Options

GNU:
-ffprefetch-loop-arrays
   If supported by the target machine, generate instructions to prefetch memory to improve the performance of loops that access large arrays.

PGI:
-Mprefetch[=option:n] -Mnoprefetch
   Add (don’t add) prefetch instructions for those processors that support them (Pentium 4, Opteron); -Mprefetch is default on Opteron; -Mnoprefetch is default on other processors.

INTEL:
-03 Enable -02 optimizations and in addition, enable more aggressive optimizations such as loop and memory access transformation, and prefetching.

TEST IF WHAT THEY CLAIM TO DO IS WHAT THEY ACTUALLY DO

21/02/08
program sum
  integer, parameter :: N=5000, ntimes=20000
  real :: x(N), y(N)
  real :: begin, end
  real :: rtime(2)
  common/saver/x, y
  !DIR CACHE_ALIGN
  call random_number(x)
  call random_number(y)
  begin = dtime(rtime)
  do it=1, ntimes
    do i=1, N
      y(i) = x(i) + i
    end do
    if (mod(it, 1000).eq.0) print *, x(1), y(1)
  end do
  end = dtime(rtime)
  print *, ' my loop time (s) is ', end
  flop = (N*ntimes)/end*1.0e-6
  print *, ' loop runs at ', flop, ' MFLOP'
end program sum

loop runs at 68.96552 MFLOP
Loop Unrolling

- Loop unrolling is an optimization technique which can be applied to loops which perform calculations on array elements.
- Consists of replicating the body of the loop so that calculations are performed on several array elements during each iteration.
- Reason for unrolling is to take advantage of pipelined functional units. Consecutive elements of the arrays can be in the functional unit simultaneously.
- Programmer usually does not have to unroll loops “by hand” -- compiler options and directives are usually available. Performing unrolling via directives and/or options is preferable.
  - Code is more portable to other systems
  - Code is self-documenting and easier to read
- Loops with small trip counts or data dependencies should not be unrolled!
Loop Unrolling (con’t)

Advantages:
- Can be done automatically by compiler.
- Makes use of pipelined functional units.
- Makes use of multiple pipelined functional units on superscalar processors.
- Compiler may be able to pre-load array operands into registers, hiding load latencies.

Disadvantages:
- Limited number of FP registers
  - Pentium III: 8
  - Pentium 4: 8
  - Athlon: 8
  - Opteron: 4
  - Itanium: 128(!)
Loop Unrolling Example

- Normal loop

  do i=1,N
    a(i)=b(i)+x*c(i)
  enddo

- Manually unrolled loop

  do i=1,N,4
    a(i)=b(i)+x*c(i)
    a(i+1)=b(i+1)+x*c(i+1)
    a(i+2)=b(i+2)+x*c(i+2)
    a(i+3)=b(i+3)+x*c(i+3)
  enddo
Loop Unrolling Compiler Options

**GNU compilers:**
- `-funroll-loops` Enable loop unrolling
- `-funroll-all-loops` Unroll all loops; not recommended

**PGI compilers:**
- `-Munroll` Enable loop unrolling
- `-Munroll=c:N` Unroll loops with trip counts of at least \( N \)
- `-Munroll=n:M` Unroll loops up to \( M \) times

**Intel compilers:**
- `-unroll` Enable loop unrolling
- `-unrollM` Unroll loops up to \( M \) times

TEST IF WHAT THEY CLAIM TO DO IS WHAT THEY ACTUALLY DO
program dirunroll
  integer, parameter :: N=1000000
  real, dimension (N) :: a, b, c
  real :: begin, end
  real, dimension (2) :: rtime
  common/saver/a, b, c
  call random_number(b)
  call random_number(c)
  x=2.5
  begin=dtime(rtime)
!DIR$ UNROLL 4
  do i=1,N
    a(i)=b(i)+x*c(i)
  end do
  end=dtime(rtime)
  print *, 'my loop time (s) is ', (end)
  flop=(2.0*N)/(end)*1.0e-6
  print *, 'loop runs at ', flop, ' MFLOP'
  print *, a(1), b(1), c(1)
end
Loop Fusion and Fission

Fusion: Merge multiple loops into one

Fission: Split one loop into multiple loops
Example of Loop Fusion

Potential for Fusion: dependent operations in separate loops

Advantage:
• Re-usage of array B()

Disadvantages:
• In total 4 arrays now contend for cache space
• More registers needed
Example of Loop Fission

Potential for Fission: independent operations in a single loop

**Advantage:**
- First loop can be scheduled more efficiently and be parallelised as well

**Disadvantages:**
- Less opportunity for out-of-order superscalar execution
- Additional loop created (a minor disadvantage)
Floating Point Division

- Floating point division is an expensive operation
  - Takes 22-60 CPs to complete (average about 32 CPs)
  - Usually not pipelined
  - According to the IEEE floating point standard, divisions **must** be carried out as such and not replaced with a multiplication by a reciprocal (even for division by a constant!)

- A possible optimization technique is to “relax” the IEEE requirements and replace a division with multiplication by a reciprocal. Most compilers do this automatically at higher levels of optimization.
Floating Point Division Compiler
Options

GNU:
-funsafe-math-optimizations
Allow optimizations for floating-point arithmetic that (a) assume that arguments and results are valid and (b) may violate IEEE or ANSI standards.

PGI:
--Kieee -Knoieee (default)
Perform floating-point operations in strict conformance with the IEEE 754 standard. Some optimizations are disabled with -Kieee, and a more accurate math library is used. The default -Knoieee uses faster but very slightly less accurate methods.

INTEL:
--no-prec-div (i32 and i32em)
Enables optimizations that give slightly less precise results than full IEEE division. With some optimizations, such as -xN and -xB, the compiler may change floating-point division computations into multiplication by the reciprocal of the denominator.
Floating Point Division Example

• Consider the following piece of code
  
  ```fortran
  do i=1,N
    a(i) = a(i)/x
  enddo
  ```

• On using the Intel ifort compiler, the compiler will automatically compute \((1.0/x)\) once outside the loop and replace the division in the loop with a multiplication by this value. The new multiplication loop will then be unrolled. The resulting performance is \(~93\) MFLOPS.
  
  - If the user forces IEEE conformance (ife -mp), the performance is reduced to 56 MFLOPS.
Floating Point Division With Arrays

Consider the following loop nest in which the array $A(i,j)$ is scaled by different factors stored in array $B(i)$:

\[
\begin{align*}
\text{do } & j=1,N \\
& \quad \text{do } i=1,N \\
& \quad \quad A(i,j) = A(i,j)/B(i) \\
& \quad \text{enddo} \\
& \text{enddo}
\end{align*}
\]

The compiler can do no automatic optimization to this, because $B(i)$ is not a scalar. However, you can manually do the following:

- Create a temporary array to hold the inverses of the $B(i)$ array.
- Replace the division in the inner loop with multiplication by the temporary array.
- The resulting code can be unrolled and/or pipelined.
Optimization based on Microprocessor Architectures

- Pipelined Functional Units
- Superscalar Processors
  - Processors which have multiple functional units are said to be *superscalar*.

- Instruction Set Extensions
  - Newer processors have additional instructions beyond the usual floating point add and multiply instructions:
    - SSE2/SSE3/3DNow! Etc...
    - Cat /proc/cpuinfo..
Pipelined Functional Units

- For the processors in most modern parallel machines, the circuitry on the chip which performs a given type of operation on operands in registers is known as a *functional unit*.

- Most integer and floating point functional units are *pipelined*, meaning that they can have multiple independent executions of the same instruction placed in a queue. The idea is that after an initial startup latency, the functional unit should be able to generate one result every clock period (CP).

- Each stage of a pipelined operation can be working simultaneously on different sets of operands.
Superscalar Processors

- Processors which have multiple functional units are said to be *superscalar*.
- Examples:
  - Intel Pentium 3
    - 1 Floating point unit
    - 1 MMX/SSE unit
    - 2 Integer units
    - 2 Load/store units
  - Intel Pentium 4
    - 2 Floating point
    - 1/MMX/SSE units
    - 2 Integer units
    - 2 Load/store units
Pipelining Compiler Options

GNU:
  none

PGI:
  --Mvect[]=option[,option,...] -Mnovect (default)
    Pass options to the internal vectorizer.

INTEL:
  none
Instruction Set Extensions

Newer processors have additional instructions beyond the usual floating point add and multiply instructions:

- **Intel MMX (Matrix Math eXtensions):**
  - introduced in 1997 supported by all current processors

- **Intel SSE (Streaming SIMD Extensions):**
  - introduced on the Pentium III in 1999
  - useless for scientific computation: single precision

- **Intel SSE2 (Streaming SIMD Extensions 2):**
  - introduced on the Pentium 4 in Dec 2000
Instruction Set Extensions (2)

- AMD 3DNow!:
  - introduced in 1998 (extends MMX)

- AMD 3DNow!+ (or 3DNow! Professional, or 3DNow! Athlon):
  - introduced with the Athlon (includes SSE)

- To check what you have on your machine:
  - cat /proc/cpuinfo
Instruction Set Extension Compiler

Options

GNU:
- mmmx/no-mmx
- -msse
- mno-sse
- msse2 / -mno-sse2
- msse3 / -mno-sse3
- m3dnow / -mno-3dnow

These switches enable or disable the use of built-in functions that allow direct access to the MMX, SSE, SSE2, SSE3 and 3Dnow extensions of the instruction set.

PGI:
--fastsse

Chooses generally optimal flags for a processor that supports SSE instructions (Pentium 3/4, AthlonXP/MP, Opteron) and SSE2 (Pentium 4, Opteron). Use pgf90 -fastsse -help to see the equivalent switches.

INTEL:
- arch SSE
  Optimizes for Intel Pentium 4 processors with Streaming SIMD Extensions (SSE).
- arch SSE2
  Optimizes for Intel Pentium 4 processors with Streaming SIMD Extensions 2 (SSE2).
General techniques

- Blocking/tiling
- !! Use of optimized libraries !!
Blocking for cache (tiling)

Blocking for cache is:

- An optimization that applies for datasets that do not entirely fit in the (2nd level) data cache
- A way to increase spatial locality of reference i.e. exploit full cache lines
- A way to increase temporal locality of reference i.e. Improves data re-usage

By way of example, let discuss the transpose of a matrix...

```fortran
do i=1,n
    do j=1,n
        a(i,j)=b(j,i)
    end do
end do
```
Block algorithm for transposing a matrix:

- block data size = bsize
  - mb = n/bsize
  - nb = n/bsize
- Code is a little bit more complicated if
  - MOD(n, bsize) is not zero
  - MOD(m, bsize) is not zero
Results... (Carlo Cavazzoni data)
Optimizing Matrix Multiply for Caches

• Several techniques for making this faster on modern processors
  – heavily studied
• Some optimizations done automatically by compiler, but can do much better
• In general, you should use optimized libraries (often supplied by vendor) for this and other very common linear algebra operations
  – BLAS = Basic Linear Algebra Subroutines
• Other algorithms you may want are not going to be supplied by vendor, so need to know these techniques
Summary

- Performance programming on uniprocessors requires
  - understanding of memory system
    - levels, costs, sizes
  - understanding of fine-grained parallelism in processor to produce good instruction mix
  - understanding your program

- Compilers are good at instruction level optimization and loop transformation

- User is responsible to present code in most natural way for compiler optimizations.

- The techniques work for any architecture, but choosing details depends on the architecture

- Blocking (tiling) is a basic approach that can be applied to many matrix algorithms
A long time ago a couple of guys...

Production runs '97