

Supercomputing in Plain English Stupid Compiler Tricks

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Tuesday February 12 2013







This is an experiment!

It's the nature of these kinds of videoconferences that FAILURES ARE GUARANTEED TO HAPPEN! NO PROMISES!

- So, please bear with us. Hopefully everything will work out well enough.
- If you lose your connection, you can retry the same kind of connection, or try connecting another way.
- Remember, if all else fails, you always have the toll free phone bridge to fall back on.







H.323 (Polycom etc) #1

- If you want to use H.323 videoconferencing for example, Polycom – then:
- If you AREN'T registered with the OneNet gatekeeper (which is probably the case), then:
 - Dial 164.58.250.47
 - Bring up the virtual keypad.
 - On some H.323 devices, you can bring up the virtual keypad by typing: #
 - (You may want to try without first, then with; some devices won't work with the #, but give cryptic error messages about it.)
 - When asked for the conference ID, or if there's no response, enter:
 0409
 - On most but not all H.323 devices, you indicate the end of the ID with:
 #







H.323 (Polycom etc) #2

If you want to use H.323 videoconferencing – for example, Polycom – then:

 If you ARE already registered with the OneNet gatekeeper (most institutions aren't), dial:

2500409

Many thanks to Skyler Donahue and Steven Haldeman of OneNet for providing this.







Wowza #1

You can watch from a Windows, MacOS or Linux laptop using Wowza from either of the following URLs:

http://www.onenet.net/technical-resources/video/sipe-stream/

OR

https://vcenter.njvid.net/videos/livestreams/page1/

Wowza behaves a lot like YouTube, except live.

Many thanks to Skyler Donahue and Steven Haldeman of OneNet and Bob Gerdes of Rutgers U for providing this.







Wowza #2

Wowza has been tested on multiple browsers on each of:

- Windows (7 and 8): IE, Firefox, Chrome, Opera, Safari
- MacOS X: Safari, Firefox
- Linux: Firefox, Opera

We've also successfully tested it on devices with:

- Android
- iOS

However, we make no representations on the likelihood of it working on your device, because we don't know which versions of Android or iOS it might or might not work with.









If one of the Wowza URLs fails, try switching over to the other one.

If we lose our network connection between OU and OneNet, then there may be a slight delay while we set up a direct connection to Rutgers.







Toll Free Phone Bridge

IF ALL ELSE FAILS, you can use our toll free phone bridge: 800-832-0736 * 623 2847 #

Please mute yourself and use the phone to listen.

Don't worry, we'll call out slide numbers as we go.

- Please use the phone bridge <u>ONLY</u> if you cannot connect any other way: the phone bridge can handle only 100 simultaneous connections, and we have over 350 participants.
- Many thanks to OU CIO Loretta Early for providing the toll free phone bridge.







No matter how you connect, please mute yourself, so that we cannot hear you.

- (For Wowza, you don't need to do that, because the information only goes from us to you, not from you to us.)
- At OU, we will turn off the sound on all conferencing technologies.
- That way, we won't have problems with echo cancellation.
- Of course, that means we cannot hear questions.
- So for questions, you'll need to send e-mail.







Questions via E-mail Only

Ask questions by sending e-mail to:

sipe2013@gmail.com

All questions will be read out loud and then answered out loud.







TENTATIVE Schedule

Tue Jan 29: Compilers: What the Heck is Supercomputing? Tue Jan 29: The Tyranny of the Storage Hierarchy Tue Feb 12: Instruction Level Parallelism Tue Feb 12: Stupid Compiler Tricks Tue Feb 19: Shared Memory Multithreading Tue Feb 26: Distributed Multiprocessing Tue March 5: Applications and Types of Parallelism Tue March 12: Multicore Madness Tue March 19: NO SESSION (OU's Spring Break) Tue March 26: High Throughput Computing Tue Apr 2: GPGPU: Number Crunching in Your Graphics Card Tue Apr 9: Grab Bag: Scientific Libraries, I/O Libraries, Visualization





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Supercomputing Exercises #1

Want to do the "Supercomputing in Plain English" exercises?

• The 3rd exercise will be posted soon at:

http://www.oscer.ou.edu/education/

 If you don't yet have a supercomputer account, you can get a temporary account, just for the "Supercomputing in Plain English" exercises, by sending e-mail to:

hneeman@ou.edu

Please note that this account is for doing the <u>exercises only</u>, and will be shut down at the end of the series. It's also available only to those at institutions in the USA.

 This week's Introductory exercise will teach you how to compile and run jobs on OU's big Linux cluster supercomputer, which is named Boomer.







Supercomputing Exercises #2

You'll be doing the exercises on your own (or you can work with others at your local institution if you like).

These aren't graded, but we're available for questions:

hneeman@ou.edu





Thanks for helping!

- OU IT
 - OSCER operations staff (Brandon George, Dave Akin, Brett Zimmerman, Josh Alexander, Patrick Calhoun)
 - Horst Severini, OSCER Associate Director for Remote & Heterogeneous Computing
 - Debi Gentis, OU Research IT coordinator
 - Kevin Blake, OU IT (videographer)
 - Chris Kobza, OU IT (learning technologies)
 - Mark McAvoy
- Kyle Keys, OU National Weather Center
- James Deaton, Skyler Donahue and Steven Haldeman, OneNet
- Bob Gerdes, Rutgers U
- Lisa Ison, U Kentucky
- Paul Dave, U Chicago







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From Computational Biophysics to Systems Biology, May 19-21, Norman OK

Great Plains Network Annual Meeting, May 29-31, Kansas City XSEDE2013, July 22-25, San Diego CA
IEEE Cluster 2013, Sep 23-27, Indianapolis IN
OKLAHOMA SUPERCOMPUTING SYMPOSIUM 2013, Oct 1-2, Norman OK
SC13, Nov 17-22, Denver CO





OK Supercomputing Symposium 2013



2003 Keynote: Peter Freeman NSF Computer & Information Science & Engineering Assistant Director



2004 Keynote: Sangtae Kim NSF Shared Cyberinfrastructure **Division Director**



2005 Keynote: Walt Brooks NASA Advanced Supercomputing **Division Director**



2006 Keynote: Dan Atkins Head of NSF's Office of



2007 Keynote: Jay Boisseau Director Texas Advanced Cyberinfrastructure Computing Center U. Texas Austin



2008 Keynote: José Munoz **Deputy Office** Director/ Senior Scientific Advisor NSF Office of Cyberinfrastructure



2009 Keynote: **Douglass Post Chief Scientist** US Dept of Defense **HPC** Modernization Program



2010 Keynote: Horst Simon **Deputy Director** Lawrence Berkeley National Science National Laboratory

2011 Keynote: **Barry Schneider** Program Manager Foundation

2012 Keynote:

Thom Dunning

Director

National Center for

Supercomputing

Applications



2013 Keynote to be announced! **FREE! Wed Oct 2 2013 @ OU** http://symposium2013.oscer.ou.edu/

Reception/Poster Session Tue Oct 1 2013 @ OU Symposium Wed Oct 2 2013 @ OU



Supercomputing in Plain English: Compilers Tue Feb 12 2013



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Outline

- Dependency Analysis
 - What is Dependency Analysis?
 - Control Dependencies
 - Data Dependencies
- Stupid Compiler Tricks
 - Tricks the Compiler Plays
 - Tricks You Play With the Compiler
 - Profiling





Dependency Analysis



What Is Dependency Analysis?

- **Dependency analysis** describes of how different parts of a program affect one another, and how various parts require other parts in order to operate correctly.
- A *control dependency* governs how different sequences of instructions affect each other.
- A *data dependency* governs how different pieces of data affect each other.

Much of this discussion is from references [1] and [6].







Control Dependencies

- Every program has a well-defined *flow of control* that moves from instruction to instruction to instruction.
- This flow can be affected by several kinds of operations:
 - Loops
 - Branches (if, select case/switch)
 - Function/subroutine calls
 - I/O (typically implemented as calls)

Dependencies affect **parallelization**!







Branch Dependency (F90)

 $\mathbf{y} = \mathbf{7}$

IF (x /= 0) THEN

y = 1.0 / x

END IF

Note that (x /= 0) means "x not equal to zero."

The value of \mathbf{y} depends on what the condition ($\mathbf{x} /= 0$) evaluates to:

- If the condition (x /= 0) evaluates to .TRUE., then y is set to 1.0 / x. (1 divided by x).
- Otherwise, **y** remains **7**.







Branch Dependency (C)

y = 7; if (x != 0) { y = 1.0 / x; }

Note that (x != 0) means "x not equal to zero."

The value of y depends on what the condition (x != 0) evaluates to:

- If the condition (x != 0) evaluates to true, then
 y is set to 1.0 / x (1 divided by x).
- Otherwise, **y** remains **7**.





Loop Carried Dependency (F90)

DO i = 2, length
 a(i) = a(i-1) + b(i)

END DO

Here, each iteration of the loop depends on the previous:
 iteration i=3 depends on iteration i=2,
 iteration i=4 depends on iteration i=3,
 iteration i=5 depends on iteration i=4, etc.

This is sometimes called a *loop carried dependency*.

There is no way to execute iteration i until after iteration i-1 has completed, so this loop can't be parallelized.





Loop Carried Dependency (C)

```
for (i = 1; i < length; i++) {
    a[i] = a[i-1] + b[i];
}</pre>
```

Here, each iteration of the loop depends on the previous:
 iteration i=3 depends on iteration i=2,
 iteration i=4 depends on iteration i=3,
 iteration i=5 depends on iteration i=4, etc.

This is sometimes called a *loop carried dependency*.

There is no way to execute iteration i until after iteration i-1 has completed, so this loop can't be parallelized.









Loops are the favorite control structures of High Performance Computing, because compilers know how to <u>optimize</u> their performance using instruction-level parallelism: superscalar, pipelining and vectorization can give excellent speedup.

Loop carried dependencies affect whether a loop can be parallelized, and how much.





Loop or Branch Dependency? (F)

Is this a <u>loop carried dependency</u> or a <u>branch dependency</u>?

```
DO i = 1, length
    IF (x(i) /= 0) THEN
        y(i) = 1.0 / x(i)
    END IF
END DO
```





Loop or Branch Dependency? (C)

Is this a **loop carried dependency** or a **branch dependency**?

```
for (i = 0; i < length; i++) {
    if (x[i] != 0) {
        y[i] = 1.0 / x[i];
    }
}</pre>
```







Call Dependency Example (F90)

- $\mathbf{x} = 5$
- y = myfunction(7)
- z = 22

The flow of the program is interrupted by the <u>call</u> to **myfunction**, which takes the execution to somewhere else in the program.

It's similar to a branch dependency.







Call Dependency Example (C)

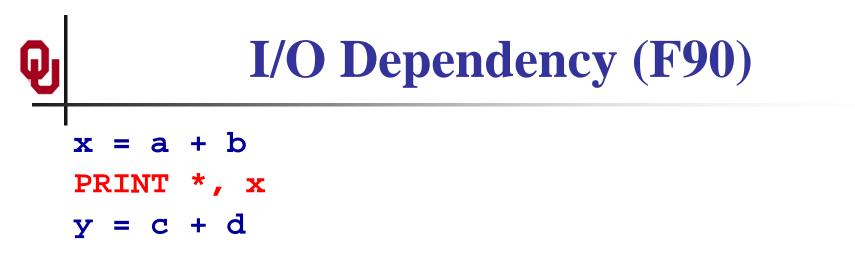
- x = 5;
- y = myfunction(7);
- z = 22;

The flow of the program is interrupted by the <u>call</u> to **myfunction**, which takes the execution to somewhere else in the program.

It's similar to a branch dependency.







Typically, I/O is implemented by hidden subroutine calls, so we can think of this as equivalent to a call dependency.







I/O Dependency (C)

x = a + b; printf("%f", x); y = c + d;

Typically, I/O is implemented by hidden subroutine calls, so we can think of this as equivalent to a call dependency.





Reductions Aren't Dependencies

```
array_sum = 0
DO i = 1, length
array_sum = array_sum + array(i)
END DO
```

A *reduction* is an operation that converts an array to a scalar.

- Other kinds of reductions: product, **.AND.**, **.OR.**, minimum, maximum, index of minimum, index of maximum, number of occurrences of a particular value, etc.
- Reductions are so common that hardware and compilers are optimized to handle them.
- Also, they aren't really dependencies, because the order in which the individual operations are performed doesn't matter.





Reductions Aren't Dependencies

```
array_sum = 0;
for (i = 0; i < length; i++) {
  array_sum = array_sum + array[i];
}
```

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- Other kinds of reductions: product, **&&**, **|**, minimum, maximum, index of minimum, index of maximum, number of occurrences of a particular value, etc.
- Reductions are so common that hardware and compilers are optimized to handle them.
- Also, they aren't really dependencies, because the order in which the individual operations are performed doesn't matter.







Data Dependencies (F90)

- "A data dependence occurs when an instruction is dependent on data from a previous instruction and therefore cannot be moved before the earlier instruction [or executed in parallel]."^[7]
- $\mathbf{a} = \mathbf{x} + \mathbf{y} + \cos(\mathbf{z})$
- b = a * c

The value of **b** depends on the value of **a**, so these two statements <u>**must**</u> be executed in order.





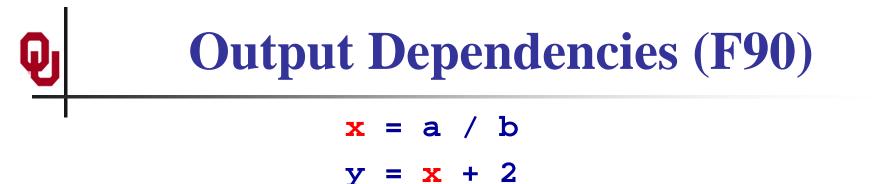


Data Dependencies (C)

- "A data dependence occurs when an instruction is dependent on data from a previous instruction and therefore cannot be moved before the earlier instruction [or executed in parallel]."^[7]
- a = x + y + cos(z);
- b = a * c;
- The value of **b** depends on the value of **a**, so these two statements <u>**must**</u> be executed in order.







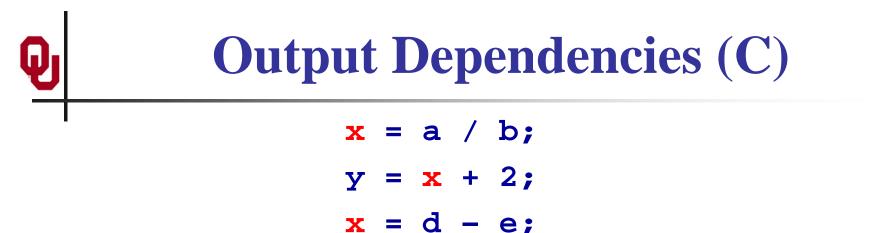
 $\mathbf{x} = \mathbf{d} - \mathbf{e}$

Notice that \mathbf{x} is assigned <u>two different values</u>, but only one of them is retained after these statements are done executing. In this context, the final value of \mathbf{x} is the "output."

Again, we are forced to execute in order.







Notice that \mathbf{x} is assigned <u>two different values</u>, but only one of them is retained after these statements are done executing. In this context, the final value of \mathbf{x} is the "output."

Again, we are forced to execute in order.







Why Does Order Matter?

- Dependencies can affect whether we can execute a particular part of the program in <u>parallel</u>.
- If we cannot execute that part of the program in parallel, then it'll be <u>SLOW</u>.





Loop Dependency Example

```
if ((dst == src1) && (dst == src2))
  for (index = 1; index < length; index++) {</pre>
    dst[index] = dst[index-1] + dst[index];
else if (dst == src1) {
  for (index = 1; index < length; index++) {
    dst[index] = dst[index-1] + src2[index];
else if (dst == src2) {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = src1[index-1] + dst[index];
else if (src1 == src2) {
  for (index = 1; index < length; index++) {
    dst[index = src1[index-1] + src1[index];
élse {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = src1[index-1] + src2[index];
                  Supercomputing in Plain English: Compilers
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```





Loop Dep Example (cont'd)

```
if ((dst == src1) && (dst == src2)) {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = dst[index-1] + dst[index];
else if (dst == src1) {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = dst[index-1] + src2[index];
else if (dst == src2) {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = src1[index-1] + dst[index];
else if (src1 == src2) {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = src1[index-1] + src1[index];
élse {
  for (index = 1; index < length; index++) {</pre>
    dst[index] = src1[index-1] + src2[index];
}
```

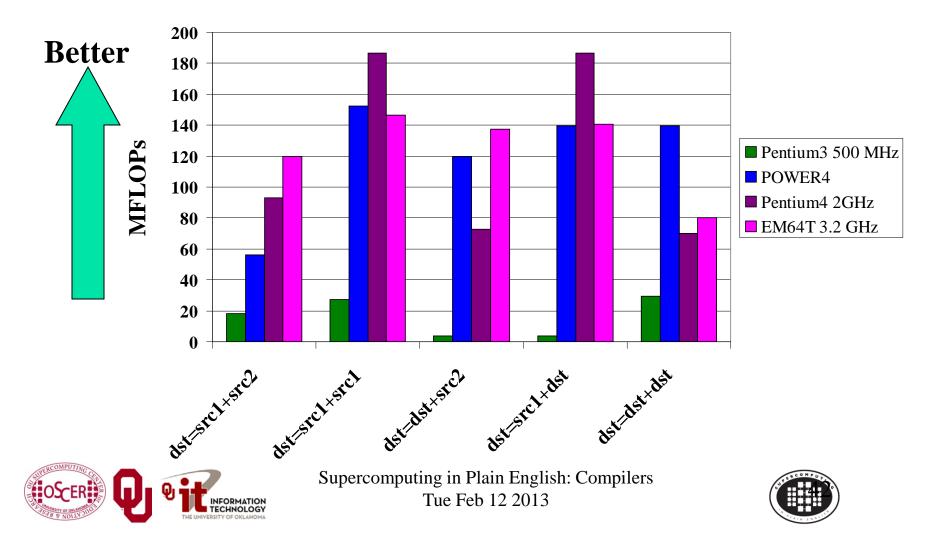
The various versions of the loop either:
do have loop carried dependencies, or
don't have loop carried dependencies.





Loop Dependency Performance

Loop Carried Dependency Performance





Stupid Compiler Tricks



Stupid Compiler Tricks

- Tricks Compilers Play
 - Scalar Optimizations
 - Loop Optimizations
 - Inlining
- Tricks You Can Play with Compilers
 - Profiling
 - Hardware counters









The people who design compilers have a lot of experience working with the languages commonly used in High Performance Computing:

- Fortran: 50+ years
- C: 40+ years
- C++: 25+ years, plus C experience

So, they've come up with clever ways to make programs run faster.





Q Tricks Compilers Play



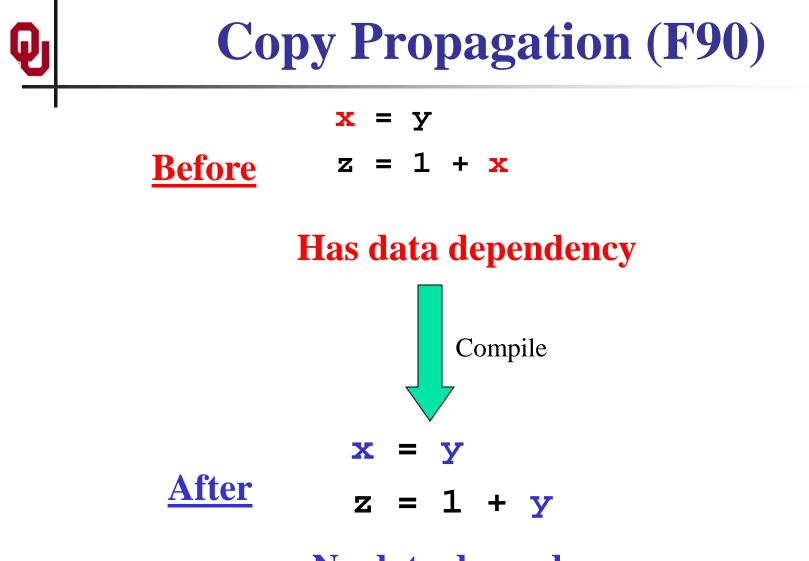
Scalar Optimizations

- Copy Propagation
- Constant Folding
- Dead Code Removal
- Strength Reduction
- Common Subexpression Elimination
- Variable Renaming
- Loop Optimizations
- Not every compiler does all of these, so it sometimes can be worth doing these by hand.

Much of this discussion is from [2] and [6].







No data dependency



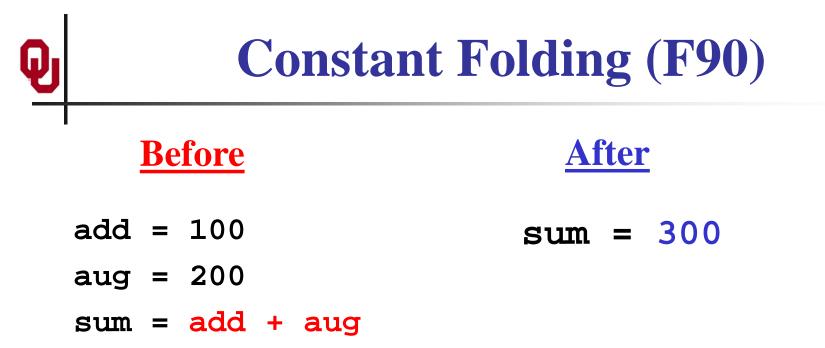


Copy Propagation (C) $\mathbf{x} = \mathbf{y};$ z = 1 + x;**Before** Has data dependency Compile X = y; After \mathbf{Z} + y;

No data dependency



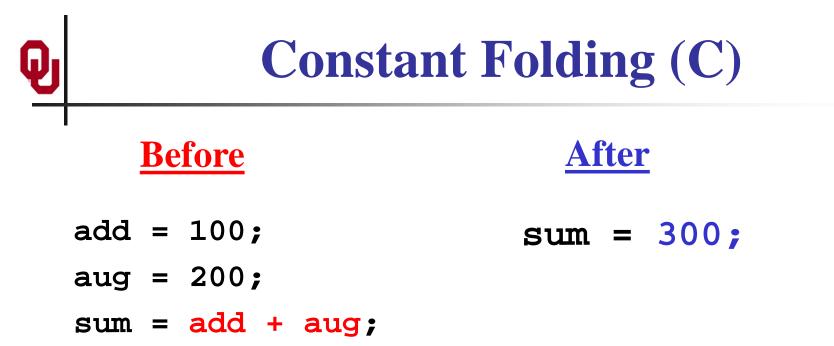




Notice that **sum** is actually the sum of two constants, so the compiler can precalculate it, eliminating the addition that otherwise would be performed at runtime.







Notice that **sum** is actually the sum of two constants, so the compiler can precalculate it, eliminating the addition that otherwise would be performed at runtime.





Q	Dead	Code Removal (F90)
	Before	After
var =	5	var = 5
PRINT	*, var	PRINT *, var
STOP		STOP
PRINT	*, var *	2

Since the last statement never executes, the compiler can eliminate it.



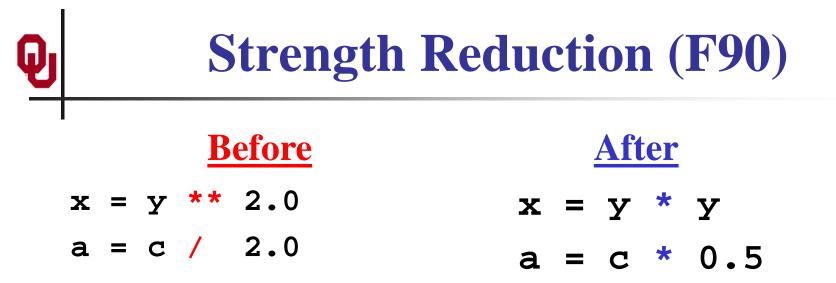


Q	Dead Code	Removal (C)	
	Before	After	
var	= 5;	var = 5;	
pri	.ntf("%d", var);	<pre>printf("%d", var);</pre>	
exi	t(-1);	exit(-1);	
pri	nf("%d", var * 2)	•	

Since the last statement never executes, the compiler can eliminate it.





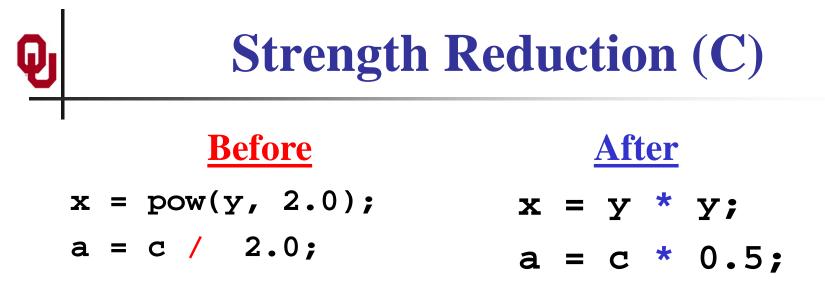


Raising one value to the power of another, or dividing, is more expensive than multiplying. If the compiler can tell that the power is a small integer, or that the denominator is a constant, it'll use multiplication instead.

Note: In Fortran, "y **** 2.0**" means "y to the power 2."







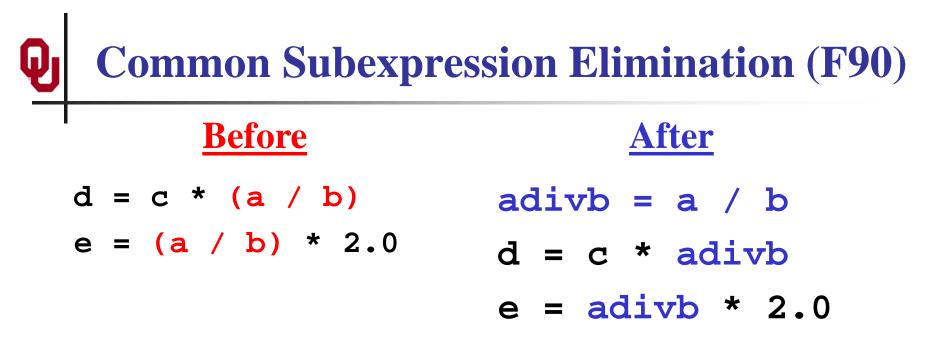
Raising one value to the power of another, or dividing, is more expensive than multiplying. If the compiler can tell that the power is a small integer, or that the denominator is a constant, it'll use multiplication instead.

Note: In C, "pow(y, 2.0)" means "y to the power 2."





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The subexpression (a / b) occurs in both assignment statements, so there's no point in calculating it twice.

This is typically only worth doing if the common subexpression is expensive to calculate.





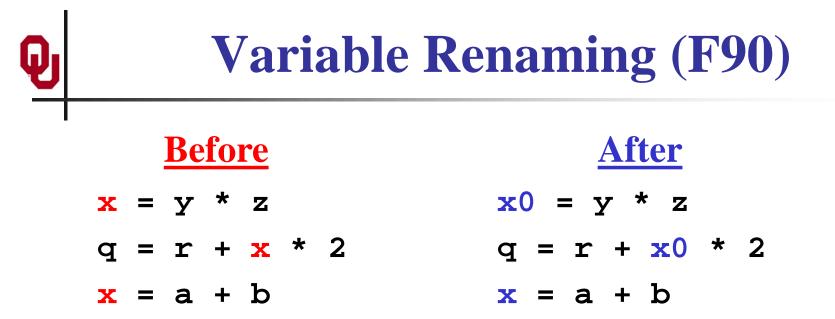
Common Subexpression Elimination (C) <u>Before</u> d = c * (a / b); e = (a / b) * 2.0; <u>After</u> adivb = a / b; d = c * adivb; e = adivb * 2.0;

The subexpression (a / b) occurs in both assignment statements, so there's no point in calculating it twice.

This is typically only worth doing if the common subexpression is expensive to calculate.



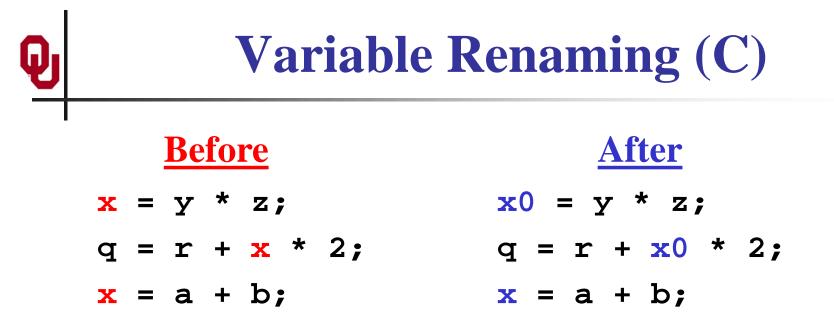




The original code has an <u>output dependency</u>, while the new code <u>doesn't</u> – but the final value of \mathbf{x} is still correct.







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Loop Optimizations

- Hoisting Loop Invariant Code
- Unswitching
- Iteration Peeling
- Index Set Splitting
- Loop Interchange
- Unrolling
- Loop Fusion
- Loop Fission
- Not every compiler does all of these, so it sometimes can be worth doing some of these by hand.

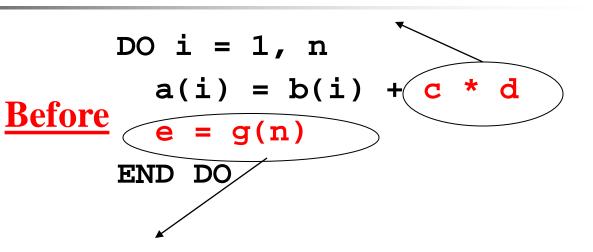
Much of this discussion is from [3] and [6].





Hoisting Loop Invariant Code (F90)

Code that doesn't change inside the loop is known as *loop invariant*. It doesn't need to be calculated over and over.



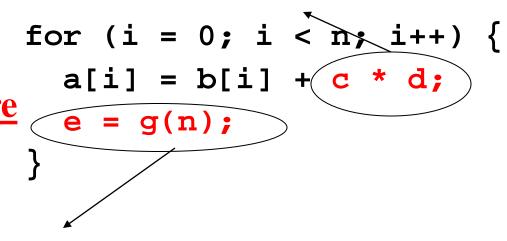
	temp = c * d
	DO $i = 1$, n
Δ fter	a(i) = b(i) + temp
	END DO
	e = g(n)





Hoisting Loop Invariant Code (C)

Code that doesn't change inside the loop is known as *loop invariant*. It doesn't need to be calculated over and over.



temp = c * d; for (i = 0; i < n; i++) { a[i] = b[i] + temp; } e = g(n);







Unswitching (F90)

```
The condition is
DO i = 1, n
  DO j = 2, n
                                         j-independent.
    IF (t(i) > 0) THEN
      a(i,j) = a(i,j) * t(i) + b(j)
    ELSE
      a(i,j) = 0.0
                                             Before
    END IF
  END DO
END DO
DO i = 1, n
                                      So, it can migrate
  IF (t(i) > 0) THEN
    DO j = 2, n
      a(i,j) = a(i,j) * t(i) + b(j) outside the j loop.
    END DO
  ELSE
    DO j = 2, n
a(i,j) = 0.0
                                              After
    END DO
  END IF
END DO
```





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Unswitching (C)

```
for (i = 0; i < n; i++)</pre>
                                                      The condition is
  for (j = 1; j < n; j++) {
    if (t[i] > 0)
                                                      j-independent.
        a[i][j] = a[i][j] * t[i] + b[j];
     }
     else {
                                                          Before
        a[i][j] = 0.0;
for (i = 0; i < n; i++) {</pre>
  if(t[i] > 0)
                                                    So, it can migrate
     for (j = 1; j < n; j++) {
    a[i][j] = a[i][j] * t[i] + b[j];</pre>
                                                    outside the j loop.
   }
                                                            After
  else {
     for (j = 1; j < n; j++) {
    a[i][j] = 0.0;</pre>
   }
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                                                                          64
```

Iteration Peeling (F90)

I	DO i = 1, n
	IF ((i == 1) .OR. (i == n)) THEN
	x(i) = y(i)
DC	ELSE
<u>Before</u>	x(i) = y(i + 1) + y(i - 1)
	END IF
	END DO

We can eliminate the IF by *peeling* the weird iterations.

	x(1) = y(1)
	DO $i = 2, n - 1$
<u>After</u>	x(i) = y(i + 1) + y(i - 1)
	END DO
	x(n) = y(n)
~ 100	





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Iteration Peeling (C)

We can eliminate the IF by *peeling* the weird iterations.







Index Set Splitting (F90)

```
DO i = 1, n

a(i) = b(i) + c(i)

IF (i > 10) THEN

d(i) = a(i) + b(i - 10)

END IF

END DO
```

```
DO i = 1, 10
    a(i) = b(i) + c(i)
END DO
DO i = 11, n
    a(i) = b(i) + c(i)
    d(i) = a(i) + b(i - 10)
END DO
```

Note that this is a generalization of **<u>peeling</u>**.



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Before

After



Index Set Splitting (C)

```
for (i = 0; i < n; i++) {</pre>
  a[i] = b[i] + c[i];
  if (i >= 10) {
    d[i] = a[i] + b[i - 10];
                                       Before
for (i = 0; i < 10; i++) {
  a[i] = b[i] + c[i];
for (i = 10; i < n; i++) {</pre>
  a[i] = b[i] + c[i];
                                        After
 d[i] = a[i] + b[i - 10];
```

Note that this is a generalization of **<u>peeling</u>**.







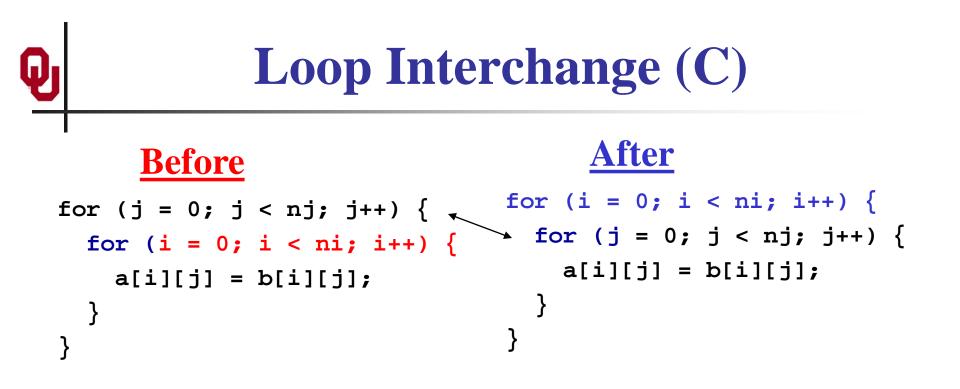
Loop Interchange (F90)

BeforeAfterDO i = 1, niDO j = 1, njDO j = 1, njDO i = 1, njDO i = 1, nia(i,j) = b(i,j)END DOEND DOEND DOEND DO

Array elements **a(i,j)** and **a(i+1,j)** are near each other in memory, while **a(i,j+1)** may be far, so it makes sense to make the **i** loop be the inner loop. (This is reversed in C, C++ and Java.)







Array elements **a**[**i**][**j**] and **a**[**i**][**j**+**1**] are near each other in memory, while **a**[**i**+**1**][**j**] may be far, so it makes sense to make the **j** loop be the inner loop. (This is reversed in Fortran.)







Unrolling (F90)

DO i = 1, n
Before
$$a(i) = a(i)+b(i)$$

END DO
DO i = 1, n, 4

DO 1 = 1, n, 4 a(i) = a(i) + b(i) a(i+1) = a(i+1) + b(i+1) a(i+2) = a(i+2) + b(i+2) a(i+3) = a(i+3) + b(i+3)END DO

You generally **shouldn't** unroll by hand.







Unrolling (C)

for (i = 0; i < n; i++) {
 Before a[i] = a[i] + b[i];
 }</pre>

for (i = 0; i < n; i += 4) {
 a[i] = a[i] + b[i];
 a[i+1] = a[i+1] + b[i+1];
 a[i+2] = a[i+2] + b[i+2];
 a[i+3] = a[i+3] + b[i+3];
}</pre>

You generally shouldn't unroll by hand.





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Why Do Compilers Unroll?

- We saw last time that a loop with a lot of operations gets better performance (up to some point), especially if there are lots of arithmetic operations but few main memory loads and stores.
- Unrolling creates multiple operations that typically load from the same, or adjacent, cache lines.
- So, an unrolled loop has more operations without increasing the memory accesses by much.
- Also, unrolling decreases the number of comparisons on the loop counter variable, and the number of branches to the top of the loop.







Loop Fusion (F90)

```
DO i = 1, n
  a(i) = b(i) + 1
END DO
DO i = 1, n
  c(i) = a(i) / 2
END DO
                             Before
DO i = 1, n
  d(i) = 1 / c(i)
END DO
DO i = 1, n
  a(i) = b(i) + 1
  c(i) = a(i) / 2
                              After
  d(i) = 1 / c(i)
END DO
```

As with unrolling, this has fewer branches. It also has fewer total memory references.







Loop Fusion (C)

```
for (i = 0; i < n; i++) {</pre>
  a[i] = b[i] + 1;
for (i = 0; i < n; i++) {</pre>
  c[i] = a[i] / 2;
                                Before
for (i = 0; i < n; i++) {</pre>
  d[i] = 1 / c[i];
for (i = 0; i < n; i++) {</pre>
  a[i] = b[i] + 1;
  c[i] = a[i] / 2;
                                After
  d[i] = 1 / c[i];
```

As with unrolling, this has fewer branches. It also has fewer total memory references.







Loop Fission (F90)

```
DO i = 1, n
  a(i) = b(i) + 1
  c(i) = a(i) / 2
                             Before
  d(i) = 1 / c(i)
END DO
DO i = 1, n
  a(i) = b(i) + 1
END DO
DO i = 1, n
  c(i) = a(i) / 2
END DO
                              After
DO i = 1, n
  d(i) = 1 / c(i)
END DO
```

Fission reduces the cache footprint and the number of operations per iteration.







Loop Fission (C)



Fission reduces the cache footprint and the number of operations per iteration.







To Fuse or to Fizz?

The question of when to perform fusion versus when to perform fission, like many many optimization questions, is highly dependent on the application, the platform and a lot of other issues that get very, very complicated.

Compilers don't always make the right choices.

That's why it's important to examine the actual behavior of the executable.







Inlining (F90)

Before	<u>After</u>
DO i = 1, n a(i) = func(i)	DO i = 1, n a(i) = i * 3
END DO	END DO
REAL FUNCTION func (x)	
func = x * 3	
END FUNCTION func	

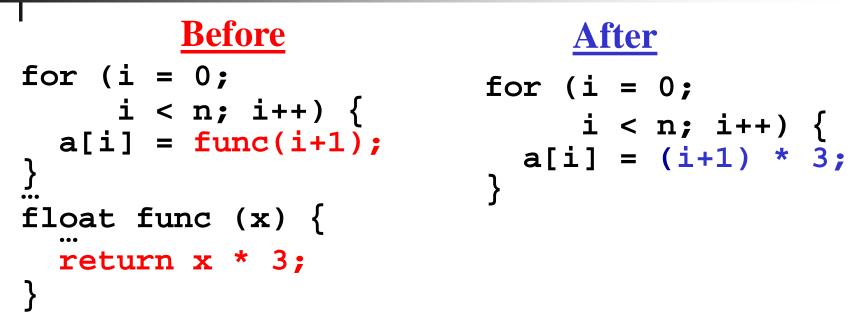
When a function or subroutine is *inlined*, its contents are transferred directly into the calling routine, eliminating the overhead of making the call.







Inlining (C)



When a function or subroutine is *inlined*, its contents are transferred directly into the calling routine, eliminating the overhead of making the call.





Tricks You Can Play with Compilers

Q



The Joy of Compiler Options

Every compiler has a different set of options that you can set. Among these are options that control single processor optimization: superscalar, pipelining, vectorization, scalar optimizations, loop optimizations, inlining and so on.







Example Compile Lines

 IBM XL xlf90 -0 -qmaxmem=-1 -qarch=auto -qtune=auto -qcache=auto -qhot
 Intel ifort -0 -march=corei7-avx -xAVX -xhost
 Portland Group f90 pgf90 -03 -tp=sandybridge
 NAG f95

nagfor -04 -Ounsafe





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What Does the Compiler Do? #1

Example: NAG nagfor compiler [4]

nagfor -O<level> source.f90

Possible levels are -00, -01, -02, -03, -04:

- -00 No optimisation. ...
- -01 Minimal quick optimisation.
- -02 Normal optimisation.
- -03 Further optimisation.
- -04 Maximal optimisation.

The man page is pretty cryptic.





What Does the Compiler Do? #2

Example: Intel **ifort** compiler ^[5]

ifort -O<level> source.f90

Possible levels are -00, -01, -02, -03:

-00 Disables all -O<n> optimizations. ...

-01 ... [E]nables optimizations for speed. ...

-02

Inlining of intrinsics.

Intra-file interprocedural optimizations, which include: inlining, constant propagation, forward substitution, routine attribute propagation, variable address-taken analysis, dead static function elimination, and removal of unreferenced variables.

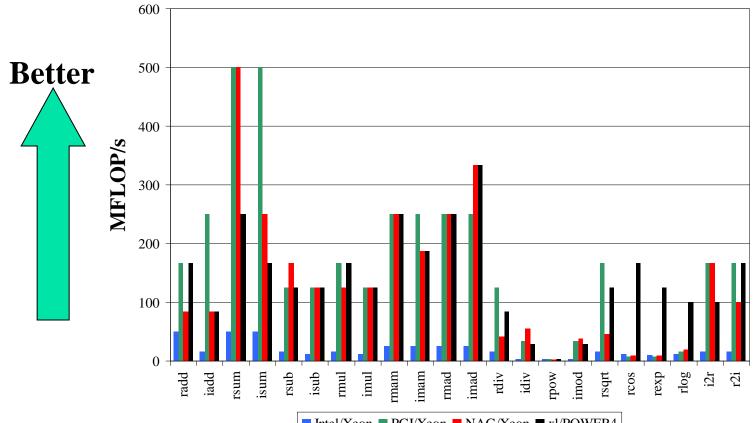
-03 Enables -02 optimizations plus more aggressive optimizations, such as prefetching, scalar replacement, and loop transformations. Enables optimizations for maximum speed, but does not guarantee higher performance unless loop and memory access transformations take place. ...





Arithmetic Operation Speeds

Ordered Arithmetic Operations

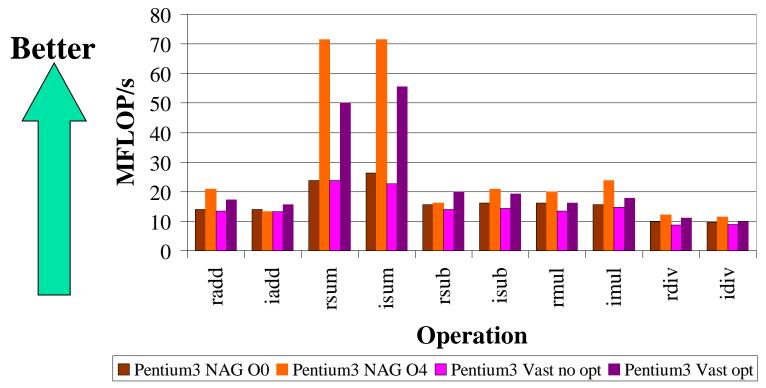


■ Intel/Xeon ■ PGI/Xeon ■ NAG/Xeon ■ xl/POWER4





Optimization Performance



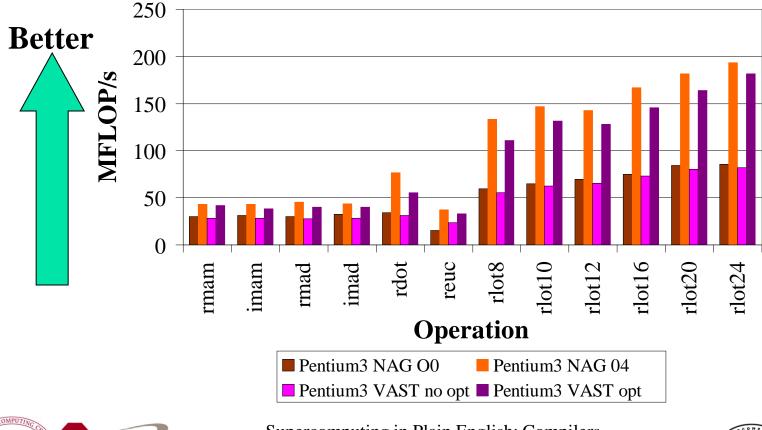






More Optimized Performance

Performance









Profiling



Profiling

Profiling means collecting data about how a program executes. The two major kinds of profiling are:

- Subroutine profiling
- Hardware timing







Subroutine Profiling

Subroutine profiling means finding out how much time is spent in each routine.

- <u>**The 90-10 Rule</u>**: Typically, a program spends 90% of its runtime in 10% of the code.</u>
- Subroutine profiling tells you what parts of the program to spend time optimizing and what parts you can ignore.
- Specifically, at regular intervals (e.g., every millisecond), the program takes note of what instruction it's currently on.







Profiling Example

On GNU compilers systems:

gcc -0 -g -pg ...

The **-g -pg** options tell the compiler to set the executable up to collect profiling information.

Running the executable generates a file named **gmon.out**, which contains the profiling information.







Profiling Example (cont'd)

When the run has completed, a file named **gmon.out** has been generated.

Then:

gprof executable

produces a list of all of the routines and how much time was spent in each.







Profiling Result

%	cumulative	self		self	total	
tim	ne seconds	seconds	calls	ms/call	ms/call	name
27.	6 52.72	52.72	480000	0.11	0.11	longwave_ [5]
24.	3 99.06	46.35	897	51.67	51.67	mpdata3_ [8]
7.	9 114.19	15.13	300	50.43	50.43	turb_ [9]
7.	2 127.94	13.75	299	45.98	45.98	turb_scalar_ [10]
4.	7 136.91	8.96	300	29.88	29.88	advect2_z_ [12]
4.	1 144.79	7.88	300	26.27	31.52	cloud_ [11]
3.	9 152.22	7.43	300	24.77	212.36	radiation_ [3]
2.	3 156.65	4.43	897	4.94	56.61	smlr_ [7]
2.	2 160.77	4.12	300	13.73	24.39	tke_full_ [13]
1.	7 163.97	3.20	300	10.66	10.66	shear_prod_ [15]
1.	5 166.79	2.82	300	9.40	9.40	rhs_ [16]
1.	4 169.53	2.74	300	9.13	9.13	advect2_xy_ [17]
1.	3 172.00	2.47	300	8.23	15.33	poisson_ [14]
1.	2 174.27	2.27	480000	0.00	0.12	long_wave_ [4]
1.	0 176.13	1.86	299	6.22	177.45	<pre>advect_scalar_ [6]</pre>
0.	9 177.94	1.81	300	6.04	6.04	buoy_ [19]

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Thanks for your attention!



Questions? www.oscer.ou.edu



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