

COMPILER OPTIMIZATIONS

What's New for Intel compilers 19.1?

Advance Support for Intel[®] Architecture – Use Intel compiler to generate optimized code for Intel Atom[®] processor through Intel[®] Xeon[®] Scalable processor and Intel[®] Xeon Phi[™] processor families

Achieve Superior Parallel Performance – Vectorize & thread your code (using OpenMP*) to take full advantage of the latest SIMD-enabled hardware, including Intel[®] Advanced Vector Extensions 512 (Intel[®] AVX-512)

What's New in C++

Initial C++20, and full C++ 17 enabled

- Enjoy advanced lambda and constant expression support
- Standards-driven parallelization for C++ developers
 Initial OpenMP* 5.0, and full OpenMP* 4.5
 support
- Modernize your code by using the latest parallelization specifications

What's New in Fortran

Substantial Fortran 2018 support

- Enjoy enhanced C-interoperability features for effective mixed language development
- Use advanced coarray features to parallelize your modern Fortran code

Initial OpenMP* 5.0, and substantial OpenMP* 4.5 support

• Customize your reduction operations by user-defined reductions



Common optimization options

	Linux*
Disable optimization	-00
Optimize for speed (no code size increase)	-01
Optimize for speed (default)	-02
High-level loop optimization	-03
Create symbols for debugging	-g
Multi-file inter-procedural optimization	-ipo
Profile guided optimization (multi-step build)	-prof-gen -prof-use
Optimize for speed across the entire program ("prototype switch") <i>fast</i> options definitions changes over time!	-fast same as: -ipo –O3 -no-prec-div –static –fp-model fast=2 -xHost)
OpenMP support	-qopenmp
Automatic parallelization	-parallel



High-Level Optimizations

Basic Optimizations with icc -O...

- -O0 no optimization; sets -g for debugging
- -O1 scalar optimizations excludes optimizations tending to increase code size
- -O2 default for icc/icpc (except with -g)

includes **auto-vectorization**; some loop transformations, e.g. unrolling, loop interchange; inlining within source file; start with this (after initial debugging at -OO)

-O3 more aggressive loop optimizations including cache blocking, loop fusion, prefetching, ... suited to applications with loops that do many floating-point calculations or process large data sets



InterProcedural Optimizations (IPO)

Multi-pass Optimization

icc -ipo

Analysis and optimization across function and/or source file boundaries, e.g.

• Function inlining; constant propagation; dependency analysis; data & code layout; etc.

2-step process:

- Compile phase objects contain intermediate representation
- "Link" phase compile and optimize over all such objects
- Seamless: linker automatically detects objects built with -ipo and their compile options
- May increase build-time and binary size
- But build can be parallelized with -ipo=n
- Entire program need not be built with IPO, just hot modules

Particularly effective for applications with many smaller functions

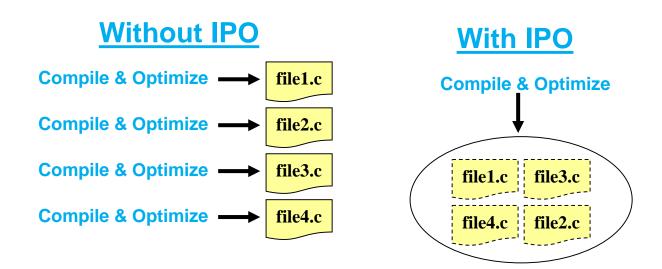
Get report on inlined functions with -qopt-report-phase=ipo



InterProcedural Optimizations

Extends optimizations across file boundaries

-ip	Only between modules of one source file
-ipo	Modules of multiple files/whole application





Profile-Guided Optimizations (PGO)

Static analysis leaves many questions open for the optimizer like:

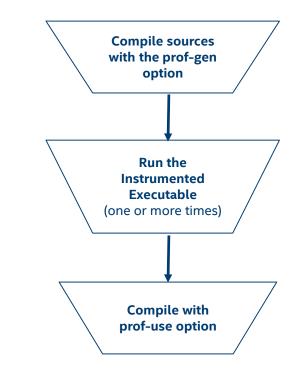
- How often is x > y
- What is the size of count
- Which code is touched how often

if (x > y)do this(); else do that();

for(i=0; i<count; ++i)
do_work();</pre>

Use execution-time feedback to guide (final) optimization Enhancements with PGO:

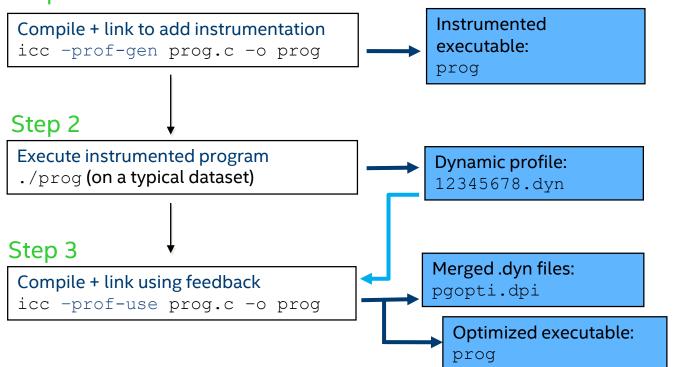
- More accurate branch prediction
- Basic block movement to improve instruction cache behavior
- Better decision of functions to inline (help IPO)
- Can optimize function ordering
- Switch-statement optimization
- Better vectorization decisions





PGO Usage: Three-Step Process

Step 1





Math Libraries

icc comes with Intel's optimized math libraries

- libimf (scalar) and libsvml (scalar & vector)
- Faster than GNU* libm
- Driver links libimf automatically, ahead of libm
- Additional functions (replace math.h by mathimf.h)

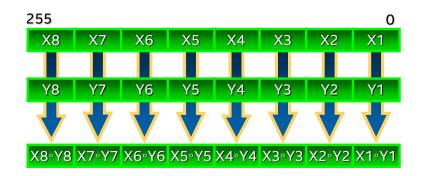
Don't link to libm explicitly!



- May give you the slower libm functions instead
- Though the Intel driver may try to prevent this
- gcc needs -lm, so it is often found in old makefiles



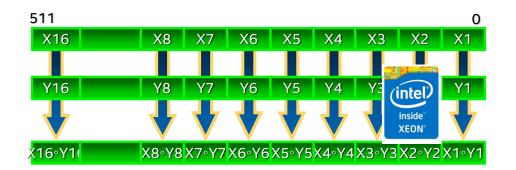
SIMD Types for Intel® Architecture



AVX

Vector size: **256 bit** Data types:

- 8, 16, 32, 64 bit integer
- 32 and 64 bit float VL: 4, 8, 16, 32



Intel® AVX-512 Vector size: 512 bit Data types:

- 8, 16, 32, 64 bit integer
- 32 and 64 bit float

VL: 8, 16, 32, 64

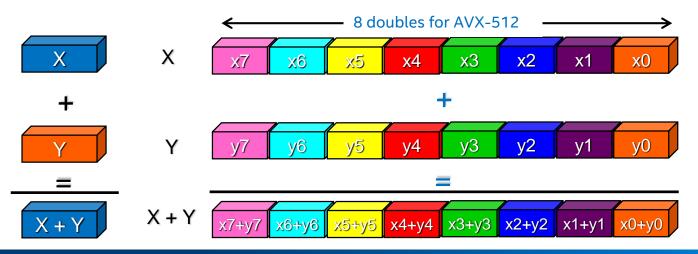


SIMD: <u>Single Instruction</u>, <u>Multiple Data</u>

for (i=0; i<n; i++) z[i] = x[i] + y[i];</pre>

- Scalar mode
 - one instruction produces one result
 - E.g. vaddss, (vaddsd)

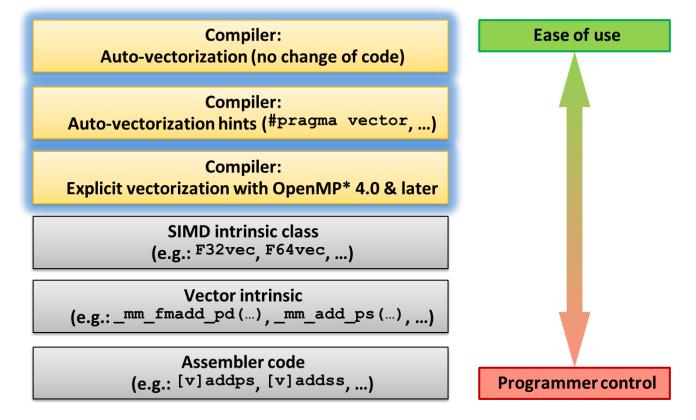
- □ Vector (SIMD) mode
 - one instruction can produce multiple results
 - E.g. vaddps, (vaddpd)



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Many ways to vectorize



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Basic Vectorization Switches I

- Requirement: -O2 or higher
- -x<feature>
 - Might enable Intel processor specific optimizations
 - Processor-check added to "main" routine: Application errors in case SIMD feature missing or non-Intel processor with appropriate/informative message
- -ax<features>
 - Multiple code paths: baseline and optimized/processor-specific
 - Optimized code paths for Intel processors defined by <features>
 - Multiple SIMD features/paths possible, e.g.: -axSSE2,AVX,AVX2
 - Baseline code path defaults to -msse2
 - The baseline code path can be modified by <u>-m<feature></u> or <u>-x<feature></u>

Basic Vectorization Switches II

- -m<feature>
 - Neither check nor specific optimizations for Intel processors:
 Application optimized for both Intel and non-Intel processors for selected SIMD feature
- Default for Linux: -msse2
 - Activated implicitly
 - Implies the need for a target processor with at least Intel[®] SSE2
- Special switch: -xHost
 - Compiler checks SIMD features of current host processor (where built on) and makes use of latest SIMD feature available
 - Code only executes on processors with same SIMD feature or later as on build host
 - As for -x<feature>, if "main" routine is built with -xHost or the final executable only runs on Intel processors



Compiler Reports – Optimization Report

- -qopt-report[=n]: tells the compiler to generate an optimization report
 - n: (Optional) Indicates the level of detail in the report. You can specify values 0 through 5. If you specify zero, no report is generated. For levels n=1 through n=5, each level includes all the information of the previous level, as well as potentially some additional information. Level 5 produces the greatest level of detail. If you do not specify n, the default is level 2, which produces a medium level of detail.
- -qopt-report-phase[=list]: specifies one or more optimizer phases for which optimization reports are generated.
 - loop: the phase for loop nest optimization
 - vec: the phase for vectorization
 - par: the phase for auto-parallelization
 - all: all optimizer phases
- -qopt-report-filter=string: specified the indicated parts of your application, and generate optimization reports for those parts of your application.

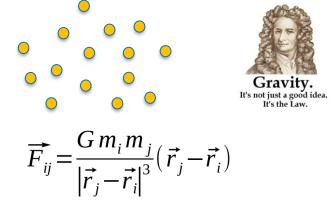


Lab: Nbody gravity simulation

Let's consider a distribution of point masses located at $r_1, ..., r_n$ and have masses $m_1, ..., m_n$

We want to calculate the position of the particles after a certain time interval using the Newton law of gravity

```
struct Particle
{
    public:
        Particle() { init();}
        void init()
        {
            pos[0] = 0.; pos[1] = 0.; pos[2] = 0.;
            vel[0] = 0.; vel[1] = 0.; vel[2] = 0.;
            acc[0] = 0.; acc[1] = 0.; acc[2] = 0.;
            mass = 0.;
        }
        real_type pos[3];
        real_type vel[3];
        real_type mass;
};
```



$$\vec{F} = m\vec{a} = m\frac{d\vec{v}}{dt} = m\frac{d^2\vec{x}}{dt^2}$$

Lab: Nbody kernel implementation

```
GSimulation.cpp:
```

```
. . .
for (i = 0; i < n; i++) {
                                                             // update acceleration
    for (j = 0; j < n; j++) {
          real type distance, dx, dy, dz;
          real type distanceSqr = 0.0;
          real type distanceInv = 0.0;
          dx = particles[j].pos[0] - particles[i].pos[0];
          dy = particles[j].pos[1] - particles[i].pos[1];
          dz = particles[j].pos[2] - particles[i].pos[2];
          distSqr = dx^*dx + dy^*dy + dz^*dz + softeningSquared;
          distInv = 1.0 / sqrt(distanceSqr);
          particles[i].acc[0] += dx * G * particles[j].mass * distInv * distInv * distInv;
          particles[i].acc[1] += ...
          particles[i].acc[2] += ...
```

. . .



- module load intel64/19.1up01
- Go to the folder nbody-demo/ver0
 - Take a look on the code to learn data structures, main loops, Flops calculation, etc.
- Type make to compile code

icpc -g -std=c++11 -02 -o nbody.x GSimulation.o main.o

- Type make run to run the test and measure the timing.
 - ./nbody.x
- If everything works fine, you should see the following results:



Run the default test case on CPU: ./nbody.x _____ Initialize Gravity Simulation nPart = 16000; nSteps = 10; dt = 0.1 dt kenergy time (s) GFlops S 0.1 26.405 5.1986 1.4281 1 0.2 313.77 5.2053 2 1.4263 3 0.3 926.56 5.3105 1.398 4 0.4 1866.4 5.2772 1.4069 5 0.5 3135.6 5.1825 1,4326 6 0.6 4737.6 5.1845 1.432 7 0.7 6676.6 5.1921 1.4299 0.8 8957.7 5.2037 1.4267 8 9 0.9 11587 5.2009 1.4275 10 14572 5.1825 1.4326 1 # Number Threads : 1 # Total Time (s) : 52.139 # Average Performance : 1.4233 +- 0.012399

- Observe how the performance changes according to the number of particles: ./nbody.x 8000
- Is the perfomance saturating?
- What is the maximum performance?
- Has the code performance issues?

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Run the default test case on CPU: ./nbody.x						
Initialize Gravity Simulation nPart = 16000; nSteps = 10; dt = 0.1						
s	dt	kenergy	time (s)	GFlops		
1 2 3 4 5 6 7 8 9 10	0.2 0.3 0.4 0.5 0.6 0.7 0.8	313.77 926.56 1866.4 3135.6 4737.6 6676.6 8957.7 11587	5.1986 5.2053 5.3105 5.2772 5.1825 5.1845 5.1921 5.2037 5.2009 5.1825	1.4263 1.398 1.4069 1.4326 1.432 1.4299		
<pre># Number Threads : 1 # Total Time (s) : 52.139 # Average Performance : 1.4233 +- 0.012399</pre>						

- Observe how the performance changes according to the number of particles: ./nbody.x 8000
- Is the perfomance saturating?
- What is the maximum performance?
- Has the code performance issues?
- Generate the compiler report: make REPORT=yes

or

icpc -g -std=c++11 -02 -qoptreport=5 -o nbody.x GSimulation.cpp main.cpp

LOOP BEGIN at GSimulation.cpp(127,20) remark #15542: loop was not vectorized: inner loop was already vectorized

```
LOOP BEGIN at GSimulation.cpp(130,5)
remark #15542: loop was not vectorized: inner
loop was already vectorized
```

.

. . . .

```
remark #15417: vectorization support: number of FP up
converts: single precision to double precision 1 [
GSimulation.cpp(143,4) ]
```

```
remark #15300: LOOP WAS VECTORIZED
    remark #15452: unmasked strided loads: 6
    remark #15475: --- begin vector cost summary
---
    remark #15476: scalar cost: 145
    remark #15477: vector cost: 79.500
    remark #15478: estimated potential speedup:
1.820
    remark #15487: type converts: 23
```

- Observe how verbose is the report looking at the file: GSimulation.optrpt
 - 2974 number of lines
- You can filter the report by running: make REPORT=yes FILTER=yes
 -qopt-report-phase=vec
 -qopt-report filter="GSimulation.cpp,125-175"
- Look at the Makefile for more details on the compiler options used
- Open the new report file and read the results

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- How can we improve the performance using only the compiler?
 - Use different compiler options and try to target the underlying architecture:
 - -xCORE-AVX2 or -xHost
 - -O3
 - -ipo
 - -prof-gen and -prof-use
 - -parallel
- Explain why some options don't bring additional speed-up. Try some more tests here: git clone <u>https://github.com/ivorobts/compiler-optimization.git</u>

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