



Erlangen Regional
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A HPM based Methodology for SIMD Code Analysis

HPCC High Performance
Computing

Instrument hot spot and measure HPM counters for 1 process running on one core with fixed frequency.

1. Ensure code is limited by instruction throughput
2. Measure executed instruction counts for the following categories:
 - FP arithmetic scalar
 - FP arithmetic SIMD (different flavors)
 - Load and store
 - Branch
 - Total
3. Measure CPI/IPC to quantify execution efficiency
4. Compute relations between counts

Do this for scalar and all SIMD variants you are interested in.
(to disable autovectorization with Intel compilers use **-no-vec**)

1. Percentages of **category instruction counts** compared to total number of instructions
2. Factor of **total and arithmetic instruction count** for SIMD variants vs scalar baseline

Outcome

- For a code where algorithmic work consists of mostly floating point computations the **instruction decomposition** should be dominated by **arithmetic floating point instructions**.
- The **SIMD factor**, means the number of lanes in a SIMD register should be visible for overall as well as arithmetic instruction counts. SIMD saves instructions but also often requires to add new instructions to enable SIMD processing.

Compare **CPI** and **runtime** for all SIMD variants vs. scalar baseline.

Outcome

- CPI quantifies superscalar execution. Instruction count times CPI gives runtime in cycles.
- The runtime factor should be equal to the **#instructions X CPI**

- Algorithmic and SIMD benefits can be quantified on an **instruction level**
- It can be quantified how well the SIMD vectorization works separated for **instruction level** and **execution side**.
- Overheads can be named, most effective together with static assembly code analysis:
 - Overhead added getting data from memory into registers
 - Overhead added by compiler for implementing programming language or programming model
 - Identify potential to save instructions

Proxy app from **MANTEVO**.

<https://mantevo.org>



- **MiniMD** – MPI + OpenMP
 - App. 3000 lines, C++
 - Extracted from LAMMPS (<http://lammps.sandia.gov>)
Molecular dynamics application code
 - Other implementations: OpenACC, OpenCL, Kokkos
 - Supports Lennard-Jones (default) and eam force calculation
 - Uses neighbour lists

- Load modules for compiler and MPI
- Edit `include_ICC.mk` for compilation options and flags

- Compile with: `$ make`
- Clean up with: `$ make clean`

- You have to `chdir` to the `./data` directory to run the code:
`$ cd ./data`

- Run the benchmark by calling:
`$../miniMD-ICC --half_neigh <0|1> -n 200`