
Solution for Project 1

Contents

1. Rosa Warm-Up	<i>(5 Points)</i>	2
1.1. exercise 1		2
1.2. exercise 2		2
1.3. exercise 3		3
1.4. exercise 4		3
1.5. exercise 5		4
2. Performance Characteristics	<i>(30 Points)</i>	4
2.1. Peak performance		4
2.2. Memory Hierarchies		5
2.3. Bandwidth: STREAM benchmark		6
2.4. Performance model: A simple roofline model		6
3. Optimize Square Matrix-Matrix Multiplication	<i>(50 Points)</i>	6
3.1. Motivation		6
3.2. Importance of data access optimization		6
3.3. Optimizing square matrix multiplication		6
4. Quality of the Report	<i>(15 Points)</i>	6

1. Rosa Warm-Up

(5 Points)

1.1. exercise 1

The **module system** is a utility that allows the user to dynamically manage their software environment on the Rosa HPC cluster and to load different compilers, libraries and applications in order to modify environment variables (like PATH, LD_LIBRARY_PATH, MANPATH, etc.) without creating conflicts between different software versions or dependencies.

As reported in the USI resource page the module system provides several commands to manage the environment.

- `module avail` – lists all available modules (on the current system)
- `module list` – lists all currently loaded modules
- `module show` – display information about
- `module load` – loads module
- `module switch` – unloads, loads
- `module rm` – unloads module
- `module purge` – unloads all loaded modules

1.2. exercise 2

The **Slurm** (Simple Linux Utility for Resource Management) is a job scheduler for Linux clusters. Main features of Slurm are:

- Job scheduling and resource management
- Framework for starting, executing, and monitoring work (jobs) on a set of allocated nodes
- Queuing management to handle multiple users and jobs

The two main components are:

- *slurmd*: the daemon that runs on each compute node responsible for launching, monitoring, and terminating jobs
- *slurmctld*: the central management daemon that manages job queues and allocates resources

Main commands:

- `srun`: submit a job for execution
- `sbatch`: submit a batch job
- `squeue`: view the status of jobs in the queue
- `scancel`: cancel a job
- `salloc`: allocate resources for an interactive job

1.3. exercise 3

Here below is a simple program in C that prints "Hello World" and the information about the system where it is executed.

```
1 #include <stdio.h>
2 #include <unistd.h>
3
4 int main(void) {
5     char hostname[256];
6     gethostname(hostname, sizeof(hostname));
7     printf("Hello - World - from - host : -%s\n", hostname);
8 }
```

Listing 1: Hello World C Program - *src/1-Rosa-warm-up/hello_world.c*

We can process the script with the following command:

```
srunk -N1 --time=00:01:00 ./hello_worldc > hello_worldc.out 2> hello_worldc.err
```

and we can see from the output that the program has been correctly compiled and executed on the cluster on node `icsnode22` from the *slim* partition.

```
Hello World from host: icsnode22
```

1.4. exercise 4

We can see the output of the command `sinfo` here below:

PARTITION	AVAIL	TIMELIMIT	NODES	STATE	NODELIST
slim*	up	2-00:00:00	7	mix	icsnode[22,27-28,32-35]
slim*	up	2-00:00:00	12	alloc	icsnode[17-21,23-26,29-31]
slim*	up	2-00:00:00	2	idle	icsnode[36-37]
lr-slim	up	30-00:00:0	1	idle	icsnode38
gpu	up	2-00:00:00	8	mix	icsnode[05-06,08-13]
gpu	up	2-00:00:00	2	idle	icsnode[14-15]
fat	up	2-00:00:00	4	idle	icsnode[01-04]
bigMem	up	2-00:00:00	2	idle	icsnode[07,15]
debug-slim	up	4:00:00	1	idle	icsnode39
debug-gpu	up	4:00:00	1	idle	icsnode16
multi-gpu	up	2-00:00:00	2	idle	icsnode[41-42]

Listing 2: Output of line command `sinfo`

As we can see nodes are divided in partitions with names that already give us some information about their characteristics. As explained in the `sbatch` guide on `slurm` website, we can use different flags on the `sbatch` command to specify the partition to use with different commands.

The flag that applies to our case is:

```
sbatch --partition=fat job_script.sh
```

Submit with `bigMem` partition to run on nodes with very large memory:

```
sbatch --partition=bigMem job_script.sh
```

Similarly, for GPU partitions:

```
sbatch --partition=gpu job_script.sh
```

For example, we can modify the script file `slurm_job_one.sh` to specify the partition with the `gpu` partition with the following line:

```
#SBATCH --partition=gpu
```

After inserting the line into the file and reprocessing the script, we can see the following result:

```

Loading gcc/13.2.0-gcc-8.5.0-5hghkwo
Loading requirement: gcc-runtime/8.5.0-gcc-8.5.0-7fyorqa
                    gmp/6.2.1-gcc-8.5.0-lrpevy5  mpfr/4.2.1-gcc-8.5.0-ybeybcx
                    mpc/1.3.1-gcc-8.5.0-cv2gjfw  zlib-ng/2.1.6-gcc-8.5.0-ztbc5xt
                    zstd/1.5.5-gcc-8.5.0-azepnn7
Currently Loaded Modulefiles:
 1) gcc-runtime/8.5.0-gcc-8.5.0-7fyorqa    5) zlib-ng/2.1.6-gcc-8.5.0-ztbc5xt
 2) gmp/6.2.1-gcc-8.5.0-lrpevy5            6) zstd/1.5.5-gcc-8.5.0-azepnn7
 3) mpfr/4.2.1-gcc-8.5.0-ybeybcx          7) gcc/13.2.0-gcc-8.5.0-5hghkwo
 4) mpc/1.3.1-gcc-8.5.0-cv2gjfw
Model name:          Intel(R) Xeon(R) CPU E5-2650 v3 @ 2.30GHz
Hello World from host: icsnode08

```

Listing 3: Output of the job script `slurm_job_one.sh` after specifying the partition

From the very last line we can assert that the job has been submitted to node `icsnode08`, and from the `sinfo` output before (2) we can see that this node belongs to the *gpu* partition instead of the *slim* partition as before (1).

1.5. exercise 5

In order to run our program on two nodes is sufficient to add the following line of our script:

```
#SBATCH --nodes=2                # Number of nodes
```

This line is already implemented in the script `slurm_job_two.sh` provided in the `src` folder. Once we process the script with the command we get the the message Submitted batch job 51103 and after a while we can check the output file `slurm_job_two-51103.out` (4) to see the result of our job.

```

Loading gcc/13.2.0-gcc-8.5.0-5hghkwo
Loading requirement: gcc-runtime/8.5.0-gcc-8.5.0-7fyorqa
                    gmp/6.2.1-gcc-8.5.0-lrpevy5  mpfr/4.2.1-gcc-8.5.0-ybeybcx
                    mpc/1.3.1-gcc-8.5.0-cv2gjfw  zlib-ng/2.1.6-gcc-8.5.0-ztbc5xt
                    zstd/1.5.5-gcc-8.5.0-azepnn7
Currently Loaded Modulefiles:
 1) gcc-runtime/8.5.0-gcc-8.5.0-7fyorqa    5) zlib-ng/2.1.6-gcc-8.5.0-ztbc5xt
 2) gmp/6.2.1-gcc-8.5.0-lrpevy5            6) zstd/1.5.5-gcc-8.5.0-azepnn7
 3) mpfr/4.2.1-gcc-8.5.0-ybeybcx          7) gcc/13.2.0-gcc-8.5.0-5hghkwo
 4) mpc/1.3.1-gcc-8.5.0-cv2gjfw
Model name:          Intel(R) Xeon(R) CPU E5-2650 v3 @ 2.30GHz
Hello World from host: icsnode22
Hello World from host: icsnode21

```

Listing 4: Output of the job script `slurm_job_two.sh`

From the output we can see that the job has been submitted to two different nodes `icsnode22` and `icsnode21` confirming that the command has been correctly processed twice on two different nodes.

2. Performance Characteristics

(30 Points)

2.1. Peak performance

Rosa's nodes are equipped with dual-socket Intel Xeon E5-2650 v3 processors, 10 cores and AVX2 units that operates with 256-bit vectors[1, 3]. Haswell micro-architecture have two 256-bit FMA instructions per cycle meaning that each core performs four double-precision at 2.30 GHz [2].

The Rosa documentation (and also the output from command `sinfo 2`) shows 42 compute nodes (icsnode01–icsnode42).[3]

The calculations for the aggregate peak throughput for the partition is written in the provided file [INTEL_XEON_E5-2650.txt] in the 02 folder and reported here below:

Intel Xeon E5–2650 v3 @ 2.30GHz:

$$\begin{aligned}
 P_{\{core\}} &= 4 \text{ lanes} \times 2 \text{ flop} \times 2 \text{ FMA} \times 2.30 \text{ GHz} = 36.8 \text{ GFlops/s} \\
 P_{\{CPU\}} &= 10 \text{ cores} \times P_{\{core\}} = 368 \text{ GFlops/s} \\
 P_{\{node\}} &= 2 \text{ sockets} \times P_{\{CPU\}} = 736 \text{ GFlops/s} \\
 P_{\{EVII\}} &= 42 \text{ nodes} \times P_{\{node\}} = 30,912 \text{ GFlops/s} \\
 &= 30.912 \text{ TFlops/s}
 \end{aligned}$$

Listing 5: Peak throughput breakdown for Intel Xeon E5-2650 v3

2.2. Memory Hierarchies

In order to study the memory hierarchy of Rosa compute nodes, I used the commands: `lscpu`, `cat /proc/meminfo`, and `hwloc-ls`. The complete output files are available in the submission in the folder 02.

The results are summarized in Table 1.

Component	Size
Main memory (total)	62 GB
Main memory (per NUMA node)	31 GB
L3 cache (shared per socket)	25 MB
L2 cache (per core)	256 KB
L1d cache (per core)	32 KB
L1i cache (per core)	32 KB

Table 1: Memory hierarchy of Rosa compute node

The graphical representation of the memory hierarchy is shown in Figure 1.

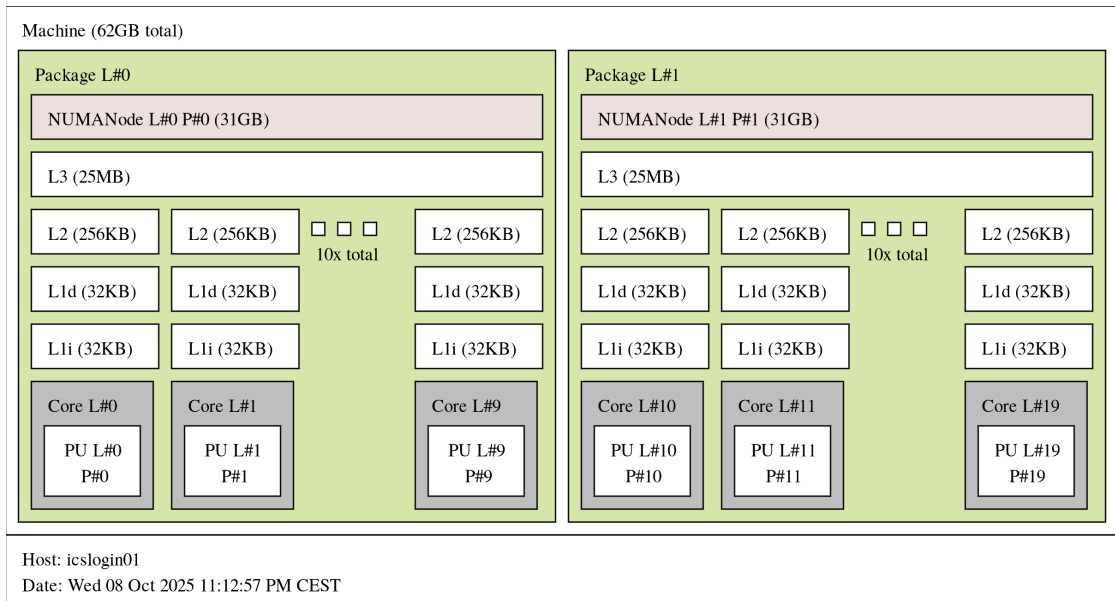


Figure 1: Graphical representation of Rosa node topology generated by command `hwloc-ls`.

2.3. Bandwidth: STREAM benchmark

2.4. Performance model: A simple roofline model

3. Optimize Square Matrix-Matrix Multiplication (50 Points)

3.1. Motivation

3.2. Importance of data access optimization

3.3. Optimizing square matrix multiplication

4. Quality of the Report (15 Points)

References

- [1] Intel Corporation. Intel® xeon® processor e5-2650 v3 (25m cache, 2.30 ghz) product specifications, 2014.
- [2] Intel Corporation. *Intel® 64 and IA-32 Architectures Optimization Reference Manual*, 2023.
- [3] Università della Svizzera italiana, Advanced Computing Lab. Rosa cluster hardware overview, 2024.