# Energy Efficient Computing

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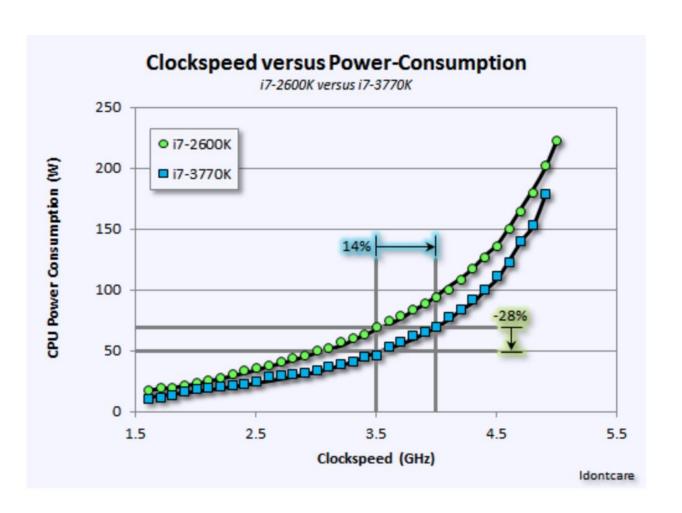


### Energy-aware focus





# Power scales with CPU Frequency

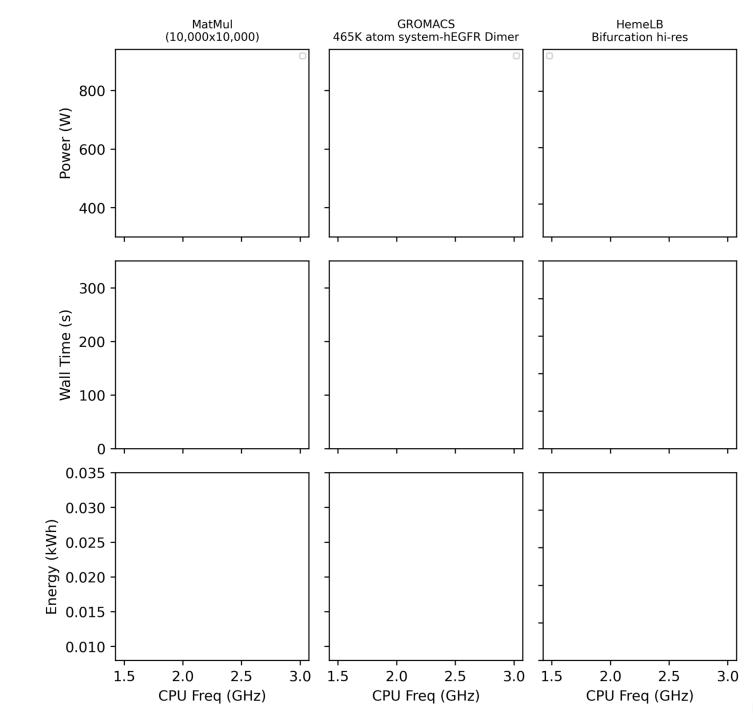


 $Power \propto Freq * Voltage^2$ 



To boost, or not too boost, that is the question

- Rome "Zen2" Released 07 August 7 2019
  - Nominal Freq (2.6 Ghz)
- Genoa "Zen4" Released 10 November 2022
  - Nominal Freq (2.4 Ghz)

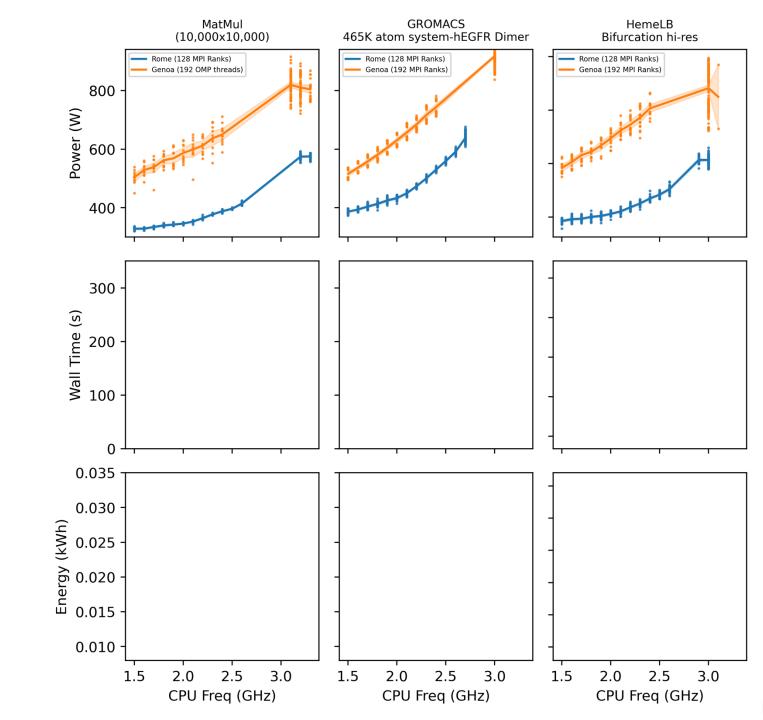


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#### Observations:

- 1. Genoa draws more power
- Each application has a different "power draw signature"

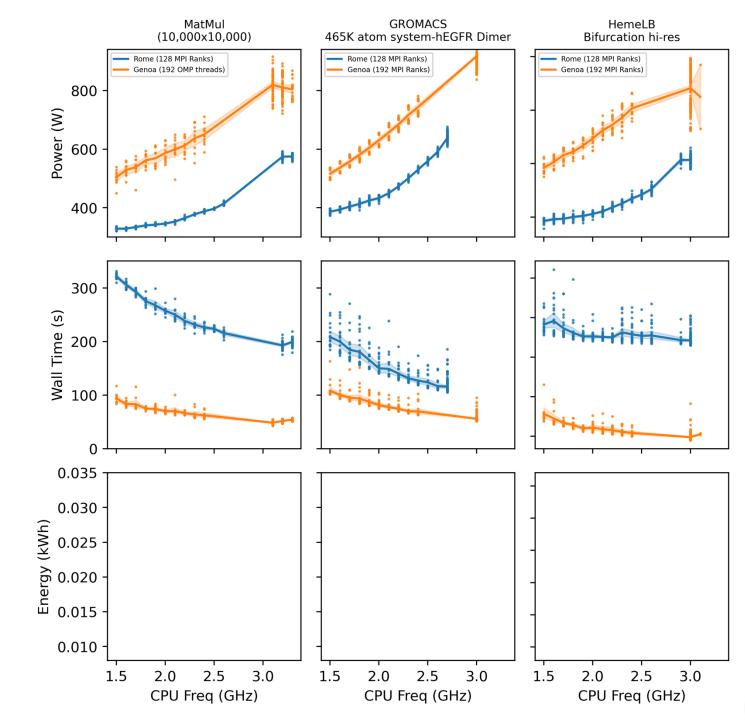


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#### **Observations:**

- 1. Genoa is more performant
  - More Logical cores to devote to problem
- In the 3 cases, Boost does not have the similar effect in performance vs power draw. i.e. its performance increase is flat as compared to its power draw

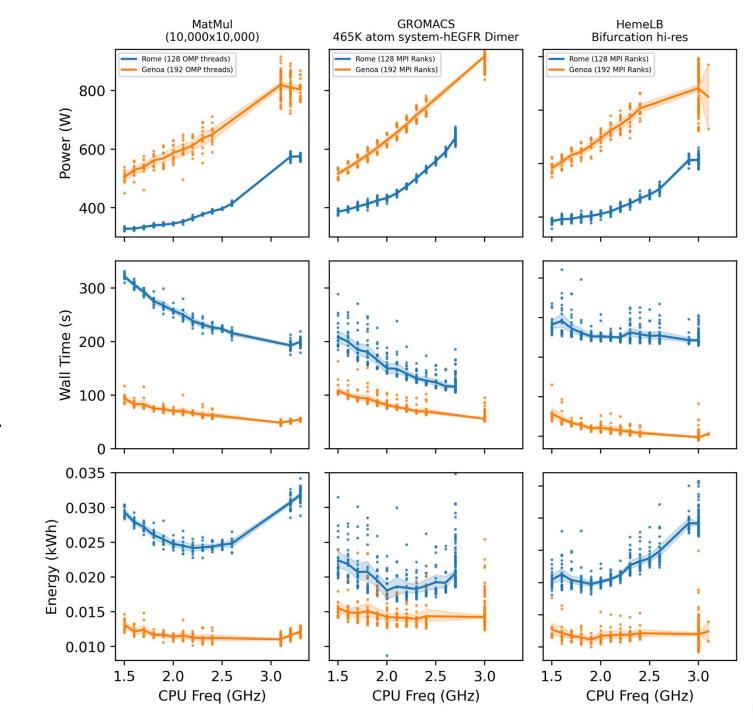


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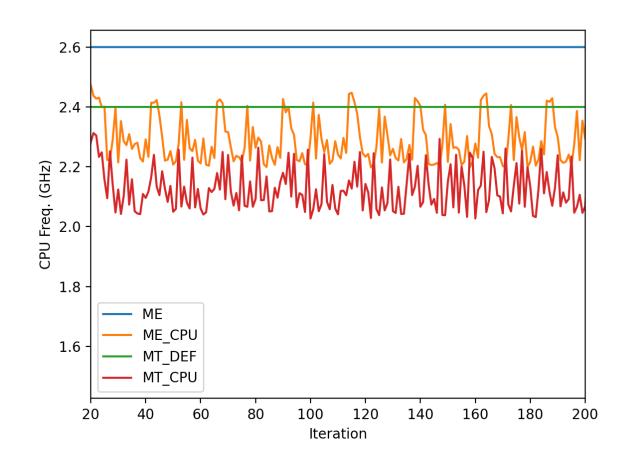
#### Observations:

- 1. Genoa is more energy efficient.
- 2. Each application has its own energy minimum.
- 3. DVFS "by hand" is HARD!!!!!!!



# Application Optimization via EAR Policies

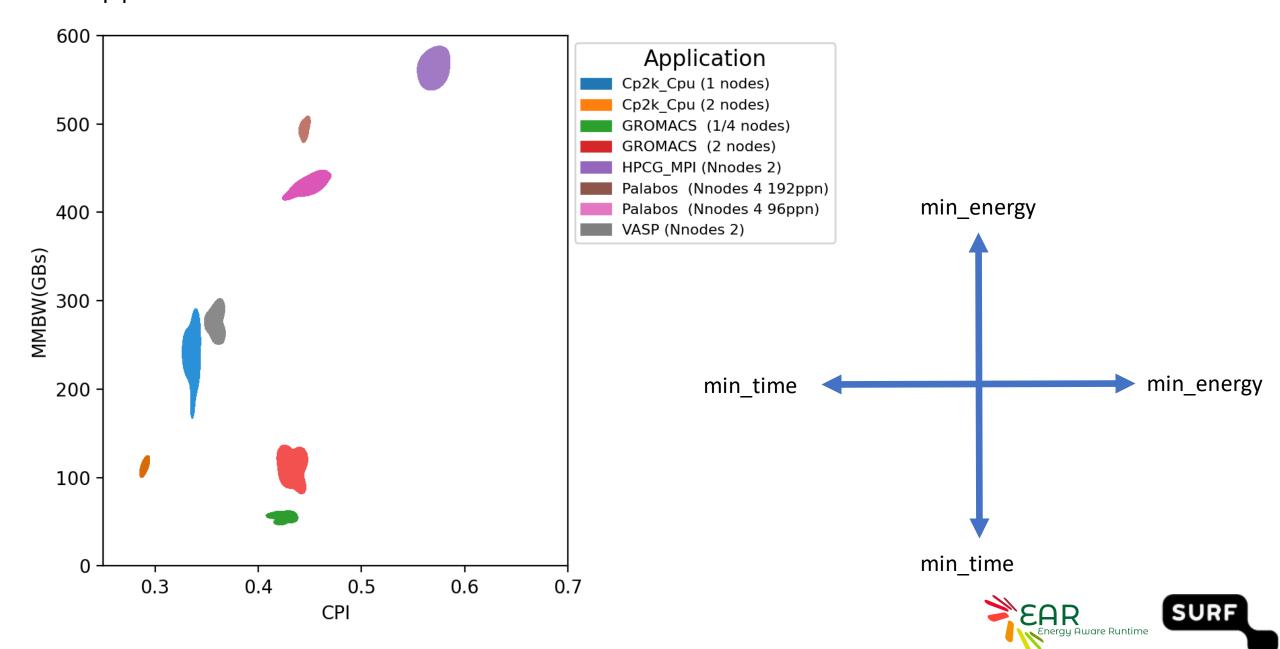
- monitoring
  - Does not affect the CPU Freq.
     Only monitors applications.
- min\_energy
  - Memory intensive applications
- min\_time
  - CPU intensive applications

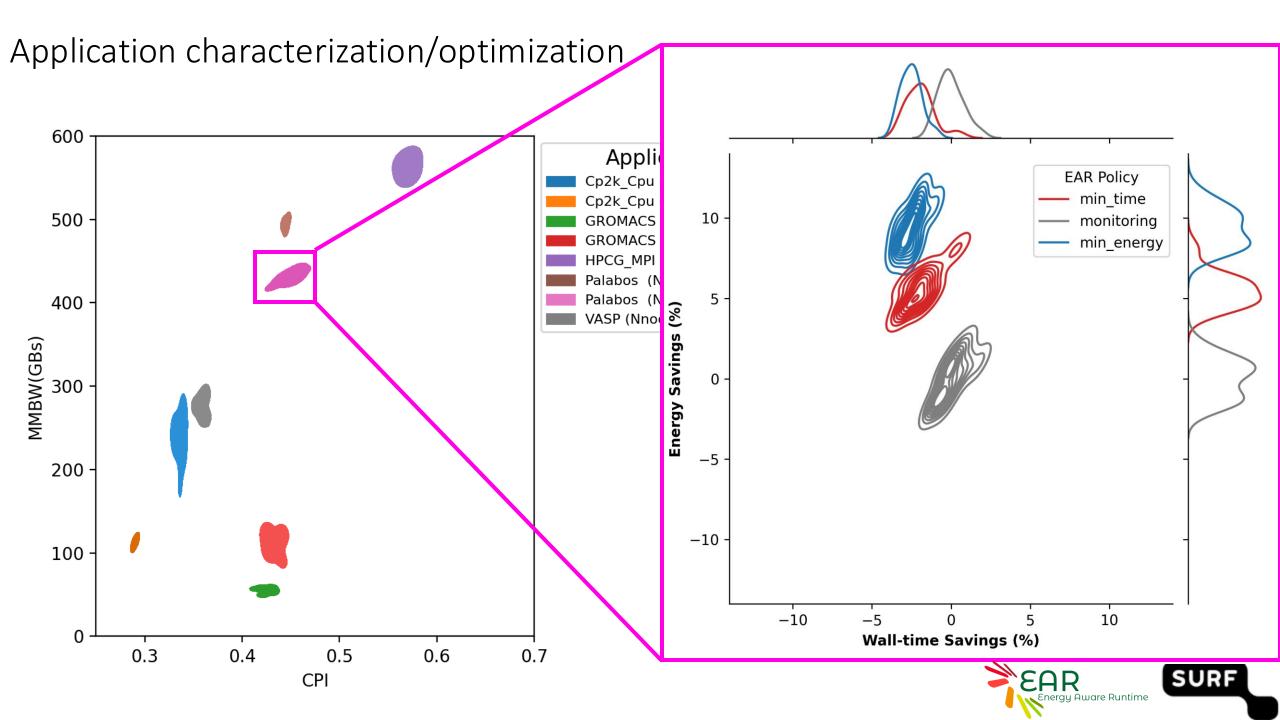






### CPU Application characterization





### Energy Aware Runtime (EAR)

Energy Aware Runtime (EAR) package provides an energy management framework for super computers.

### Its Documented and available for use!!!

https://servicedesk.surf.nl/wiki/display/WIKI/Energy



Pages / ... / Energy

#### **EAR example: GROMACS**

i In this short example we will show you how to run a simple GROMACS benchmark with EAR enabled.

GROMACS (https://www.gromacs.org)

The HECBioSim Benchmarks (https://www.hecbiosim.ac.uk/access-hpc/benchmarks)

**GROMACS** A free and open-source software suite for high-performance molecular dynamics and output analysis.

**HECBioSim benchmark suite** consists of a set of simple benchmarks for a number of popular Molecular Dynamics (MD) engines, each of which is set at a different atom count. The benchmark suite currently contains benchmarks for the AMBER, GROMACS, LAMMPS and NAMD molecular dynamics packages.

- 1 Prepare the benchmark input
- 2 Prepare the SLURM job-script
- 3 Get energy and performance metrics report from your job

#### Prepare the SLURM job-script

Notice all that is special here is Lines 10 and 11. This is how you call the EAR runtime library (EARL).

Read more detail here Energy Aware Runtime (EAR) if you want more information on how EAR works.

```
#!/bin/bash

#$SBATCH -p rome
#$SBATCH -n 128

#$SBATCH -t 00:20:00

#$SBATCH --exclusive
#$SBATCH --output=GROMACS_run.out
#$SBATCH --error=GROMACS_run.err

#$SBATCH --ear=on
#$SBATCH --ear=on
#$SBATCH --ear-policy=monitoring

module load 2022
module load foss/2022a
module load GROMACS/2021.6-foss-2022a

srun --ntasks=128 --cpus-per-task=1 gmx_mpi mdrun -s hEGFRDimer_benchmark.tpr
```

Obviously you need to submit it to the SLURM scheduler!!

So...

```
sbatch myjobscript.sh
```

# Thank you!

- Contact us
  - https://servicedesk.surf.nl/
- EAR User Tutorial on Snellius
  - https://servicedesk.surf.nl/wiki/display/WIKI/Energy+Aware+Runtime+(EAR)
- Energy Aware Runtime (EAR) --- Energy Aware Solutions
  - https://www.eas4dc.com/solutions
  - https://gitlab.bsc.es/ear\_team/ear/-/wikis/home



