

# BSC Tools Hands-On

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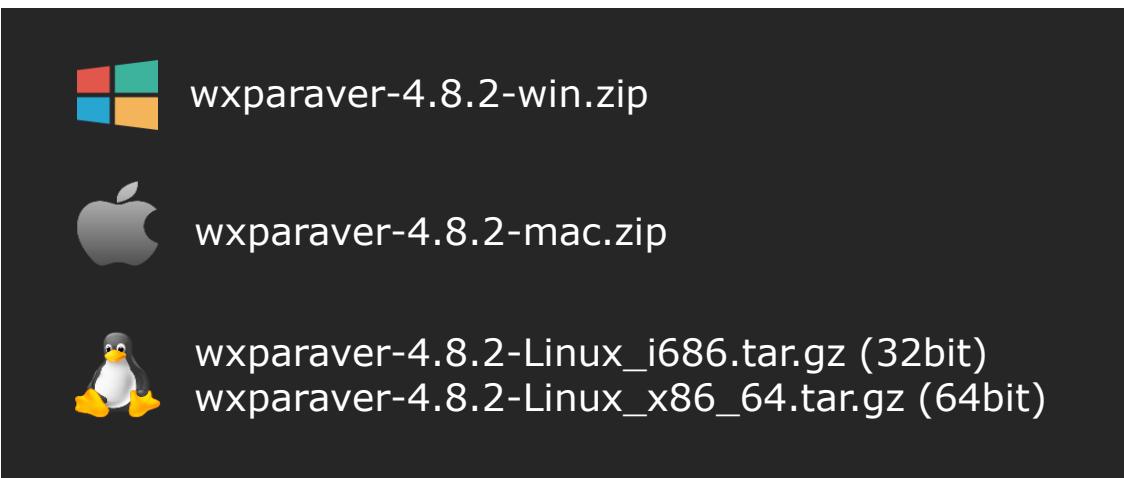
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Barcelona Supercomputing Center

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# Installing Paraver in your computer

# Install Paraver in your laptop

- Download a binary for your OS
  - <https://tools.bsc.es/downloads>



The screenshot shows the BSC Tools website's "Downloads" section. The "CORE TOOLS" section contains a link to "Get PARAVER" which is highlighted with a yellow arrow. The "PERFORMANCE ANALYTICS" section contains links to "Get CLUSTERING", "Get SPECTRAL", "Get TRACKING", and "Get BASIC ANALYSIS". The "DIMEMAS" section contains a link to "Get DIMEMAS". Each tool entry includes its version and file size.

Category	Tool	Description	Version	File Size
CORE TOOLS	EXTRAE	Instrumentation framework to generate execution traces of the most used parallel runtimes.	Version 3.4.1	2.24 MB
	PARAVER	Expressive powerful and flexible trace visualizer for post-mortem trace analysis.	Version 1.0.2	1.65 MB
	DIMEMAS	High-abstracted network simulator for message-passing programs.	Version 5.2.12	1.09 MB
	CLUSTERING	Automatically expose the main performance trends in applications' computation structure.	Version 2.6.6	2 MB
PERFORMANCE ANALYTICS	SPECTRAL	Signal processing techniques to select representative regions from Paraver traces.	Version 3.4.0	0.31 MB
	TRACKING	Analyze how the behavior of a parallel application evolves through different scenarios.	Version 2.6.5	1.9 MB
	FOLDING	Combined instrumentation and sampling for instantaneous metric evolution with low overhead.	Version 1.0.2	11.06 MB
	BASIC ANALYSIS	Framework for automatic extraction of fundamental factors for Paraver traces.	Version 0.2	10.89 MB

# Install Paraver tutorials

- Download tutorials archive
  - <https://tools.bsc.es/paraver-tutorials>

The screenshot shows a web browser displaying the BSC tools website. The URL in the address bar is `news@tools:~ > Paraver 4.7.2 avail`. The page title is "Home » Documentation » Paraver tutorials". The content area describes seven tutorials available for Paraver 4.7.2. A yellow callout box labeled "All tutorials" points to the first tutorial in the list: "Paraver introduction (MPI)". Another yellow callout box at the bottom points to the download links for ".tar.gz format" and ".zip format".

These seven tutorials can be opened with wxParaver versions newer than 4.3.0, and you'll be able to follow the steps within the tool. To install them, download and untar the package and follow the instructions of the Help/Tutorial option on the Paraver main window. Following there is a list of available tutorials:

- Paraver introduction (MPI)
- Dimemas introduction
- Introduction to Paraver and Dimemas methodology
- Methodology
- Tutorial on HydroC analysis (MPI, Dimemas, CUDA)
- Trace preparation
- Trace alignment tutorial.

If you prefer, you can download all of them together in a single package:

- [.tar.gz format \(127 Mb\)](#)
- [.zip format \(127 Mb\)](#)

## Install Paraver (III)

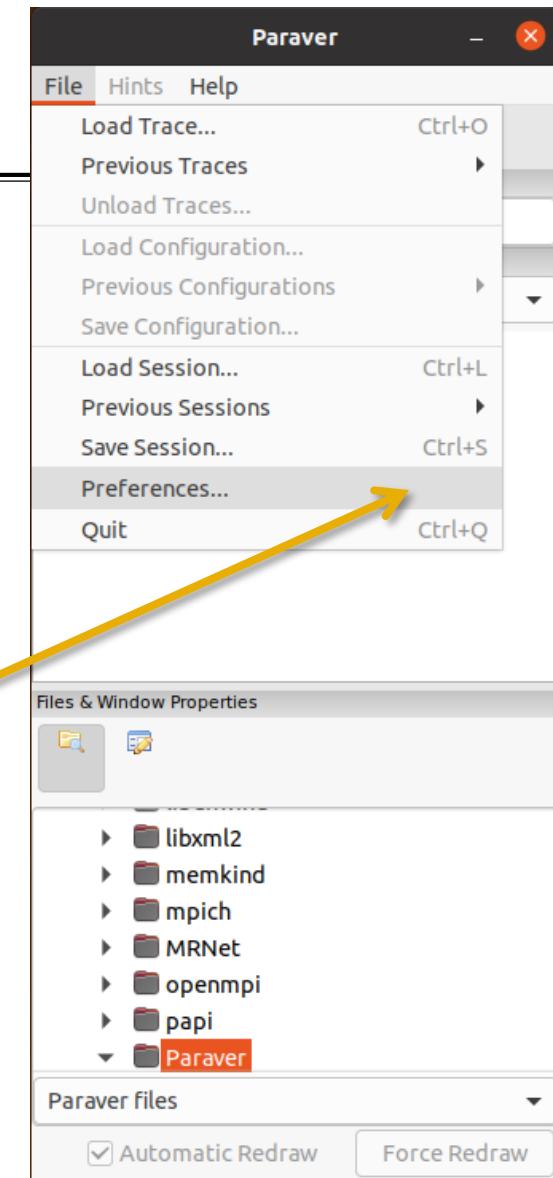
- Uncompress downloaded packages
- Rename the folders:
  - wxparaver-4.8.2-\* → paraver
  - paraver-tutorials-20150526 → tutorials
- Start Paraver locally:
  - Linux: Run the command:

```
laptop$ paraver/bin/wxparaver
```

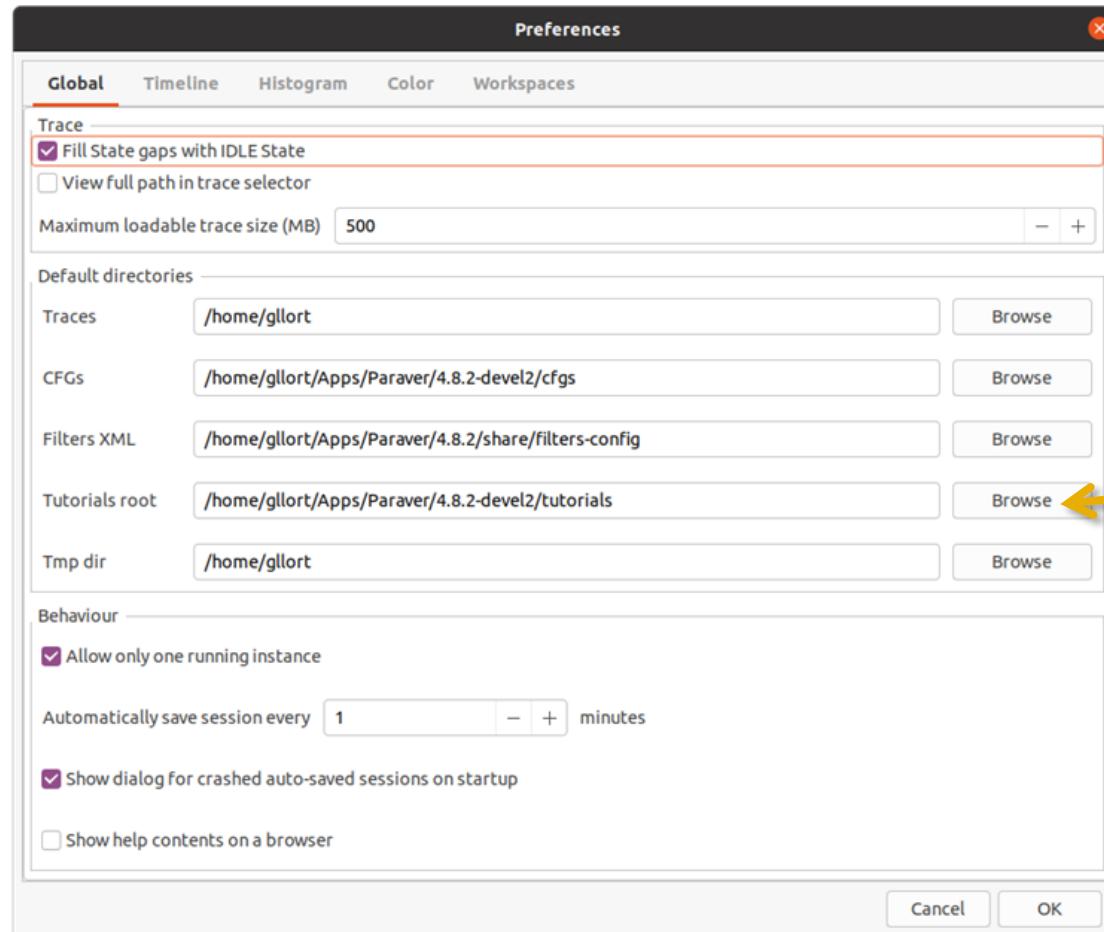
- Windows: Double-click on paraver/wxparaver.exe
  - MAC: Double click on paraver/wxparaver.app
- Any issue? Start Paraver in HAWK:

```
hawk> module load bsc_tools
hawk> wxparaver &
```

- Open File → Preferences



## Install Paraver (IV)



- Setup the “Tutorials root” pointing to your folder “tutorials”

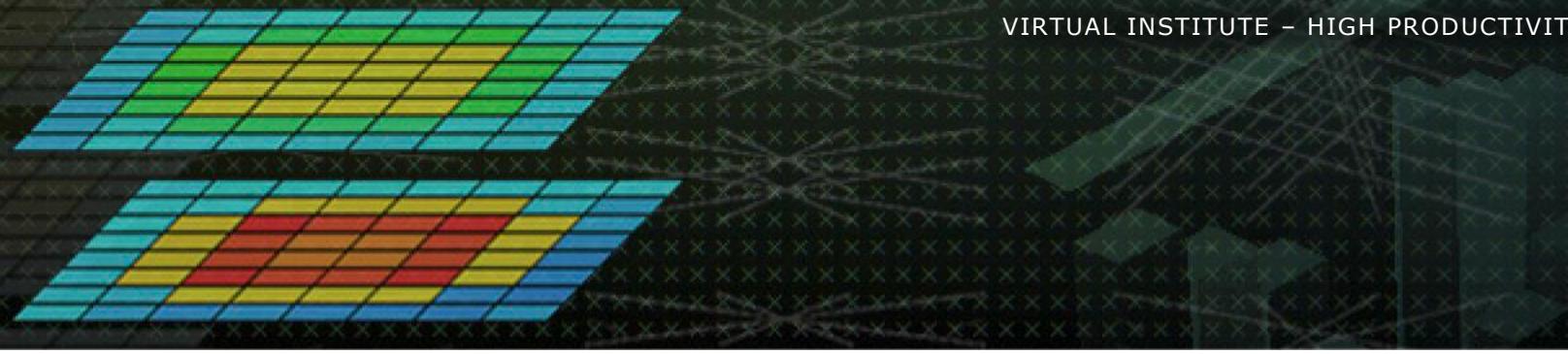
Click Browse and select your folder “tutorials”

## Install Paraver (IV)

- Check tutorials are properly installed



- Follow these tutorials by clicking on the hyperlinks and reading the explanations. When you click on a link, one or more views will open.

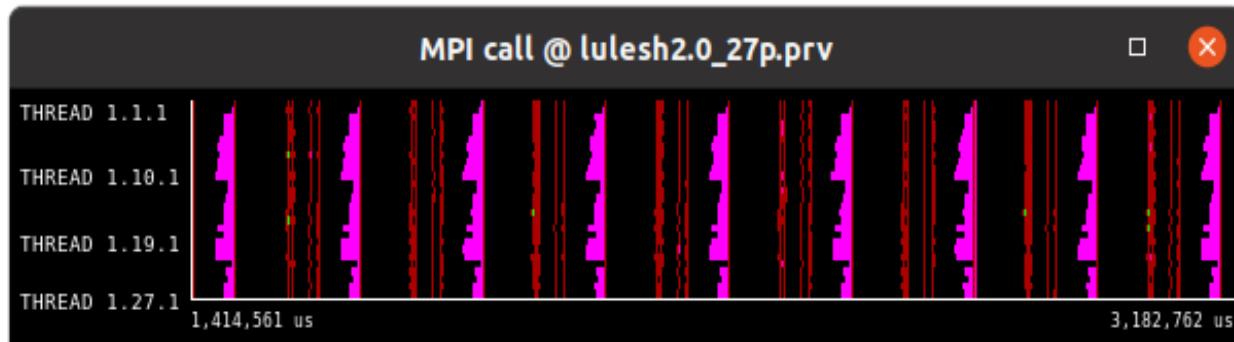


# Brief introduction to Paraver

## 3 main views of Paraver (I)

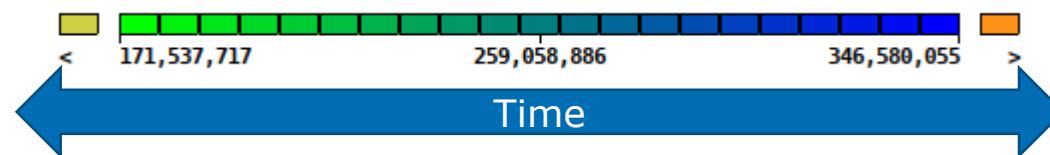
### ▪ Timeline

↑ Processes (and threads)



**Code color**  
(e.g. 1 color for each MPI call)

- Outside MPI
- MPI\_Isend
- MPI\_Irecv
- MPI\_Wait
- MPI\_Waitall
- MPI\_BARRIER
- MPI\_Reduce
- MPI\_Allreduce
- MPI\_Comm\_rank
- MPI\_Finalize

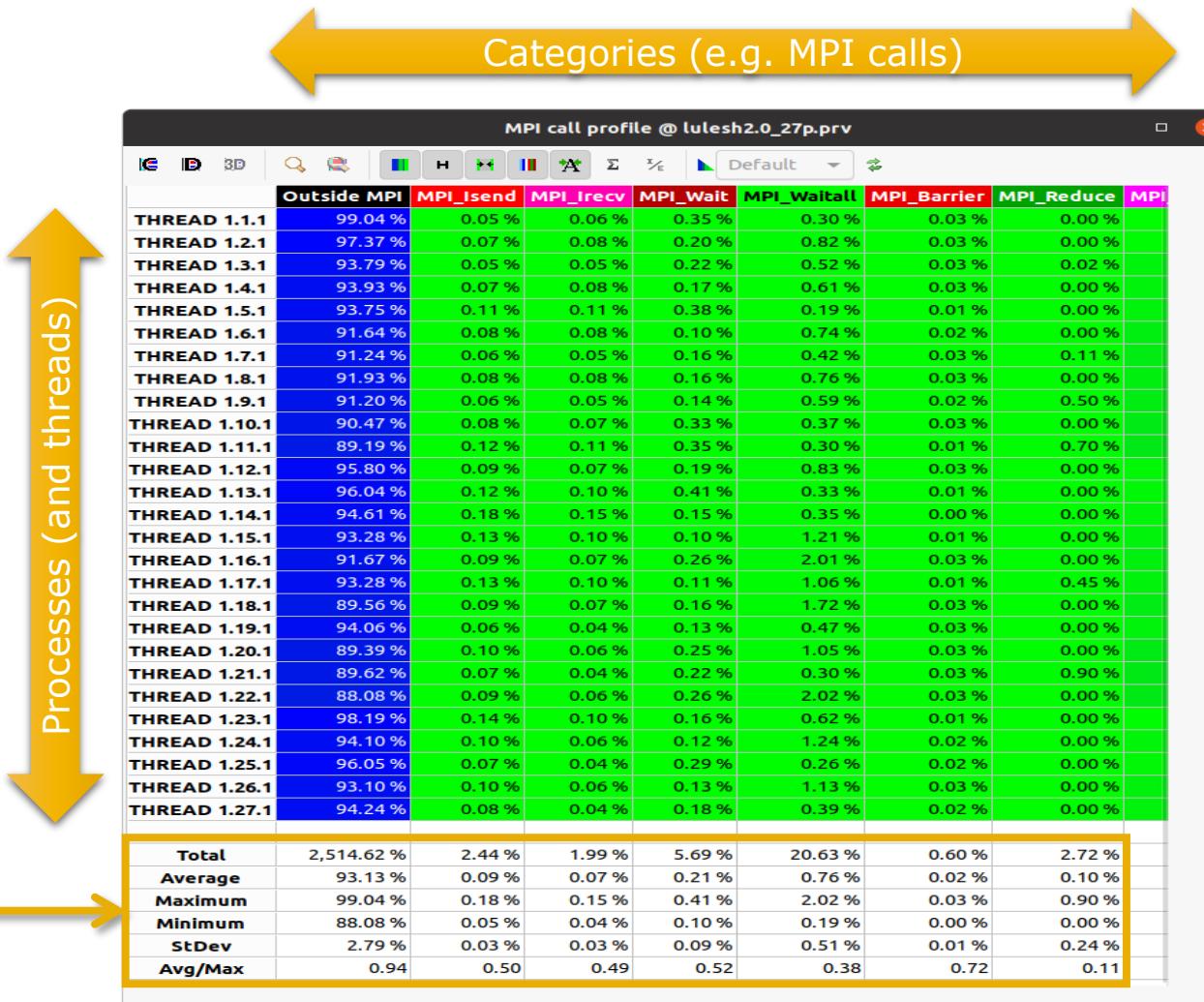


**Gradient color**  
(e.g. from low  
#instructions to  
high  
#instructions)

← Time →

## 3 main views of Paraver (II)

- Table (Profile)



The table can display a variety of statistics (e.g. % of time, # of calls, etc.) with gradient coloring showing from low values to high values

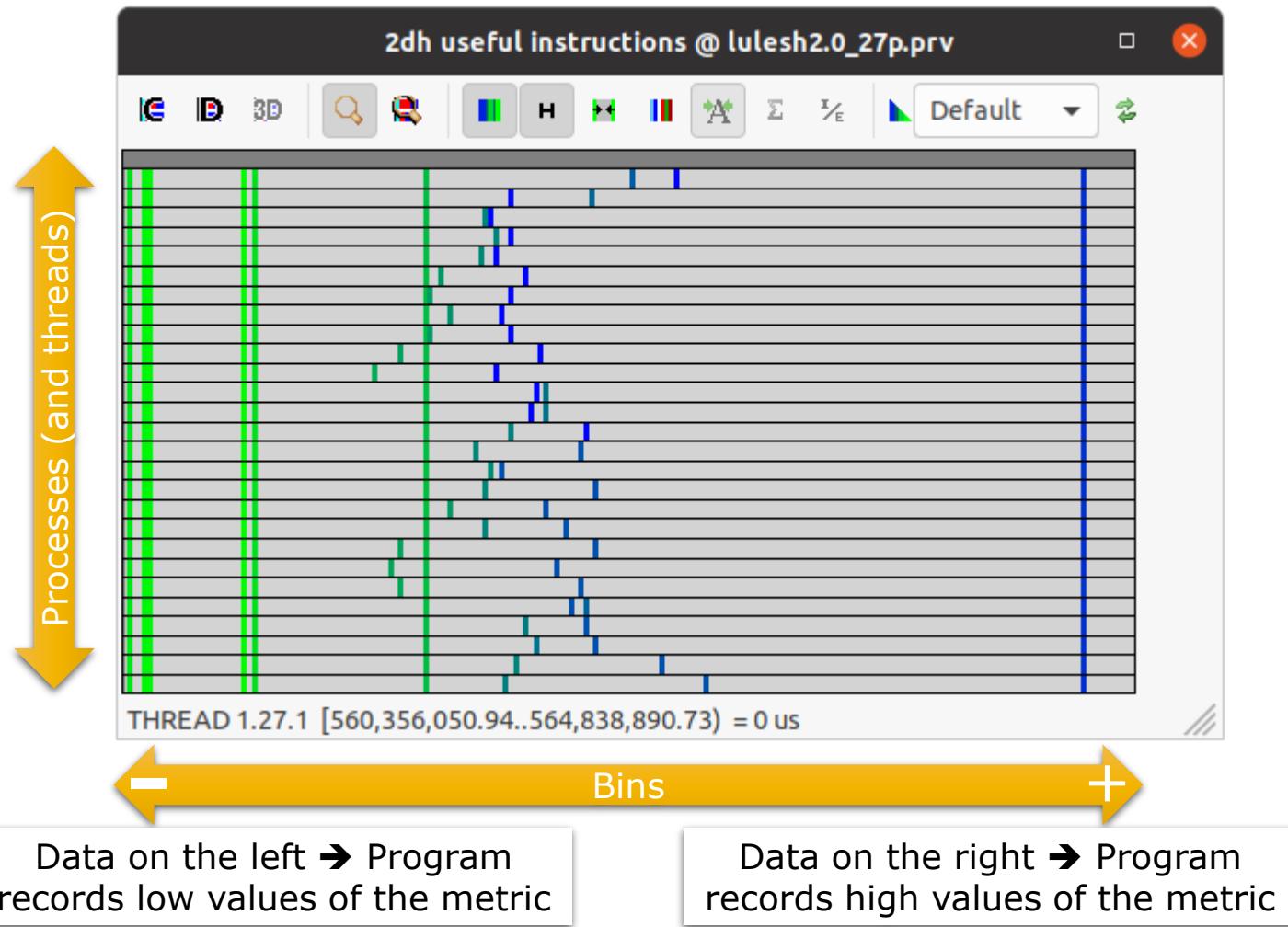
## 3 main views of Paraver (III)

### ▪ Histogram

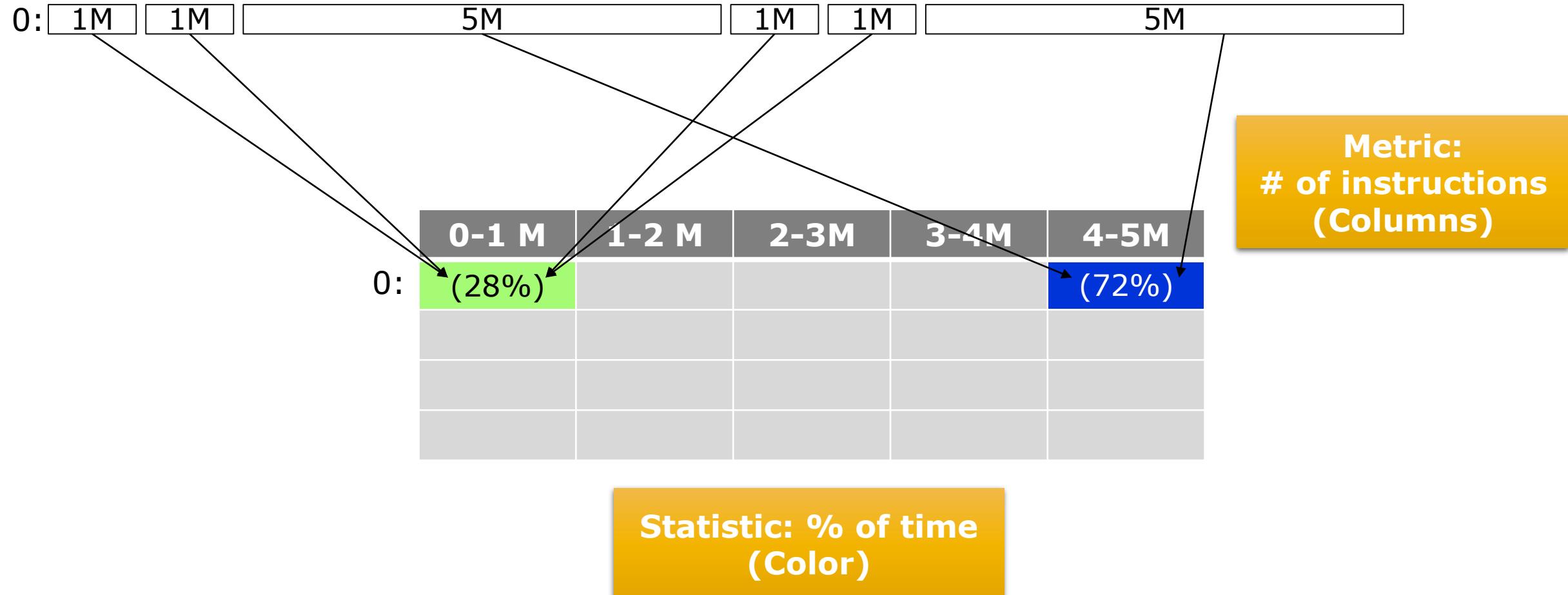
Displays continuous metrics (e.g. **instructions executed**, duration of computations, bytes sent/received, etc. )

Gradient color represents if the selected statistic (% of time, # of calls, etc.) for each behavior is **high** or **low**

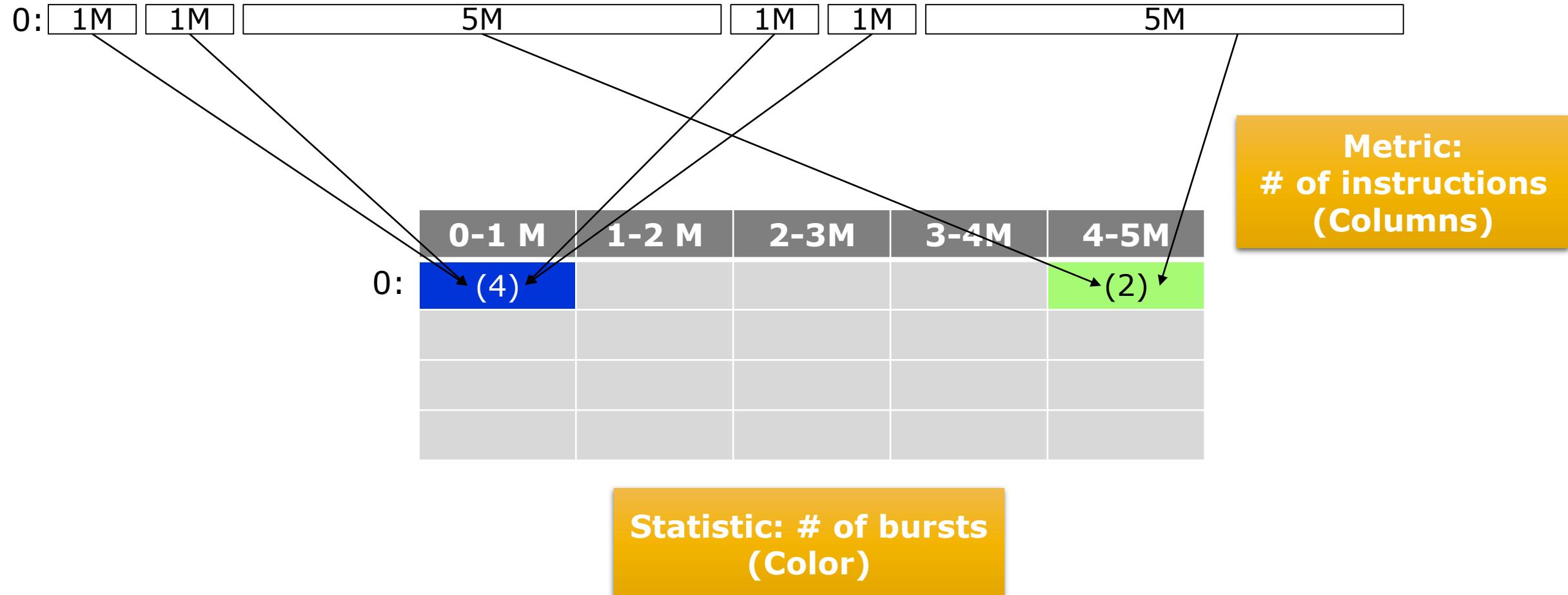
**General tip:** straight lines are good (all processes show same behavior), while variabilities usually indicate imbalances



## Histogram: position, metric and statistic ... wait, what!?



## Histogram: position, color, metric and statistic ... wait, what!?



## Histogram: position, color, metric and statistic ... wait, what!?

0:	1M	1M	5M	1M	1M	5M
1:	1M	1M	5M	1M	1M	5M
2:	1M	1M	5M	1M	1M	5M
3:	1M	1M	5M	1M	1M	5M

	0-1 M	1-2 M	2-3M	3-4M	4-5M
0:	(4)				(2)
1:	(4)				(2)
2:	(4)				(2)
3:	(4)				(2)

**Metric:**  
**# of instructions**  
**(Columns)**

**Statistic: # of bursts**  
**(Color)**

## Histogram: position, color, metric and statistic ... wait, what!?

0:	1M	1M	5M	1M	1M	5M
1:	1M	1M	5M	1M	1M	5M
2:	2M	2M	2M	2M	2M	2M
3:	1M	1M	5M	1M	1M	5M

	0-1 M	1-2 M	2-3M	3-4M	4-5M
0:	(4)				(2)
1:	(4)				(2)
2:			(7)		(2)
3:	(4)				(2)

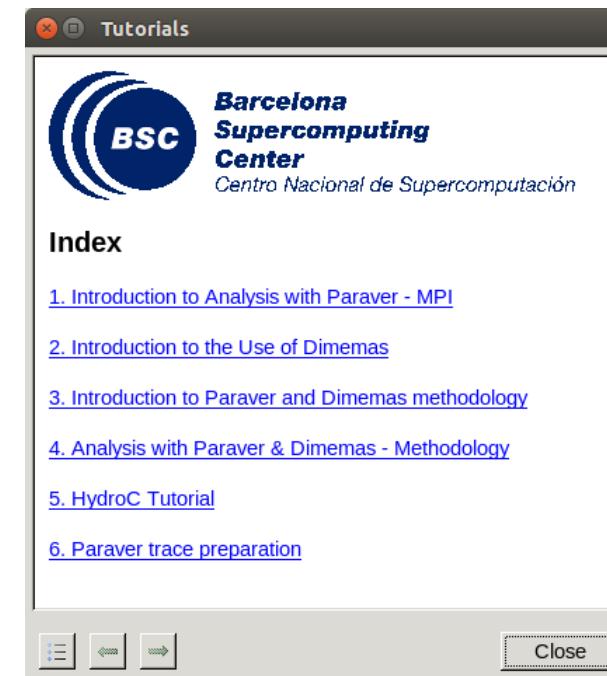
**Metric:**  
**# of instructions**  
**(Columns)**

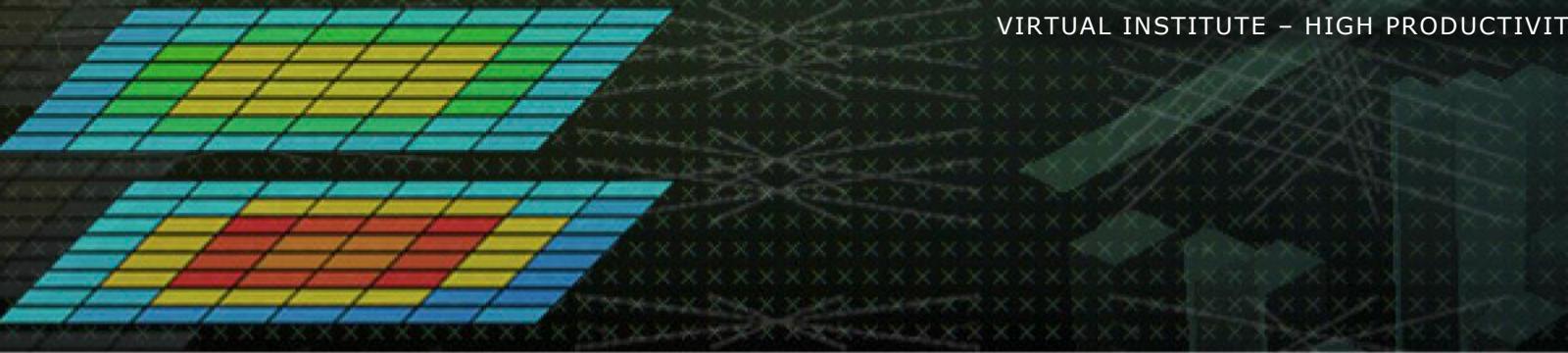
**Statistic: # of bursts**  
**(Color)**

## First steps with Paraver

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- Follow tutorial number...
  - 1 → Explains basic navigation with the tool
  - 3 → Basic analysis methodology (first 4 bullets, Clustering and Dimemas part not covered today!)
  - 5 → Analysis methodology applied to a real application





# Getting a trace with Extrae

## Extrae features

- Platforms
  - Intel, Cray, BlueGene, MIC, ARM, Android, Fujitsu Sparc ...
- Parallel programming models
  - MPI, OpenMP, pthreads, OmpSs, CUDA, OpenCL, Java, Python ...
- Performance Counters
  - Using PAPI interface
- Link to source code
  - Callstack at MPI routines
  - OpenMP outlined routines
  - Selected user functions (Dyninst)
- Periodic sampling
- User events (Extrae API)

No need to  
recompile  
nor relink!

## Extrae Overheads

---

	Average values
Recording 1 Event	150 – 200ns
1 Event + PAPI hardware counters	750 – 1000ns
1 Event + Callstack (1 level)	1µs
1 Event + Callstack (6 levels)	2µs

## How does Extrae work?

---

- Symbol substitution through LD\_PRELOAD
  - Specific libraries for each combination of runtimes
    - MPI
    - OpenMP
    - OpenMP+MPI
    - ...
- Dynamic instrumentation
  - Based on Dyninst (developed by U.Wisconsin / U.Maryland)
    - Instrumentation in memory
    - Binary rewriting
- Alternatives
  - Static link (i.e., PMPI, Extrae API)



Recommended

## Extrae on HAWK

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- Log-in to HAWK:

```
laptop$ ssh -Y <USER>@hawk.hww.hlr.de
```

- Extrae is available via modules for 2 MPI versions... choose yours!

- Cray MPT (default)

```
hawk$ module load mpt
hawk$ module load extrae
```

- OpenMPI

```
hawk$ module load openmpi
hawk$ module load extrae
```

## Getting your first trace

---

- Copy this folder to your `$HOME` and you are ready to follow this hands-on tutorial

```
hawk$ cp -r /lustre/cray/ws9/2/ws/hpcjgrac-tw35/BSC/tools-material $HOME
```

- Provided folder **tools-material** in `/lustre/cray/ws9/2/ws/hpcjgrac-tw35/BSC` contains:
  - Test application compiled for OpenMPI (`lulesh2.0_openmpi`)
  - Jobscripts to execute and trace (`job.pbs`, `trace.sh`)
  - Configuration of the tracing tool (`extrae.xml`)
  - Already generated tracefiles (`traces/*.{pcf,prv,raw}`)
  - Clustering analysis configuration file (`cluster.xml`)
  - A copy of this slides

## Using Extrae in 3 steps

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### 1. Adapt your job submission scripts

### 2. Configure what to trace

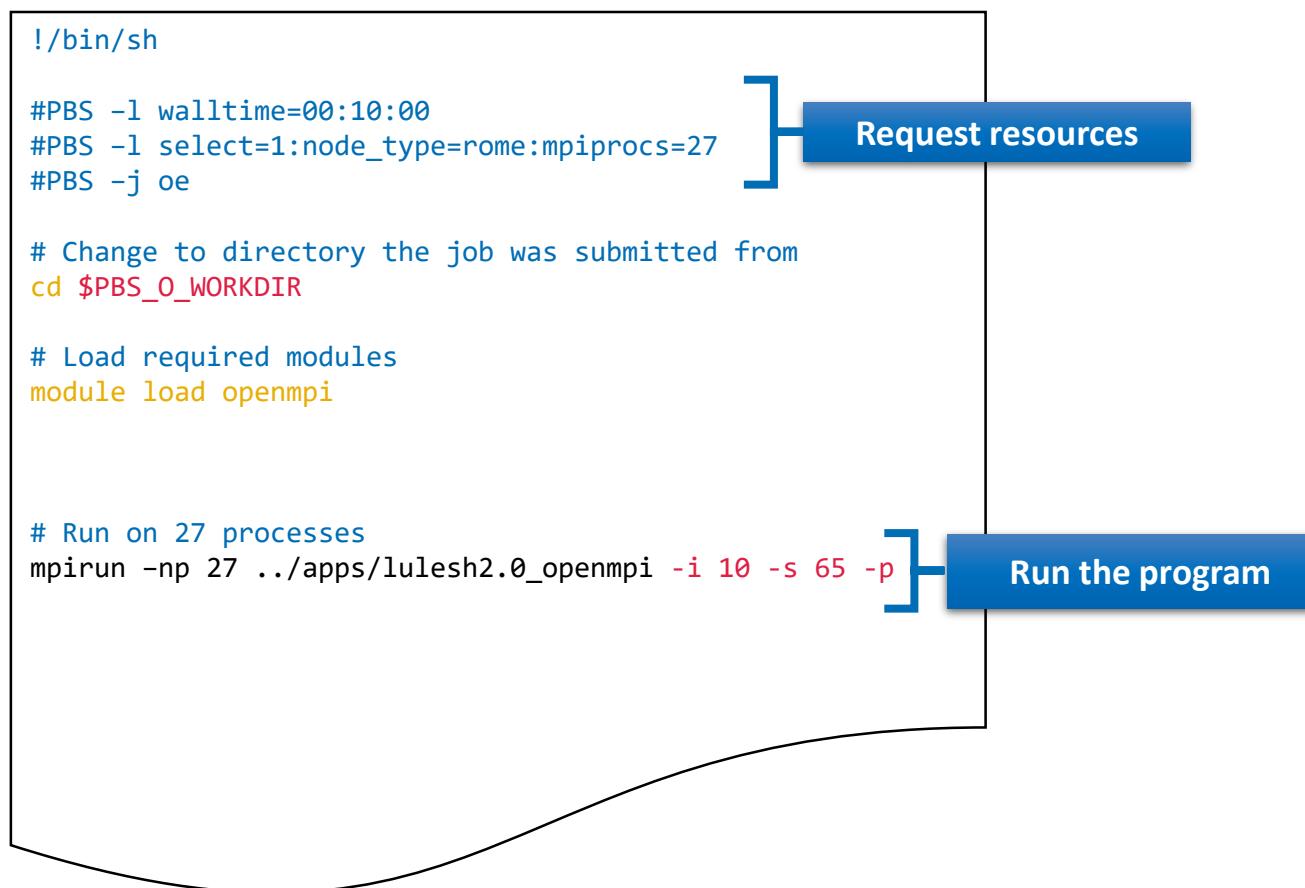
- XML configuration file
- More example configurations at `$EXTRAE_HOME/share/example`

### 3. Run it!

- For further reference check the **Extrae User Guide**:
  - <https://tools.bsc.es/doc/html/extrae>
  - Also distributed with Extrae at `$EXTRAE_HOME/share/doc`

## Step 1: Adapt the job script to load Extrae

- Example of a standard jobscrip (without tracing)



## Step 1: Adapt the job script to load Extrae

- Jobscript modified to load Extrae (extrae/job.pbs)

```
#!/bin/sh

#PBS -l walltime=00:10:00
#PBS -l select=1:node_type=rome:mpiprocs=27
#PBS -j oe

# Change to directory the job was submitted from
cd $PBS_O_WORKDIR

# Load required modules
module load openmpi

export TRACE_NAME=lulesh2.0_openmpi_27p.prv

# Run on 27 processes
mpirun -np 27 ./trace.sh ./apps/lulesh2.0_openmpi
-i 10 -s 65 -p
```

Specify name of output trace (optional)

Run with Extrae

## Step 1: Adapt the job script to load Extrae

- Tracing launcher helper script (extrae/trace.sh)

```
#!/bin/sh

#PBS -l walltime=00:10:00
#PBS -l select=1:node_type=rome:mpiprocs=27
#PBS -j oe

# Change to directory the job was submitted from
cd $PBS_O_WORKDIR

# Load required modules
module load openmpi

export TRACE_NAME=lulesh2.0_openmpi_27p.prv

# Run on 27 processes
mpirun -np 27 ./trace.sh ./apps/lulesh2.0_openmpi
-i 10 -s 65 -p
```

```
#!/bin/sh

# Extrae configuration file
export EXTRAE_CONFIG_FILE=./extrae.xml What to trace?

# Select tracing library (depends on runtime & lang)

# For MPI/C apps
export LD_PRELOAD=${EXTRAE_HOME}/lib/libmpitrace.so
# For MPI/Fortran apps
#export LD_PRELOAD=${EXTRAE_HOME}/lib/libmpitracef.so
...
# Run the application
$*
```

Choose a tracing library depending on the app type (see next slide)

## Step 1: Which tracing library?

---

- Choose depending on the application type

Library	Serial	MPI	OpenMP	pthread	CUDA
libseqtrace	✓				
libmpitrace[f] <sup>1</sup>		✓			
libomptrace			✓		
libpttrace				✓	
libcudatrace					✓
libompitrace[f] <sup>1</sup>		✓	✓		
libptmpitrace[f] <sup>1</sup>		✓		✓	
libcudampitrace[f] <sup>1</sup>	✓				✓

<sup>1</sup> add suffix “f” in Fortran codes

## Step 2: Extract XML configuration – most relevant settings

```
<mpi enabled="yes">
  <counters enabled="yes" /> ←
</mpi>

<openmp enabled="yes">
  <locks enabled="no" />
  <counters enabled="yes" />
</openmp>

<pthread enabled="no">
  <locks enabled="no" />
  <counters enabled="yes" />
</pthread>

<callers enabled="yes">
  <mpi enabled="yes">1-3</mpi> ←
  <sampling enabled="no">1-5</sampling>
</callers>
```

Instrument the MPI calls  
(What's the program doing?)

Instrument the call-stack  
(Where in my code?)

## Step 2: Extract XML configuration – most relevant settings (II)

```
<counters enabled="yes">
  <cpu enabled="yes" starting-set-distribution="1">
    <set enabled="yes" domain="all" changeat-time="5s">
      PAPI_TOT_INS,PAPI_TOT_CYC,PAPI_L1_DCM,PAPI_L2_DCM
    </set>
    <set enabled="yes" domain="all" changeat-time="5s">
      PAPI_TOT_INS,PAPI_TOT_CYC,PAPI_FP_OPS
    </set>
  </cpu>
  <network enabled="no"/>
  <resource-usage enabled="no"/>
  <memory-usage enabled="no"/>
</counters>
```

Select which HW counters are measured  
(How's the machine doing?)

> papi\_avail

← List of available counters (PAPI)

> papi\_best\_set

← Make groups of compatible counters (Extracte)

```
<buffer enabled="yes">
  <size enabled="yes">5000000</size>
  <circular enabled="no" />
</buffer>

<sampling enabled="no" type="default" period="50m" variability="10m"/>

<merge enabled="yes" synchronization="default" tree-fan-out="16" max-memory="512" joint-states="yes" keep-mpits="yes" sort-addresses="yes" overwrite="yes" $TRACE_NAME$>
</merge>
```

Extracte buffer size  
(Flush/memory trade-off)

Additional sampling  
(Want more details?)

Automatic post-processing to generate the Paraver trace

## Step 3: Run it!

---

- Submit your job as usual

```
hawk$ qsub -q R_tw job.pbs
```

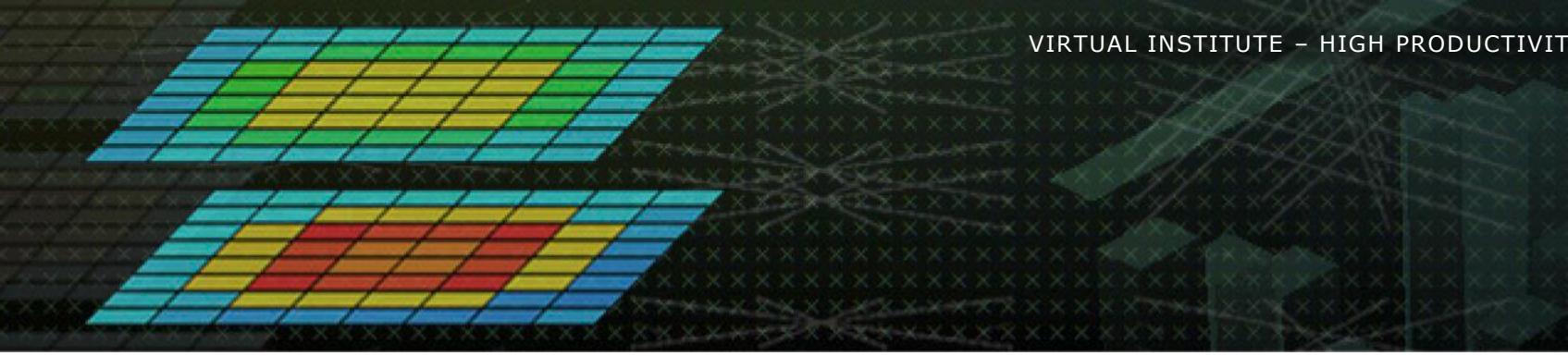
## All done! Check your resulting trace

---

- Once finished (check with “`qstat -u $USER`”) you will have the trace (3 files):

```
hawk$ ls -l
...
lulesh2.0_openmpi_27p.pcf
lulesh2.0_openmpi_27p.prv
lulesh2.0_openmpi_27p.row
```

- Any trouble? There's a trace already generated under the “traces” folder
- Now let's look into it!



# Analysing a trace with Paraver

## First steps of analysis

---

- Copy the trace to your computer (scp, WinSCP, etc.)

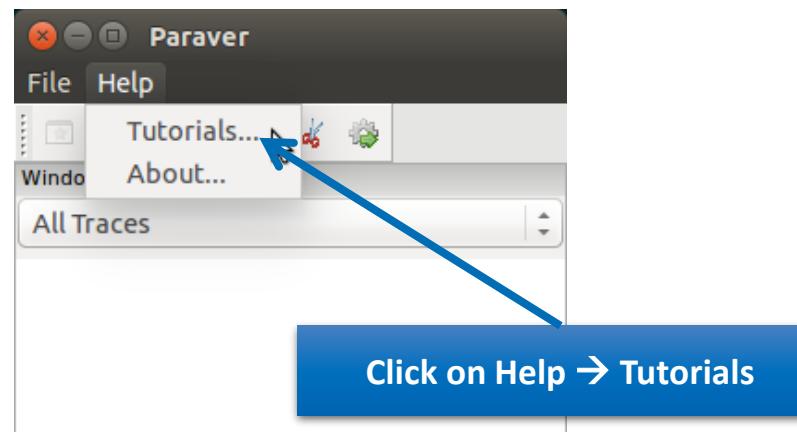
```
laptop$ scp <USER>@hawk.hww.hlr.de:tools-material/extrاء/*.{prv,pcf,row} .
```

- Load the trace with Paraver



## First steps of analysis

- Follow Tutorial #3
  - Introduction to Paraver and Dimemas methodology



# Measure the parallel efficiency

- Click on “mpi\_stats.cfg”
  - Check the **Average** for the column “Outside MPI”

Tutorials

The first question to answer when analyzing a parallel code is "how efficient does it run?". The efficiency of a parallel program can be defined based on two aspects: the parallelization efficiency and the efficiency obtained in the execution of the serial regions. These two metrics would be the first checks on the proposed methodology.

- To measure the parallel efficiency load the configuration file [cfgs/mpi\\_mpi\\_stats.cfg](#) This configuration pops up a table with %time that every thread spends in every MPI call. Look at the global statistics at the bottom of the outside mpi column. Entry **Average** represents the application parallel efficiency, entry **Avg/Max** represents the global load balance and entry **Maximum** represents the communication efficiency. If any of those values are lower than 85% is recommended to look at the corresponding metric in detail. Open the control window to identify the phases and iterations of the code.
- To measure the computation time distribution load the configuration file [cfgs/general/2dh\\_usefulduration.cfg](#) This configuration pops up a histogram of the duration for the computation regions. The computation regions are delimited by the exit from an MPI call and the entry to the next call. If the histogram does not show vertical lines, it indicates the computation time may be not balanced. Open the control window to look at the time distribution and visually correlate both views.
- To measure the computational load (instructions) distribution load the configuration file [cfgs/papi/2dh\\_useful\\_instructions.cfg](#) This configuration pops up a histogram of the instructions for the computation regions. The computation regions are delimited by the exit from an MPI call and the entry to the next call. If the histogram doesn't show vertical lines, it indicates the distribution of the instructions may be not balanced. Open the control window to look at the time distribution and correlate both views.
- To measure the serial regions performance look at the TPC timeline

**Parallel efficiency (Avg)**

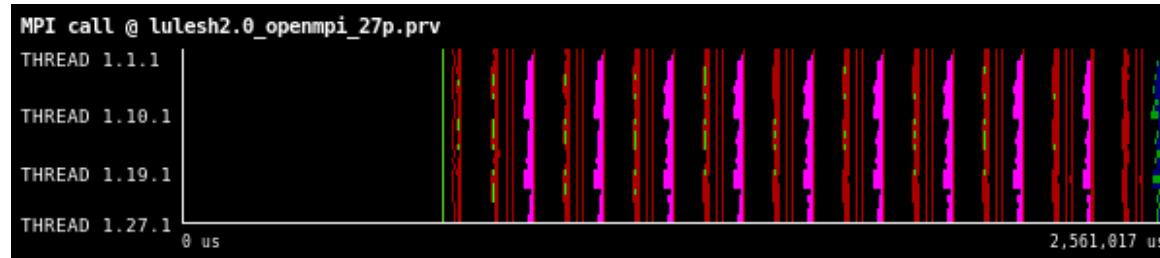
**Comm efficiency (Max)**

**Load balance (Avg/Max)**

Close

	THREAD 1.17.1	93.37 %	0.38 %	0.30 %	0.24 %	0.22 %	0.01 %	0.34 %
THREAD 1.18.1	91.20 %	0.27 %	0.20 %	0.15 %	0.92 %	0.01 %	0.00 %	
THREAD 1.19.1	92.87 %	0.18 %	0.14 %	0.13 %	0.80 %	0.05 %	0.00 %	
THREAD 1.20.1	86.08 %	0.28 %	0.20 %	0.18 %	4.02 %	0.04 %	0.00 %	
THREAD 1.21.1	85.98 %	0.19 %	0.13 %	2.41 %	1.18 %	0.06 %	0.78 %	
THREAD 1.22.1	86.42 %	0.28 %	0.19 %	0.62 %	2.85 %	0.03 %	0.00 %	
THREAD 1.23.1	93.78 %	0.42 %	0.29 %	0.43 %	2.82 %	0.03 %	0.00 %	
THREAD 1.24.1	90.60 %	0.30 %	0.18 %	0.99 %	3.40 %	0.03 %	0.00 %	
THREAD 1.25.1	94.65 %	0.20 %	0.12 %	0.63 %	0.70 %	0.03 %	0.00 %	
THREAD 1.26.1	94.19 %	0.32 %	0.17 %	0.13 %	0.66 %	0.01 %	0.00 %	
THREAD 1.27.1	93.14 %	0.21 %	0.11 %	0.10 %	1.42 %	0.04 %	0.00 %	
Total	2,454.20 %	7.19 %	6.20 %	15.50 %	52.66 %	0.79 %	2.44 %	
Average	90.90 %	0.27 %	0.23 %	0.57 %	1.95 %	0.03 %	0.09 %	
Maximum	97.91 %	0.54 %	0.48 %	2.94 %	4.02 %	0.06 %	0.78 %	
Minimum	85.98 %	0.13 %	0.11 %	0.10 %	0.22 %	0.00 %	0.00 %	
StDev	3.16 %	0.09 %	0.08 %	0.65 %	1.21 %	0.01 %	0.20 %	
Avg/Max	0.93	0.49	0.48	0.20	0.49	0.51	0.12	

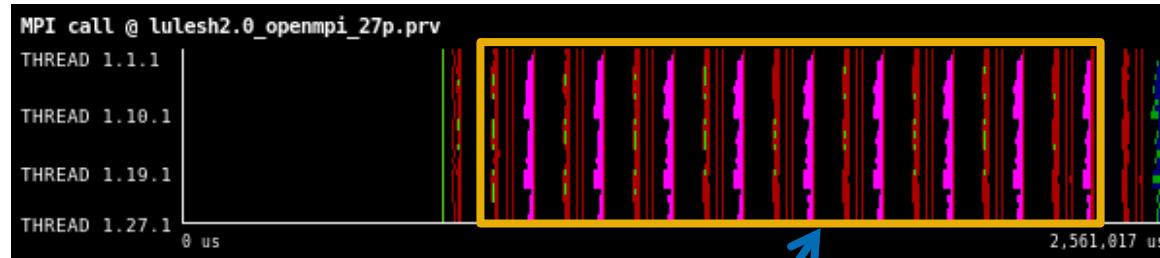
## Focus on the iterative part



Click on  
“Open Control  
Window”

	93.37 %	0.38 %	0.30 %	0.24 %	0.22 %	0.01 %	0.34 %
THREAD 1.17.1	93.37 %	0.38 %	0.30 %	0.24 %	0.22 %	0.01 %	0.34 %
THREAD 1.18.1	91.20 %	0.27 %	0.20 %	0.15 %	0.92 %	0.01 %	0.00 %
THREAD 1.19.1	92.87 %	0.18 %	0.14 %	0.13 %	0.80 %	0.05 %	0.00 %
THREAD 1.20.1	86.08 %	0.28 %	0.20 %	0.18 %	4.02 %	0.04 %	0.00 %
THREAD 1.21.1	85.98 %	0.19 %	0.13 %	2.41 %	1.18 %	0.06 %	0.78 %
THREAD 1.22.1	86.42 %	0.28 %	0.19 %	0.62 %	2.85 %	0.03 %	0.00 %
THREAD 1.23.1	93.78 %	0.42 %	0.29 %	0.43 %	2.82 %	0.03 %	0.00 %
THREAD 1.24.1	90.60 %	0.30 %	0.18 %	0.99 %	3.40 %	0.03 %	0.00 %
THREAD 1.25.1	94.65 %	0.20 %	0.12 %	0.63 %	0.70 %	0.03 %	0.00 %
THREAD 1.26.1	94.19 %	0.32 %	0.17 %	0.13 %	0.66 %	0.01 %	0.00 %
THREAD 1.27.1	93.14 %	0.21 %	0.11 %	0.10 %	1.42 %	0.04 %	0.00 %
Total	2,454.20 %	7.19 %	6.20 %	15.50 %	52.66 %	0.79 %	2.44 %
Average	90.90 %	0.27 %	0.23 %	0.57 %	1.95 %	0.03 %	0.09 %
Maximum	97.91 %	0.54 %	0.48 %	2.94 %	4.02 %	0.06 %	0.78 %
Minimum	85.98 %	0.13 %	0.11 %	0.10 %	0.22 %	0.00 %	0.00 %
StDev	3.16 %	0.09 %	0.08 %	0.65 %	1.21 %	0.01 %	0.20 %
Avg/Max	0.93	0.49	0.48	0.20	0.49	0.51	0.12

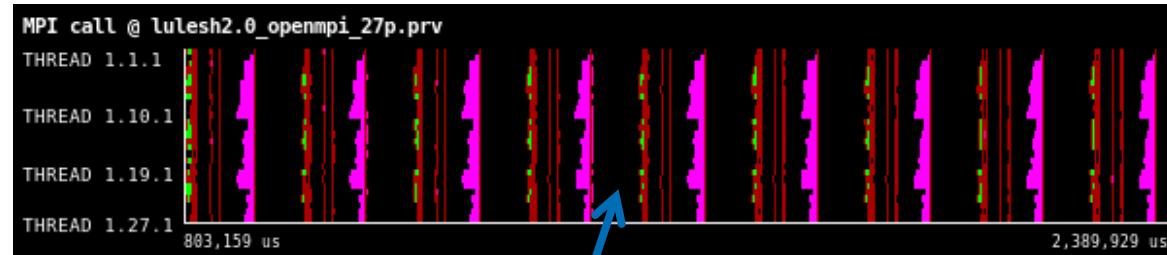
## Focus on the iterative part



Drag & drop on this area  
to zoom on the iterative  
region

MPI call profile @ lulesh2.0_openmpi_27p.prv							
	Default						
THREAD 1.17.1	93.37 %	0.38 %	0.30 %	0.24 %	0.22 %	0.01 %	0.34 %
THREAD 1.18.1	91.20 %	0.27 %	0.20 %	0.15 %	0.92 %	0.01 %	0.00 %
THREAD 1.19.1	92.87 %	0.18 %	0.14 %	0.13 %	0.80 %	0.05 %	0.00 %
THREAD 1.20.1	86.08 %	0.28 %	0.20 %	0.18 %	4.02 %	0.04 %	0.00 %
THREAD 1.21.1	85.98 %	0.19 %	0.13 %	2.41 %	1.18 %	0.06 %	0.78 %
THREAD 1.22.1	86.42 %	0.28 %	0.19 %	0.62 %	2.85 %	0.03 %	0.00 %
THREAD 1.23.1	93.78 %	0.42 %	0.29 %	0.43 %	2.82 %	0.03 %	0.00 %
THREAD 1.24.1	90.60 %	0.30 %	0.18 %	0.99 %	3.40 %	0.03 %	0.00 %
THREAD 1.25.1	94.65 %	0.20 %	0.12 %	0.63 %	0.70 %	0.03 %	0.00 %
THREAD 1.26.1	94.19 %	0.32 %	0.17 %	0.13 %	0.66 %	0.01 %	0.00 %
THREAD 1.27.1	93.14 %	0.21 %	0.11 %	0.10 %	1.42 %	0.04 %	0.00 %
Total	2,454.20 %	7.19 %	6.20 %	15.50 %	52.66 %	0.79 %	2.44 %
Average	90.90 %	0.27 %	0.23 %	0.57 %	1.95 %	0.03 %	0.09 %
Maximum	97.91 %	0.54 %	0.48 %	2.94 %	4.02 %	0.06 %	0.78 %
Minimum	85.98 %	0.13 %	0.11 %	0.10 %	0.22 %	0.00 %	0.00 %
StDev	3.16 %	0.09 %	0.08 %	0.65 %	1.21 %	0.01 %	0.20 %
Avg/Max	0.93	0.49	0.48	0.20	0.49	0.51	0.12

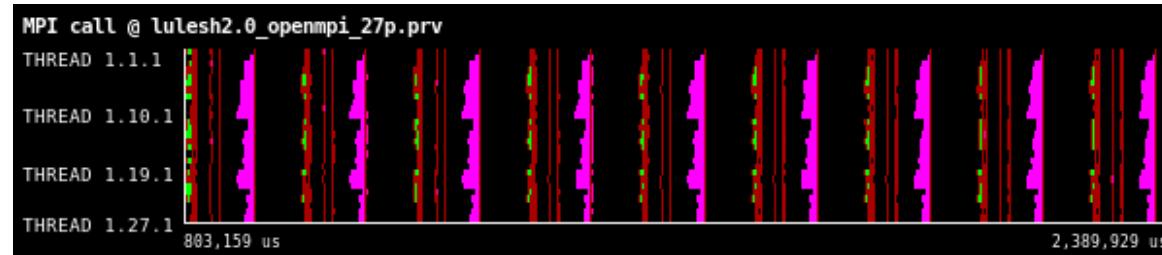
## Recalculate efficiency of iterative region



**Right click →  
Copy**

MPI call profile @ lulesh2.0_openmpi_27p.prv							
	Default						
<b>THREAD 1.17.1</b>	93.37 %	0.38 %	0.30 %	0.24 %	0.22 %	0.01 %	0.34 %
<b>THREAD 1.18.1</b>	91.20 %	0.27 %	0.20 %	0.15 %	0.92 %	0.01 %	0.00 %
<b>THREAD 1.19.1</b>	92.87 %	0.18 %	0.14 %	0.13 %	0.80 %	0.05 %	0.00 %
<b>THREAD 1.20.1</b>	86.08 %	0.28 %	0.20 %	0.18 %	4.02 %	0.04 %	0.00 %
<b>THREAD 1.21.1</b>	85.98 %	0.19 %	0.13 %	2.41 %	1.18 %	0.06 %	0.78 %
<b>THREAD 1.22.1</b>	86.42 %	0.28 %	0.19 %	0.62 %	2.85 %	0.03 %	0.00 %
<b>THREAD 1.23.1</b>	93.78 %	0.42 %	0.29 %	0.43 %	2.82 %	0.03 %	0.00 %
<b>THREAD 1.24.1</b>	90.60 %	0.30 %	0.18 %	0.99 %	3.40 %	0.03 %	0.00 %
<b>THREAD 1.25.1</b>	94.65 %	0.20 %	0.12 %	0.63 %	0.70 %	0.03 %	0.00 %
<b>THREAD 1.26.1</b>	94.19 %	0.32 %	0.17 %	0.13 %	0.66 %	0.01 %	0.00 %
<b>THREAD 1.27.1</b>	93.14 %	0.21 %	0.11 %	0.10 %	1.42 %	0.04 %	0.00 %
<b>Total</b>	2,454.20 %	7.19 %	6.20 %	15.50 %	52.66 %	0.79 %	2.44 %
<b>Average</b>	90.90 %	0.27 %	0.23 %	0.57 %	1.95 %	0.03 %	0.09 %
<b>Maximum</b>	97.91 %	0.54 %	0.48 %	2.94 %	4.02 %	0.06 %	0.78 %
<b>Minimum</b>	85.98 %	0.13 %	0.11 %	0.10 %	0.22 %	0.00 %	0.00 %
<b>StDev</b>	3.16 %	0.09 %	0.08 %	0.65 %	1.21 %	0.01 %	0.20 %
<b>Avg/Max</b>	0.93	0.49	0.48	0.20	0.49	0.51	0.12

## Recalculate efficiency of iterative region



**Right click →  
Paste → Time**

MPI call profile @ lulesh2.0_openmpi_27p.prv							
	Default						
<b>THREAD 1.17.1</b>	93.37 %	0.38 %	0.30 %	0.24 %	0.22 %	0.01 %	0.34 %
<b>THREAD 1.18.1</b>	91.20 %	0.27 %	0.20 %	0.15 %	0.92 %	0.01 %	0.00 %
<b>THREAD 1.19.1</b>	92.87 %	0.18 %	0.14 %	0.13 %	0.80 %	0.05 %	0.00 %
<b>THREAD 1.20.1</b>	86.08 %	0.28 %	0.20 %	0.18 %	4.02 %	0.04 %	0.00 %
<b>THREAD 1.21.1</b>	85.98 %	0.19 %	0.13 %	2.41 %	1.18 %	0.06 %	0.78 %
<b>THREAD 1.22.1</b>	86.42 %	0.28 %	0.19 %	0.62 %	2.85 %	0.03 %	0.00 %
<b>THREAD 1.23.1</b>	93.78 %	0.72 %	0.29 %	0.43 %	2.82 %	0.03 %	0.00 %
<b>THREAD 1.24.1</b>	90.60 %	0.30 %	0.18 %	0.99 %	3.40 %	0.03 %	0.00 %
<b>THREAD 1.25.1</b>	94.65 %	0.20 %	0.12 %	0.63 %	0.70 %	0.03 %	0.00 %
<b>THREAD 1.26.1</b>	94.19 %	0.32 %	0.17 %	0.13 %	0.66 %	0.01 %	0.00 %
<b>THREAD 1.27.1</b>	93.14 %	0.21 %	0.11 %	0.10 %	1.42 %	0.04 %	0.00 %
<b>Total</b>	2,454.20 %	7.19 %	6.20 %	15.50 %	52.66 %	0.79 %	2.44 %
<b>Average</b>	90.90 %	0.27 %	0.23 %	0.57 %	1.95 %	0.03 %	0.09 %
<b>Maximum</b>	97.91 %	0.54 %	0.48 %	2.94 %	4.02 %	0.06 %	0.78 %
<b>Minimum</b>	85.98 %	0.13 %	0.11 %	0.10 %	0.22 %	0.00 %	0.00 %
<b>StDev</b>	3.16 %	0.09 %	0.08 %	0.65 %	1.21 %	0.01 %	0.20 %
<b>Avg/Max</b>	0.93	0.49	0.48	0.20	0.49	0.51	0.12

## Efficiency of iterative region

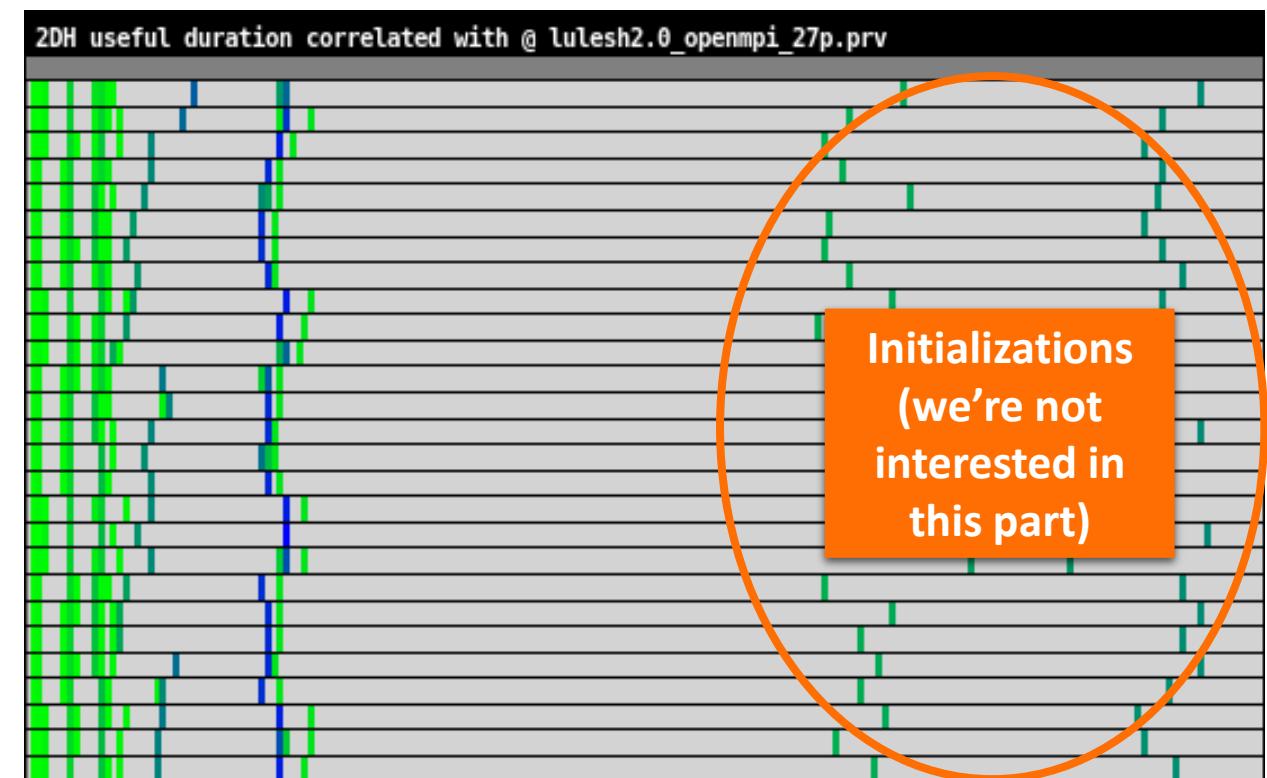
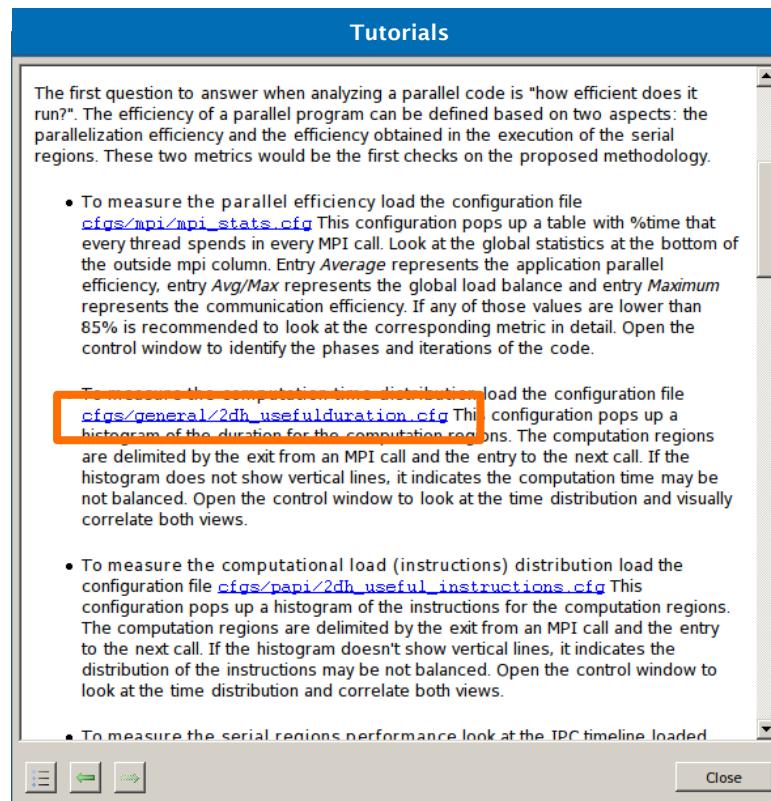
- 3 numbers to quickly describe the efficiency of your code
  - Parallel efficiency → % of time my program is computing (100% is perfect)
  - Comm efficiency → At least 1 process can finish all communications in 100 - Maximum % of the program's time (100% is perfect)
  - Load balance → Ratio of slow/fast processes (1 is perfectly balanced)
  - Any value below 85% (0.85)?
    - Pay attention there



MPI call profile @ lulesh2.0_openmpi_27p.prv								
	Default							
THREAD 1.17.1	90.61 %	0.52 %	0.41 %	0.33 %	0.31 %	7.65 %	0.16	
THREAD 1.18.1	87.65 %	0.36 %	0.27 %	0.21 %	1.13 %	10.20 %	0.17	
THREAD 1.19.1	89.90 %	0.24 %	0.19 %	0.18 %	1.16 %	8.14 %	0.18	
THREAD 1.20.1	80.40 %	0.38 %	0.27 %	0.25 %	5.54 %	12.97 %	0.19	
THREAD 1.21.1	80.14 %	0.26 %	0.17 %	3.31 %	1.68 %	14.28 %	0.17	
THREAD 1.22.1	80.85 %	0.37 %	0.26 %	0.89 %	3.89 %	13.58 %	0.17	
THREAD 1.23.1	91.44 %	0.57 %	0.38 %	0.62 %	3.89 %	2.94 %	0.16	
THREAD 1.24.1	86.87 %	0.41 %	0.25 %	1.38 %	4.71 %	6.23 %	0.17	
THREAD 1.25.1	92.48 %	0.27 %	0.16 %	0.92 %	0.93 %	5.06 %	0.18	
THREAD 1.26.1	91.76 %	0.44 %	0.23 %	0.17 %	0.96 %	6.25 %	0.18	
THREAD 1.27.1	90.23 %	0.29 %	0.14 %	0.12 %	2.03 %	7.01 %	0.18	
Total	2,355.04 %	9.64 %	8.43 %	21.58 %	72.49 %	228.18 %	4.64	
Average	87.22 %	0.36 %	0.31 %	0.80 %	2.68 %	8.45 %	0.17	
Maximum	97.63 %	0.73 %	0.66 %	4.04 %	5.54 %	14.37 %	0.19	
Minimum	80.14 %	0.18 %	0.14 %	0.12 %	0.31 %	0.03 %	0.16	
StDev	4.51 %	0.13 %	0.12 %	0.89 %	1.66 %	3.68 %	0.01	
Avg/Max	0.89	0.49	0.47	0.20	0.48	0.59	0.	

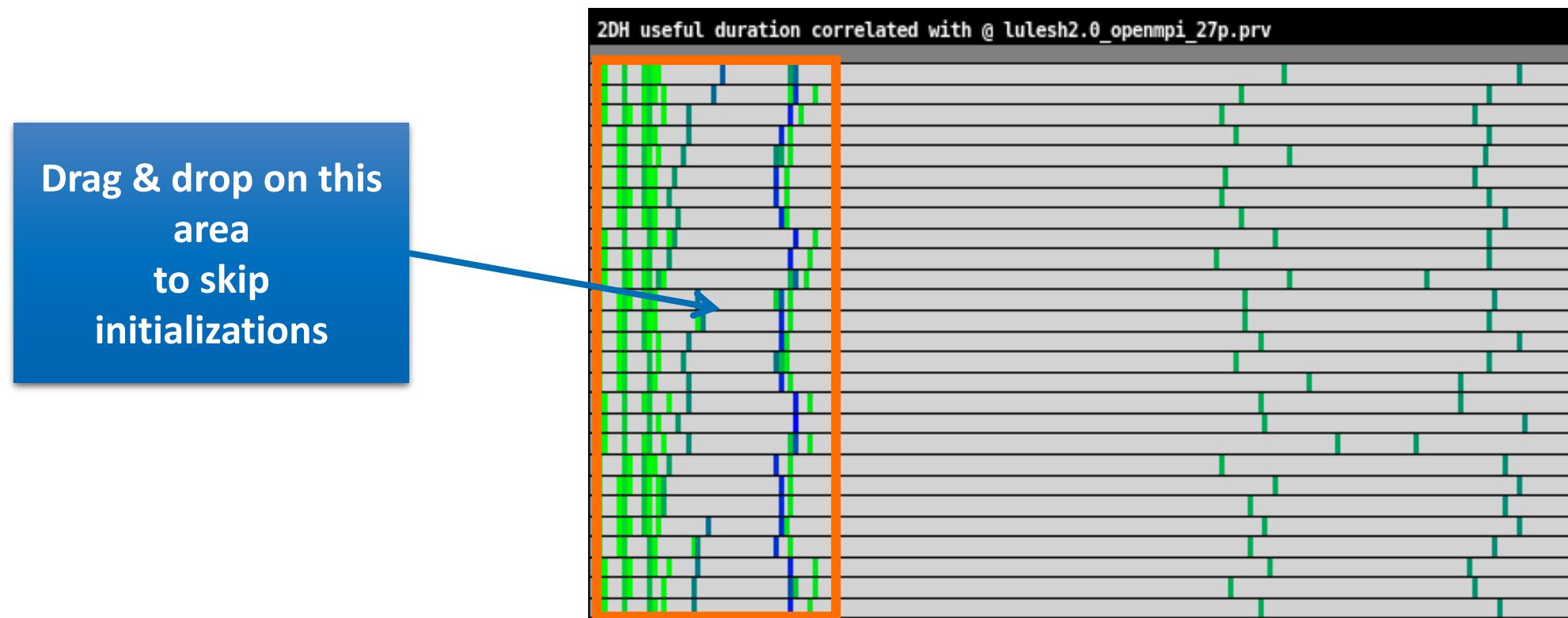
## Computation time distribution

- Click on “2dh\_usefulduration.cfg” (2nd link) → Shows **time computing**



## Focus on the iterative part

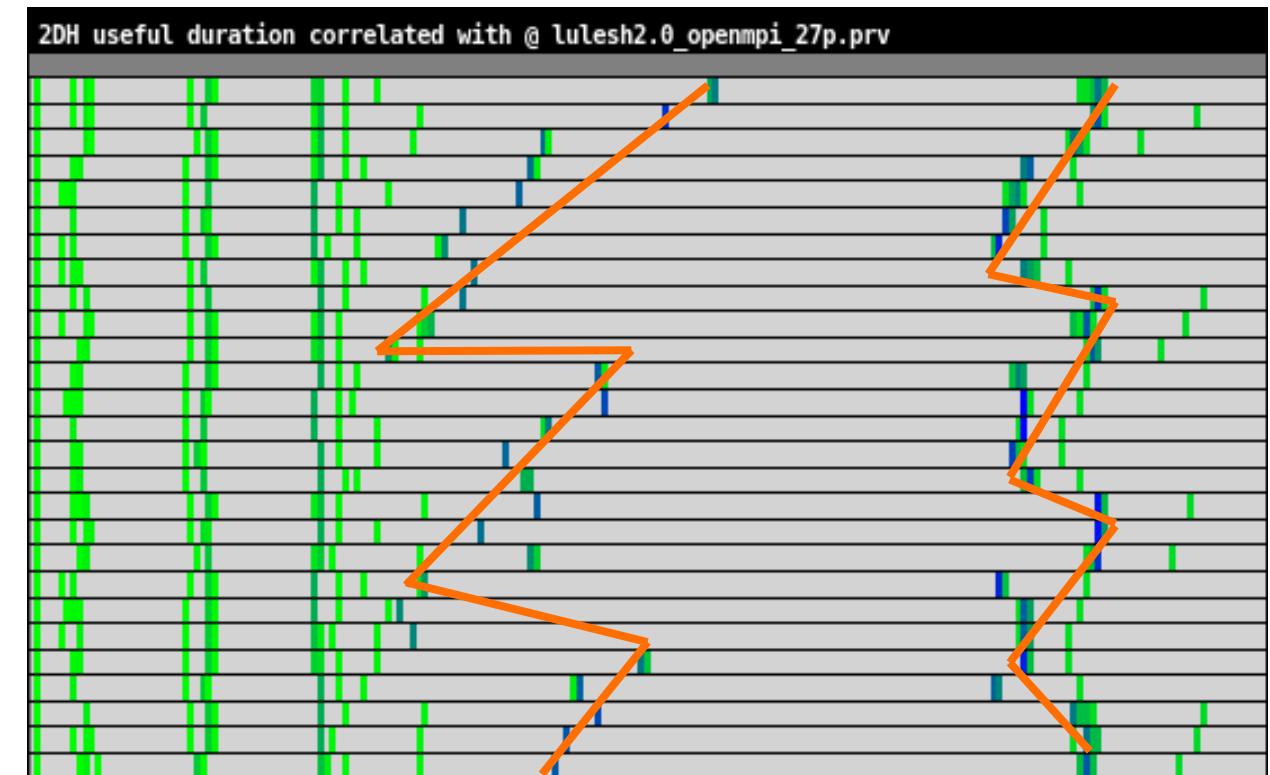
- Click on “2dh\_usefulduration.cfg” (2nd link) → Shows **time computing**



## Computation time distribution

- Click on “2dh\_usefulduration.cfg” (2nd link) → Shows **time computing**

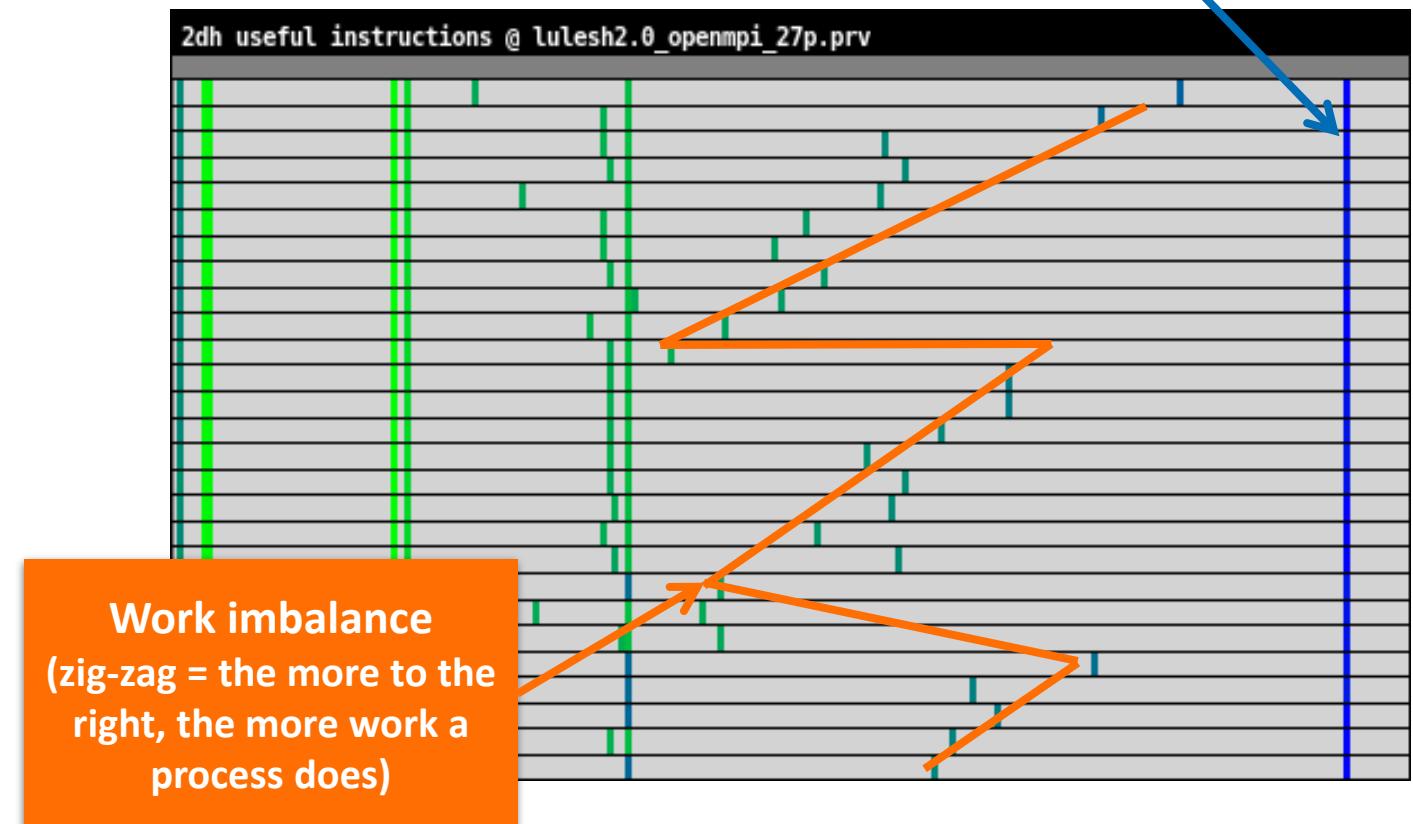
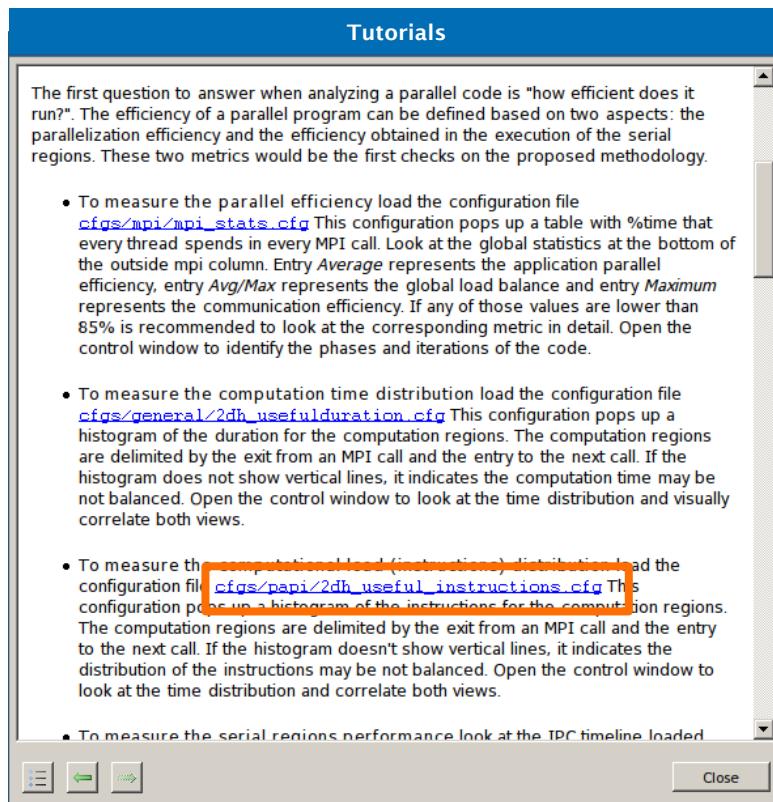
Duration imbalance  
(zig-zag = the more to  
the right, the more time  
a process takes)



# Computation load distribution

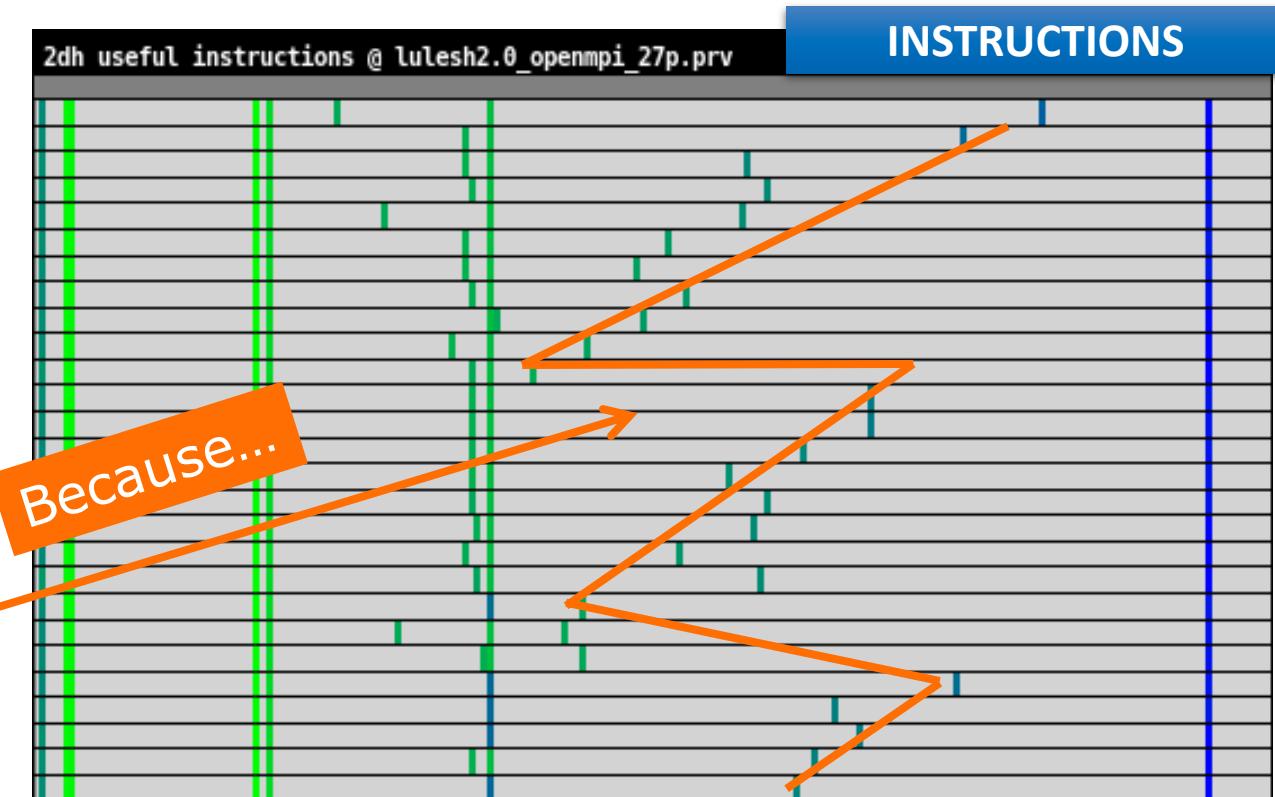
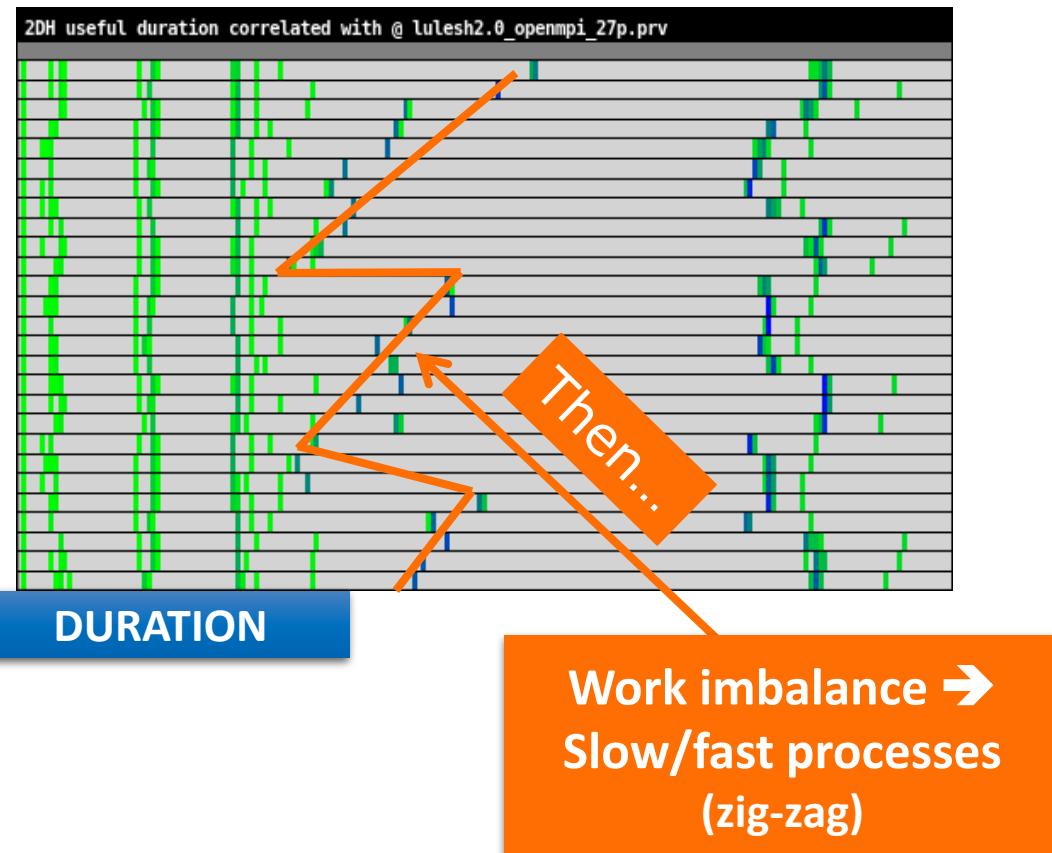
- Click on “2dh\_useful\_instructions.cfg” (3rd link) → Shows **amount of work**

Perfect work distribution in other parts of the program (straight line)



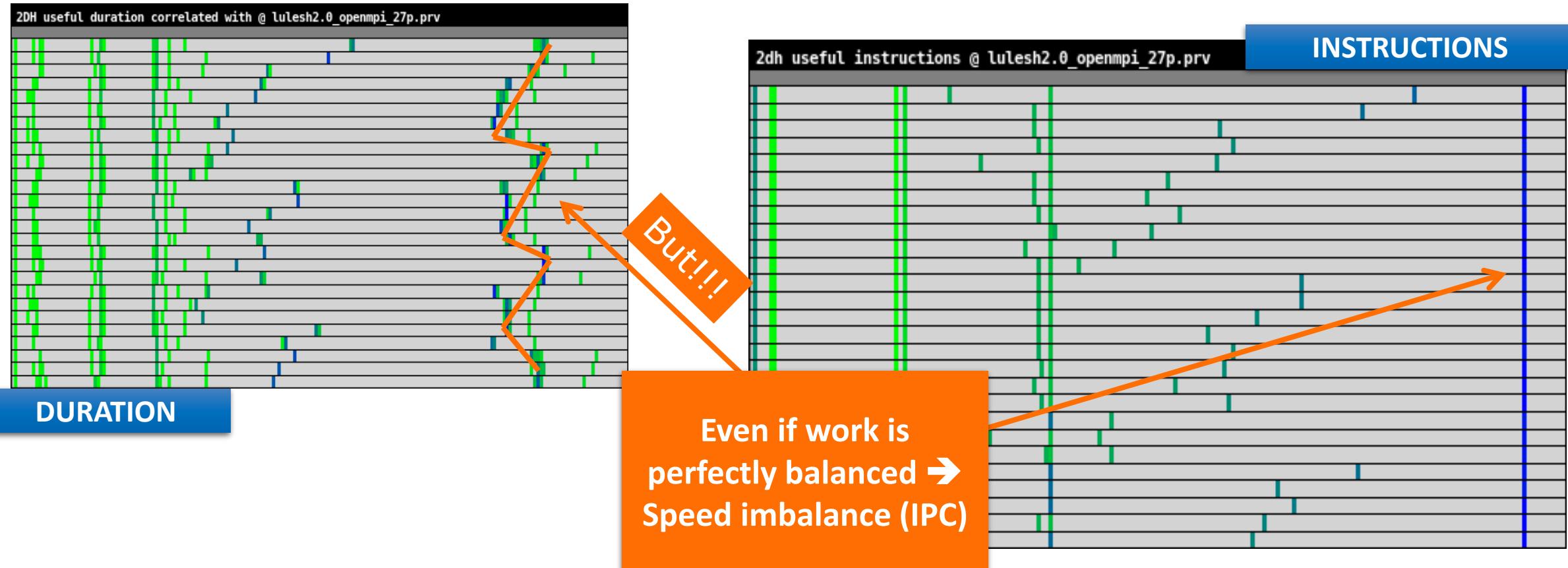
## Computation load distribution

- Comparing the two histograms → **Similar shapes** → Work distribution determines time computing



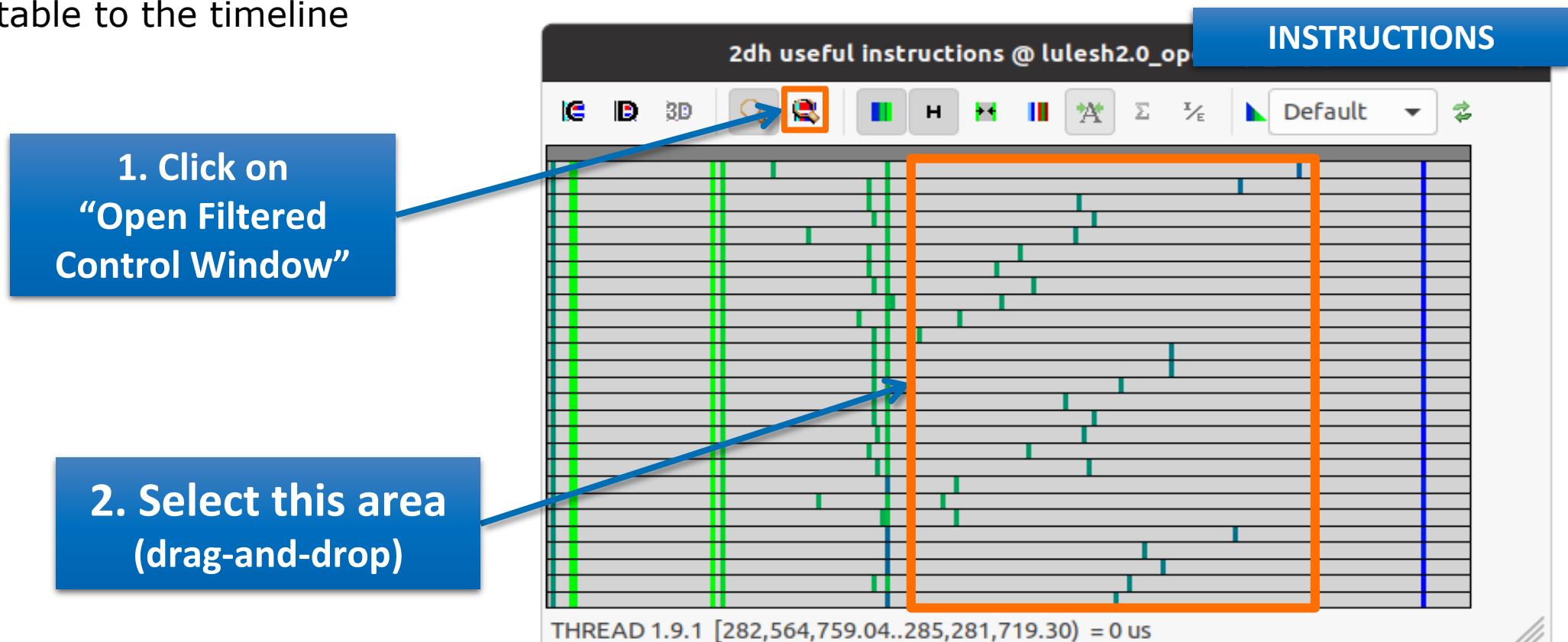
## Computation load distribution

- Comparing two histograms → **Not similar shapes** → Balanced workload executes at variable speed



## Where does this happen?

- Go from the table to the timeline



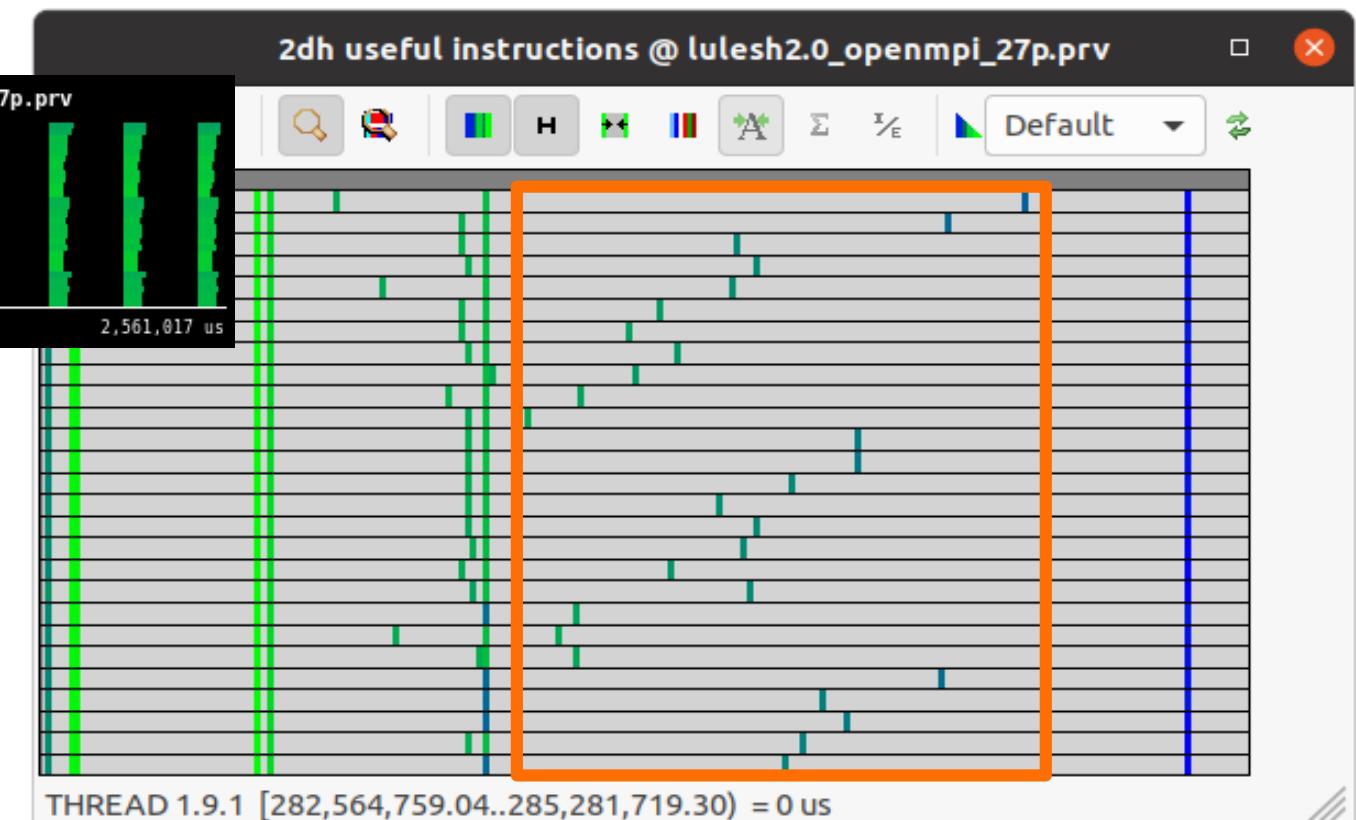
## Where does this happen?

- Go from the table to the timeline



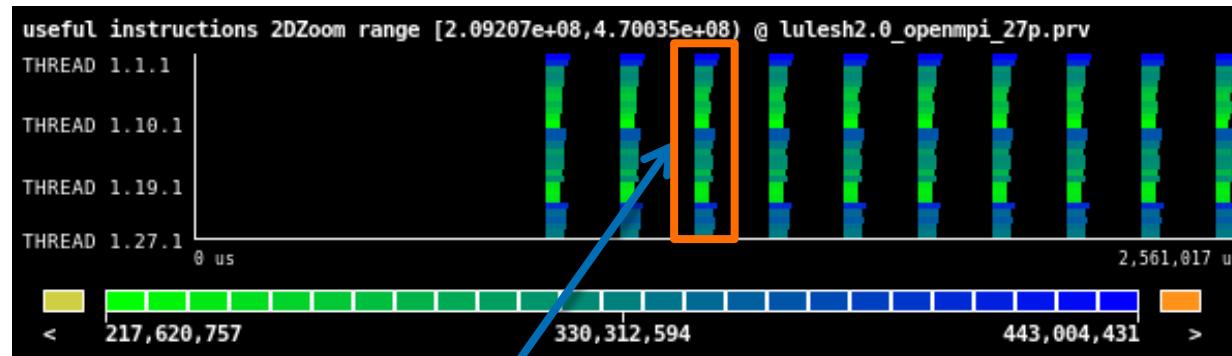
Clicking here always rescales the gradient. Same as...

Right click →  
Fit Semantic Scale → Fit Both  
(do this when colors are too similar)



## Where does this happen?

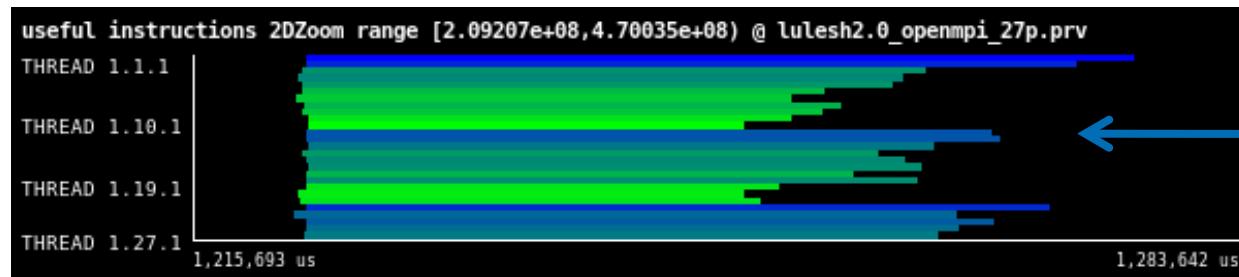
- **Slow** & **Fast** at the same time? → Imbalance



Zoom into  
1 of the iterations  
(by drag-and-dropping)

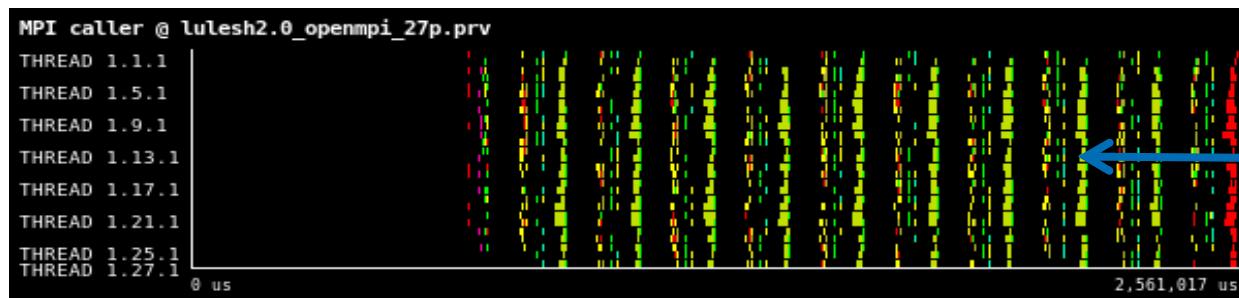
## Where does this happen?

- **Slow** & **Fast** at the same time? → Imbalance



1. Right click →  
Copy

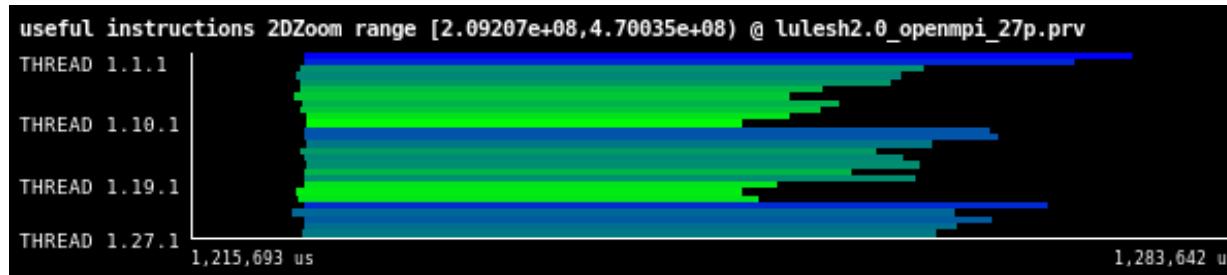
- Hints → Call stack references → Caller function



2. Right click →  
Paste → Time

## Where does this happen?

- **Slow** & **Fast** at the same time? → Imbalance

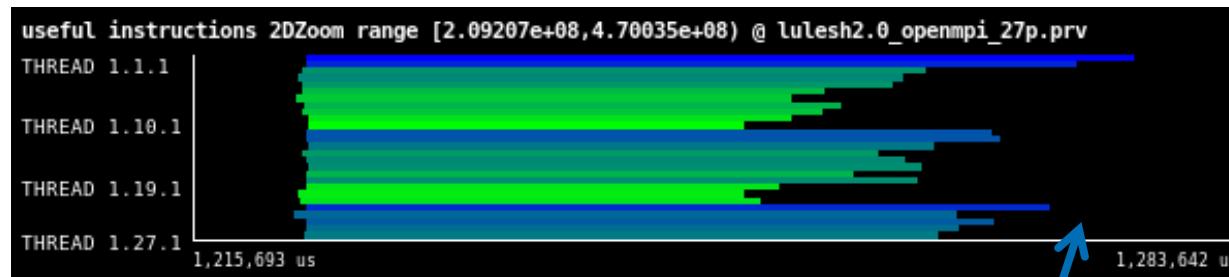


- Hints → Call stack references → Caller function

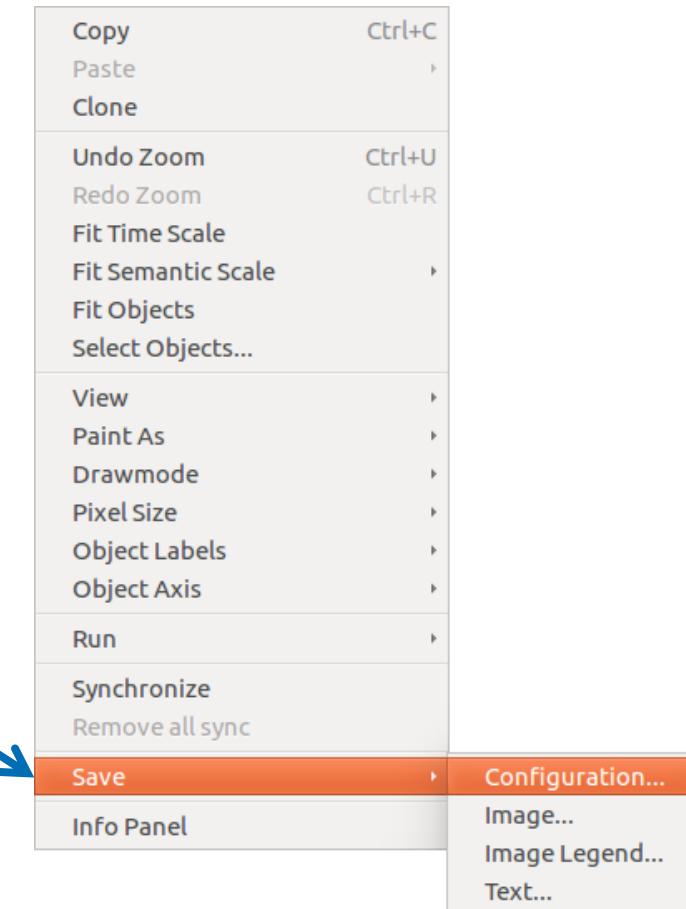


User code with  
imbalanced workload  
between **CommMonoQ** &  
**TimeIncrement**  
(you can hover the mouse to see the  
legend over the colors)

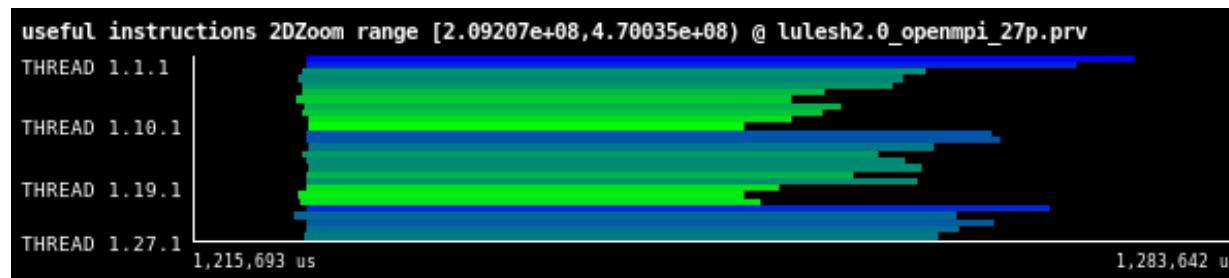
## Save CFG's (method 1)



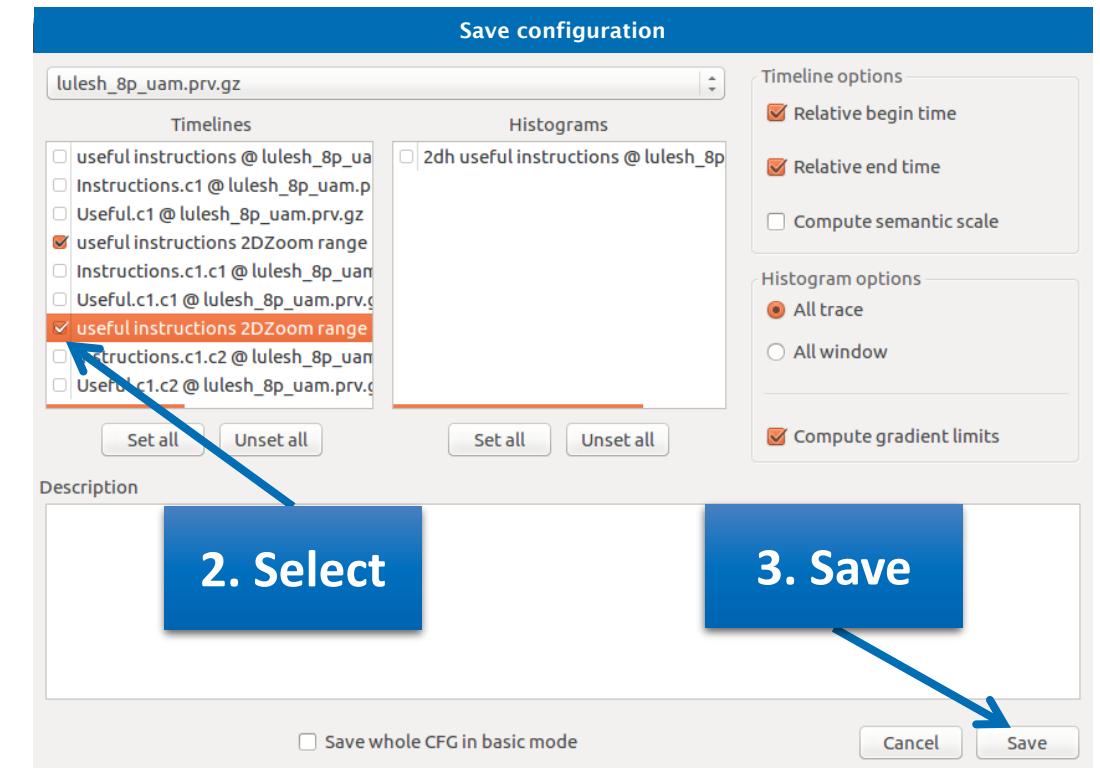
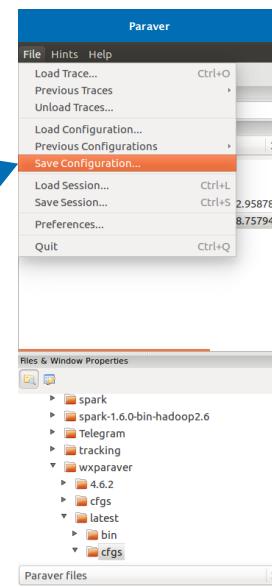
Right click on  
timeline



## Save CFG's (method 2)

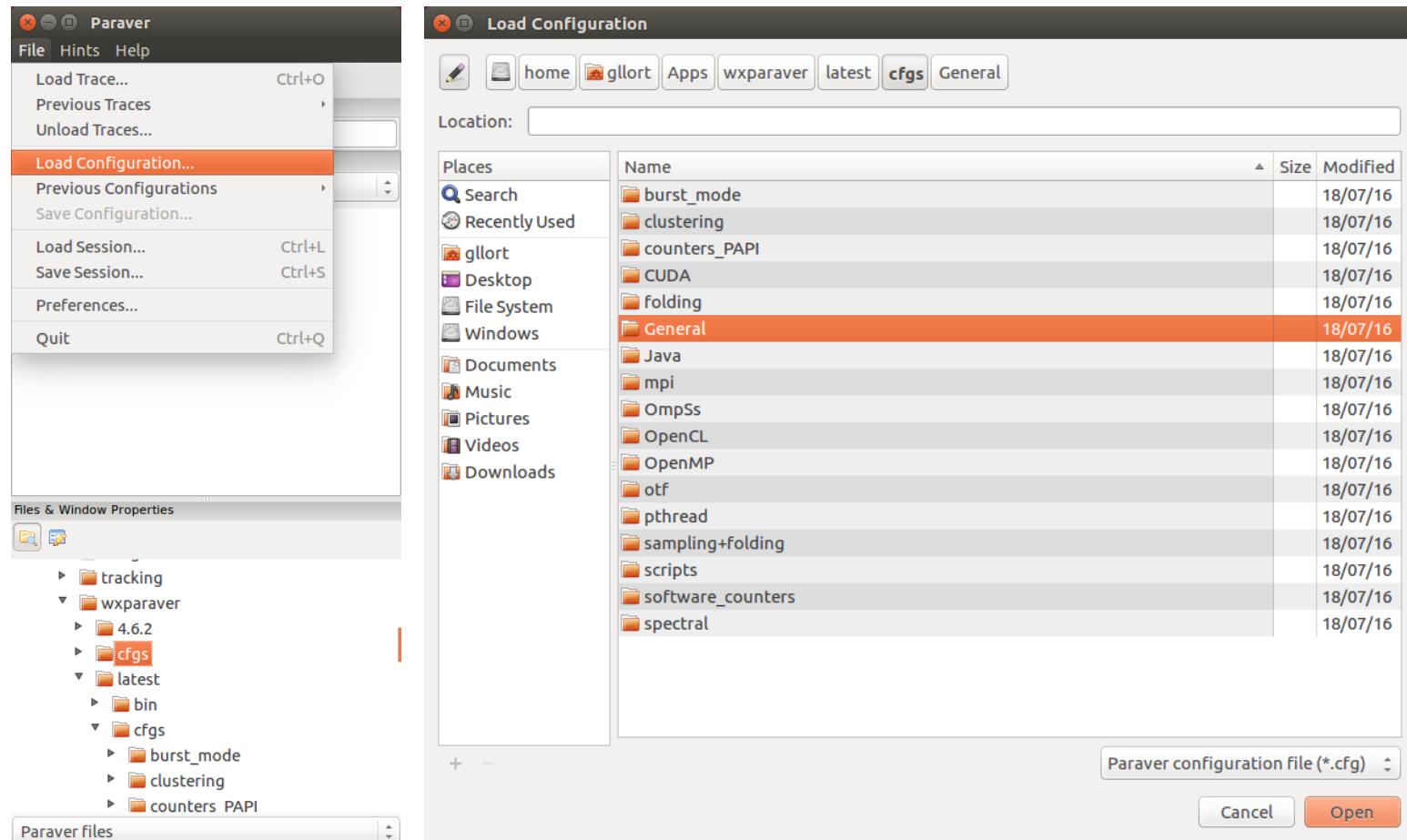


1. File → Save Configuration



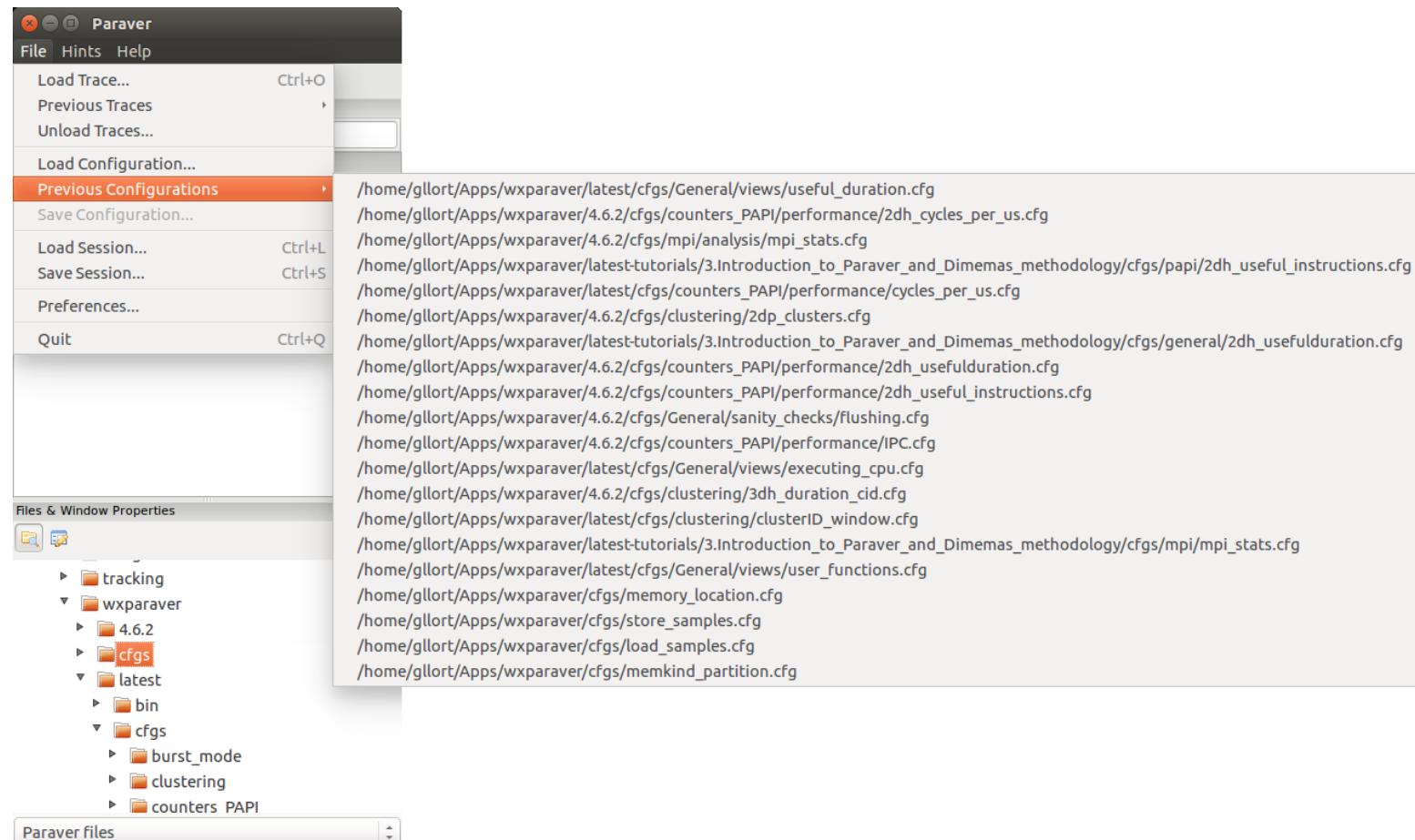
## CFG's distribution

- Paraver comes with many included CFG's: File → Load Configuration → Apply any CFG to any trace!



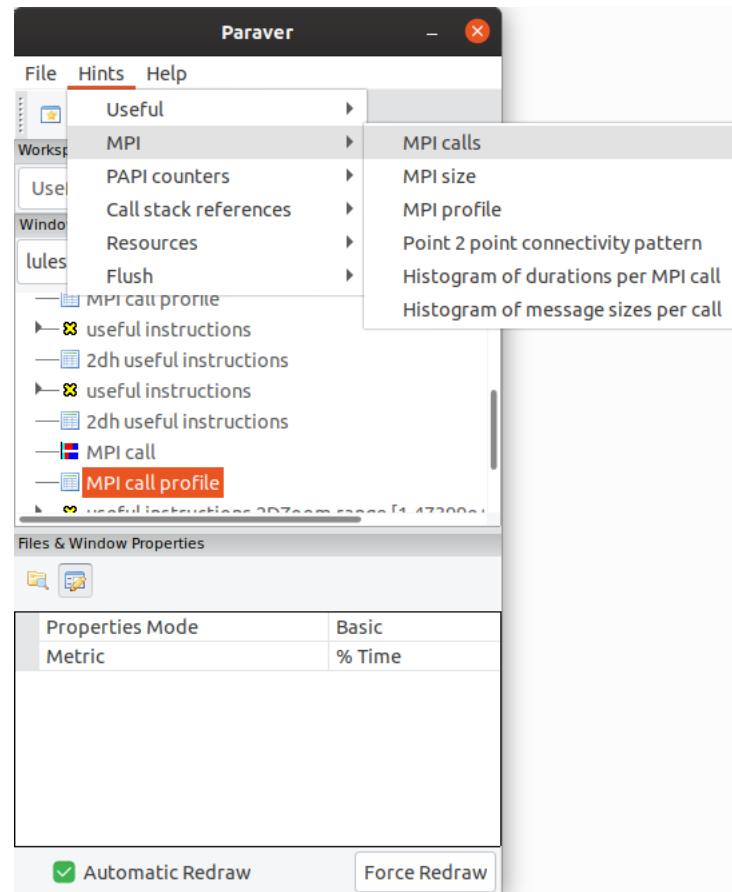
## CFG's distribution

- Recently opened CFG's are always at hand: File → Previous Configurations



## Hints: a good place to start!

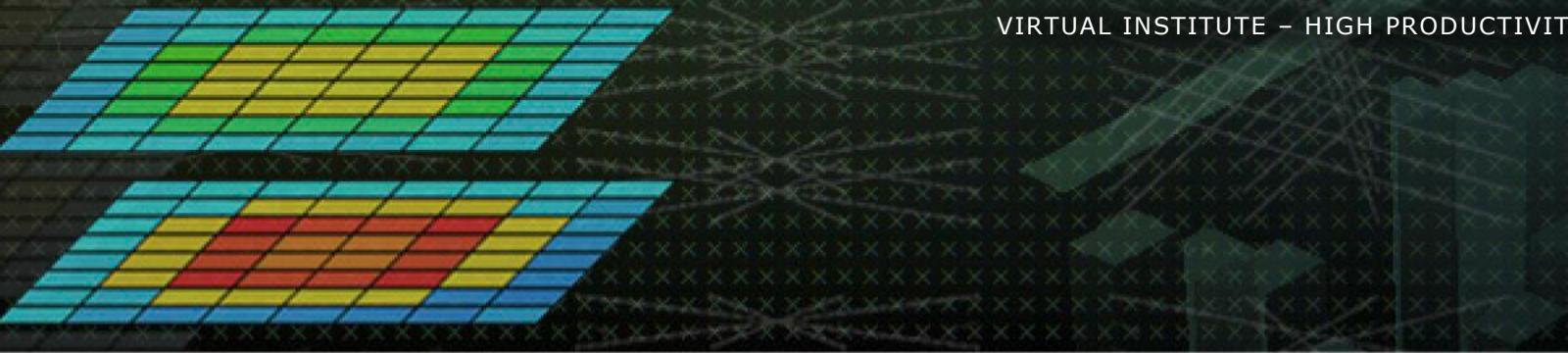
- Paraver suggests CFG's based on the contents of the trace



## Do it on your code!

---

- Follow guidelines from slides 7-15 to your own code to get a trace
  - There are more examples of tracing scripts for different programming models under  
\$EXRAE\_HOME/share/examples
- Follow guidelines from slides 17-35 to conduct an initial analysis
  - The usual suspects:
    - Parallel Efficiency is low? Load balance issues?
    - Imbalances in the durations of computations?
    - Are these caused by work imbalance?



# Cluster-based analysis

## Use clustering analysis

---

- Run clustering

```
hawk> cd $HOME/tools-material/clustering  
hawk> cat ./run_clustering.sh ← Have a look, it's just a wrapper for the tool BurstClustering  
hawk> ./run_clustering.sh
```

- If you didn't get your own trace, you can edit `run_clustering.sh` and change the input trace to use a prepared one from:

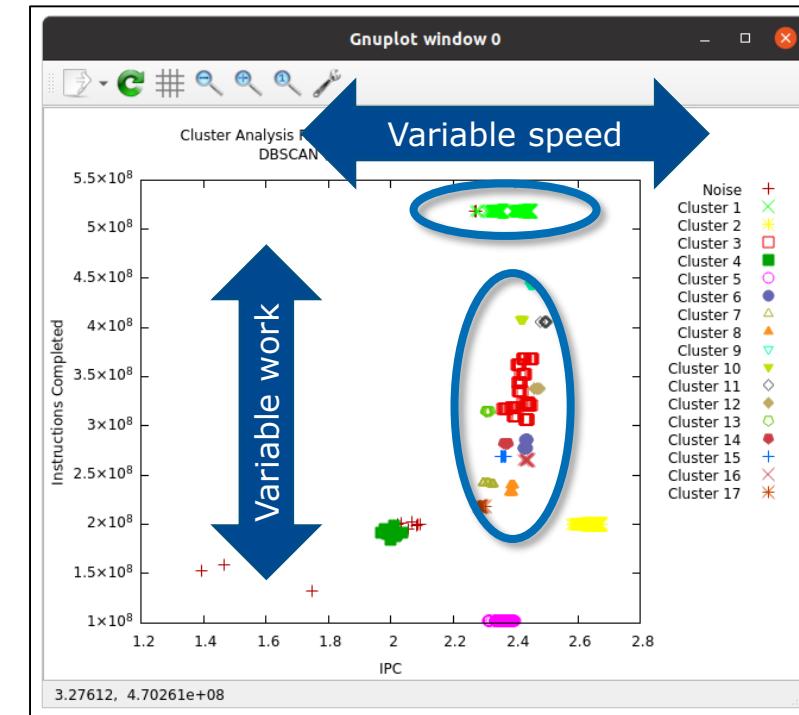
```
-i ../traces/lulesh2.0_openmpi_27p.prv
```

## Cluster-based analysis

- Check the resulting scatter plot

```
hawk> gnuplot lulesh2.0_openmpi_27p_clustered.IPC.PAPI_TOT_INS.gnuplot
```

- Identify main computing trends
- Work (Y) vs. Speed (X)
- Look at the clusters shape
  - Variability in both axes indicate **potential imbalances**



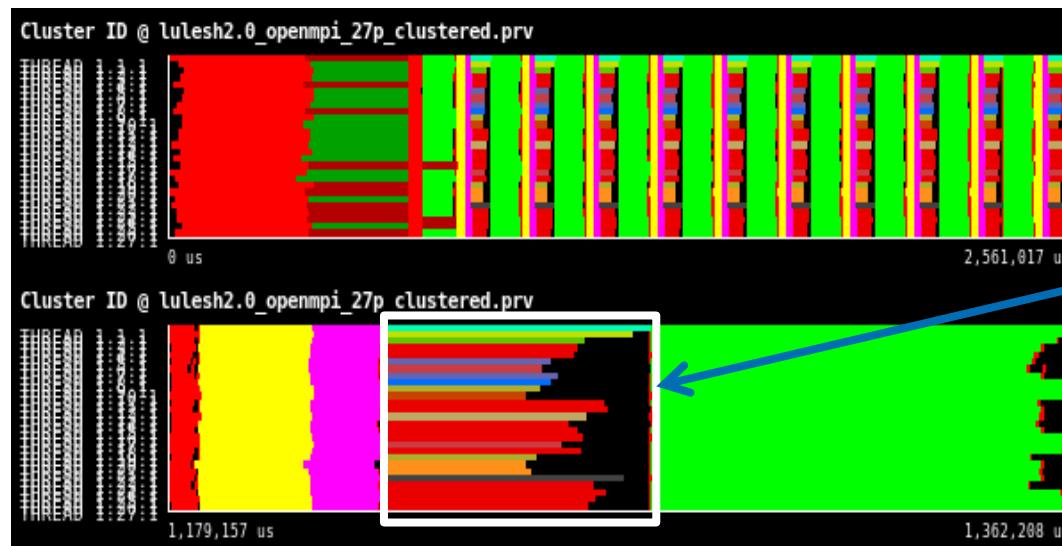
## Correlating scatter plot and time distribution

- Copy the clustered trace to your laptop and open it with Paraver:

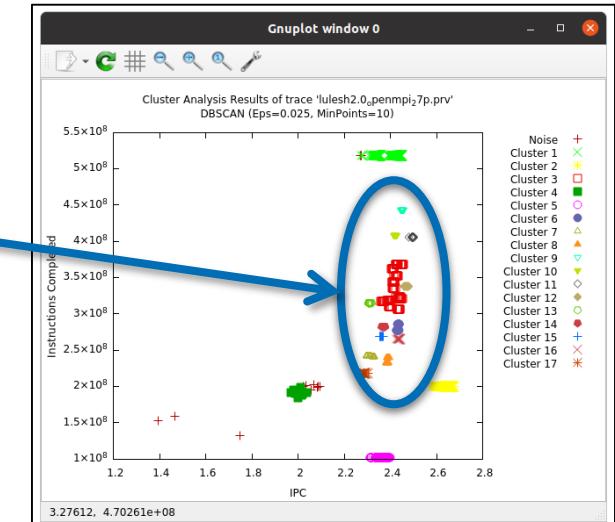
```
laptop> $HOME/paraver/bin/wxparaver <path-to>/lulesh2.0_openmpi_27p_clustered.prv
```

- Display the distribution of clusters over time

- File → Load configuration → \$HOME/paraver/cfgs/clustering/clusterID\_window.cfg



Variable work /  
speed  
+  
Simultaneously  
@ different  
processes  
=  
Imbalances



# BSC Tools Hands-On

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(tools@bsc.es)  
Barcelona Supercomputing Center

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