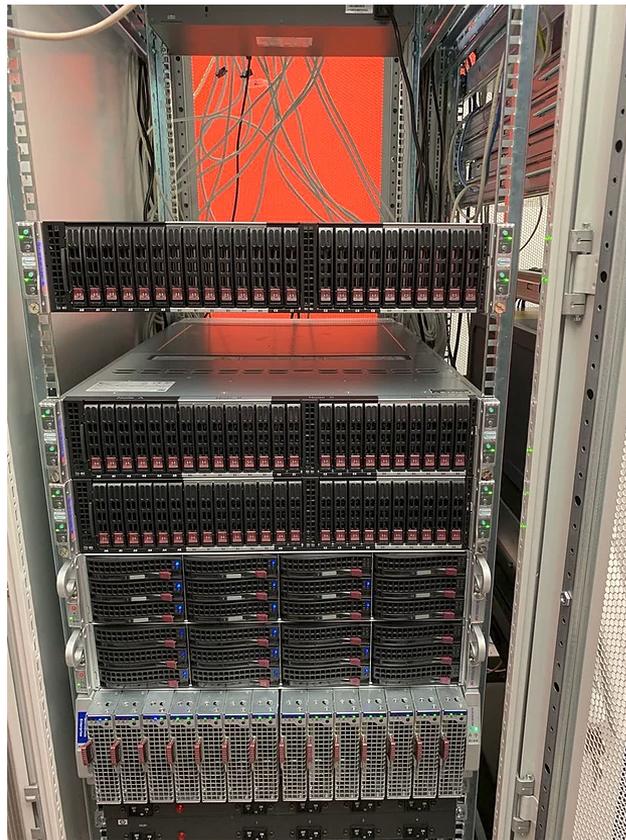


# ZTF computer cluster



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# Development

- 1) 2016 UKF Lazić 335k HRK
- 2) 2017 UKF Lazić 70k HRK
- 3) 2017 HrZZ Franchini 100k HRK
- 4) 2019 CEMS Lončarić 170k HRK
- 5) 2019 Sinergija ZTF 400k HRK
- 6) 2020 PZS Lončarić 340k HRK
- 7) 2021 HrZZ Grisanti 300k HRK
- 8) 2021 Hrzz Lončarić 330k HRK

# Significance

- The only capital equipment of ZTF
- Total over 2M HRK
- Enables competitive research in theoretical physics, especially in condensed matter
- Potential for commercial activities

# Technical details - hardware

- 2 management/storage nodes, RAID 10 storage 40TB, RAID 5 storage 18 TB
- 28 nodes: 8 Intel Xeon cores + 64 GB RAM
- 8 nodes: 48 AMD Epyc cores + 256 GB RAM
- 4 nodes: 64 AMD Epyc + 512 GB RAM
- 4 nodes: 64 AMD Epyc + 256 GB RAM
- 1 GPU server for ML
- 10 G network between all nodes and storage

# Technical details - disks

- /home (slow) and /storage (fast) available from everywhere
- /scratch (very fast) on some nodes
- No backup!

# Technical details - software

- OpenHPC
- Centos 8
- SLURM
- Environment Module System
- Anaconda for Python
- Intel and GNU compilers

# How to use

- Send me e-mail with request for opening the account
- Choose username and password
- Login via ssh `username@10.133.24.18`  
(available only from IRB internal network)
- Prepare submission script, software, inputs
- Run job, collect data and enjoy

# Submission script example - serial

```
#!/bin/bash
#SBATCH --partition=serial
#SBATCH --job-name=test
#SBATCH --cpus-per-task=1
#SBATCH --mem=2gb
#SBATCH --nodes=1
#SBATCH --ntasks-per-node=1
#SBATCH --output=%x-%j.out
#SBATCH --error=%x-%j.err
```

```
module load Anaconda3
```

```
python test.py > log
```

# Submission script example - parallel

```
#!/bin/bash
#SBATCH --partition=normal
#SBATCH --job-name=test
#SBATCH --cpus-per-task=1
#SBATCH --mem=100gb
#SBATCH --nodes=1
#SBATCH --ntasks-per-node=24
#SBATCH --output=%x-%j.out
#SBATCH --error=%x-%j.err

module purge
module load QE/6.8
export MKL_CBWR="AVX2"
export I_MPI_FABRICS=shm:ofi
ulimit -s unlimited
mpiexec.hydra -bootstrap slurm -n $SLURM_NTASKS pw.x -i scf.in > scf.out
```

# Typical use

- Submit the script to SLURM with:  
    `sbatch script.sh`
- Check status with:  
    `squeue`
- Copy results to your PC or backup:  
    `scp ...`

# Support

- No support, try:
  - 1) Google
  - 2) Ask colleague with more experience how to google
  - 99) Ask me, but understand that I manage the cluster in my free time