Running parallel jobs on the Linux-Cluster

On all HPC systems at LRZ, the SLURM scheduler is used to execute parallel jobs. This document describes usage, policies and resources available for submission and management of such jobs.

- Examples, Policies and Commands
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Examples, Policies and Commands

| Examples | provides example job scripts which cover the most common usage patterns. |
| Policies | provides information about the policies, such as memory limits, run time limits etc. |
| SLURM Workload Manager | lists SLURM commands and options, and explains them, making appropriate recommendations where necessary. Provides hints for aborting jobs. |

Detailed Instructions (for Beginners)

All parallel programs in the parallel segments of the cluster must be started up using either

- an interactive SLURM shell
- a SLURM batch script

In order to access the SLURM infrastructure described here, please first log in to a login node of the cluster as described in the introduction.

This document provides information on how to configure, submit and execute SLURM jobs, as well as information about batch processing policies. In particular, please be aware that misuse of the resources described here can result in the invalidation of the violating account. In particular, all parallel runs should always use either a salloc shell (for testing) or a scripted SLURM job.

Interactive SLURM shell for parallel testing

For performing program testing and short runs the following sequence of commands can be used: First, salloc is invoked to reserve the needed resources. Then, mpiexec can be used to start up a program on these resources.

<table>
<thead>
<tr>
<th>Command</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>salloc</td>
<td>Use at LRZ the srun command only for pure OpenMP Jobs. Please invoke this command from lxlogin5,6,7.</td>
</tr>
<tr>
<td>export OMP_NUM_THREADS=28</td>
<td></td>
</tr>
<tr>
<td>srun ./myprog.exe</td>
<td></td>
</tr>
<tr>
<td>exit</td>
<td></td>
</tr>
<tr>
<td>mpiexec ./myprog.exe</td>
<td></td>
</tr>
<tr>
<td>exit</td>
<td></td>
</tr>
<tr>
<td>salloc --ntasks=56</td>
<td>Start an MPP mode Intel MPI program using two 28-way nodes on the CooLMUC-2 cluster</td>
</tr>
<tr>
<td>mpiexec ./myprog.exe</td>
<td>Please invoke this command from lxlogin5,6,7.</td>
</tr>
<tr>
<td>exit</td>
<td></td>
</tr>
<tr>
<td>salloc --ntasks=8 --cpus-per-task=7</td>
<td>Start a hybrid mode Intel MPI program on the CooLMUC-2 cluster using 8 MPI tasks, with 7 OpenMP threads per task (2 nodes will be needed).</td>
</tr>
<tr>
<td>export OMP_NUM_THREADS=7</td>
<td>Please invoke this command from lxlogin5,6,7. Note that currently there are 28 cores per node available on CooLMUC-2, but no hyperthreading.</td>
</tr>
<tr>
<td>mpiexec -n 8 ./myprog.exe</td>
<td></td>
</tr>
<tr>
<td>exit</td>
<td></td>
</tr>
</tbody>
</table>
salloc --nodes=2 --tasks-per-node=3
export OMP_NUM_THREADS=8
mpiexec -n 6 ./myprog.exe
exit

by default, a SLURM shell generated via salloc will run for **15 minutes**. This interval can be extended to the partition maximum by specifying a suitable `--time=hh:mm:ss` argument. Also, the `--partition` option can be used to explicitly specify a desired partition, but we advise against doing so because the different login nodes have different environments that are incompatible.

**Notes and Warnings:**

- Only application/commands which are started with mpiexec are executed on the allocated nodes. All other commands will still be executed on the login node. This might block the login node for other users. A workaround would be to start memory or time consuming commands with "mpiexec -n 1", even if they are serial, optionally packing them into a script and starting it with mpiexec. Try "mpiexec -n 2 hostname" and compare the output with that of just typing "hostname".

- SLURM also has its own `srun` command which you can try to use.

- Once the allocation expires, the program will be signalled and killed; further programs can not be started. Please issue the exit command and start a new allocation.

## Batch Jobs

This type of execution method should be used for all production runs. A step-by-step recipe for the simplest type of parallel job is given, illustrating the use of the SLURM commands for users of the bash shell. See the documentation section at the end for pointers to more complex setups.

### Step 1: Edit a job script

The following script is assumed to be stored in the file `myjob.cmd`.

```
#!/bin/bash

#SBATCH -o /home/hpc/<group>/<user>/myjob.%j.%N.out
#SBATCH -D /home/hpc/<group>/<user>/mydir
#SBATCH -J <job_name>
#SBATCH --clusters=mpp2
#SBATCH --get-user-env
#SBATCH --ntasks=64
#SBATCH --cpus-per-task=2
#SBATCH --mail-type=end
#SBATCH --mail-user=<email_address>@<domain>

(Placeholder) standard output and error go there. Note that the directory where the output file is placed must exist before the job starts, and the full path name must be specified (no environment variable!). The %j encodes the job ID into the output file. The %N encodes the master node of the job and should be added since job IDs from different SLURM clusters may be the same.

(directory used by script as starting point (working directory))

(Placeholder) name of job (not more than 10 characters please)

The name "mpp2" specifies the cluster to be used - here the CoolMUC-2 MPP cluster.

Set user environment properly

Number of MPI tasks assigned to job. By default, SLURM will start as many MPI tasks per node as there are virtual CPUs in the node.

One the nodes where hyperthreading is enabled you need to this. Hyperthreading is enabled on:

```
ivymuc
```

Send an e-mail at job completion

(Placeholder) e-mail address (don't forget, and please enter a valid address!)
#SBATCH --export=NONE
Do not export the environment of the submitting shell into the job; while SLURM allows to also use ALL here, this is strongly discouraged, because the submission environment is very likely to be inconsistent with the environment required for execution of the job.

#SBATCH --time=08:00:00
maximum run time is 8 hours 0 minutes 0 seconds; this may be increased up to the queue limit

source /etc/profile.d/modules.sh
initialize module system

module load gsl # ... etc
load any required environment modules (may be needed if program is linked against shared libraries). "gsl" of course is only a placeholder.

mpiexec ./my_mpi_prog.exe
start MPI executable. The MPI variant used depends on the loaded module set; non-MPI programs may fail to start up - please consult the example jobs or the software-specific documentation for other startup mechanisms.

This script essentially looks like a bash script. However, there are specially marked comment lines ("control sequences"), which have a special meaning in the SLURM context explained on the right hand of the above table. The entries marked "Placeholder" must be suitably modified to have valid user-specific values.

For this script, the environment of the submitting shell will not be exported to the job's environment; the job will start 64 MPI tasks on as many cores.

Step 2: Submission procedure

The job script is submitted to the queue via the command

sbatch myjob.cmd

At submission time the control sequences are evaluated and stored in the queuing database, and the script is copied into an internal directory for later execution. If the command was executed successfully, the Job ID will be returned as follows:

Submitted batch job 65648.

It is a good idea to note down your Job ID's, for example to provide to LRZ HPC support as information if anything goes wrong. The submission command can also contain control sequences. For example,

sbatch --time=12:00:00 myjob.cmd

would override the setting inside the script, forcing it to run 12 instead of 8 hours.

Step 3: Checking the status of a job

Once submitted, the job will be queued for some time, depending on how many jobs are presently submitted. Eventually, more or less after previously submitted jobs have completed, the job will be started on one or more of the systems determined by its resource requirements.

The status of the job can be queried with the squeue --clusters=[all | cluster_name] command, which will give an output like

<table>
<thead>
<tr>
<th>CLUSTER: mpp1</th>
</tr>
</thead>
<tbody>
<tr>
<td>JOBID</td>
</tr>
<tr>
<td>65646</td>
</tr>
<tr>
<td>65647</td>
</tr>
<tr>
<td>65648</td>
</tr>
</tbody>
</table>

(assuming mpp1 is specified as the clusters argument) indicating that the job is queued. Once the job is running, the output would indicate the state to be "R" (=running), and would also list the host(s) it was running on. For jobs that have not yet started, the --start option, applied to kbd>squeue, will provide an estimate (!) for the starting time. The sinfo --clusters=[all | cluster_name] command prints out an overview of the status of all clusters or a particular clusters in the SLURM configuration.

Inspection and modification of jobs

Queued jobs can be inspected for their characteristics via the command

scontrol --clusters=<cluster_name> show jobid=<job ID>

which will print out a list of "Keyword=Value" pairs which characterize the job. As long as a job is waiting in the queue, it is possible to modify at least some of these; for example, the command

scontrol --clusters=<cluster_name> update jobid=65648 TimeLimit=04:00:00

would change the run time limit of the above-mentioned example job from 8 hours to 4 hours.
Deleting jobs from the queue

To forcibly remove a job from SLURM, the command

```
scancel --clusters=<cluster_name> <JOB_ID>
```

can be used. Please do not forget to specify the cluster! The scancel (1) man page provides further information on the use of this command.