









Available File Systems				
<ul> <li>The PC<sup>2</sup> provides three types of file systems:</li> </ul>				
<ul> <li>A shared file system located on an external filer. This file system is available on all clusters.</li> </ul>				
<ul> <li>Cluster local file systems. They provide fast access (typically realized by a parallel file system).</li> </ul>				
<ul> <li>Node local disks.</li> </ul>				
<ul> <li>Based on this file systems, every user can access the following directories:</li> </ul>				
Environment Variable	Purpose			
\$HOME	Login home, small data			
\$PC2GROUPS/HPC-LCO-SIMON	Group related data	Files needed for exercises		
\$PC2WORK	Temporary, data center wide	working data		
\$PC2SCRATCH	Temporary, system local working data			
\$CCS_TMPDIR	Temporary, node local data			
More details see				
https://wikis.upi-paderborn.de/pc2doc/PC2-FileSystems				
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Executing Programs on Arminius
Set Environment <ul> <li>module add intel/compiler</li> </ul>
Compile OpenMP program <ul> <li>icc –openmp –o pi pi.c</li> </ul>
Workload Management System (OpenCCS) <ul> <li>ccsinfo -a</li> <li>ccsinfo -s</li> <li>ccsinfo -smine</li> </ul>
Job Submission (example) <ul> <li>ccsalloc -Igroup=HPC-LCO-SIMON -t 5m -c 1 pi 1000000000</li> <li>ccsalloc -I -g HPC-LCO-SIMONres=rset=1:ncpus=6:ompthreads=6,place=free ./pi 1000000000</li> </ul>
Explicit Job Termination <ul> <li>ccskill <reqid></reqid></li> </ul>
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Executing MPI-Programs	
<ul> <li>MPI-Job Submission</li> <li>24 MPI-processes each associated with one core ccsalloc -I -c 24 ompi/hello_mpi</li> <li>2 MPI-processes each placed on a separate node ccsalloc -I -t 4m -n 2 ompi/hello_mpi</li> <li>8 MPI-processes each with 3 threads ccsalloc -I -t 5mres=rset=8:ncpus=3:mpiprocs=1:ompthreads=3:ib=true,place=pack / ompi/hello_mpi</li> <li>ccsalloc -I -t 5mres=rset=2:ncpus=12:mpiprocs=4:ompthreads=3:ib=true,place=excl / ompi/hello_mpi</li> <li>ccsalloc -I -t 5mres=rset=4:ncpus=6:mpiprocs=2:ompthreads=3:ib=true,place=free / ompi/hello_mpi</li> </ul>	
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